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Exploration Associates

CA71221

M4 Relief Road - Stage 2
Preliminary Chemical Investigation

150006
May 2000

Client:
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London
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W1P 6BQ

Engineer:
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150006

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APPENDIX

General Notes

1. INTRODUCTION

On the instruction of Ove Arup & Partners a preliminary chemical investigation was carried out by Exploration Associates for the proposed M4 Relief Road - Stage 2. The instruction to proceed with the investigation was received by a letter Ove Arup & Partners, dated 11 January 2000, reference 53600/69.

This factual report provides a description of the site, a summary of the investigation procedures adopted and presents the findings of the site works together, with results of the subsequent in situ and contaminant testing.

The work was carried out in general accordance with the relevant British Standards and the enclosed general notes.

2. THE SITE

The site comprised ten separate areas, described as follows:

Site 1 - Church Road: *Part of Parc Golf Centre immediately south west of the LG factory; identified as a former inert landfill site. Located at National Grid reference ST 273 839.*

Site 2 - Green Lane: *Field located on Wentlooge Levels, immediately south of the Great Western Railway line; identified as a former inert landfill site. Located at National Grid Reference ST 291 834.*

Site 3 - Newport Docks: *Area north of Alexandra South Dock, east of the River Ebbw and west of the River Usk. Located at National Grid Reference points ST 309 848 and ST 320 854.*

Site 4 - Docks East of Usk (Eastern Wharf): *Plot of land on north eastern bank of the River Usk. Located at National Grid Reference ST 326 855.*

Site 5 - Docks East of Usk (Marshall's Yard: *Works off Corporation Road, north east of River Usk. Located at National Grid Reference ST 328 855.*

Site 6 - Docks East of Usk (Birdport): *Plot of land immediately to the south west of the dead-end of Corporation Road and north of the container terminal. Located at National Grid Reference ST 332 854.*

Site 7 - Solutia Chemical Works: *Plot of land south of the Solutia Chemical Works (formerly Monsanto). Located at National Grid Reference ST 337 852.*

Site 8 - *Llanwern Research Laboratories: Fields off Meadows Road, south of Leeway Industrial Estate, Pye Corner. Located at National Grid Reference ST 346 854.*

Site 9 - *Llanwern Tippings: Fields on Caldicot Levels immediately south of Llanwern Steelworks. Located between National Grid Reference points ST 385 854 and ST 395 856.*

Site 10 - *Magor Depot: Depot immediately east of the brewery, south of Junction 23 of the M4. Located at National Grid Reference ST 419 876.*

3. FIELDWORK

3.1 General

The fieldwork was carried out during the period 25 January to the 16 February 2000 and generally comprised Window Sampling, the excavation of Trial Pits and in situ testing.

The fieldwork was carried out in general accordance with BS 5930: 1981.

3.2 Window Sampling

Thirty three 80mm, 60mm and 40mm nominal diameter boreholes were drilled using a combination of a hand held Window Sampler and a Geotool Tracked Rig, to depths of between 0.40 metres and 5.00 metres below existing ground level.

The depths and descriptive details of the strata encountered , comments on the groundwater conditions, details of samples taken and drilling progress are shown in the Exploratory Hole Records, Enclosure A.

During the course of the window sampling large disturbed samples (bulk) and small disturbed samples (jar) were obtained at regular intervals for identification and descriptive purposes, and to facilitate laboratory testing. On encountering groundwater, levels were recorded during 5 minute intervals for a 20 minute period of monitoring.

The locations of the boreholes are shown on the Exploratory Hole Location plan, Enclosure D.

3.3 Trial Pits

Twenty nine pits were excavated using a wheeled hydraulic excavator to depths of between 0.75 metres and 3.20 metres below existing ground level. The details of which are presented on the Trial Pit Records, Enclosure A.

3.4 In Situ Testing

During the course of the site works, in situ water testing was carried out for the following; temperature, pH value and conductivity. The results of which are presented in Enclosure B.

3.5 Survey

On completion of the fieldwork, a survey was undertaken by John Vincent Surveys under the supervision of Exploration Associates. The levels and co-ordinates are presented on the Exploratory Hole Records, Enclosure A.

4. LABORATORY TESTING

Laboratory testing schedules were prepared by Ove Arup & Partners . The resultant contaminant testing was undertaken TES Bretby laboratories, the results of which are presented in Enclosure C.

5. GROUND CONDITIONS

5.1 Published Geology

Reference to the British Geological Survey 1:50,000 Scale Series, Drift Map 249, Newport; indicates the site to be underlain by superficial deposits of Marine Alluvium. The solid geology of the area is shown to Mercia Mudstone of Triassic age unconformably overlying St Maughans Group and Raglan Marl of Devonian age.

5.2 Strata Encountered

For full details of the strata encountered, reference should be made to the Exploratory Hole Records. The exploratory holes generally confirmed the published geology, however a laterally impersistant layer of Made Ground was encountered at most locations.

Groundwater

Groundwater was encountered in some of the exploratory holes.

For full details of groundwater encountered, reference should be made to the individual Exploratory Hole Records.

It should be noted that groundwater levels may be subject to variations due to changes in climatic and other conditions and may therefore at times be different to those measured during the fieldworks.

FOR AND ON BEHALF OF EXPLORATION ASSOCIATES

A.Figgis
B.Sc(Hons)
ENGINEERING GEOLOGIST

A handwritten signature in dark ink, appearing to read 'R.H. Griffiths', with a long horizontal stroke extending to the right.

R.H. Griffiths
B.Sc(Hons), M.Sc., F.G.S., C.Geol
PRINCIPAL GEOTECHNICAL ENGINEER

EXPLORATION ASSOCIATES
AF/RHG/150006/Rev. 2/MAY 2000

REFERENCES

BS5930(1999). Code of Practice for Site Investigations.BSI

BS1377(1990). Method of Test for Soil for Civil Engineering Purposes.BSI.

British Geological Survey. 1:50000 Scale. Sheet 249, Newport.

Ordnance Survey of Great Britain. Sheet 171, Cardiff, Newport & surrounding area.

ENCLOSURE A
Exploratory Hole Records

Symbol Sheet

Window Sampling Records

CH1 to CH4, CH9 to CH33,
CH55, CH56, CH15A &
CH19A

Trial Pit Records

CH5 to CH8, CH34 to
CH54,CH47A,CH52A, CH53A
& CH54A

KEY TO SYMBOLS ON EXPLORATORY HOLE RECORDS

All linear dimensions are in metres or millimetres

DESCRIPTIONS

** : Drillers Description

SAMPLES

U () : Undisturbed 102mm diameter sample, () denotes number of blows to drive sampler
 U ()F, U ()P : F - not recovered, P - partially recovered
 U38 : Undisturbed 38mm diameter sample
 P(F),(P) : Piston sample, F - not recovered, P - partially recovered
 B : Bulk sample - disturbed
 D : Jar Sample - disturbed
 W : Water Sample
 CBR : California Bearing Ratio mould sample
 G : Gas Sample and depth of hole at time of sampling

CORE RECOVERY AND ROCK QUALITY

TCR : Total Core Recovery %
 SCR : Solid Core Recovery %
 RQD : Rock Quality Designation %
 FI : Fracture Index (discontinuities per metre) NI - not intact, NR - not recordable, NA - not applicable

GROUNDWATER

∇ : Groundwater strike
 $\bar{y}$: Groundwater level after standing period
 Date/Water : Date of shift (day/month)/Depth to water at end of previous shift shown above the date and depth to water at beginning of shift given below the date.

IN SITU TESTING

S : Standard Penetration Test - split barrel sampler
 C : Standard Penetration Test - solid 60° cone
 V(H)(R) : Vane Test (Hand) (R) demonstrates remoulded strength
 K(F), (C), (R), (P) : Permeability Test (falling, constant or rising head, packer)
 PT : Pressuremeter Test
 HP : Hand Penetrometer Test

MEASURED PROPERTIES

N : Standard Penetration Test - blows required to drive 300mm after seating drive
 $\frac{x}{y}$: Denotes x blows for y mm within the Standard Penetration Test
 $\frac{x^*}{y}$: Denotes x blows for y mm within the seating drive
 c_u : Undrained Shear Strength (kN/m²)
 CBR : California Bearing Ratio

ROTARY DRILLING SIZES

Index Letter	NOMINAL DIAMETER (mm)	
	Borehole	Core
N	75	54
H	99	76
P	120	92
S	146	113

EXPLORATORY HOLE SYMBOLS

Project

Contract

150006



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Figure

Sampling					Strata				
Depth	Type	Casing Depth	Date/ Water	SPT N (Cu)	Description	Depth (Thickness)	Level	Legend	
0.20-0.45	T		25/01 0000 DRY		Brown sandy slightly clayey TOPSOIL	G.L.	10.41		
0.45-0.90	B				Soft to firm brown slightly sandy very silty CLAY with occasional to some fine to coarse rounded to sub-rounded gravel of sandstone	0.20	10.21		
0.90-1.40	T				Soft to firm red brown sandy very silty CLAY with some rounded to sub-rounded fine to coarse gravel of sandstone	0.45	9.96		
1.45-1.90	B					(1.45)			
1.90-2.38	T					1.90	8.51		
2.38-2.70	T				Moderately compact brown sandy slightly silty clayey sub-rounded fine to coarse GRAVEL of sandstone	(0.48)			
2.70		NIL	DRY		Firm red brown very silty clayey highly weathered MUDSTONE	2.38	8.03		
					 End of Borehole.	2.70	7.71		
Equipment: Geotool R780 Tracked Rig					Groundwater		Sealed		
Borehole Dia (mm) Casing Dia (mm)					No. Struck Behaviour		Ground Level 10.41 m OD		
80 to 1.00m					No groundwater encountered		Coordinates 327310.67		
60 to 2.70m							183955.65 mE		
							mN		
							Drilled by DC		
							Logged by DC		
							Checked by AF		
Remarks Refusal at 2.70 (Stiff)									
See key sheet and appendices for explanations.									
Borehole Record					Project			Contract	
					M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners			150006	
Exploration Associates								Borehole	
								CH1(1 of 1)	

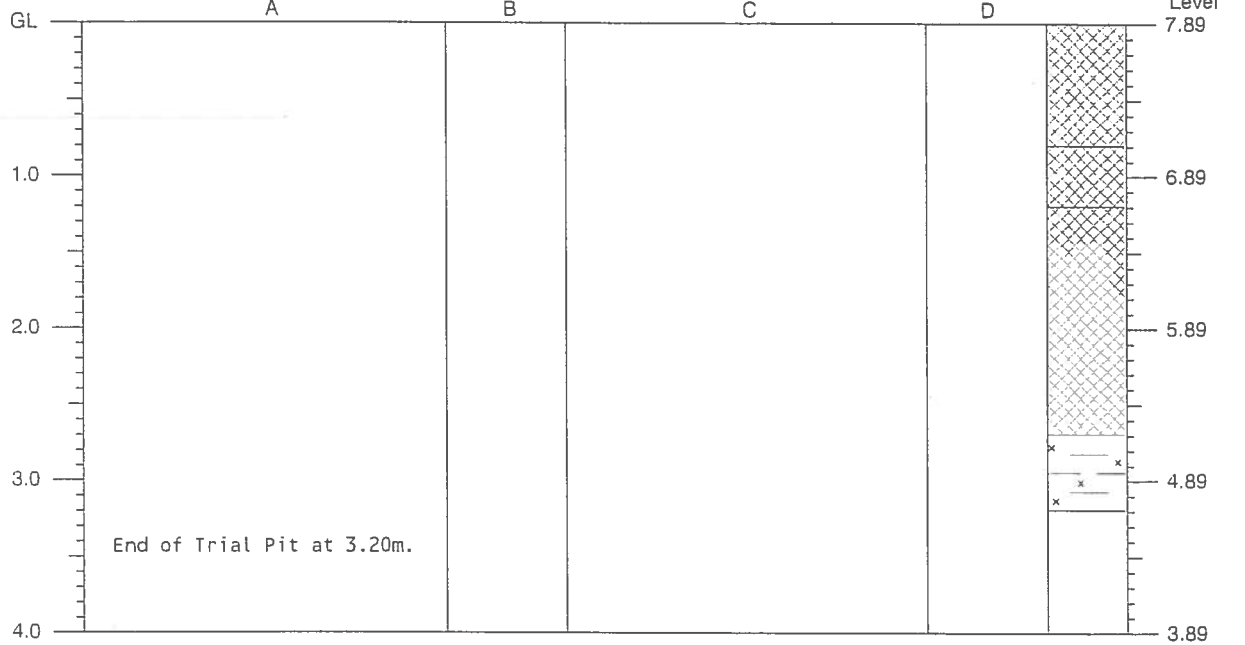
Sampling					Strata				
Depth	Type	Casing Depth	Date/ Water	SPT N (Cu)	Description	Depth (Thickness)	Level	Legend	
0.20-0.75	T		25/01 0000 DRY		Brown sandy clayey TOPSOIL	G.L.	10.78		
0.75-0.94	B				Firm orange brown sandy very silty CLAY with occasional to some rounded to sub-rounded gravel of sandstone	0.20	10.58		
0.94-1.12	T					(0.74)			
1.12		NIL	DRY		Compact orange brown very sandy slightly silty sub-angular to sub-rounded fine to coarse GRAVEL of sandstone	0.94 1.12	9.84 9.66		
					 End of Borehole.				
Equipment: Geotool R780 Tracked Rig					Groundwater		Ground Level		
Borehole Dia (mm) Casing Dia (mm)					No. Struck Behaviour Sealed		Coordinates		
80 to 1.00m 60 to 1.12m					No groundwater encountered		10.78 m OD 327349.63 mE 183940.41 mN		
							Drilled by DC Logged by DC Checked by AF		
Remarks Refusal at 1.12 (Could not penetrate gravels)									
See key sheet and appendices for explanations.									
Borehole Record					Project		Contract		
					M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners		150006		
Exploration Associates							Borehole		
							CH2(1 of 1)		

Sampling					Strata			
Depth	Type	Casing Depth	Date/ Water	SPT N (Cu)	Description	Depth (Thickness)	Level	Legend
0.10-0.35	T		25/01 2000 DRY		Brown sandy slightly silt clayey TOPSOIL	G.L. 0.10	10.86 10.76	
0.35-0.85	TO				MADE GROUND: Moderately compact brown sandy very silty slightly clayey angular to sub-angular fine to coarse gravel of sandstone, slag, ash and some refuse of plastic.	0.35	10.51	
0.85-1.17	B					(0.82)		
1.17-1.67	T				Firm brown locally orange brown sandy very silty CLAY with some sub-angular to sub-rounded fine to coarse gravel of sandstone and mudstone with some pockets of orange brown very silty fine to coarse sand.	1.17	9.69	
1.67		NIL	DRY		Compact red brown and green grey mottled dark grey very sandy slightly silty angular to sub-angular fine to coarse GRAVEL of sandstone and mudstone	(0.50) 1.67	9.19	
					 End of Borehole.			
Equipment: Geotool R780 Tracked Rig					Groundwater			
Borehole Dia (mm) 80 to 1.00m Casing Dia (mm) 60 to 1.67m					No. Struck Behaviour Sealed		Ground Level Coordinates 10.86 m OD 327381.55 183942.57 mE mN	
					No groundwater encountered		Drilled by DC Logged by DC Checked by AF	
Remarks Refusal at 1.67 (Unable to penetrate gravels) See key sheet and appendices for explanations.								
Borehole Record					Project		Contract	
					M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners		150006	
							Borehole	
							CH3(1 of 1)	

Sampling					Strata			
Depth	Type	Casing Depth	Date/ Water	SPT N (Cu)	Description	Depth (Thickness)	Level	Legend
0.20-0.60	T		25/01 0000 DRY		Brown sandy slightly silty slightly clayey TOPSOIL	G.L. 0.20	10.91 10.71	
0.60-1.00	TO				Firm brown slightly sandy very silty CLAY with occasional sub-rounded fine to coarse gravel of sandstone and mudstone with occasional rootlets	(0.40) 0.60	10.31	
1.00-1.50 1.00-2.00	B T				Firm red brown mottled grey slightly sandy very silty slightly organic CLAY with some rounded to sub-rounded fine to coarse gravel of mudstone and sandstone	(0.40) 1.00 (1.00)	9.91	
2.17		NIL	DRY		Firm red brown mottled green grey slightly sandy very silty CLAY	2.00 2.17	8.91 8.74	
					Compact brown very sandy sub-rounded fine to coarse GRAVEL of sandstone			
					 End of Borehole.			
Equipment: Geotool R780 Tracked Rig					Groundwater No. Struck Behaviour	Sealed	Ground Level Coordinates	10.91 m OD 327375.64 183925.14 mE mN
Borehole Dia (mm) Casing Dia (mm) 80 to 1.00m 60 to 2.17m					No groundwater encountered		Drilled by Logged by Checked by	DC DC AF
Remarks Refusal At 2.17m (could not penetrate gravels) See key sheet and appendices for explanations.								
Borehole Record					Project		Contract	
Exploration Associates					M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners		150006	
							Borehole CH4(1 of 1)	

Dimensions : 2.75 x 1.45

Orientation : N



Strata			Samples and Tests		
Depth (m)	No.	Description	Depth (m)	Type	Results
0.00-0.80	1	MADE GROUND: Loose red brown sandy very silty clayey angular to sub-angular fine to coarse gravel and cobbles of concrete, brick, sandstone and limestone with some refuse of plastic 5%.	0.00-0.80	BTO	
0.80-1.20	2	MADE GROUND: Loose grey sandy very silty clayey angular to sub-angular fine to coarse gravel and cobbles of brick, concrete, slag, ash and sandstone with some boulders (<550mm) of reinforced concrete (strong creosote odour) from wooden sleepers at 1.00m).	0.80-1.20	BTO	
1.20-2.70	3	MADE GROUND: Soft to firm red brown and grey sandy very silty clay with some angular to sub-angular fine to coarse gravel of slag, brick, ash, sandstone and mudstone with occasional cobbles of brick and concrete and many wooden sleepers encountered at 1.90m with some refuse of plastic 5%.	1.20-1.70	BTO	
2.70-3.20	4	Firm grey mottled brown very silty organic CLAY.	2.70-3.20	BTO	
			2.70	W	

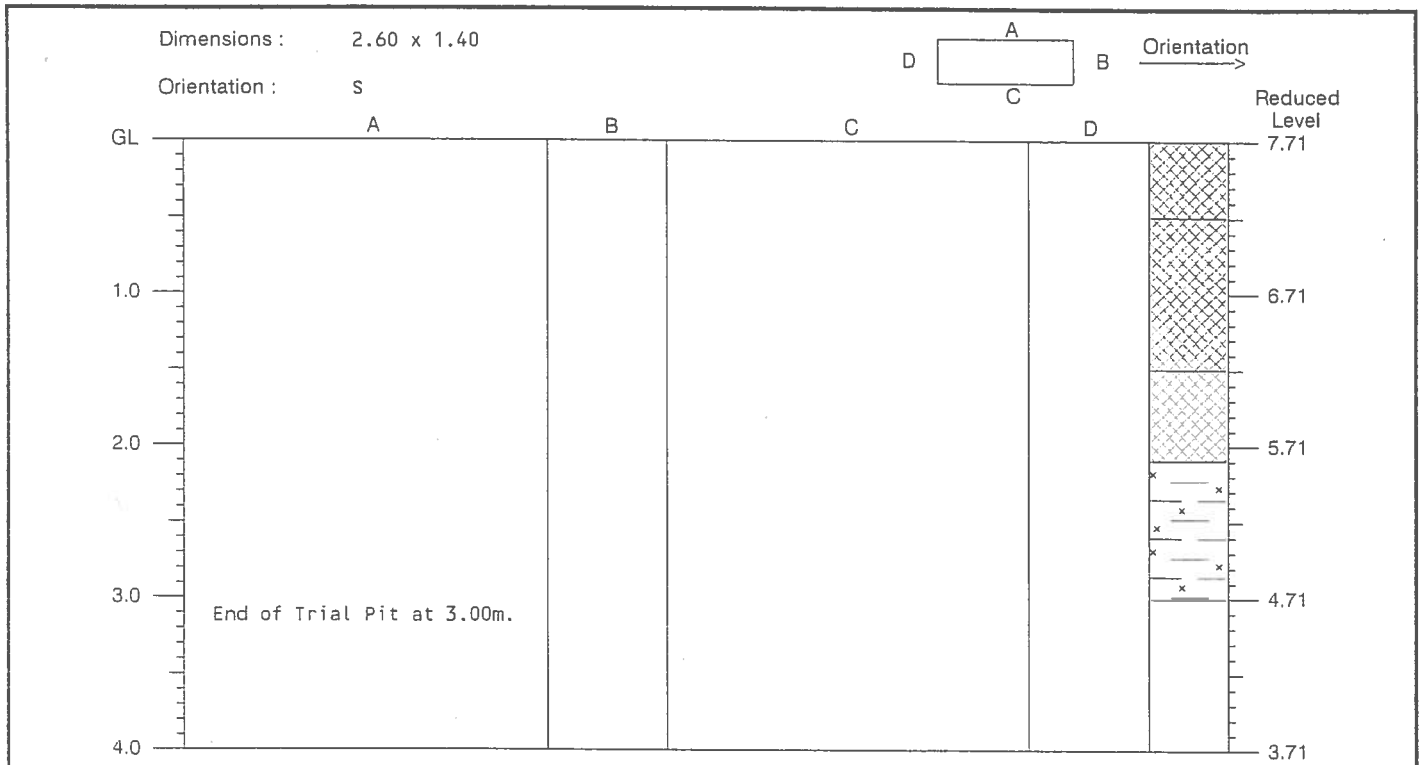
Date of Excavation 31/01/00 Equipment JCB 3CX Stability Stable	Groundwater No. Struck Behaviour 1 2.70 medium flow	Ground Level 7.89 m OD Coordinates 329050.29 mE 183417.98 mN Logged by DC Checked by AF
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Remarks

See key sheet and appendices for explanations.

Form 2/0

Trial Pit Record	Project M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners	Contract 150006
 Exploration Associates		Trial Pit CH5



Strata			Samples and Tests		
Depth (m)	No.	Description	Depth (m)	Type	Results
0.00-0.50	1	MADE GROUND: Firm reddish brown sandy very silty clay with much sub-angular to sub-rounded fine to coarse gravel and cobbles of brick, slag, ash, sandstone, siltstone and limestone.	0.00-0.50	BTO	
0.50-1.50	2	MADE GROUND: Moderately compact grey very sandy slightly silty slightly clayey sub-angular to sub-rounded fine to coarse gravel and cobbles of concrete, brick, sandstone, slate, slag, ash and siltstone with some sub-angular boulder (<500mm) of reinforced concrete with occasional pockets of red brown very silty sandy clayey marl and some refuse of pipe 1%, plastic 1% and wood 10%.	0.50-1.00	BTO	
1.50-2.10	3	MADE GROUND: Firm red brown sandy very silty clay with some angular to sub-angular fine to coarse gravel of brick, sandstone and concrete with much wood.	1.50-2.00	BTO	
2.10-3.00	4	Firm grey mottled blue grey and brown slightly sandy very silty organic CLAY.	2.10-2.70	BTO W	

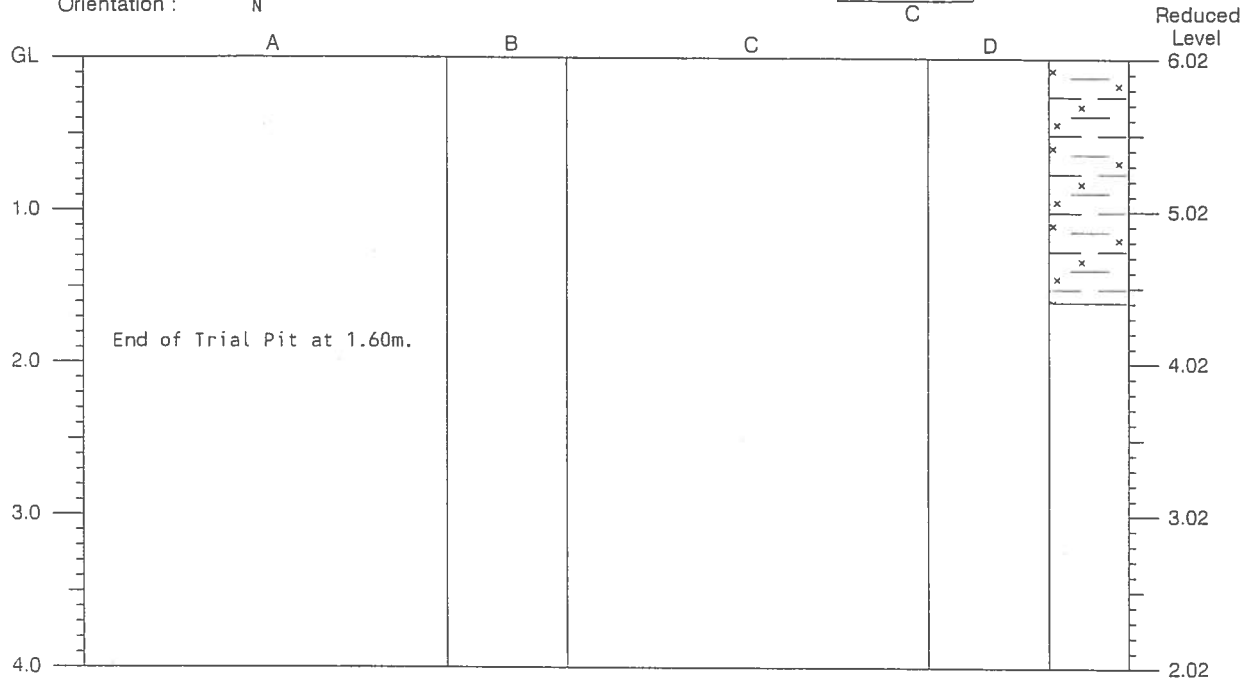
Date of Excavation 31/01/00 Equipment JCB 3CX Stability Stable	Groundwater No. Struck Behaviour 1 1.80 slight seepage	Ground Level 7.71 m OD Coordinates 329064.57 mE 183391.17 mN Logged by DC Checked by AF
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Remarks Slight water seepage at 1.80m (insufficient amount for water sampling). Water sample taken in rear adjacent to CH6 (known as "Site No. 2" rear sample) See key sheet and appendices for explanations.	Project M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners	Contract 150006 Trial Pit CH6
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Trial Pit Record 	Form 2/0
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Dimensions : 2.10 x 1.20

Orientation : N



Strata

Samples and Tests

Depth (m)	No.	Description	Depth (m)	Type	Results
0.00-1.60	1	Soft to firm grey mottled brown very slightly organic CLAY.	0.00-0.50 0.50-1.00 1.00-1.60	BTO B	

Date of Excavation 31/01/00
Equipment Jcb 3CX
Stability Stable

Groundwater
No. Struck Behaviour
1 1.50 slight seepage

Ground Level 6.02 m OD
Coordinates 329130.33 mE
183458.91 mN

Logged by DC
Checked by AF

Remarks Slight seepage at 1.50m (insufficient amount for water sampling).

See key sheet
and appendices
for explanations.

Form 2/0

Trial Pit Record

Project

Contract 150006

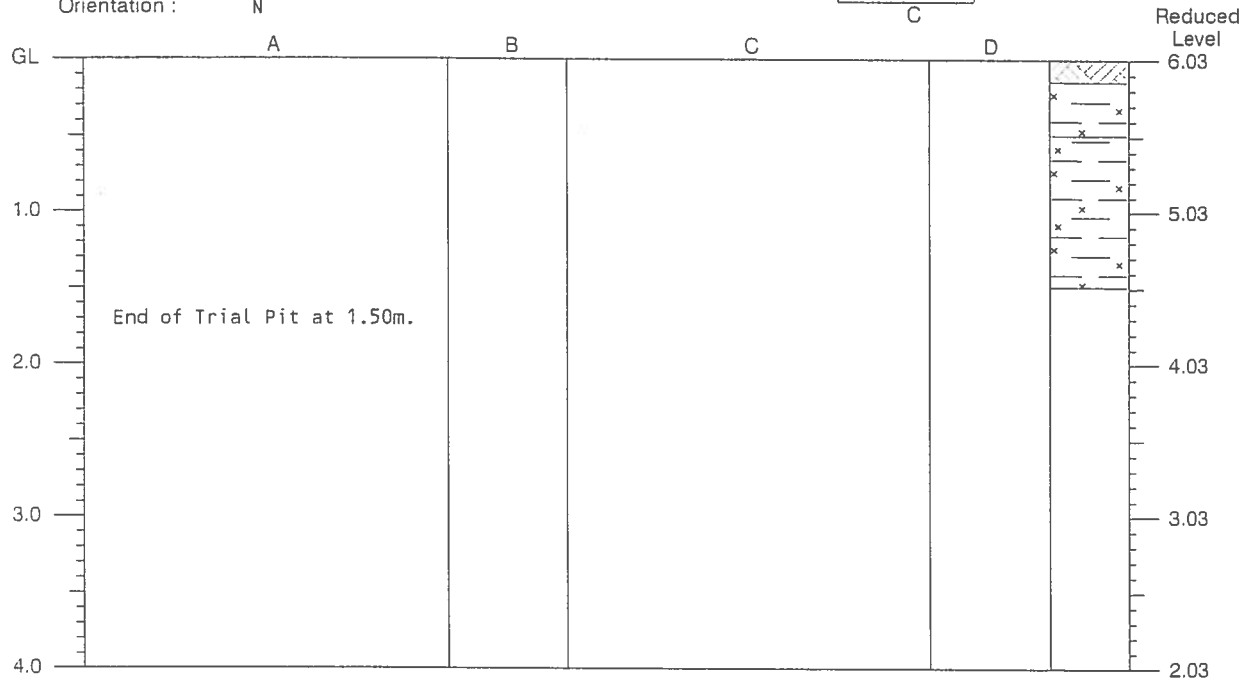
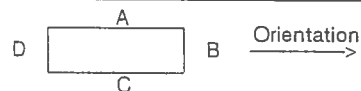
 **Exploration Associates**

M4 Relief Road - Stage 2. Preliminary
Chemical Investigation
Ove Arup & Partners

Trial Pit CH7

Dimensions : 2.00 x 1.10

Orientation : N



Strata			Samples and Tests		
Depth (m)	No.	Description	Depth (m)	Type	Results
0.00-0.15	1	Brown sandy clayey TOPSOIL.			
0.15-0.50	2	Grey mottled brown very silty organic CLAY.	0.15	B	
0.50-1.50	3	Blue grey mottled brown very silty CLAY.	1.00-1.50	BTO	
			1.50	W	
Date of Excavation 31/01/00 Equipment JCB 3CX Stability Stable			Groundwater No. Struck Behaviour 1 1.50 medium flow		
			Ground Level 6.03 m OD Coordinates 329178.75 mE 183519.18 mN Logged by DC Checked by AF		

Remarks

See key sheet and appendices for explanations.

Form 2/0

 Exploration Associates	Project M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners	Contract 150006
		Trial Pit CH8

Sampling					Strata				
Depth	Type	Casing Depth	Date/ Water	SPT N (Cu)	Description	Depth (Thickness)	Level	Legend	
0.00-0.50	TO		07/02 0000 DRY		MADE GROUND: Moderately compact brown grey and red brown sandy very silty clayey angular to sub-angular fine to coarse gravel of coal, brick, concrete, slag, sandstone and limestone with some cobbles of sub-rounded sandstone Firm brown mottled grey very silty CLAY no recovery 2.50m - 3.00m End of Borehole.	G.L.	9.80		
0.50-1.00	B			(1.50)					
1.00-1.50	TO								
1.50-2.00	TO			1.50		8.30			
2.00-2.50	B			(1.50)					
3.00		NIL	DRY		3.00	6.80			
Equipment: Hand Held Window Sampler Borehole Dia (mm) 80 to 1.00m Casing Dia (mm) 60 to 3.00m					Groundwater No. Struck Behaviour Sealed No groundwater encountered		Ground Level Coordinates 9.80 m OD 331020.20 mE 184756.13 mN Drilled by DC Logged by DC Checked by AF		
Remarks See key sheet and appendices for explanations.									
Borehole Record					Project M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners		Contract 150006 Borehole CH9(1 of 1)		
Exploration Associates									

Sampling					Strata			
Depth	Type	Casing Depth	Date/ Water	SPT N (Cu)	Description	Depth (Thickness)	Level	Legend
0.40-1.00	TO		07/02 2020 DRY		MADE GROUND: Weak grey concrete	G.L. (0.40)	10.29	
1.00-1.40	TO				MADE GROUND: Moderately compact dark grey and red brown sandy very silty clayey sub-angular to sub-rounded fine to coarse gravel of slag, ash, brick, concrete and limestone with occasional sub-rounded cobbles of sandstone.	0.40	9.89	
1.40-2.00	TO					(1.60)		
2.00		NIL	DRY		 End of Borehole.	2.00	8.29	
Equipment: Hand Held Window Sampler Borehole Dia (mm) Casing Dia (mm) 202 to 0.40m 80 to 1.40m 60 to 2.00m					Groundwater No. Struck Behaviour Sealed No groundwater encountered		Ground Level 10.29 m OD Coordinates 331076.30 mE 184823.93 mN Drilled by DC Logged by DC Checked by AF	
Remarks Concrete core 202mm ø to a depth of 0.40m (1hr) Refusal at 2.00m (possible cobbles?) See key sheet and appendices for explanations.								
Borehole Record Exploration Associates					Project M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners		Contract 150006 Borehole CH10(1 of 1)	

Sampling					Strata				
Depth	Type	Casing Depth	Date/ Water	SPT N (Cu)	Description	Depth (Thickness)	Level	Legend	
0.35-0.85	TO		07/02 0000 DRY		MADE GROUND: Black tarmac	G.L.	10.10		
0.85-1.15	B				MADE GROUND: Moderately compact grey slightly sandy angular to sub-angular fine to coarse gravel of siltstone and limestone(sub-base)	0.15 0.35	9.95 9.75		
1.15-1.65	B				MADE GROUND: Firm red brown slightly sandy very silty clay with much angular to sub-angular fine to coarse gravel of slag,ash, coal,sandstone and brick	(1.30)			
1.65-2.15	TO					1.65	8.45		
					MADE GROUND: Soft grey very silty clayey fine to coarse sand with occasional angular to sub-angular fine to coarse gravel of slag and ash (Strong fuel odour) no recovery 2.15m - 3.00m	(1.35)			
3.00		NIL	DRY			3.00	7.10		
					End of Borehole.				
Equipment: Hand Held Window Sampler					Groundwater No. Struck Behaviour	Sealed	Ground Level Coordinates	10.10 m OD 331177.34 184920.35	mE mN
Borehole Dia (mm) 202 to 0.15m 80 to 1.15m 60 to 3.00m		Casing Dia (mm)		No groundwater encountered		Drilled by Logged by Checked by		DC DC AF	
Remarks Concrete Core 202mm ϕ to a depth of 0.15m (1hr) See key sheet and appendices for explanations.									
Borehole Record					Project M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners		Contract 150006		
 Exploration Associates							Borehole CH11(1 of 1)		

Sampling					Strata			
Depth	Type	Casing Depth	Date/ Water	SPT N (Cu)	Description	Depth (Thickness)	Level	Legend
0.17-0.67	TO		07/02 2000 DRY		MADE GROUND: Black tarmac	G.L. 0.17	9.66 9.49	
0.67-1.37	TO				MADE GROUND: Moderately compact dark grey black very silty slightly clayey very sandy angular to sub-angular fine to coarse gravel of slag, ash, sandstone, brick and limestone with occasional cobbles of sandstone	(1.20)		
1.37-1.70	TO					1.37	8.29	
1.70-2.00 1.35	T W				MADE GROUND: Soft to firm grey very silty clay with occasional to some sub-rounded to sub-angular fine to coarse gravel of slag and sandstone	1.70	7.96	
					Soft to firm grey very silty organic CLAY no recovery 2.00m - 3.00m	(1.30)		
3.00		NIL	1.35		 End of Borehole.	3.00	6.66	
Equipment: Hand Held Window Sampler Borehole Dia (mm) Casing Dia (mm) 202 to 0.17m 80 to 1.17m 60 to 3.00m					Groundwater No. Struck Behaviour Sealed 1 1.35 Rose to 1.35m in 20 mins 1.35/1.35/1.35 in 5/10/15 mins		Ground Level 9.66 m OD Coordinates 331249.31 mE 184935.22 mN Drilled by DC Logged by DC Checked by AF	
Remarks Concrete core 202mm ϕ to a depth of 0.17 (1hr) Only Sufficient amount of water for 1no litre sample (mineral) See key sheet and appendices for explanations.								
Borehole Record Exploration Associates					Project M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners		Contract 150006 Borehole CH12(1 of 1)	

Sampling					Strata				
Depth	Type	Casing Depth	Date/ Water	SPT N (Cu)	Description	Depth (Thickness)	Level	Legend	
0.00-0.40	TO		28/01 2000 DRY		MADE GROUND: Moderately compact brown slightly silty sandy sub-angular to sub-rounded fine to coarse gravel of chert, mudstone, brick and sandstone with occasional refuse of plastic	G.L.	9.74		
0.40-0.77	TO					(0.40)			
0.77-1.07	TO					0.40 0.47	9.34 9.27		
1.07		NIL	DRY		MADE GROUND: Concrete.	0.77 1.07	8.97 8.67		
					MADE GROUND: Soft to firm red brown slightly sandy very silty clay with occasional to some sub-angular fine to coarse gravel of chert, sandstone, coal, slag, ash and brick with some cobbles of sub-rounded cobbles of sandstone				
					MADE GROUND: Soft to firm dark grey black very silty clay with occasional to some sub-rounded to sub-angular fine to coarse gravel of chert, coal, ash, brick and sandstone End of Borehole.				
Equipment: Hand Held Window Sampler					Groundwater No. Struck Behaviour No groundwater encountered	Sealed	Ground Level 9.74 m OD Coordinates 331289.61 mE 184976.66 mN		
Borehole Dia (mm) 80 to 1.00m Casing Dia (mm)							Drilled by DC Logged by DC Checked by AF		
Remarks Hole refusal at 1.07m (possible cobbles?) See key sheet and appendices for explanations.									
Borehole Record  Exploration Associates					Project M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners		Contract 150006		
							Borehole CH13(1 of 1)		

Sampling					Strata				
Depth	Type	Casing Depth	Date/ Water	SPT N (Cu)	Description	Depth (Thickness)	Level	Legend	
0.16-0.50	TO		04/02 2000 DRY		MADE GROUND: Strong grey intact re-inforced concrete.	G.L.	9.10		
0.50-1.30	B				MADE GROUND: Compact dark grey black very silty sandy angular to sub-angular fine to coarse gravel and cobbles of slag, ash, coal, sandstone and limestone	0.16	8.94		
		NIL	DRY			(1.14)			
1.30						1.30	7.80		
					End of Borehole.				
Equipment: Hand held window sampler					Groundwater				
Borehole Dia (mm) 202 to 0.16m 80 to 1.30m					No. Struck Behaviour	Sealed	Ground Level Coordinates	9.10 m OD 331384.08 185053.24	mE mN
Casing Dia (mm)					No groundwater encountered				
Remarks					Concrete core 202mm ϕ to a depth of 0.16 (1hr) Refusal at 1.30 (Possible cobbles?)				
See key sheet and appendices for explanations.					Form 1/0				
Borehole Record					Project		Contract		
 Exploration Associates					M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners		150006		
							Borehole CH14(1 of 1)		

Sampling					Strata					
Depth	Type	Casing Depth	Date/ Water	SPT N (Cu)	Description	Depth (Thickness)	Level	Legend		
0.30-0.80	TO		27/01 2000 DRY		MADE GROUND: Black tarmac	G.L.	9.21			
					MADE GROUND: Light grey reinforced concrete.	0.10	9.11			
						0.20	9.01			
						0.30	8.91			
0.80		NIL	DRY		MADE GROUND: Compact red brown very sandy angular to sub-angular fine to coarse gravel of limestone (sub-base)	(0.50)				
					MADE GROUND: Compact dark grey very silty sandy very clayey angular to sub-angular fine to coarse gravel of sandstone, slag and ash	0.80	8.41			
					 End of Borehole.					
Equipment: Hand Held Window Sampler					Groundwater No. Struck Behaviour	Sealed	Ground Level	9.21 m OD		
Borehole Dia (mm) Casing Dia (mm) 202 to 0.20m 80 to 0.80m					No groundwater encountered		Coordinates	331417.03 185090.44	mE mN	
Remarks					Drilled by DC Logged by DC Checked by AF					
Hand Dug pit to 0.20m (1/2hr) Refusal at 0.80m (possible cobbles?)										
See key sheet and appendices for explanations.										
Borehole Record					Project		Contract			
					M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners		150006			
 Exploration Associates							Borehole			
							CH15(1 of 1)			

Sampling					Strata				
Depth	Type	Casing Depth	Date/ Water	SPT N (Cu)	Description	Depth (Thickness)	Level	Legend	
0.17-0.67	TO		28/01 2000 DRY		MADE GROUND: Strong light grey intact re-inforced concrete.	G.L. 0.17	9.24 9.07		
0.67-0.87	B				MADE GROUND: Compact dark grey sandy very silty very clayey angular to sub-angular fine to coarse gravel of sandstone,slag,ash,brick and limestone	(0.70) 0.87			
0.87		NIL	DRY		 End of Borehole.		8.37		
Equipment: Hand Held Window Sampler					Groundwater No. Struck Behaviour		Sealed	Ground Level Coordinates	9.24 m OD 331415.85 185091.46 mE mN
Borehole Dia (mm) 202 to 0.17m 80 to 0.87m					Casing Dia (mm) No groundwater encountered		Drilled by DC Logged by DC Checked by AF		
Remarks					Concrete core 202mm ϕ to a depth of 0.17 (1½hrs) Refusal at 0.87 (possible cobbles?)				
See key sheet and appendices for explanations.					Form 1/0				
Borehole Record					Project		Contract		
					M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners		150006		
 Exploration Associates							Borehole		
							CH15A(1 of 1)		

Sampling					Strata				
Depth	Type	Casing Depth	Date/ Water	SPT N (Cu)	Description	Depth (Thickness)	Level	Legend	
0.20-0.70	TO		28/01 2000 DRY		MADE GROUND: Light grey strong intact re-inforced concrete	G.L. 0.20	8.69 8.49		
0.70-1.30	B				MADE GROUND: Moderately compact dark grey very silty clayey slightly sandy angular to sub-angular fine to coarse gravel of slag, slate, mudstone, ash, coal, sandstone and limestone				
1.30-1.90	B					(2.20)			
1.90-2.40	TO								
2.40-2.70	T	NIL	DRY		Soft grey very silty CLAY	2.40	6.29		
2.70					 End of Borehole.	2.70	5.99		
Equipment: Hand Held Window Sampler					Groundwater No. Struck Behaviour		Sealed	Ground Level Coordinates	8.69 m OD 331453.57 185126.62 mE mN
Borehole Dia (mm) Casing Dia (mm) 202 to 0.20m 80 to 1.20m 60 to 2.70m					No groundwater encountered		Drilled by Logged by Checked by	DC DC AF	
Remarks Concrete core 202mm ϕ to a depth of 0.20 (1½hrs) See key sheet and appendices for explanations.									
Borehole Record					Project M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners		Contract 150006		
 Exploration Associates							Borehole CH16(1 of 1)		

Sampling					Strata				
Depth	Type	Casing Depth	Date/ Water	SPT N (Cu)	Description	Depth (Thickness)	Level	Legend	
0.05-0.60	TO		27/01 2000 DRY		Brown slightly silty very sandy TOPSOIL	G.L. 0.05 (0.55)	8.79 8.74		
0.60-1.00	TO			MADE GROUND: Dark brown and red brown sandy very silty very clayey angular to sub-angular fine to coarse gravel of slag,brick,coal,ash, chert and sandstone	0.60 (0.70)	8.19			
1.30-1.90	TO			MADE GROUND: Red brown sandy very silty clay with much angular to sub-angular fine to coarse gravel of brick,coal,slag,ash and sandstone	1.30 (0.60)	7.49			
1.90-2.40 1.00-1.30	TO B			MADE GROUND: Loose dark grey very silty sandy angular to sub-angular fine to coarse gravel of slag,ash,coal and sandstone	1.90 (1.10)	6.89			
3.00		NIL	DRY	Soft to firm grey very silty CLAY	3.00	5.79			
									
					End of Borehole.				
Equipment: Hand Held Window Sampler					Groundwater		Ground Level		
Borehole Dia (mm) Casing Dia (mm)					No. Struck Behaviour Sealed		Coordinates		
80 to 1.00m 60 to 3.00m					No groundwater encountered		8.79 m OD 331500.85 185119.04 mE mN		
Remarks					Drilled by		DC		
See key sheet and appendices for explanations.					Logged by		DC		
					Checked by		AF		
Borehole Record					Project		Contract		
					M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners		150006		
Exploration Associates							Borehole		
							CH17(1 of 1)		

Sampling					Strata				
Depth	Type	Casing Depth	Date/Water	SPT N (Cu)	Description	Depth (Thickness)	Level	Legend	
0.20-0.50	TO		27/01 2000 DRY		MADE GROUND: Moderately compact red brown very sandy silty angular to sub-angular fine to coarse gravel of limestone and slag	G.L. 0.20	10.02 9.82		
0.50-1.00	TO					0.50	9.52		
1.00-1.60	B				MADE GROUND: Firm brown locally mottled orange brown sandy very silty clay with occasional to some sub-rounded fine to coarse gravel of sandstone	(1.10)			
1.60-2.20	TO				MADE GROUND: Moderately compact dark grey black very sandy very silty angular to sub-angular fine to coarse fine to coarse gravel of slag, ash, coal, chert and sandstone	1.60	8.42		
					MADE GROUND: Loose brown silty fine to coarse sand with occasional angular to sub-angular fine to coarse gravel of slag, ash, coal and sandstone no recovery 2.20m - 3.20m	(2.40)			
3.20-4.00	T								
4.00-4.50	T					4.00	6.02		
4.50-5.00	T				Soft to firm grey very silty CLAY with occasional sub-rounded fine to coarse gravel of sandstone	(0.50)			
						4.50	5.52		
5.00		NIL	DRY		Firm red brown very silty CLAY with occasional to some sub-angular fine to medium gravel of mudstone End of Borehole.	(0.50) 5.00	5.02		
Equipment: Hand Held Window Sampler					Groundwater				
Borehole Dia (mm) Casing Dia (mm)					No. Struck Behaviour Sealed		Ground Level Coordinates 10.02 m OD mE		
80 to 1.00m 60 to 3.00m 40 to 5.00m					No groundwater encountered		185145.01 mN		
Remarks					Drilled by DC		Checked by AF		
See key sheet and appendices for explanations.									
Borehole Record					Project		Contract		
Exploration Associates					M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners		150006		
							Borehole		
							CH18(1 of 1)		

Sampling					Strata				
Depth	Type	Casing Depth	Date/ Water	SPT N (Cu)	Description	Depth (Thickness)	Level	Legend	
0.20-0.75	TO		27/01 2000 DRY		MADE GROUND: Black tarmac	G.L.	10.32		
0.75-2.00	B				MADE GROUND: Loose dark grey black very silty sandy angular to sub-angular fine to coarse gravel of slag,ash,brick,sandstone and limestone	0.20	10.12		
						(1.80)			
2.00-2.50	TO					2.00	8.32		
2.50-3.00	TO				Soft to firm grey slightly sandy very clayey SILT	(1.00)			
3.00		NIL	DRY			3.00	7.32		
					End of Borehole.				
Equipment: Hand Held Window Sampler					Groundwater		Ground Level		
Borehole Dia (mm) 80 to 1.00m 60 to 3.00m					No. Struck Behaviour Sealed		Coordinates 10.32 m OD 331668.12 mE 185238.85 mN		
Casing Dia (mm)					No groundwater encountered		Drilled by DC Logged by DC Checked by AF		
Remarks Hand dug pit 0.00m - 0.20m (1/2hr)									
See key sheet and appendices for explanations.									
Borehole Record					Project		Contract		
Exploration Associates					M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners		150006		
							Borehole CH19(1 of 1)		

Sampling					Strata				
Depth	Type	Casing Depth	Date/ Water	SPT N (Cu)	Description	Depth (Thickness)	Level	Legend	
0.00-0.30	TO		27/01 2000 DRY		MADE GROUND: Loose yellow brown slightly silty fine to coarse sand	G.L.	10.32		
0.30-0.80	TO				MADE GROUND: Loose dark brown sandy very silty angular to sub-angular fine to coarse gravel of brick,slag,ash and sandstone	0.30 (0.50)	10.02		
1.00-1.50	TO				MADE GROUND: Soft brown very sandy clayey silt	0.80 (1.20)	9.52		
1.50-2.00	B								
2.00-2.50	TO				MADE GROUND: Soft grey very sandy clayey silt.	2.00 (0.50)	8.32		
2.50		NIL	DRY			2.50	7.82		
					End of Borehole.				
Equipment: Hand held Window Sampler					Groundwater				
Borehole Dia (mm) 80 to 1.00m 60 to 2.50m					No. Struck Behaviour	Sealed	Ground Level Coordinates	10.32 m OD 331745 185300	mE mN
Casing Dia (mm)					No groundwater encountered				
Remarks					Refusal at 2.50m (Gravels)				
See key sheet and appendices for explanations.					Form 1/0				
Borehole Record					Project		Contract		
 Exploration Associates					M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners		150006		
							Borehole CH19A(1 of 1)		

Sampling					Strata				
Depth	Type	Casing Depth	Date/ Water	SPT N (Cu)	Description	Depth (Thickness)	Level	Legend	
0.00-0.50	TO		26/01 2000 DRY		MADE GROUND: Compact dark brown very sandy silty angular to sub-angular fine to coarse gravel of slag, ash, sandstone and limestone Soft to firm grey very silty slightly organic CLAY with occasional fine to medium sub-rounded gravel of sandstone no recovery 2.20m - 3.00m	G.L.	8.85		
0.50-1.00	B					(1.00)			
1.00-1.50	TO					1.00	7.85		
1.50-2.20	TO					(2.00)			
3.00		NIL	DRY			3.00	5.85		
					End of Borehole.				
Equipment: Hand Held Window Sampler Borehole Dia (mm) 80 to 1.00m 60 to 3.00m					Groundwater No. Struck Behaviour Sealed No groundwater encountered		Ground Level 8.85 m OD Coordinates 331885.47 185314.89 mE mN Drilled by DC Logged by DC Checked by AF		
Remarks See key sheet and appendices for explanations.									
Borehole Record 					Project M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners		Contract 150006 Borehole CH20(1 of 1)		

Sampling					Strata			
Depth	Type	Casing Depth	Date/ Water	SPT N (Cu)	Description	Depth (Thickness)	Level	Legend
0.00-0.50	TO		26/01 2000 DRY		MADE GROUND: Compact dark brown and grey very sandy silty angular to sub-angular fine to coarse gravel of slag and ash	G.L.	8.47	
0.50-0.75	B					(0.75)		
0.75-1.50	TO					0.75	7.72	
1.50-2.00	B							
2.00-2.50	B					(2.25)		
2.50-3.00	TO							
3.00		NIL	DRY		 End of Borehole.	3.00	5.47	
Equipment: Hand Held Window Sampler Borehole Dia (mm) Casing Dia (mm) 80 to 1.00m 60 to 3.00m					Groundwater No. Struck Behaviour Sealed No groundwater encountered			
Ground Level Coordinates 8.47 m OD 331943.00 mE 185338.48 mN					Drilled by DC Logged by DC Checked by			
Remarks See key sheet and appendices for explanations.								
Borehole Record Exploration Associates					Project M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners		Contract 150006 Borehole CH21(1 of 1)	

Sampling					Strata			
Depth	Type	Casing Depth	Date/ Water	SPT N (Cu)	Description	Depth (Thickness)	Level	Legend
0.30-0.60	TO		26/01 2000 DRY		MADE GROUND: Black tarmac	G.L.	8.37	
0.60-0.75	T				MADE GROUND: Loose brown very sandy angular fine to coarse gravel of limestone (sub-base)	0.15	8.22	
0.75-0.95	T					0.30	8.07	
1.00-1.50	TO				MADE GROUND: Loose dark brown sandy very silty angular to sub-angular fine to coarse gravel of slag, ash, sandstone and limestone	0.60	7.77	
1.50-2.00	B					0.75	7.62	
2.00-2.50	B				MADE GROUND: Brown and grey very silty fine to medium sand	0.95	7.42	
2.50-3.00	TO				MADE GROUND: Loose dark grey sandy very silty angular to sub-angular fine to coarse gravel of slag and ash	(2.00)		
3.00		NIL	DRY		Soft to firm grey locally mottled brown very silty CLAY	2.95	5.42	
								
					End of Borehole.			
Equipment: Hand Held Window Sampler					Groundwater			
Borehole Dia (mm) 80 to 1.00m 60 to 3.00m					No. Struck Behaviour		Sealed	
Casing Dia (mm)					No groundwater encountered		Ground Level Coordinates 8.37 m OD 332018.57 mE 185383.50 mN	
Remarks Breakout (1/2hr)					Drilled by DC		Logged by DC	
					Checked by AF			
See key sheet and appendices for explanations.								
Borehole Record					Project		Form 1/0	
Exploration Associates					M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners		Contract 150006	
					Borehole		CH22(1 of 1)	

Sampling					Strata				
Depth	Type	Casing Depth	Date/ Water	SPT N (Cu)	Description	Depth (Thickness)	Level	Legend	
0.16-0.36 0.36-0.46 0.46-1.00	T T TO		26/01 2000 DRY		MADE GROUND: Light grey intact re-inforced concrete	G.L. 0.16 0.36 0.46	8.47 8.31 8.11 8.01		
1.00-1.50	B				MADE GROUND: Loose dark brown very sandy slightly silty angular to sub-angular fine to coarse gravel of slag,ash and limestone				
1.50-2.00	B				MADE GROUND: Loose light brown very sandy sub-angular to sub-rounded fine to coarse gravel of slag,ash,slate,brick and sandstone				
2.00-2.50	B				Soft to firm grey locally mottled brown very silty CLAY	(2.61)			
2.50-3.07	TO								
3.07		NIL	DRY			3.07	5.40		
					End of Borehole.				
Equipment: Hand Held Window Sampler					Groundwater				
Borehole Dia (mm) 202 to 0.16m 80 to 1.16m 60 to 3.07m					No. Struck Behaviour	Sealed	Ground Level Coordinates	8.47 m OD 332057.74 185390.67	mE mN
Casing Dia (mm)					No groundwater encountered				
Remarks					Concrete core 202mmø to a depth of 0.16m (1½hrs)				
See key sheet and appendices for explanations.									
Borehole Record					Project		Contract		
Exploration Associates					M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners		150006		
					Borehole		CH23(1 of 1)		

Sampling					Strata			
Depth	Type	Casing Depth	Date/ Water	SPT N (Cu)	Description	Depth (Thickness)	Level	Legend
0.00-0.50	TO		08/02 2000 DRY		MADE GROUND: Compact brown slightly clayey silty fine to coarse sand with occasional to some sub-angular fine to coarse gravel of brick, chert, coal, slag and ash with occasional sub-angular cobbles of sandstone Firm grey brown slightly organic slightly sandy very silty CLAY Soft to firm grey mottled brown very silty CLAY no recovery 2.50m - 3.00m End of Borehole.	G.L.	7.41	
0.50-0.75	B					(0.75)		
0.75-1.00	TO					0.75	6.66	
1.00-1.50	TO					(1.25)		
1.50-2.00	B					2.00	5.41	
2.00-2.50	TO				(1.00)			
3.00		NIL	DRY			3.00	4.41	
Equipment: Hand Held Window Sampler Borehole Dia (mm) Casing Dia (mm) 80 to 1.00m 60 to 3.00m					Groundwater No. Struck Behaviour Sealed No groundwater encountered		Ground Level Coordinates 7.41 m OD 332622.32 mE 185515.17 mN Drilled by DC Logged by DC Checked by AF	
Remarks See key sheet and appendices for explanations.								
Borehole Record Exploration Associates					Project M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners		Contract 150006 Borehole CH24(1 of 1)	

Sampling					Strata			
Depth	Type	Casing Depth	Date/ Water	SPT N (Cu)	Description	Depth (Thickness)	Level	Legend
0.00-0.50	TO		05/02 2000 DRY		MADE GROUND: Compact brown and grey sandy very silty very clayey sub-rounded to sub-angular fine to coarse gravel of brick,slag, sandstone and limestone Firm to stiff grey mottled brown very silty CLAY becoming soft to firm below 1.80m.	G.L.	7.61	
0.50-0.80	B					(0.80)		
0.80-1.10	TO					0.80	6.81	
1.10-1.80	B							
1.80-2.30	TO					(2.20)		
3.00		NIL	DRY			3.00	4.61	
					End of Borehole.			
Equipment: Hand Held Window Sampler Borehole Dia (mm) 80 to 1.00m Casing Dia (mm) 60 to 3.00m					Groundwater No. Struck Behaviour Sealed No groundwater encountered		Ground Level 7.61 m OD Coordinates 332754.50 185473.49 mE mN Drilled by DC Logged by DC Checked by AF	
Remarks See key sheet and appendices for explanations.								
Borehole Record					Project M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners		Contract 150006 Borehole CH25(1 of 1)	
Exploration Associates								

Sampling					Strata			
Depth	Type	Casing Depth	Date/ Water	SPT N (Cu)	Description	Depth (Thickness)	Level	Legend
0.30-0.65	TO		08/02 0000 DRY		MADE GROUND: Strong grey intact re-inforced concrete.	G.L.	7.19	
0.65-1.00	T				MADE GROUND: Moderately compact grey slightly sandy silty slightly clayey sub-angular to angular fine to coarse gravel of sandstone, brick and quartz with some cobbles of sandstone	0.30	6.89	
1.00		NIL	0.65			(0.70)		
					End of Borehole.	1.00	6.19	
Equipment: Hand Held Window Sampler Borehole Dia (mm) 80 to 1.00m Casing Dia (mm)					Groundwater No. Struck Behaviour Sealed 1 0.65 Rose to 0.65m in 20 mins .65/.65/.65 in 5/10/15 mins		Ground Level Coordinates 7.19 m OD 332824.91 185489.35 mE mN Drilled by DC Logged by DC Checked by AF	
Remarks Concrete core 202mmØ to a depth of 0.30 (1hr) Refusal at 1.00m (Possible cobbles??) Insufficient amount of water for same See key sheet and appendices for explanations.								
Borehole Record Exploration Associates					Project M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners		Contract 150006 Borehole CH26(1 of 1)	

Sampling					Strata			
Depth	Type	Casing Depth	Date/ Water	SPT N (Cu)	Description	Depth (Thickness)	Level	Legend
			08/02 0000 DRY		MADE GROUND: Strong grey re-inforced concrete.	G.L.	6.78	
0.55-0.90	TO				MADE GROUND: Moderately compact brown sandy angular to sub-angular fine to coarse gravel of limestone and brick	0.25	6.53	
1.10-1.50	TO				MADE GROUND: Moderately compact dark grey black very silty slightly clayey fine sand with occasional to some sub-angular fine to coarse gravel of sandstone and slag	0.90 1.10	5.88 5.68	
2.00-2.50	TO				MADE GROUND: Firm grey very silty clay with occasional sub-angular fine to medium gravel of coal and ash	(1.90)		
3.00		NIL	DRY		Firm grey very silty CLAY from 2.00 Becoming soft to firm	3.00	3.78	
								
					End of Borehole.			
Equipment: Hand Held Window Sampler					Groundwater			
Borehole Dia (mm) 80 to 1.00m 60 to 3.00m					No. Struck Behaviour		Sealed	
Casing Dia (mm)					No groundwater encountered		Ground Level 6.78 m OD Coordinates 332877.57 185472.83 mE mN	
Remarks					Concrete core 202mmØ to a depth of 0.25 (1.5 hrs)		Drilled by DC Logged by DC Checked by AF	
See key sheet and appendices for explanations.					Form 1/0			
Borehole Record					Project		Contract	
Exploration Associates					M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners		150006	
					Borehole		CH27(1 of 1)	

Sampling					Strata			
Depth	Type	Casing Depth	Date/Water	SPT N (Cu)	Description	Depth (Thickness)	Level	Legend
0.27-0.72	TO		08/02 2000 DRY		MADE GROUND: Strong grey intact re-inforced concrete	G.L.	7.89	
						0.27	7.62	
1.40-1.90	B				MADE GROUND: Compact brown slightly clayey very silty fine to medium sand with some sub-angular to sub-rounded fine to coarse gravel of quartz, sandstone, mudstone, coal, slag and ash with occasional cobbles of slag	(1.13)		
1.90-2.50	TO				MADE GROUND: Moderately compact slightly clayey slightly silty fine to medium sand with some to much angular to sub-angular fine to coarse gravel of sandstone	1.40	6.49	
						(0.50)		
2.50-3.00	TO				Soft to firm grey mottled brown slightly sandy very silty CLAY	1.90	5.99	
					From 2.30 becoming firm to stiff	(1.10)		
3.00		NIL	DRY			3.00	4.89	
					End of Borehole.			

Equipment: Hand Held Window Sampler		Groundwater		Sealed	Ground Level Coordinates	7.89 m OD 332926.36 185443.98	mE mN
Borehole Dia (mm)	Casing Dia (mm)	No. Struck	Behaviour				
202 to 0.27m		No groundwater encountered			Drilled by	DC	
80 to 1.27m					Logged by	DC	
60 to 3.00m					Checked by	AF	

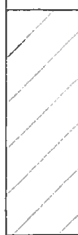
Remarks Concrete core 202mmØ to a depth of 0.27m (1hr)

See key sheet and appendices for explanations.

Form 1/0

Borehole Record		Project M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners	Contract	150006
 Exploration Associates			Borehole	CH28(1 of 1)

Sampling					Strata				
Depth	Type	Casing Depth	Date/ Water	SPT N (Cu)	Description	Depth (Thickness)	Level	Legend	
0.20-0.30	TO		08/02 0000 DRY		MADE GROUND: Strong grey intact re-inforced concrete.	G.L.	8.04		
0.40-0.50	TO					0.20	7.84		
		NIL	DRY		MADE GROUND: Moderately compact grey sandy sub-angular fine to coarse gravel of limestone (sub-base)	0.30	7.74		
0.50					MADE GROUND: Strong to weak grey concrete.	0.40	7.64		
					MADE GROUND: Compact dark grey slightly silty slightly clayey sandy angular to sub-angular fine to medium gravel of sandstone, slag and limestone	0.50	7.54		
					 End of Borehole.				
Equipment: Hand Held Window Sampler					Groundwater				
Borehole Dia (mm) Casing Dia (mm)					No. Struck Behaviour	Sealed	Ground Level	8.04 m OD	
202 to 0.20m					No groundwater encountered		Coordinates	332956.17	mE
80 to 0.30m								185427.60	mN
100 to 0.40m							Drilled by	DC	
							Logged by	DC	
							Checked by	AF	
Remarks Refusal at 0.50m (possible cobbles?)									
See key sheet and appendices for explanations.									
Borehole Record					Project		Contract		
Exploration Associates					M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners		150006		
							Borehole		
							CH29(1 of 1)		

Sampling					Strata				
Depth	Type	Casing Depth	Date/ Water	SPT N (Cu)	Description	Depth (Thickness)	Level	Legend	
0.00-0.50	TO		08/02 0000 DRY		MADE GROUND: Compact brown slightly clayey slightly silty fine sand with much angular to sub-angular fine to coarse gravel of chert, brick, sandstone and slag with occasional cobbles of sandstone from 0.75m becoming moderately compact End of Borehole.	G.L.	8.23	 	
0.50-1.00	B			(1.50)					
1.00-1.50	TO								
1.50		NIL	DRY	1.50		6.73			
Equipment: Hand Held Window Sampler					Groundwater No. Struck Behaviour Sealed No groundwater encountered		Ground Level 8.23 m OD Coordinates 333037.64 mE 185405.50 mN		
Borehole Dia (mm) 80 to 1.00m Casing Dia (mm) 60 to 1.50m							Drilled by DC Logged by DC Checked by AF		
Remarks Hole refusal at 1.50m (possible wooden sleepers?) See key sheet and appendices for explanations.									
Borehole Record					Project		Contract		
 Exploration Associates					M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners		150006		
							Borehole CH30(1 of 1)		

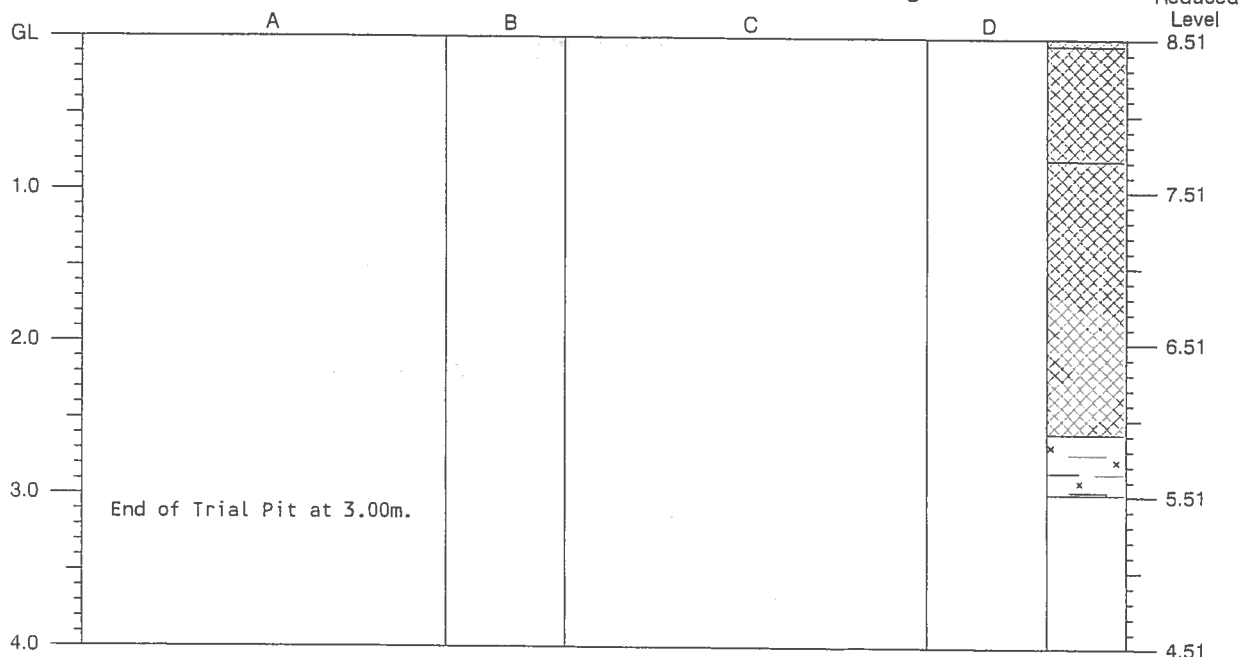
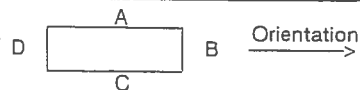
Sampling					Strata				
Depth	Type	Casing Depth	Date/ Water	SPT N (Cu)	Description	Depth (Thickness)	Level	Legend	
0.00-0.40	T		08/02 0000 DRY DRY		MADE GROUND: Very compact grey black slightly silty sandy slightly clayey sub-angular fine to coarse gravel of slag, ash, sandstone and coal with occasional cobbles of sandstone End of Borehole.	G.L. (0.40)	7.96		
0.40		NIL				0.40	7.56		
Equipment: Hand Held Window Sampler Borehole Dia (mm) 80 to 0.40m Casing Dia (mm)					Groundwater No. Struck Behaviour Sealed No groundwater encountered		Ground Level 7.96 m OD Coordinates 333206.82 mE 185368.33 mN Drilled by DC Logged by DC Checked by AF		
Remarks Hole refusal at 0.40m (could not penetrate gravels)									
See key sheet and appendices for explanations.									
Borehole Record  Exploration Associates					Project M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners			Contract 150006 Borehole CH31(1 of 1)	

Sampling					Strata				
Depth	Type	Casing Depth	Date/ Water	SPT N (Cu)	Description	Depth (Thickness)	Level	Legend	
0.00-0.35	T		08/02 0000 DRY		MADE GROUND: Moderately compact red brown slightly sandy slightly silty slightly clayey angular to sub-angular fine to coarse gravel of sandstone, limestone and brick with occasional to some cobbles of sandstone MADE GROUND: Firm to stiff brown slightly sandy very silty clay with occasional sub-angular fine to coarse gravel of coal and ash End of Borehole.	G.L.	8.13		
0.35-1.00 0.10	TO W					0.35	7.78		
1.00		NIL	0.10			(0.65) 1.00	7.13		
Equipment: Hand Held Window Sampler Borehole Dia (mm) 80 to 1.00m Casing Dia (mm)					Groundwater No. Struck Behaviour Sealed 1 0.10 Rose to 0.10m in 20 mins .10/.10/.10 in 5/10/15 mins		Ground Level Coordinates 8.13 m OD 333267.59 mE 185318.45 mN Drilled by DC Logged by DC Checked by AF		
Remarks Refusal at 1.00m (possible wooden sleepers?) See key sheet and appendices for explanations.									
Borehole Record Exploration Associates					Project M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners		Contract 150006 Borehole CH32(1 of 1)		

Sampling					Strata				
Depth	Type	Casing Depth	Date/ Water	SPT N (Cu)	Description	Depth (Thickness)	Level	Legend	
0.20-0.50	TO		08/02 0000 DRY		MADE GROUND: Stiff brown mottled grey very silty organic clay with a little sub-rounded fine to coarse gravel of brick and slag and with much root material	G.L.	8.28		
0.50-1.00	B					(1.20)			
1.50-2.00	TO				Firm to stiff blue grey mottled brown very silty organic CLAY with some root material	1.20	7.08		
2.00-2.50	B					(1.30)			
2.50-3.00	TO				Firm to stiff brown mottled grey very silty organic CLAY with much root material	2.50	5.78		
3.00		NIL	DRY			(0.50)			
						3.00	5.28		
					End of Borehole.				
Equipment: Hand Held Window Sampler					Groundwater				
Borehole Dia (mm) 80 to 1.00m 60 to 3.00m					No. Struck Behaviour		Sealed	Ground Level Coordinates 8.28 m OD 333473.63 185296.04 mE mN	
Casing Dia (mm)					No groundwater encountered		Drilled by DC Logged by DC Checked by AF		
Remarks CH33 changed from trial pit to window sample due to unsuitable access for JCB.									
See key sheet and appendices for explanations.									
Borehole Record					Project		Contract 150006		
Exploration Associates					M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners		Borehole CH33(1 of 1)		

Dimensions : 2.60 x 1.40

Orientation : W



Strata

Samples and Tests

Depth (m)	No.	Description	Depth (m)	Type	Results
0.00-0.05	1	MADE GROUND: Loose brown sandy angular to sub-angular fine to coarse gravel of limestone (hardcore).			
0.05-0.80	2	MADE GROUND: Firm brown mottled grey very silty clay with occasional sub-angular to sub-rounded fine to coarse gravel of siltstone with much root material.	0.15-0.50	BTO	
0.80-2.60	3	MADE GROUND: Firm to stiff grey mottled brown very silty clay with occasional to some root material.at 1.00m a little sub-angular fine to coarse gravel of slagat 2.00m some pockets of brown fibrous peatat 2.40m much root material	1.40-1.60	BTO	
2.60-3.00	4	Firm to stiff brown mottled blue grey very silty CLAY with some root material	2.60-2.70	BTO	

Date of Excavation 01/02/00
Equipment JCB 3CX
Stability Stable

Groundwater
No. Struck Behaviour
Not encountered during excavation

Ground Level 8.51 m OD
Coordinates 333500.60 mE
185283.99 mN

Logged by DC
Checked by AF

Remarks

See key sheet
and appendices
for explanations.

Form 2/0

Trial Pit Record

Project

Contract 150006

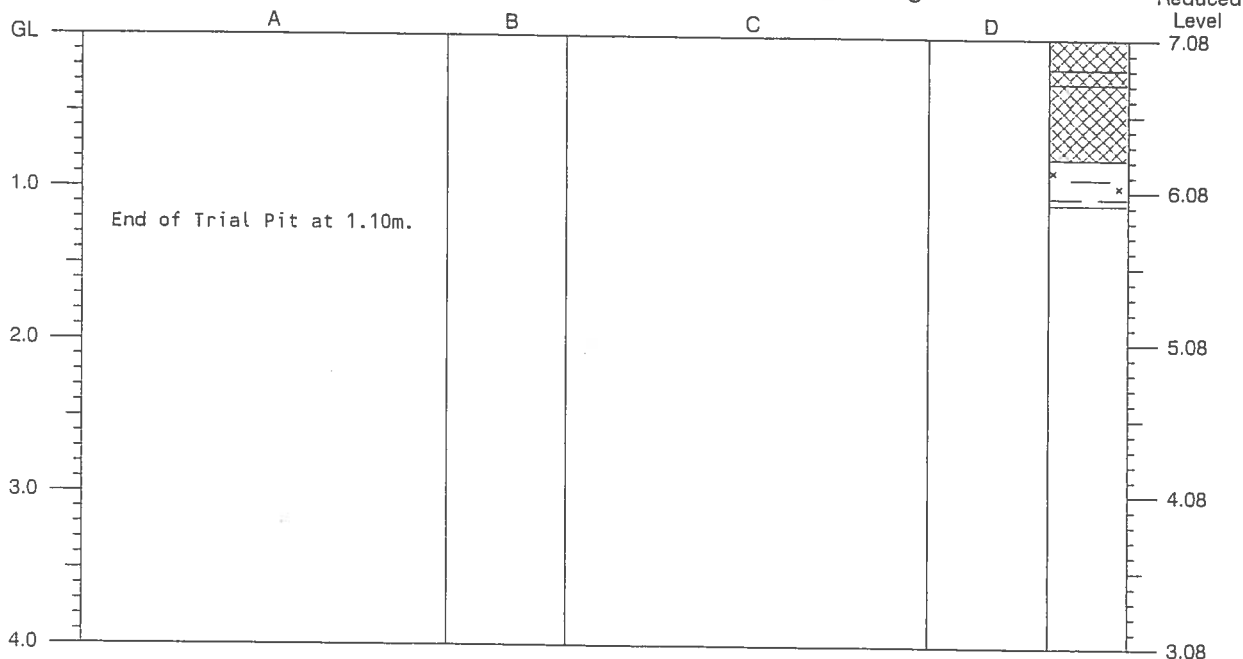
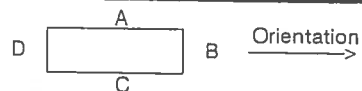
 Exploration Associates

M4 Relief Road - Stage 2. Preliminary
Chemical Investigation
Ove Arup & Partners

Trial Pit CH34

Dimensions : 2.60 x 1.30

Orientation : W



Strata

Samples and Tests

Depth (m)	No.	Description	Depth (m)	Type	Results
0.00-0.20	1	MADE GROUND: Moderately compact grey and brown slightly silty sand sub-angular to sub-rounded fine to coarse gravel and cobbles of slag, ash, concrete, brick, limestone (strong acidic odour).	0.00-0.20	BTO	
0.20-0.30	2	MADE GROUND: Moderately compact brown sandy angular to sub-angular fine to coarse gravel of limestone (sub-base).	0.20-0.30	BTO	
0.30-0.80	3	MADE GROUND: Moderately compact grey sandy very silty very clayey sub-angular fine to coarse gravel and cobbles of concrete, slag, ash, limestone and siltstone with some boulders (<440mm) of slag and concrete and some metal piping (5%).	0.50-0.70	BTO	
0.80-1.10	4	Stiff blue grey mottled brown very silty organic CLAY with much root material.	0.80-1.10	W BTO	

Date of Excavation 01/02/00
Equipment JCB 3CX
Stability Stable

Groundwater
No. Struck Behaviour
1 0.80 medium inflow

Ground Level 7.08 m OD
Coordinates 333540.67 mE
185256.56 mN

Logged by DC
Checked by AF

Remarks

See key sheet
and appendices
for explanations.

Form 2/0

Trial Pit Record

Project

Contract

150006



Exploration Associates

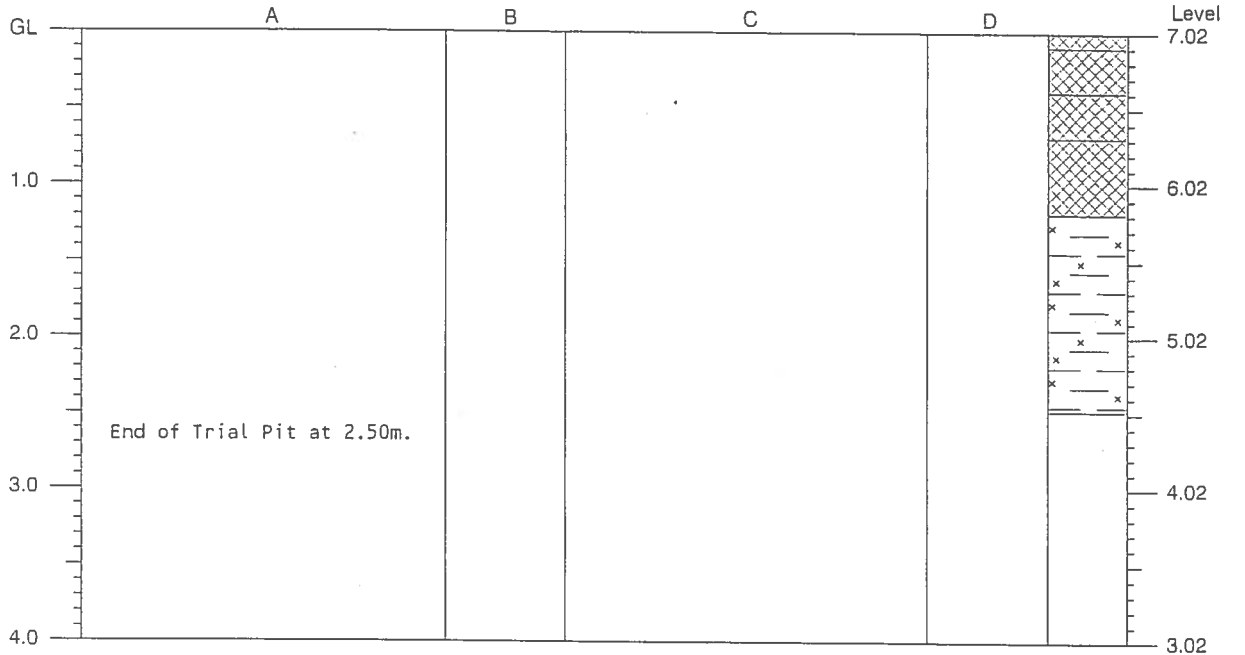
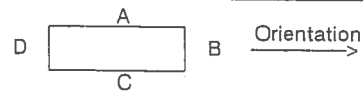
M4 Relief Road - Stage 2. Preliminary
Chemical Investigation
Ove Arup & Partners

Trial Pit

CH35

Dimensions : 1.00 x 1.80

Orientation : S



Strata			Samples and Tests		
Depth (m)	No.	Description	Depth (m)	Type	Results
0.00-0.10	1	MADE GROUND: Loose dark grey sandy slightly silty angular to sub-angular fine to coarse gravel of limestone and sandstone.	0.00-0.10	BTO	
0.10-0.40	2	MADE GROUND: Loose brown sandy angular to sub-angular fine to coarse gravel and cobbles of limestone (sub-base).	0.10-0.40	BTO	
0.40-0.70	3	Firm grey very silty organic CLAY	0.70-1.00	BTO	
0.70-1.20	4	Firm to stiff grey and brown very silty organic CLAY with much root material.	1.50-1.70	BTO	
1.20-2.50	5	Firm brown mottled grey very silty CLAY with some root material.from 1.90m becoming soft to firm			

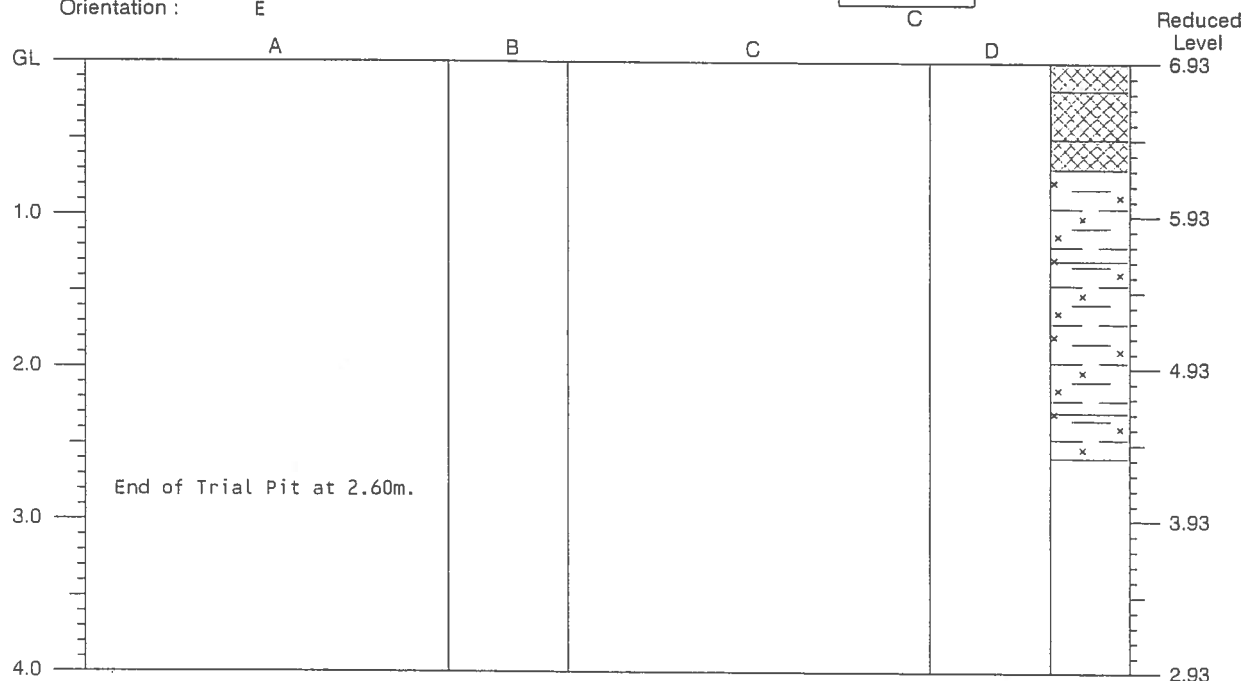
Date of Excavation 01/02/00 Equipment JCB 3CX Stability Stable	Groundwater No. Struck Behaviour Not encountered during excavation	Ground Level 7.02 m OD Coordinates 333550.07 mE 185225.06 mN Logged by DC Checked by AF
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Remarks See key sheet and appendices for explanations.		
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Trial Pit Record  Exploration Associates	Project M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners	Contract 150006 Trial Pit CH36
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Dimensions : 1.00 x 2.00

Orientation : E



Strata			Samples and Tests		
Depth (m)	No.	Description	Depth (m)	Type	Results
0.00-0.18	1	MADE GROUND: Loose dark grey sandy angular to sub-angular fine to coarse gravel of limestone (hardcore).	0.00-0.18	BTO	
0.18-0.50	2	MADE GROUND: Brown and grey slightly silty slightly clayey angular to sub-angular cobbles and boulders (<880mm) of brick and concrete with some refuse of plastic (1%).			
0.50-0.70	3	MADE GROUND: Firm blue grey very silty clay with many sub-angular fine to coarse gravel and cobbles of brick and much root material and refuse of plastic (1%).			
0.70-1.30	4	Firm to stiff blue grey mottled brown very silty organic CLAY.	1.20-1.30	BTO	
1.30-2.30	5	Firm stiff brown mottled blue grey very silty CLAY with much root material.	1.50-2.00	BTO	
2.30-2.60	6	Soft to firm blue grey very silty CLAY with occasional root material.	2.30-2.60	BTO	

Date of Excavation 01/02/00 Equipment JCB 3CX Stability Stable		Groundwater No. Struck Behaviour 1 2.00 slight seepage		Ground Level 6.93 m OD Coordinates 333588.39 mE 185235.73 mN Logged by DC Checked by Af	
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Remarks

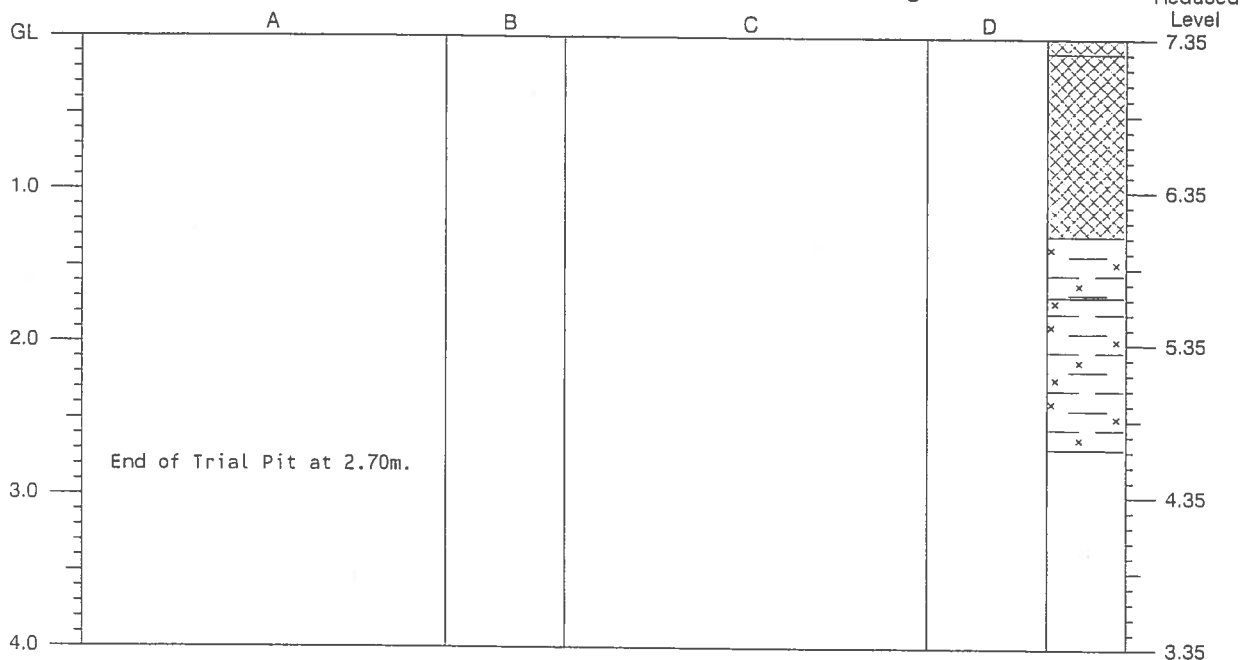
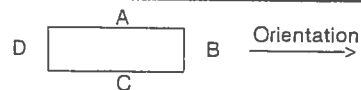
See key sheet and appendices for explanations.

Form 2/0

Trial Pit Record	Project M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners	Contract 150006
 Exploration Associates		Trial Pit CH37

Dimensions : 1.10 x 1.90

Orientation : S



Strata

Samples and Tests

Depth (m)	No.	Description	Depth (m)	Type	Results
0.00-0.10	1	MADE GROUND: Loose dark grey sandy angular to sub-angular fine to coarse gravel of limestone and siltstone (hardcore).			
0.10-1.30	2	MADE GROUND: Brown sandy sub-angular cobbles and boulders (>900mm) of concrete and brick with much refuse of wood (5%), metal (5%) and plastic (1%).	0.10-0.80	BTO	
1.30-1.70	3	Firm to stiff blue grey mottled brown very silty CLAY with much root material.	1.30-1.70	BTO	
1.70-2.70	4	Soft to firm brown mottled blue grey very silty CLAY with much root material.	2.50-2.70	BTO	

Date of Excavation 01/02/00
Equipment JCB 3CX
Stability Stable

Groundwater
No. Struck Behaviour
Not encountered during excavation

Ground Level 7.35 m OD
Coordinates 333599.73 mE
185219.05 mN

Logged by DC
Checked by AF

Remarks

See key sheet
and appendices
for explanations.

Form 2/0

Trial Pit Record

Project

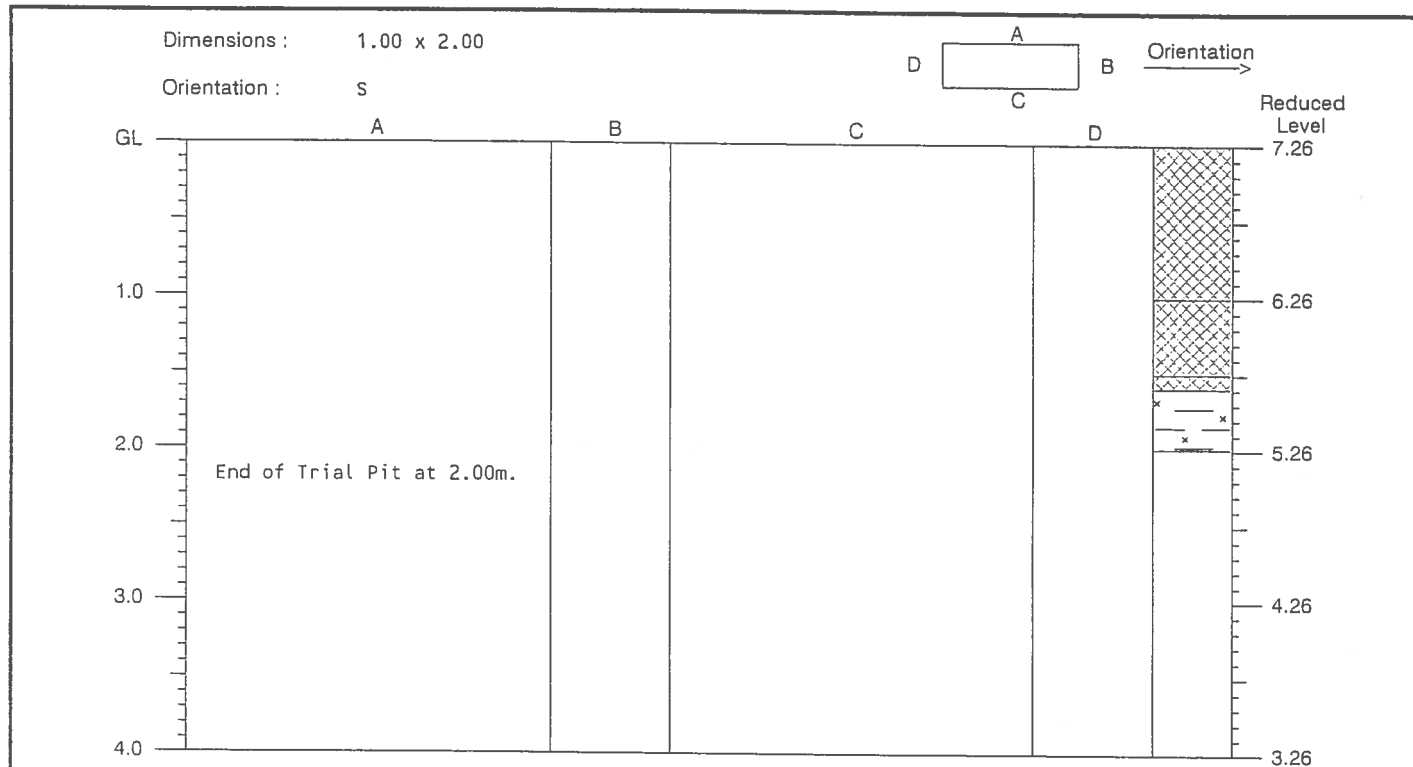
Contract 150006



Exploration Associates

M4 Relief Road - Stage 2. Preliminary
Chemical Investigation
Ove Arup & Partners

Trial Pit CH38



Strata			Samples and Tests		
Depth (m)	No.	Description	Depth (m)	Type	Results
0.00-1.00	1	MADE GROUND: Soft to firm brown very silty sandy clay with much sub-rounded to sub-angular fine to coarse gravel and cobbles of concrete, brick, sandstone and boulders (<560mm) of concrete and occasional refuse of plastic (1%).	0.00-0.60	BTO	
1.00-1.50	2	MADE GROUND: Moderately compact brown sandy very silty clayey sub-angular to sub-rounded fine to coarse gravel and cobbles of concrete, brick, slate, sandstone and slag with some boulders (<500mm) of concrete and refuse of wood (5%) and metal (5%).	1.00-1.30	BTO	
1.50-1.60	3	MADE GROUND: Concrete.	1.60-2.00	BTO	
1.60-2.00	4	Firm to stiff blue grey mottled brown very silty organic CLAY with much root material.			

Date of Excavation 02/02/00 Equipment JCB 3CX Stability Stable	Groundwater No. Struck Behaviour 1 1.40 slight seepage	Ground Level 7.26 m OD Coordinates 333612.01 mE 185209.37 mN Logged by DC Checked by AF
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Remarks

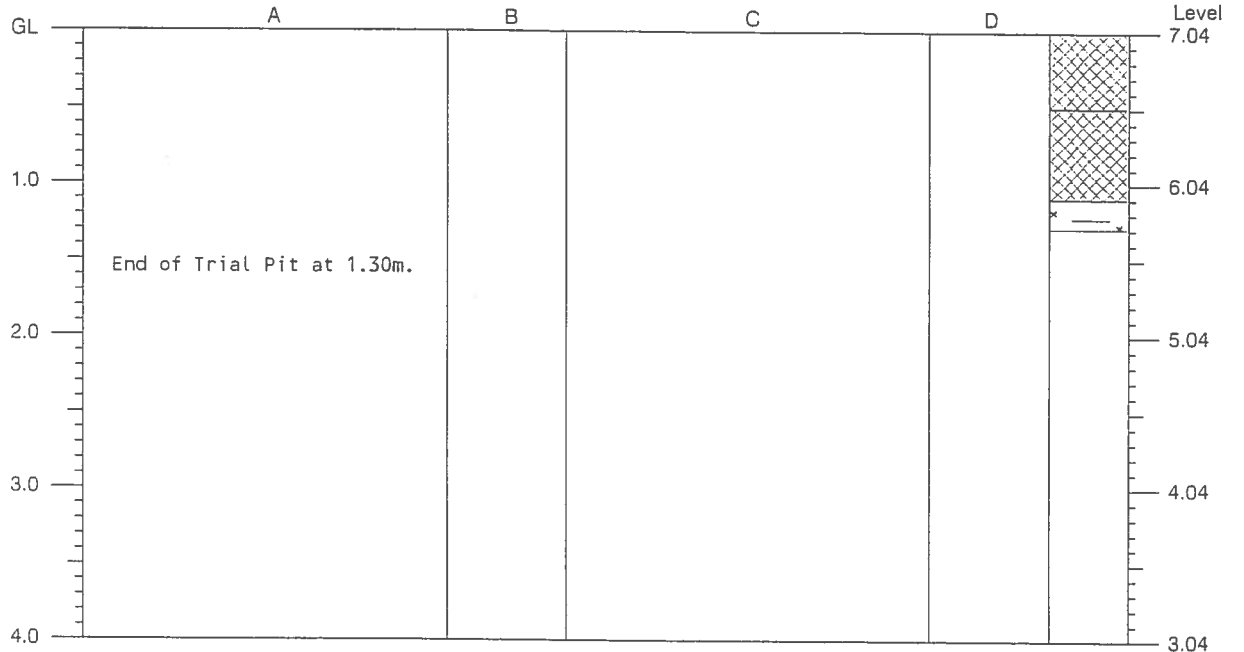
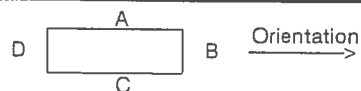
See key sheet and appendices for explanations.

Form 2/0

Trial Pit Record Exploration Associates	Project M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners	Contract 150006 Trial Pit CH39
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Dimensions : 1.20 x 2.50

Orientation : S



Strata			Samples and Tests		
Depth (m)	No.	Description	Depth (m)	Type	Results
0.00-0.50	1	MADE GROUND: Loose brown sandy angular to sub-angular fine to coarse gravel of limestone (hardcore).at 0.50m a layer of "terram".	0.00-0.50	BTO	
0.50-1.10	2	MADE GROUND: Moderately compact brown and grey sandy angular to sub-angular fine to coarse gravel and cobbles of concrete, brick, slag, sandstone and limestone with many boulders (<1400mm) of concrete and some refuse of plastic (5%) and metal (10%).	0.50-0.80	BTO	
1.10-1.30	3	Firm to stiff blue grey mottled brown very silty organic CLAY.	1.10-1.30	BTO	
			1.10	W	

Date of Excavation 02/02/00
Equipment JCB 3CX
Stability Stable

Groundwater
No. Struck Behaviour
1 1.10 fast strike

Ground Level 7.04 m OD
Coordinates 333641.12 mE
185239.97 mN

Logged by DC
Checked by AF

Remarks

See key sheet
and appendices
for explanations.

Form 2/0

Trial Pit Record

Project

M4 Relief Road - Stage 2. Preliminary
Chemical Investigation
Ove Arup & Partners

Contract

150006

Trial Pit

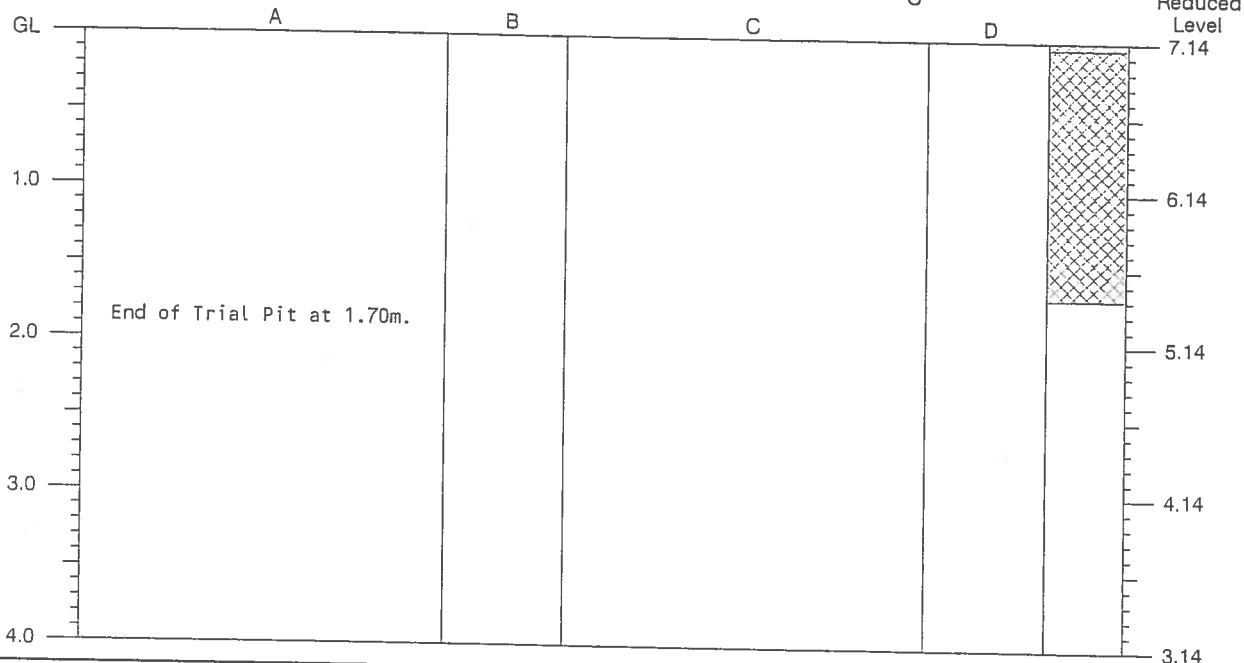
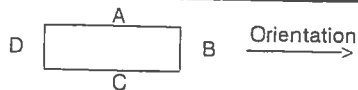
CH40



Exploration Associates

Dimensions :

Orientation :



Strata

Samples and Tests

Depth (m)	No.	Description	Depth (m)	Type	Results
0.00-0.05	1	Brown sandy TOPSOIL.			
0.05-1.70	2	MADE GROUND: Brown and grey sandy very silty clayey sub-angular fine to coarse gravel and cobbles of concrete, brick and sandstone and many boulders (<900mm) of concrete "bollards".from 1.40m becoming dark grey ("oil film" noticed on water in pit and on spoil excavated).	0.05-0.40 1.10 1.40	BTO W BTO	

Date of Excavation 01/02/00
Equipment JCB 3CX
Stability Stable

Groundwater
No. Struck Behaviour
1 1.10 medium inflow

Ground Level 7.14 m OD
Coordinates 333625.63 mE
185205.04 mN

Logged by DC
Checked by AF

Remarks

See key sheet
and appendices
for explanations.

Form 2/0

Trial Pit Record

Project

Contract

150006



Exploration Associates

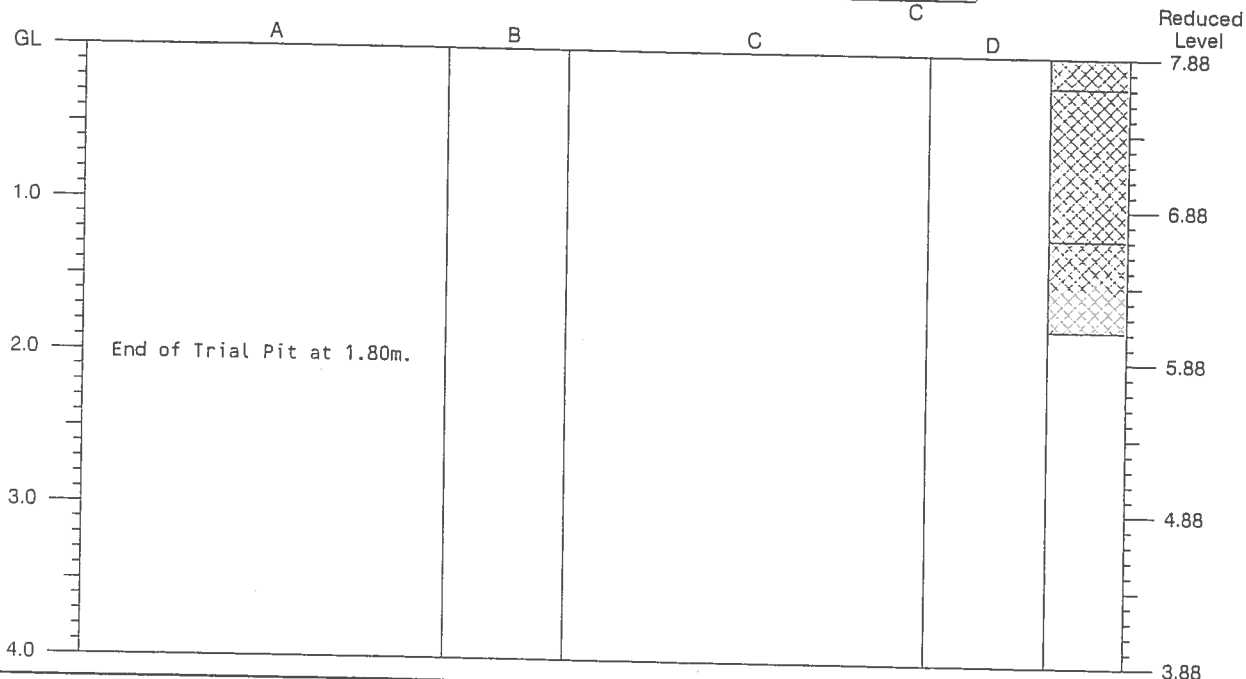
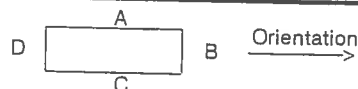
M4 Relief Road - Stage 2. Preliminary
Chemical Investigation
Ove Arup & Partners

Trial Pit

CH41

Dimensions :

Orientation :



Strata

Samples and Tests

Depth (m)	No.	Description	Depth (m)	Type	Results
0.00-0.20	1	Brown sandy very silty clayey TOPSOIL.			
0.20-1.20	2	MADE GROUND: Loose brown sandy very silty very clayey sub-angular to sub-rounded fine to coarse gravel and cobbles of brick, concrete, slag and ash.	0.20-0.70	BTO	
1.20-1.80	3	MADE GROUND: Moderately compact brown slightly sandy cobbles and boulders (<1000mm) of concrete, brick and slag with some refuse of plastic (1%) and metal (5%).	1.20-1.30	BTO	—

Date of Excavation 01/02/00
Equipment JCB 3CX
Stability Stable

Groundwater
No. Struck Behaviour
1 1.60 medium inflow

Ground Level 7.88 m OD
Coordinates 333681.23 mE
185236.03 mN

Logged by DC
Checked by Af

Remarks

See key sheet
and appendices
for explanations.

Form 2/0

Trial Pit Record

Project

Contract

150006

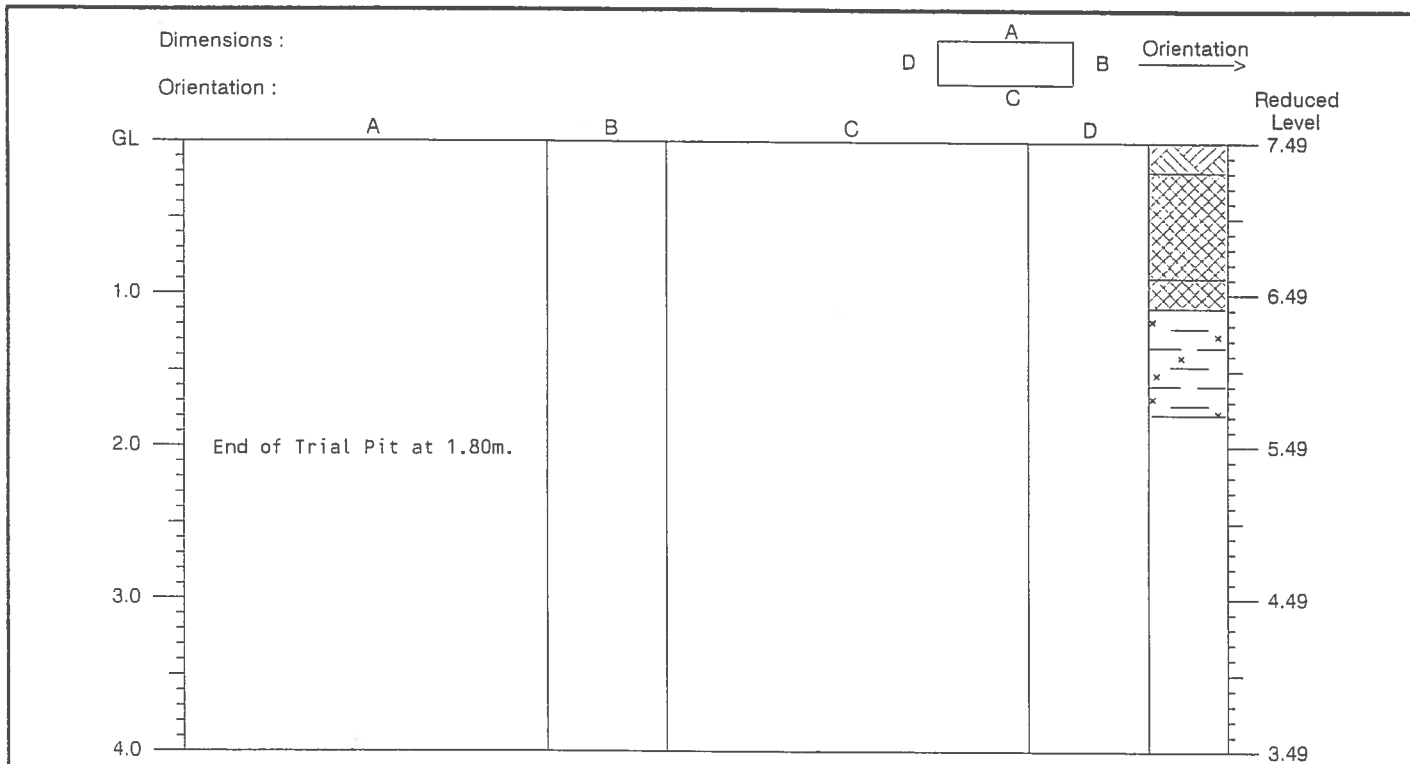


Exploration Associates

M4 Relief Road - Stage 2. Preliminary
Chemical Investigation
Ove Arup & Partners

Trial Pit

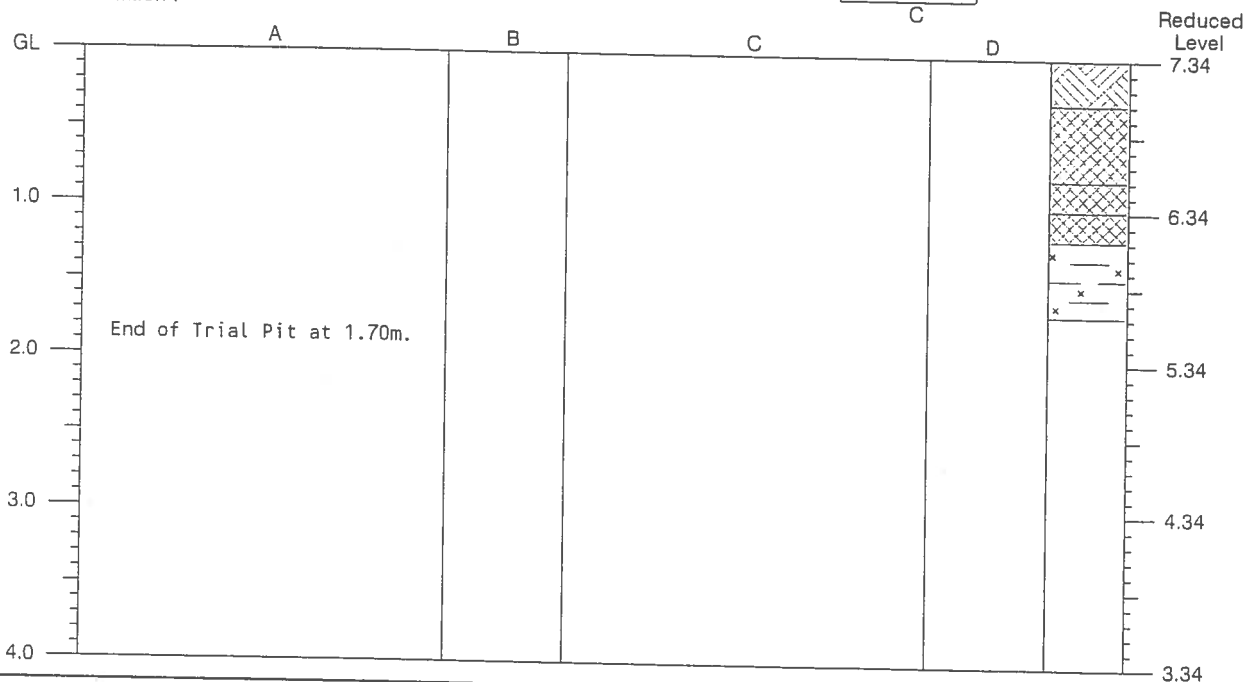
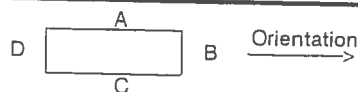
CH42



Strata			Samples and Tests		
Depth (m)	No.	Description	Depth (m)	Type	Results
0.00-0.20	1	Brown sandy very silty clayey TOPSOIL.			
0.20-0.90	2	MADE GROUND: Firm stiff blue grey and brown very silty clay with much sub-rounded to sub-angular fine to coarse gravel and cobbles of concrete, brick, sandstone and slag and with some wooden sleepers.	0.20-0.70	BTO	
0.90-1.10	3	MADE GROUND: Compact brown and grey slightly sandy cobbles and boulders (<1200) of concrete and brick.			
1.10-1.80	4	Firm to stiff blue grey mottled brown very silty slightly organic CLAY with much root material.	1.30-1.80	BTO	
Date of Excavation 02/02/00		Groundwater		Ground Level 7.49 m OD	
Equipment JCB 3CX		No. Struck Behaviour		Coordinates 333709.24 mE	
Stability Stable		Not encountered during excavation		185218.15 mN	
				Logged by DC	
				Checked by AF	
Remarks See key sheet and appendices for explanations.					
Trial Pit Record		Project		Contract 150006	
		M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners		Trial Pit CH43	
 Exploration Associates					

Dimensions :

Orientation :

**Strata****Samples and Tests**

Depth (m)	No.	Description	Depth (m)	Type	Results
0.00-0.30	1	Brown very silty clayey TOPSOIL.			
0.30-0.80	2	MADE GROUND: Loose grey and brown sandy very silty clayey angular to sub-angular fine to coarse gravel and cobbles of brick, concrete, slag, ash and sandstone with some <400mm of concrete.	0.30-0.80	BTO	
0.80-1.00	3	MADE GROUND: Compact brown and grey cobbles and boulders (<450mm) of concrete and brick with some refuse of metal (5%).			
1.00-1.20	4	MADE GROUND: Moderately compact brown and grey sandy sub-angular to sub-rounded fine to coarse gravel and cobbles of concrete, brick and sandstone.	1.00-1.20	BTO	
1.20-1.70	5	Soft to firm blue grey mottled brown very silty organic CLAY with some root material.	1.20-1.50	BTO	

Date of Excavation 02/02/00

Equipment JCB 3CX

Stability Stable

Groundwater

No. Struck Behaviour

1 1.70 slight seepage

Ground Level 7.34 m OD

Coordinates 333740.12 mE

185222.28 mN

Logged by DC

Checked by AF

RemarksSee key sheet
and appendices
for explanations.

Form 2/0

Trial Pit Record**Project**M4 Relief Road - Stage 2. Preliminary
Chemical Investigation
Ove Arup & Partners**Contract**

150006

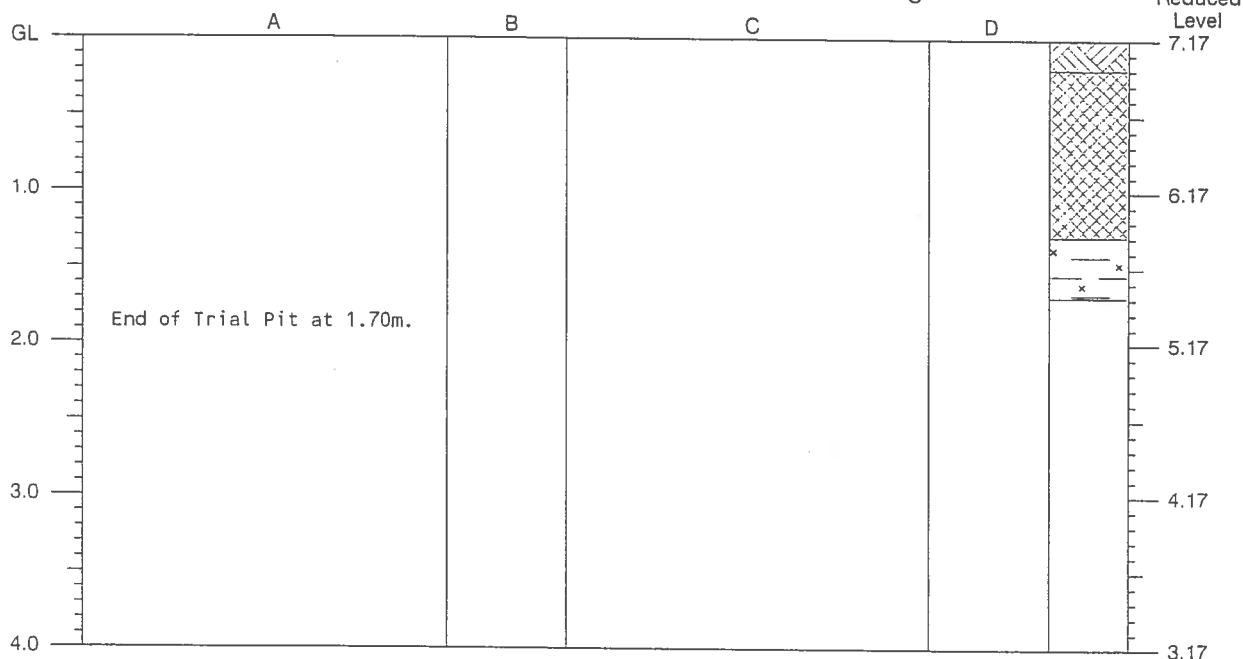
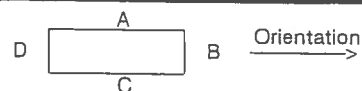
Trial Pit

CH44

**Exploration Associates**

Dimensions :

Orientation :

**Strata****Samples and Tests**

Depth (m)	No.	Description	Depth (m)	Type	Results
0.00-0.20	1	Brown very silty slightly clayey TOPSOIL.			
0.20-1.30	2	MADE GROUND: Moderately compact brown and grey sandy angular to sub-angular fine to coarse gravel and cobbles of concrete, brick and sandstone with some boulders (<560mm) of concrete and much refuse of metal (5%).	0.20-0.50	BTO	
1.30-1.70	3	Firm to stiff brown mottled orange brown and blue grey very silty organic CLAY.	1.30-1.70	BTO	
			1.30	W	

Date of Excavation 02/02/00
 Equipment JCB 3CX
 Stability Stable

Groundwater
 No. Struck Behaviour
 1 1.30 medium inflow

Ground Level 7.17 m OD
 Coordinates 333784.67 mE
 185215.33 mN

Logged by DC
 Checked by AF

Remarks

See key sheet
 and appendices
 for explanations.

Form 2/0

Trial Pit Record**Project**

M4 Relief Road - Stage 2. Preliminary
 Chemical Investigation
 Ove Arup & Partners

Contract

150006

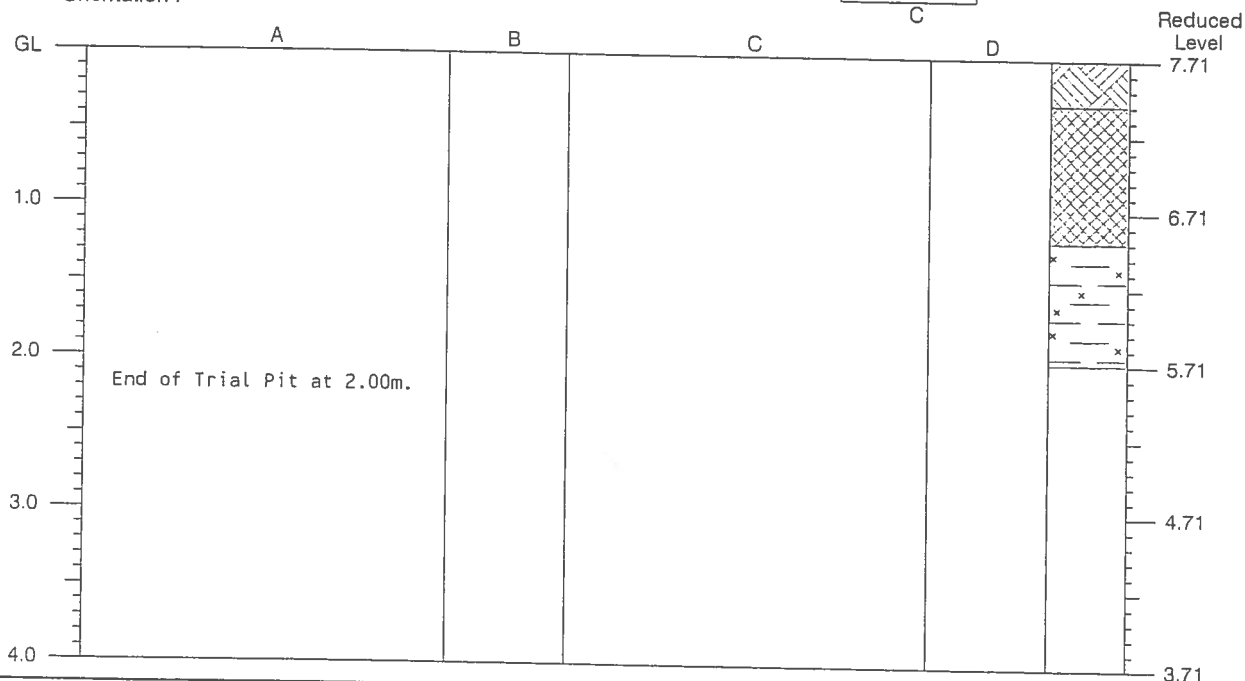
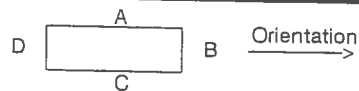
Trial Pit

CH45

**Exploration Associates**

Dimensions :

Orientation :

**Strata****Samples and Tests**

Depth (m)	No.	Description	Depth (m)	Type	Results
0.00-0.30	1	Brown very silty clayey TOPSOIL with a little angular to sub-angular fine to coarse gravel of sandstone.			
0.30-1.20	2	MADE GROUND: Loose brown sandy very silty very clayey sub-rounded to sub-angular fine to coarse gravel and cobbles of concrete, brick, slag, ash, slate, sandstone and limestone with occasional boulders (<650mm) of concrete and some refuse of plastic (1%) and metal (5%).	0.50-0.90	BTO	
1.20-2.00	3	Firm to stiff blue grey very silty organic CLAY with much root material.	1.20-1.70	BTO	

Date of Excavation 02/02/00

Equipment JCB 3CX

Stability Stable

Groundwater

No. Struck Behaviour

Not encountered during excavation

Ground Level 7.71 m OD

Coordinates 333757.41 mE
185226.90 mNLogged by DC
Checked by AF**Remarks**See key sheet
and appendices
for explanations.

Form 2/0

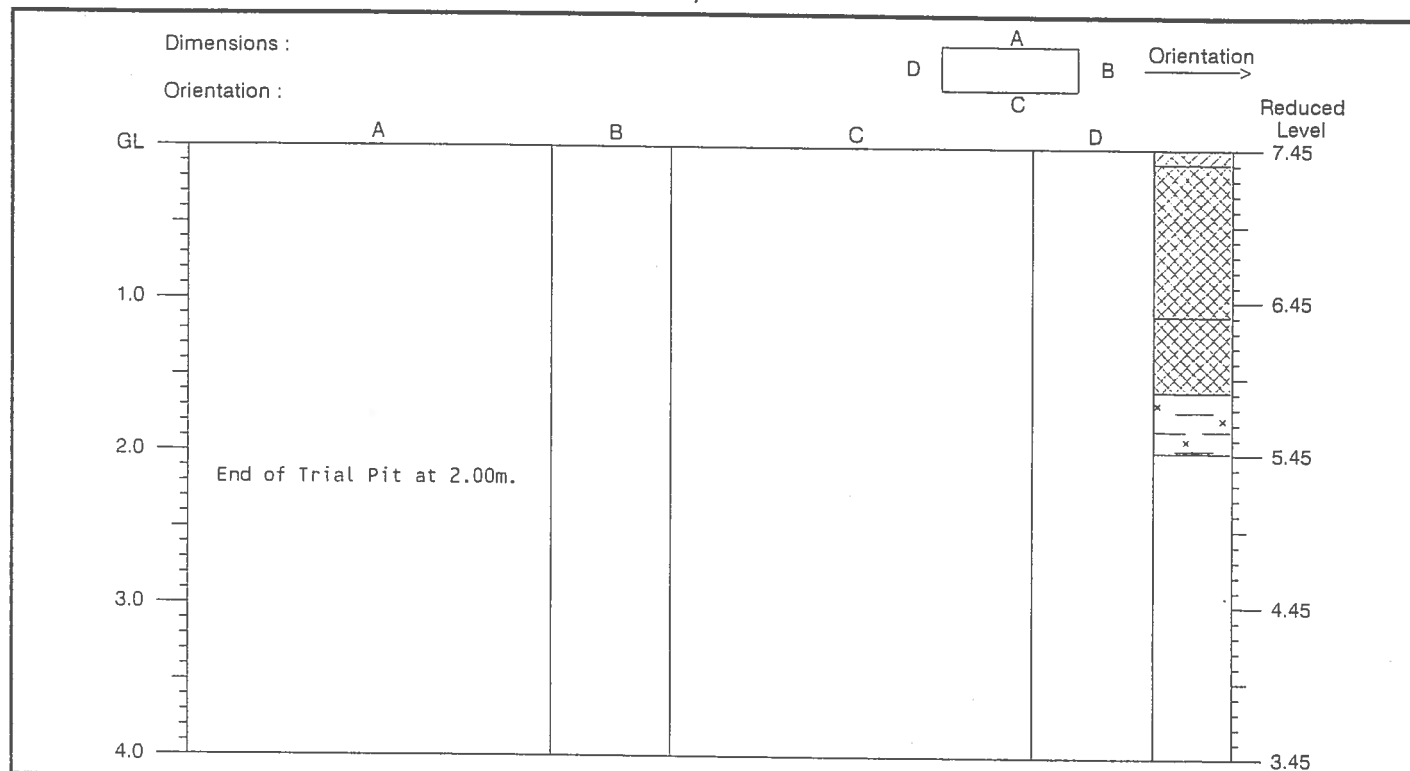
Trial Pit Record**Project**M4 Relief Road - Stage 2. Preliminary
Chemical Investigation
Ove Arup & Partners**Contract**

150006

Trial Pit

CH46

**Exploration Associates**



Strata			Samples and Tests		
Depth (m)	No.	Description	Depth (m)	Type	Results
0.00-0.10	1	Brown sandy very silty slightly clayey TOPSOIL.			
0.10-1.10	2	MADE GROUND: Moderately compact brown sandy very silty very clayey angular to sub-angular fine to coarse gravel of concrete, brick, slag, ash and sandstone with some boulders (<1080mm) of concrete and some refuse of plastic (1%), metal (5%) and wood (1%).	0.20-0.60	BTO	
1.10-1.60	3	MADE GROUND: Firm to stiff blue grey very silty slightly organic clay with a little angular to sub-angular fine to coarse gravel of concrete and brick.	1.10-1.60	BTO	
1.60-2.00	4	Firm to stiff brown very silty organic CLAY.	1.60-2.00	BTO	

Date of Excavation 02/02/00 Equipment JCB 3CX Stability Stable	Groundwater No. Struck Behaviour Not encountered during excavation	Ground Level 7.45 m OD Coordinates 333783.57 mE 185246.27 mN Logged by DC Checked by Af
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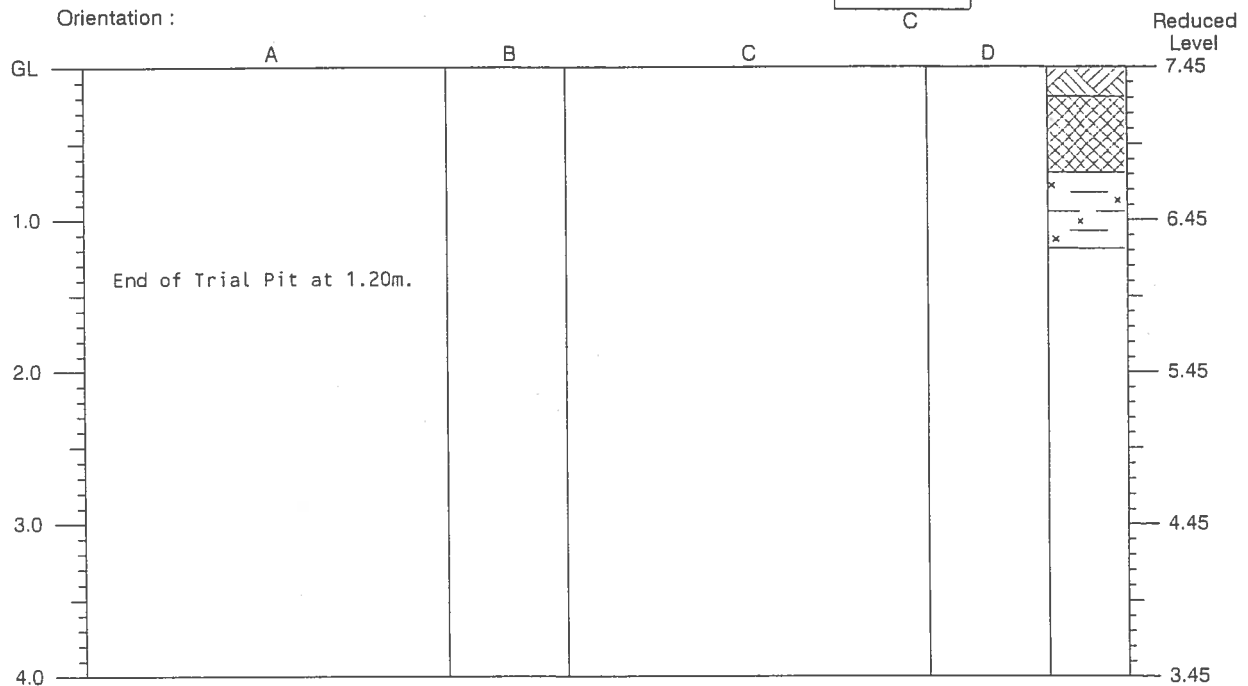
Remarks

See key sheet and appendices for explanations.

Trial Pit Record 	Project M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners	Contract 150006
		Trial Pit CH47

Dimensions :

Orientation :

**Strata****Samples and Tests**

Depth (m)	No.	Description	Depth (m)	Type	Results
0.00-0.20	1	Brown very silty clayey TOPSOIL.			
0.20-0.70	2	MADE GROUND: Soft to firm brown very silty clay with occasional refuse of plastic (1%).	0.20-0.70	BTO	
0.70-1.20	3	Firm to stiff blue grey very silty slightly organic CLAY.	0.70-1.20	BTO	

Date of Excavation 02/02/00
 Equipment JCB 3CX
 Stability Stable

Groundwater
 No. Struck Behaviour
 1 1.20 slight seepage

Ground Level 7.45 m OD
 Coordinates 333807.57 mE
 185246.27 mN

Logged by DC
 Checked by AF

Remarks

See key sheet
 and appendices
 for explanations.

Form 2/0

Trial Pit Record**Project**

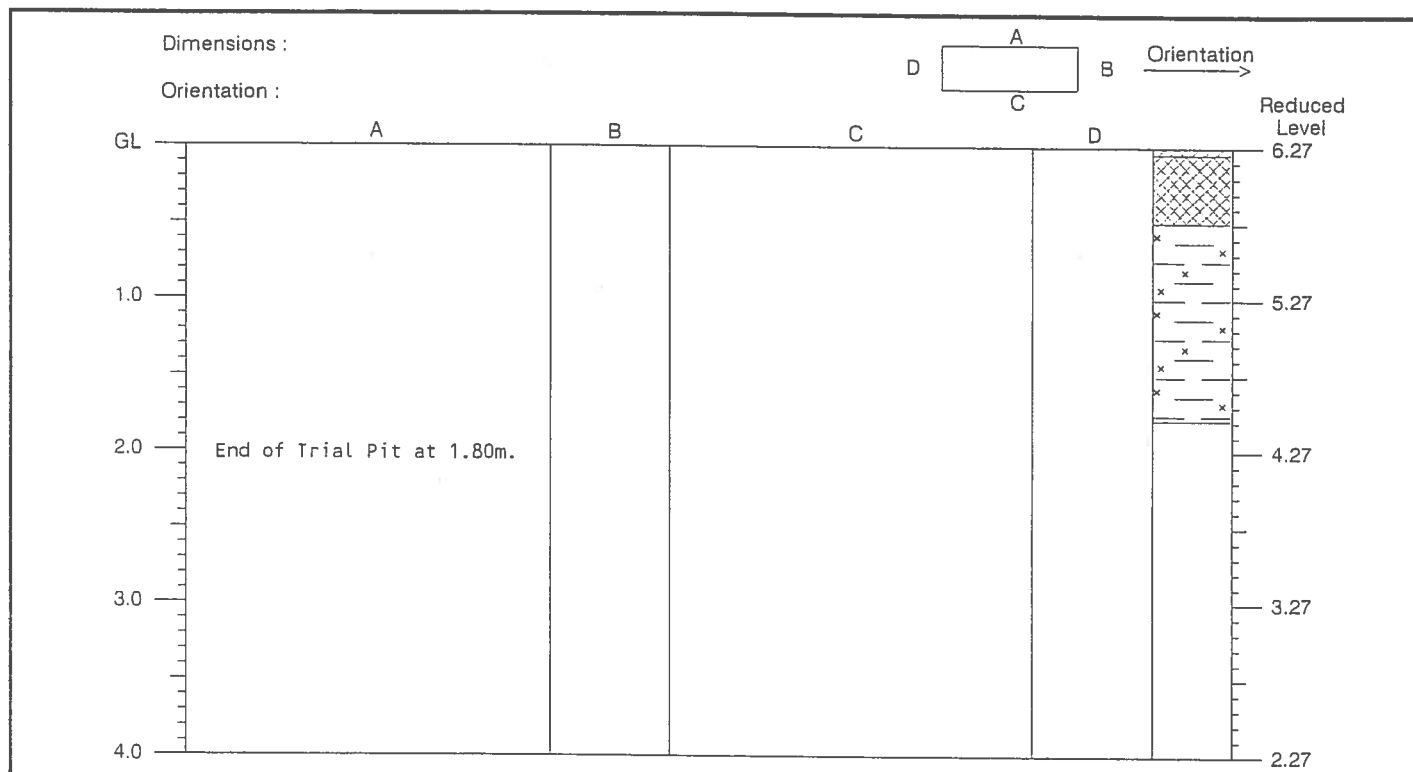
M4 Relief Road - Stage 2. Preliminary
 Chemical Investigation
 Ove Arup & Partners

Contract

150006

**Exploration Associates****Trial Pit**

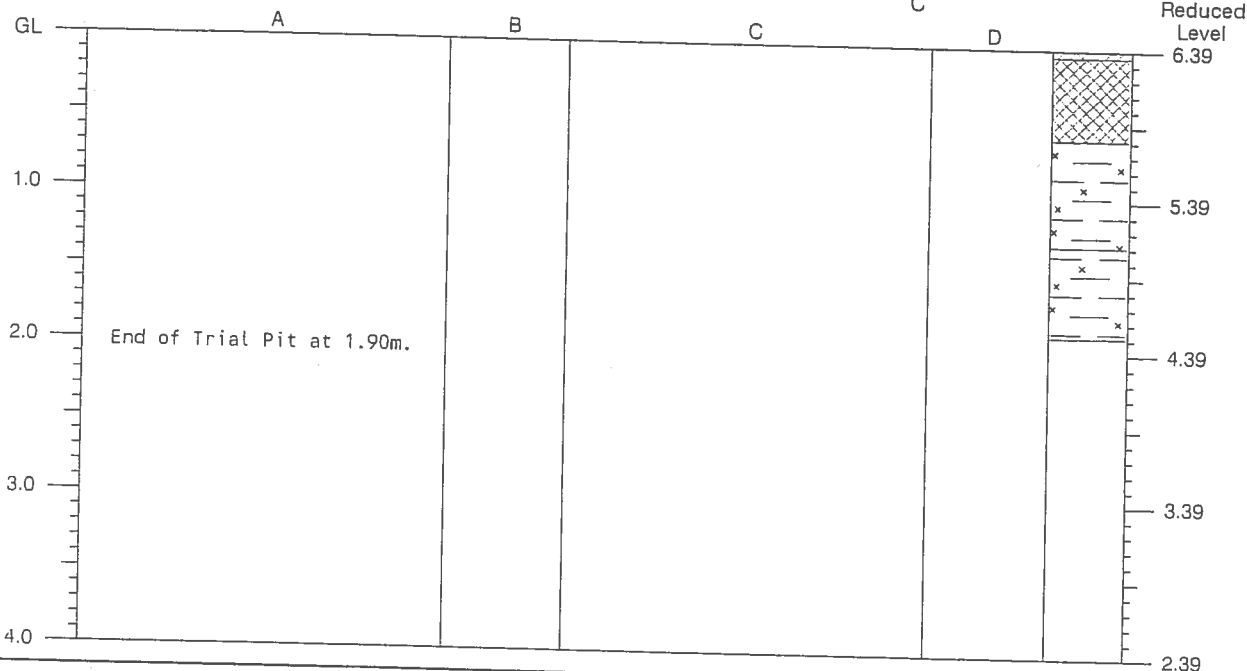
CH47A



Strata			Samples and Tests		
Depth (m)	No.	Description	Depth (m)	Type	Results
0.00-0.05	1	Brown very silty clayey TOPSOIL.			
0.05-0.50	2	MADE GROUND: Very compact brown and grey sandy sub-angular to sub-rounded fine to coarse gravel and cobbles of slag, ash, limestone and sandstone with some boulders (<400mm) of slag.	0.05-0.50	BTO	
0.50-1.80	3	Firm to stiff blue grey mottled brown very silty slightly organic CLAY with occasional some root material.from 1.50m becoming soft to firm.	1.00-1.20 1.50-1.80	BTO BTO	
Date of Excavation 04/02/00 Equipment JCB 3CX Stability Stable			Groundwater No. Struck Behaviour Not encountered during excavation		
			Ground Level 6.27 m OD Coordinates 333562.09 mE 185439.87 mN Logged by DC Checked by AF		
Remarks See key sheet and appendices for explanations.					
Trial Pit Record		Project		Contract	
		M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners		150006	
Exploration Associates				Trial Pit	
				CH48	

Dimensions :

Orientation :



Strata

Strata			Samples and Tests		
Depth (m)	No.	Description	Depth (m)	Type	Results
0.00-0.05	1	Brown very silty clayey TOPSOIL.			
0.05-0.60	2	MADE GROUND: Moderately compact brown very sandy sub-angular to sub-rounded fine to coarse gravel and cobbles of slag, ash and brick with occasional boulders (<600mm) of slag.	0.05-0.60	BTO	
0.60-1.30	3	Firm to stiff brown slightly mottled grey and orange brown very silty organic CLAY with much root material.	0.60-1.10	BTO	
1.30-1.90	4	Firm grey mottled brown very silty CLAY with occasional root material.	1.50-1.90	BTO	

Date of Excavation 04/02/00
Equipment JCB 3CX
Stability Stable

Groundwater
No. Struck Behaviour
Not encountered during excavation

Ground Level 6.39 m OD
Coordinates 334598.04 mE
185425.61 mN

Logged by DC
Checked by AF

Remarks

See key sheet
and appendices
for explanations.

Form 2/0

Trial Pit Record

Project

M4 Relief Road - Stage 2. Preliminary
Chemical Investigation
Ove Arup & Partners

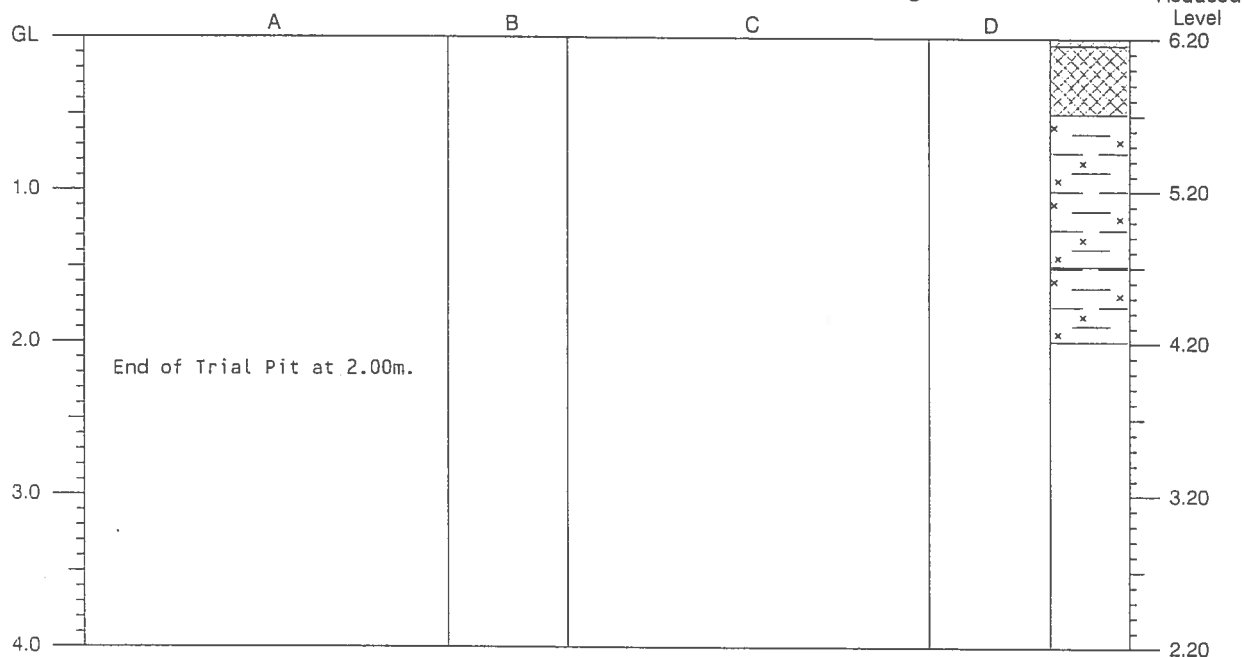
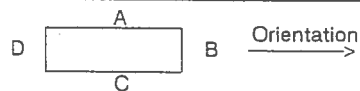
Contract 150006

Trial Pit CH49

 Exploration Associates

Dimensions :

Orientation :



Strata

Samples and Tests

Depth (m)	No.	Description	Depth (m)	Type	Results
0.00-0.05	1	Brown very silty clayey TOPSOIL.			
0.05-0.50	2	MADE GROUND: Moderately compact brown very sandy sub-angular to sub-rounded fine to coarse gravel and cobbles of slag, ash and siltstone with occasional boulders (<400mm) of slag.	0.05-0.50	BTO	
0.50-1.50	3	Firm brown mottled blue grey very silty slightly organic CLAY with much root material.	0.50-1.00	BTO	
1.50-2.00	4	Firm grey mottled brown very silty organic CLAY with occasional root material.	1.50-2.00	BTO	

Date of Excavation 04/02/00

Equipment JCB 3CX

Stability Stable

Groundwater

No. Struck Behaviour

1 0.50 slight seepage

Ground Level 6.20 m OD

Coordinates 334718.53 mE

185421.00 mN

Logged by DC

Checked by AF

Remarks

See key sheet
and appendices
for explanations.

Form 2/0

Trial Pit Record

Project

M4 Relief Road - Stage 2. Preliminary
Chemical Investigation
Ove Arup & Partners

Contract

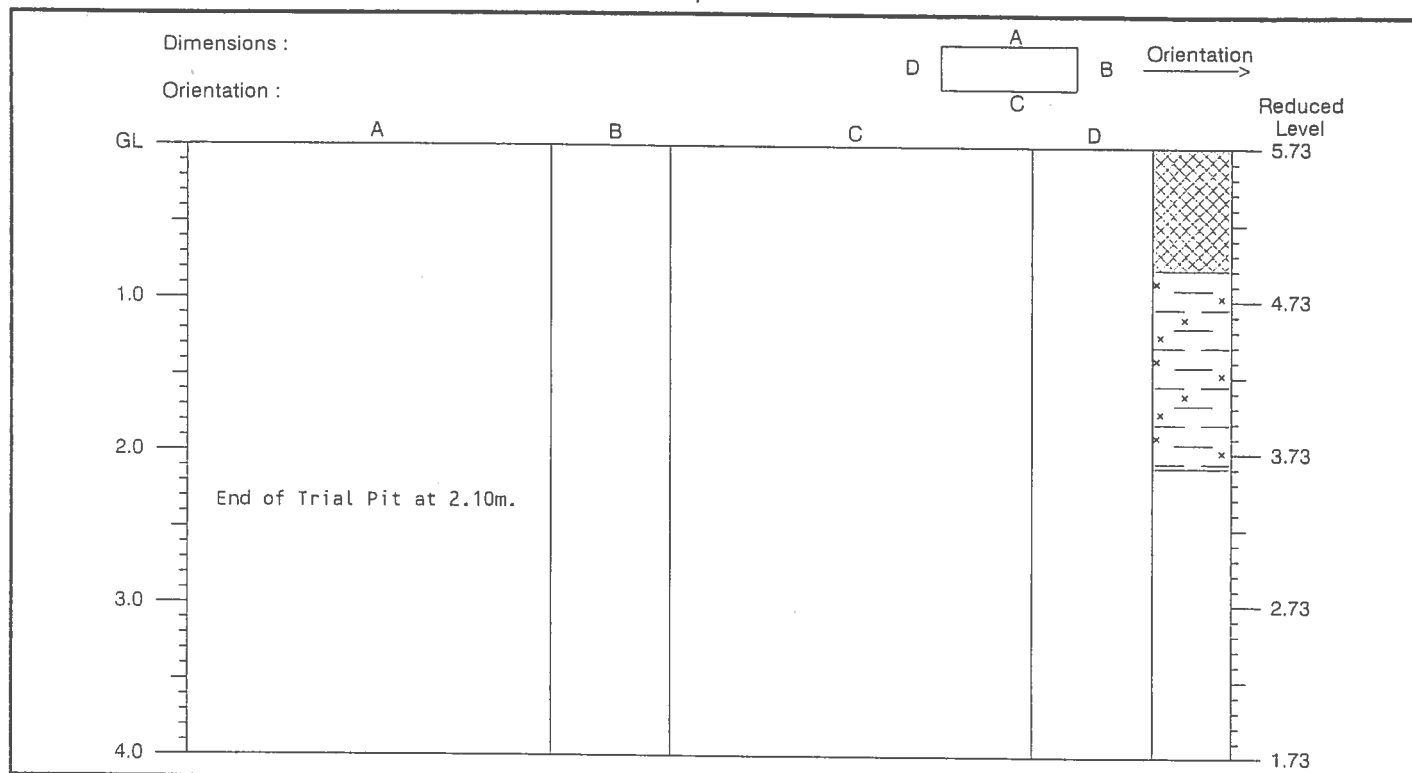
150006



Exploration Associates

Trial Pit

CH50



Strata			Samples and Tests		
Depth (m)	No.	Description	Depth (m)	Type	Results
0.00-0.80	1	MADE GROUND: Loose brown sandy slightly silty sub-angular to sub-rounded fine to coarse gravel and cobbles of slag, ash and brick.	0.00-0.50	BTO	
0.80-2.10	2	Firm to stiff blue grey mottled green grey and brown very silty organic CLAY with much root material.from 1.90m becoming soft to firm.	0.80-1.10	BTO	
			1.10-1.90	BTO	
			2.10-0.80	W	
Date of Excavation 03/02/00		Groundwater	Ground Level 5.73 m OD		
Equipment JCB 3CX		No. Struck Behaviour	Coordinates 338603.12 mE		
Stability Stable		1 0.80 medium inflow	185406.50 mN		
			Logged by DC		
			Checked by AF		
Remarks See key sheet and appendices for explanations.					
Trial Pit Record		Project	Contract		
		M4 Relief Road - Stage 2. Preliminary Chemical Investigation	150006		
		Ove Arup & Partners	Trial Pit		
			CH51		

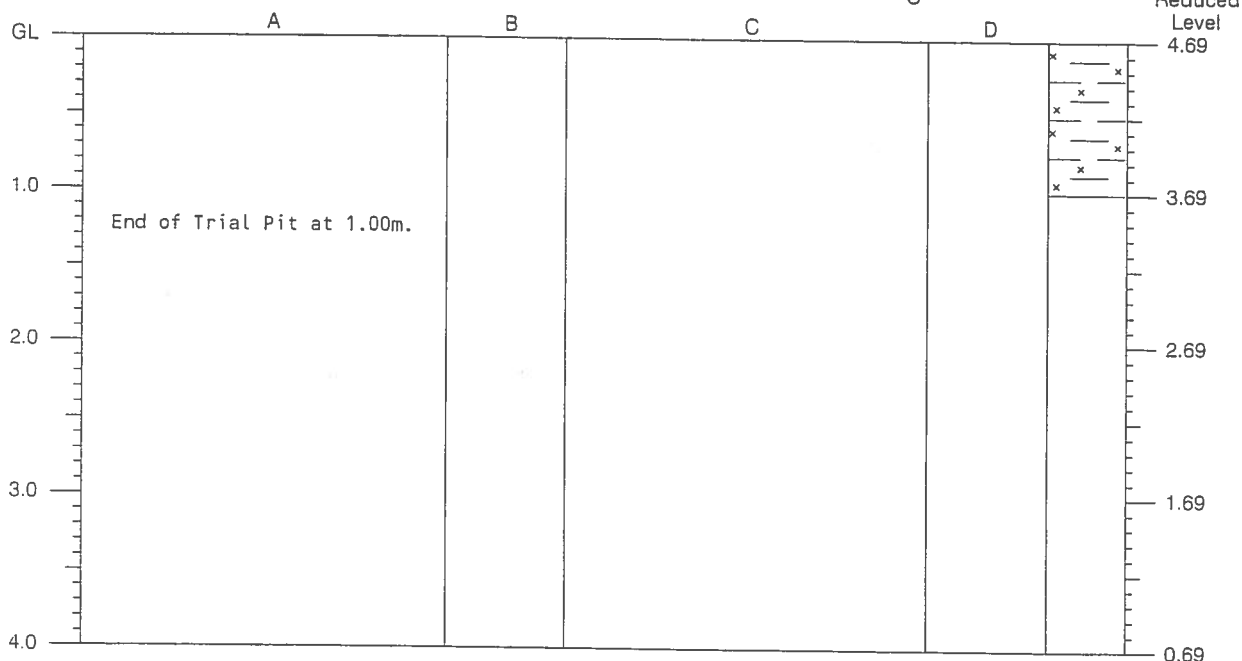
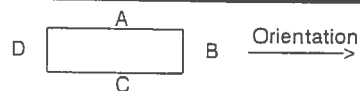


Exploration Associates

Form 2/0

Dimensions :

Orientation :



Strata

Samples and Tests

Depth (m)	No.	Description	Depth (m)	Type	Results
0.00-1.00	1	Soft to firm grey mottled brown and dark grey very silty organic CLAY with much root material.	0.50-1.00	BTO	
			0.00	W	

Date of Excavation 03/02/00

Equipment JCB 3CX

Stability Stable

Groundwater

No. Struck Behaviour

Surface groundwater present

Ground Level 4.69 m OD

Coordinates 338945.93 mE
185445.65 mNLogged by DC
Checked by AF

Remarks Pit stopped at 1.00m (surface water filled trial pit).

See key sheet
and appendices
for explanations.

Form 2/0

Trial Pit Record

Project

M4 Relief Road - Stage 2. Preliminary
Chemical Investigation
Ove Arup & Partners

Contract 150006

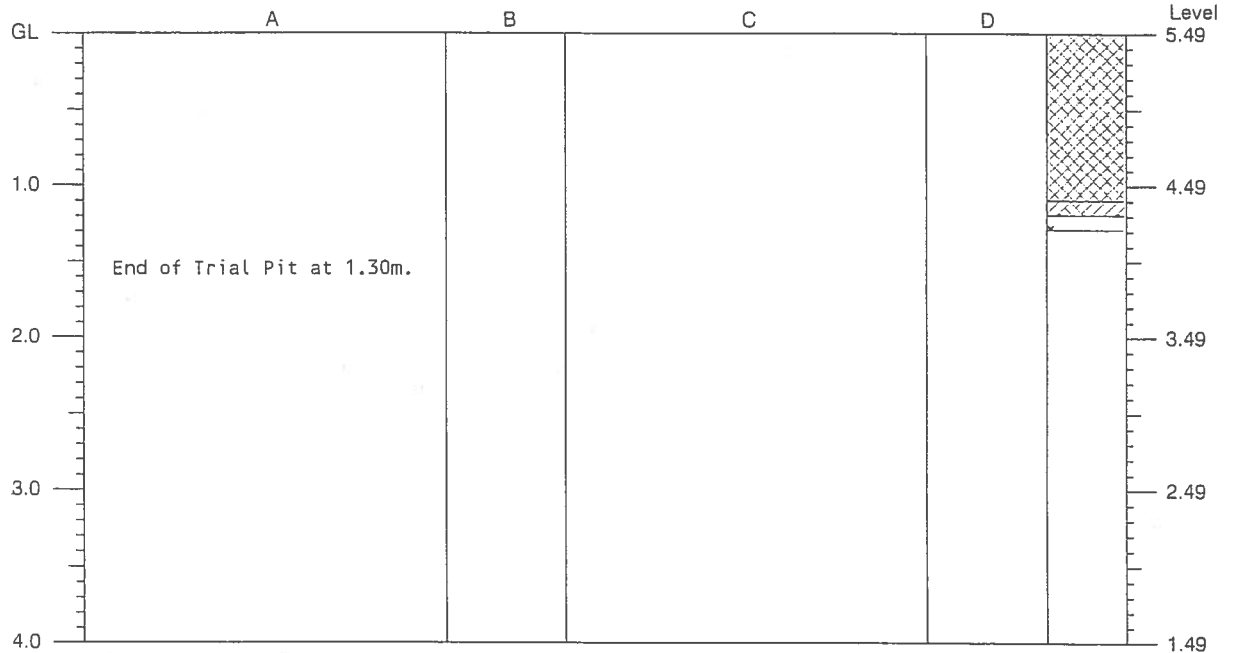
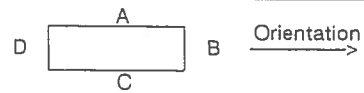
Trial Pit CH52



Exploration Associates

Dimensions : 1.80 x 1.00

Orientation : S



Strata

Samples and Tests

Depth (m)	No.	Description	Depth (m)	Type	Results
0.00-1.10	1	MADE GROUND: Moderately compact grey sandy slightly silty sub-angular to sub-rounded fine to coarse gravel and cobbles of slag, ash, siltstone and limestone	0.30 0.80	BTO W	
1.10-1.20	2	TOPSOIL			
1.20-1.30	3	Firm to stiff grey brown very silty CLAY with much root material	1.20	BTO	

Date of Excavation 16/02/00

Equipment JCB 3CX

Stability Stable

Groundwater

No. Struck Behaviour
1 0.80 Fast

Ground Level 5.49 m OD

Coordinates 338946.51 mE
185449.82 mN

Logged by DC

Checked by AF

Remarks

See key sheet
and appendices
for explanations.

Form 2/0

Trial Pit Record

Project

M4 Relief Road - Stage 2. Preliminary
Chemical Investigation
Ove Arup & Partners

Contract

150006

Trial Pit

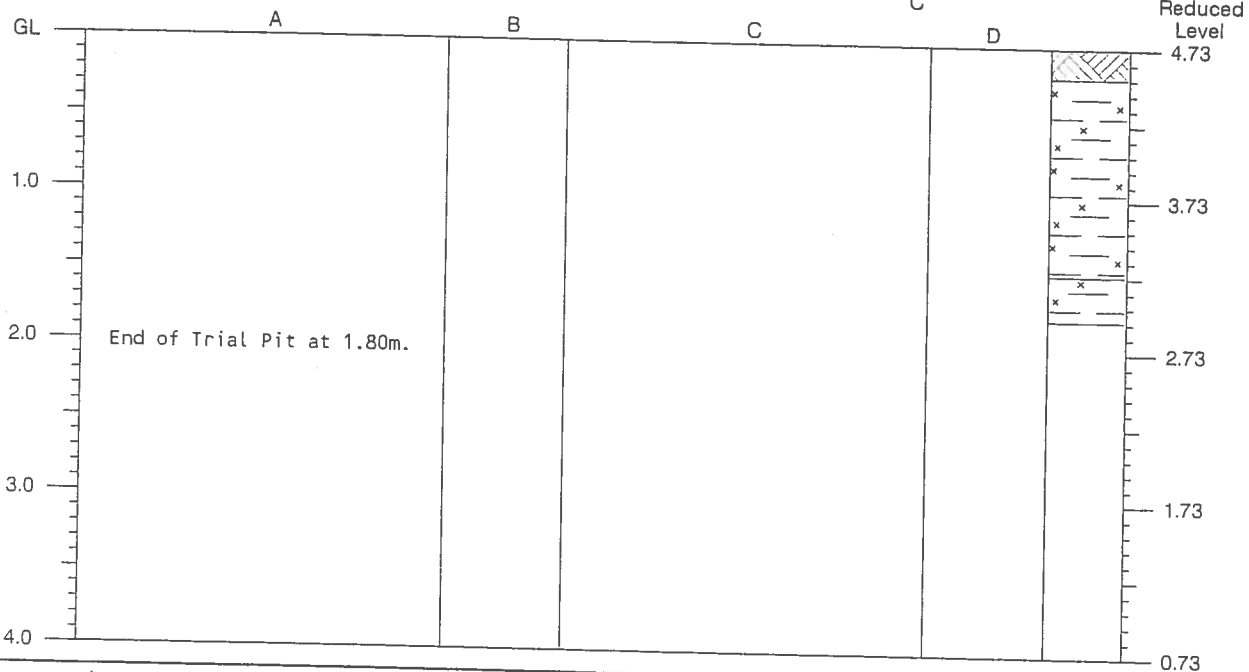
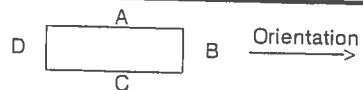
CH52A



Exploration Associates

Dimensions :

Orientation :



Strata

Samples and Tests

Depth (m)	No.	Description	Depth (m)	Type	Results
0.00-0.20	1	Brown very silty clayey TOPSOIL.			
0.20-1.50	2	Firm to stiff grey mottled brown very silty CLAY with much root material.	0.40-0.50	BTO	
1.50-1.80	3	Soft to firm blue grey mottled brown and dark grey very silty CLAY with much organic and root material.	1.50-1.80	W BTO	

Date of Excavation 03/02/00
Equipment JCB 3CX
Stability Stable

Groundwater
No. Struck Behaviour
1 1.50 medium flow

Ground Level 4.73 m OD
Coordinates 339261.29 mE
185485.43 mN

Logged by DC
Checked by AF

Remarks Water sample taken in rear adjacent to CH53 known as "size No. 9 rear sample".

See key sheet
and appendices
for explanations.

Form 2/0

Trial Pit Record

Project

Contract 150006

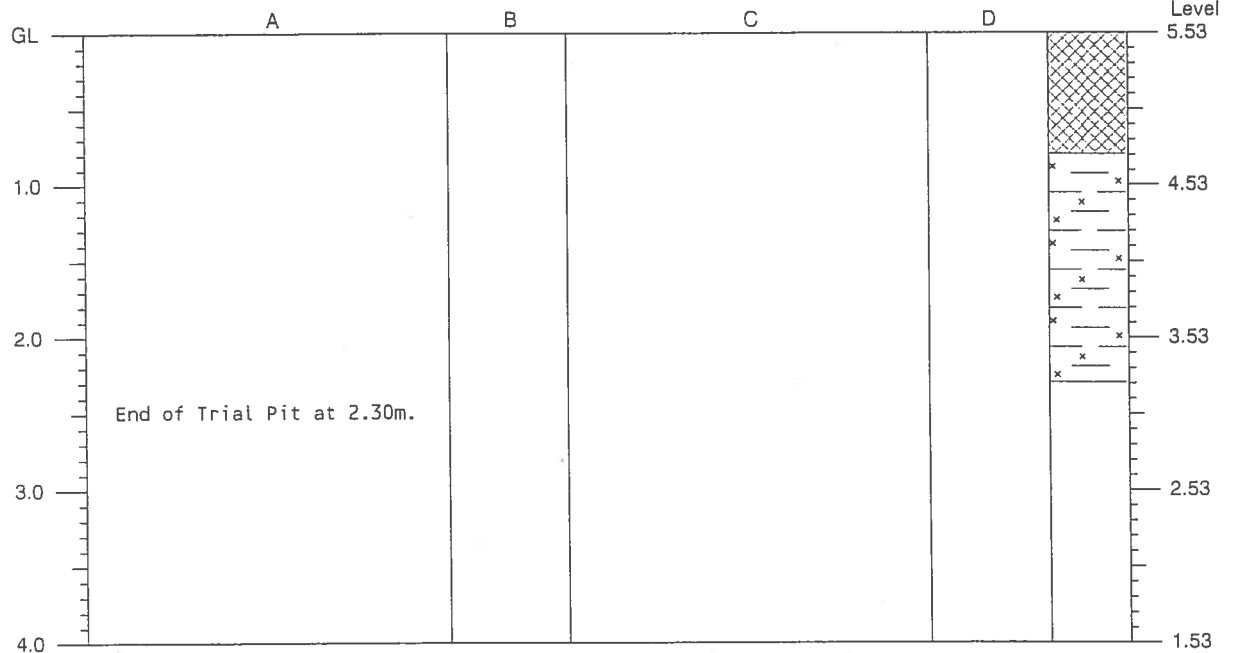
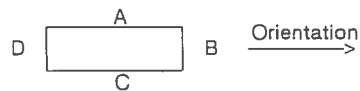
 **Exploration Associates**

M4 Relief Road - Stage 2. Preliminary
Chemical Investigation
Ove Arup & Partners

Trial Pit CH53

Dimensions : 1.90 x 1.10

Orientation : W



Strata

Samples and Tests

Depth (m)	No.	Description	Depth (m)	Type	Results
0.00-0.80	1	MADE GROUND: Loose grey sandy slightly silty sub-angular to sub-rounded fine to coarse gravel and cobbles of slag,ash,brick,siltstone and limestone	0.40	BTO	
0.80-2.30	2	Firm to stiff grey brown very silty CLAY with much root materialfrom 1.60m becoming soft to firm	1.10	BTO	

Date of Excavation 16/02/00
Equipment JCB 3CX
Stability Stable

Groundwater
No. Struck Behaviour
Not encountered during excavation

Ground Level 5.53 m OD
Coordinates 339262.37 mE
185482.23 mN

Logged by DC
Checked by AF

Remarks

See key sheet
and appendices
for explanations.

Form 2/0

Trial Pit Record

Project

M4 Relief Road - Stage 2. Preliminary
Chemical Investigation
Ove Arup & Partners

Contract

150006

Trial Pit

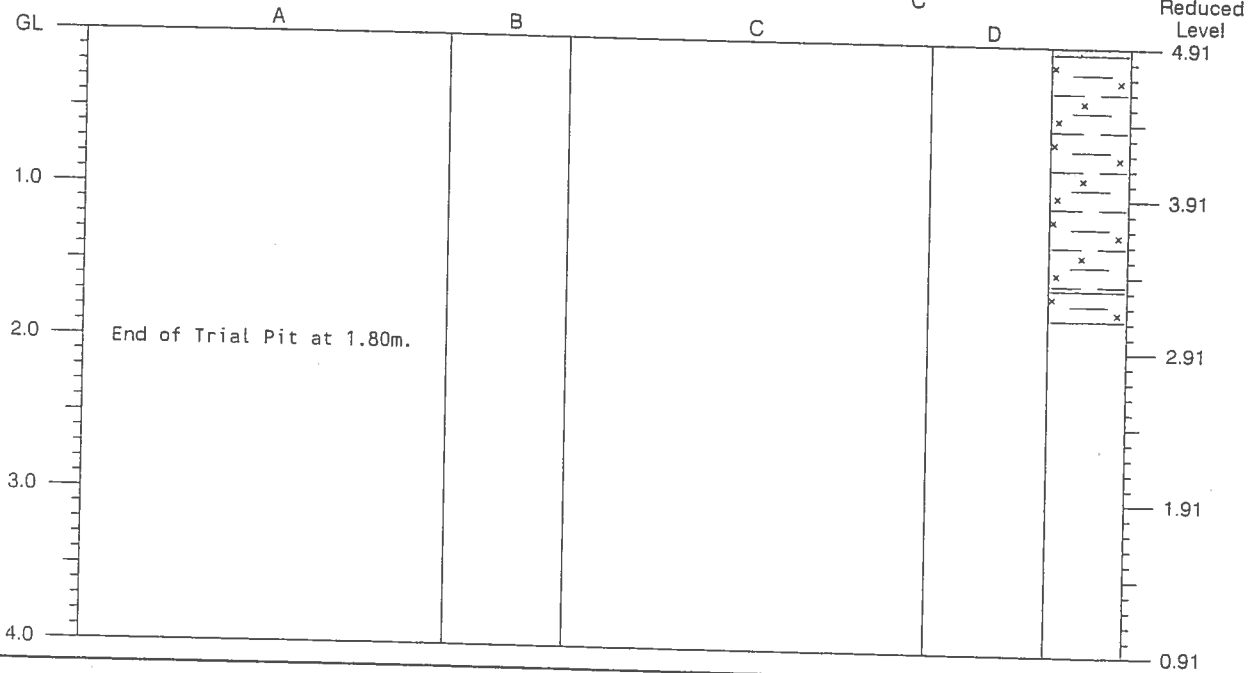
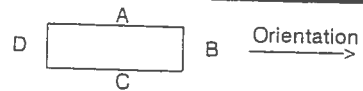
CH53A



Exploration Associates

Dimensions :

Orientation :



Strata

Depth (m)	No.	Description	Samples and Tests		
			Depth (m)	Type	Results
0.00-0.05	1	Brown very silty clayey TOPSOIL.			
0.05-1.60	2	Firm to stiff brown and grey very silty organic CLAY with much root material.from 1.10m becoming soft to firm.	0.05-0.50	BTO	
1.60-1.80	3	Soft to firm grey mottled brown and dark grey very silty organic CLAY with much root material.	1.60-1.80	BTO	

Date of Excavation 03/02/00

Equipment JCB 3CX

Stability Stable

Groundwater

No. Struck Behaviour
1 1.60 slight water seepage

Ground Level 4.91 m OD
Coordinates 339514.60 mE
185546.01 mN

Logged by DC
Checked by AF

Remarks

See key sheet
and appendices
for explanations.

Form 2/0

Trial Pit Record

Project

M4 Relief Road - Stage 2. Preliminary
Chemical Investigation
Ove Arup & Partners

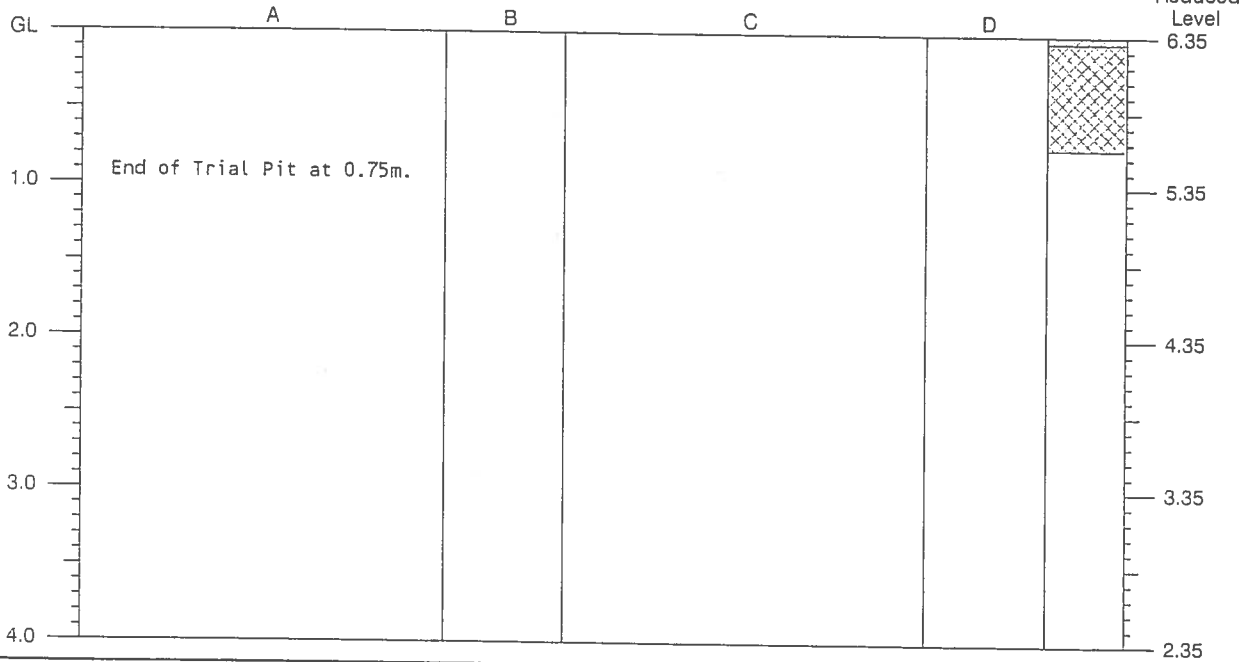
Contract 150006

Trial Pit CH54

 Exploration Associates

Dimensions : 0.50 x 0.50

Orientation : N/A



Strata

Samples and Tests

Depth (m)	No.	Description	Depth (m)	Type	Results
0.00-0.05	1	Brown clayey very silty TOPSOIL			
0.05-0.75	2	MADE GROUND: Firm to stiff grey brown very silty CLAY with occasional to some root material	0.40	BTO	

Date of Excavation 16/02/00
Equipment Hand Dug
Stability Stable

Groundwater
No. Struck Behaviour
Not encountered during excavation

Ground Level 6.35 m OD
Coordinates 339512.35 mE
185556.72 mN

Logged by DC
Checked by AF

Remarks CH54A Excavated by hand due to unsuitable access for the JCB 3CX

See key sheet
and appendices
for explanations.

Form 2/0

Trial Pit Record

Project

Contract 150006

 **Exploration Associates**

M4 Relief Road - Stage 2. Preliminary
Chemical Investigation
Ove Arup & Partners

Trial Pit CH54A

Sampling					Strata				
Depth	Type	Casing Depth	Date/ Water	SPT N (Cu)	Description	Depth (Thickness)	Level	Legend	
0.00-0.25	T		05/02		MADE GROUND: Loose brown very silty sandy sub-angular to sub-rounded fine to coarse gravel of slag, ash and sandstone with much root material	G.L.	21.45		
0.25-0.75	TO		0000			0.25	21.20		
0.75-1.00	B				Purple and green grey very silty highly weathered MUDSTONE	(1.45)			
1.00-1.70	TO								
1.70-1.75	T	NIL	DRY			1.70	19.75		
1.75					Red brown silty weathered MUDSTONE End of Borehole.	1.75	19.70		
Equipment: Hand Held Window Sampler					Groundwater No. Struck Behaviour Sealed		Ground Level 21.45 m. OD Coordinates 341851.79 mE 187531.14 mN		
Borehole Dia (mm) Casing Dia (mm) 80 to 1.00m 60 to 1.75m					No groundwater encountered		Drilled by DC Logged by DC Checked by AF		
Remarks Refusal at 1.75m See key sheet and appendices for explanations.									
Borehole Record					Project M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners		Contract 150006		
 Exploration Associates							Borehole CH55(1 of 1)		

Form 1/0

Sampling						Strata						
Depth	Type	Casing Depth	Date/ Water	SPT N (Cu)	Description	Depth (Thickness)		Level	Legend			
0.00-0.40	T		05/02 0000 DRY		Soft to firm brown very silty slightly sandy CLAY with occasional to some root material	G.L. (0.40)		20.64				
0.40-0.60	T	NIL	DRY		Red and yellow brown weathered MUDSTONE	0.40 0.60		20.24 20.04				
0.60					 End of Borehole.							
Equipment: Hand Held Window Sampler					Groundwater							
Borehole Dia (mm) Casing Dia (mm) 80 to 0.60m					No. Struck Behaviour	Sealed	Ground Level Coordinates 20.64 m OD 341890.96 mE 187529.33 mN					
					No groundwater encountered		Drilled by DC Logged by DC Checked by AF					
Remarks Refusal at 0.60m See key sheet and appendices for explanations.												
Borehole Record					Project			Contract				
Exploration Associates					M4 Relief Road - Stage 2. Preliminary Chemical Investigation Ove Arup & Partners			150006				
								Borehole CH56(1 of 1)				

ENCLOSURE B

In Situ Testing

On Site Water Testing

1

Water Testing

HOLE	TEMP. deg. C	pH	COND. µs
Site no7	7.9	6.58	721
ch52a	8.0	7.14	621
ch40	7.1	8.19	515
ch45	7.2	8.24	480
ch17	7.0	7.92	627
ch53	8.8	6.92	1590
ch52a	9.1	7.63	467
site no9	8.9	7.28	734
ch51	8.5	10.20	452
ch9	7.9	6.94	814
ch8	7.4	7.02	796
site no2	7.0	6.98	521
ch33	7.2	7.99	624
ch41	7.0	8.02	547
ch12	7.2	6.42	627
ch32	7.6	7.41	728

On Site Water Testing	Project M4 Relief Road - Stage 2,Preliminary Chemical Investigation Ove Arup & partners	Contract 150006
Exploration Associates		Figure No WT1

ENCLOSURE C

Laboratory Test Results

Contaminant Test Results

Soils	CL/000368, CL/000484, CL/000542, CL/000543 & CL/000581
Leachates	CL/000368, CL/000484, CL/000542, & CL/000543
Waters	W/EXR/000694, W/EXR/000771 & W/EXR/000800



TEST REPORT SOIL SAMPLE ANALYSIS



1252
1411

TES Report No. CL/000368

Site: M4 Relief Road

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
Mid Glamorgan
CF31 1JZ

The 14 samples described in this report were scheduled for analysis by TES Bretby on Wednesday, 2 February 2000. The analysis was completed by Wednesday, 16 February 2000.

Tests marked as 'not UKAS accredited' and any opinions or interpretations expressed herein are outside the scope of any UKAS accreditation held by TES Bretby laboratories.

The following tables are contained in this report:

Table 1 Sample Descriptions
Table 2 Main Analysis Results

On behalf of
TES Bretby : J. Hannah
J Hannah Project Co-ordinator

Date of Issue: 16/02/00

Tests marked 'not UKAS accredited' in this report are not included in the UKAS Accreditation Schedule for our laboratory.

TES Bretby accepts no responsibility for the sampling related to the above results

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TES Bretby
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TEST REPORT

SOIL SAMPLE ANALYSIS

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
Mid Glamorgan
CF31 1JZ

1252
1411

TES Report No. 000368

Site: M4 Relief Road

Page No	Assessed Area	ID No. EFS/CL	Sample	Depth(m) from to	Description
1	M4 Relief Road	0002699	CH/ 1/	0.20	CH1 0.20
1	M4 Relief Road	0002700	CH/ 2/	0.20	CH2 0.20
1	M4 Relief Road	0002701	CH/ 3/	0.10	CH3 0.10
1	M4 Relief Road	0002703	CH/ 3/	1.17	CH3 1.17
2	M4 Relief Road	0002704	CH/ 4/	0.60	CH4 0.60
2	M4 Relief Road	0002706	CH/ 20/	0.00	CH20 0.00
2	M4 Relief Road	0002710	CH/ 20/	1.00	CH20 1.00
2	M4 Relief Road	0002707	CH/ 21/	0.00	CH21 0.00
3	M4 Relief Road	0002711	CH/ 21/	2.50	CH21 2.50
3	M4 Relief Road	0002708	CH/ 22/	0.30	CH22 0.30
3	M4 Relief Road	0002709	CH/ 22/	0.75	CH22 0.75
3	M4 Relief Road	0002712	CH/ 22/	1.00	CH22 1.00
4	M4 Relief Road	0002702	CH/ 23/	0.16	CH23 0.16
4	M4 Relief Road	0002705	CH/ 23/	0.46	CH23 0.46

Date of Issue: 16/02/00

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TES Bretby accepts no responsibility for the sampling related to the above results

TES Bretby
Report 000368
Table 1
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TEST REPORT

SOIL SAMPLE ANALYSIS

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
Mid Glamorgan
CF31 1JZ

TES Report No. 000368

Site: M4 Relief Road

Customer reference	CH 1	CH 2	CH 3	CH 3
Depth (m)	0.20	0.20	0.10	1.17
Date logged	02/02/00	02/02/00	02/02/00	02/02/00
TES Bretby ID Number	CL/0002699	CL/0002700	CL/0002701	CL/0002703

UKAS accredited	Test No.	CL/0002699	CL/0002700	CL/0002701	CL/0002703
Chromium (total)	ICPSSS11	19	17	17	15
CN- (total)	ICTSCN28	<1	<1	<1	<1
Copper	ICPSSS11	6	8	8	7
Lead	ICPSSS11	30	17	30	14
Nickel	ICPSSS11	14	15	8	19
Oil FTIR	FTIRSW			1920	
PAH (screening)	PAHSCUV	<10	<10	502	<10
pH units	WSLM3	6.4	7.4	8.2	7.2
Phenol Index	WSLM4	<0.5	<0.5	<0.5	<0.5
SO4-- (acid sol)	ICPACS58	307	77	1020	26
TPH FTIR (AR)	FTIRSW			1790	
Zinc	ICPSSS11	47	47	54	41

not UKAS accredited		CL/0002699	CL/0002700	CL/0002701	CL/0002703
Arsenic (Ms)		6.40	6.70	5.40	5.00
Boron.		<0.5	<0.5	<0.5	<0.5
Cadmium (Ms)		0.12	0.05	0.28	0.09
Mercury (Ms)		<0.01	<0.01	<0.01	<0.01
Selenium (MS)				0.88	<0.50

Results expressed as mg/kg Air Dried unless stated otherwise

SO4 Analysis not conducted in accordance with BS1377

AR denotes analysis conducted on As Received Sample

Date of Issue: 16/02/00

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TES Bretby
Report 000368
Table 2
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TEST REPORT

SOIL SAMPLE ANALYSIS

1252
1411

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
Mid Glamorgan
CF31 1JZ

TES Report No. 000368

Site: M4 Relief Road

Customer reference	CH 4	CH 20	CH 20	CH 21
Depth (m)	0.60	0.00	1.00	0.00
Date logged	02/02/00	02/02/00	02/02/00	02/02/00
TES Bretby ID Number	CL/0002704	CL/0002706	CL/0002710	CL/0002707

UKAS accredited	Test No.	CL/0002704	CL/0002706	CL/0002710	CL/0002707
Chromium (total)	ICPSSS11	21	33	27	42
CN- (total)	ICTSCN28	<1	<1	<1	<1
Copper	ICPSSS11	30	315	26	164
Lead	ICPSSS11	32	82	34	102
Nickel	ICPSSS11	16	32	24	24
Oil FTIR	FTIRSW		112		172
PAH (screening)	PAHSCUV	<10	43	14	44
pH units	WSLM3	7.2	8.1	8.1	9.2
Phenol Index	WSLM4	<0.5	<0.5	<0.5	<0.5
SO4-- (acid sol)	ICPACS58	122	712	632	1080
TPH FTIR (AR)	FTIRSW		112		79
Zinc	ICPSSS11	43	272	103	116

not UKAS accredited		CL/0002704	CL/0002706	CL/0002710	CL/0002707
Arsenic (Ms)		4.00	17	11.00	17
Boron.		<0.5	<0.5	1.0	0.9
Cadmium (Ms)		0.04	0.27	0.11	0.17
Mercury (Ms)		<0.01	0.24	<0.01	1.76
Selenium (MS)		<0.50	1.55	0.90	1.60

Results expressed as mg/kg Air Dried unless stated otherwise

SO4 Analysis not conducted in accordance with BS1377

AR denotes analysis conducted on As Received Sample

Date of Issue: 16/02/00

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TEST REPORT

SOIL SAMPLE ANALYSIS

1252
1411

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
Mid Glamorgan
CF31 1JZ

TES Report No. 000368

Site: M4 Relief Road

Customer reference	CH 21	CH 22	CH 22	CH 22
Depth (m)	2.50	0.30	0.75	1.00
Date logged	02/02/00	02/02/00	02/02/00	02/02/00
TES Bretby ID Number	CL/0002711	CL/0002708	CL/0002709	CL/0002712

UKAS accredited	Test No.	CL/0002711	CL/0002708	CL/0002709	CL/0002712
Chromium (total)	ICPSSS11	39	27	15	34
CN- (total)	ICTSCN28	<1	1	<1	1
Copper	ICPSSS11	22	490	26	18
Lead	ICPSSS11	43	426	36	38
Nickel	ICPSSS11	31	44	12	28
Oil FTIR	FTIRSW		1120	87	
PAH (screening)	PAHSCUV	301	288	84	13
pH units	WSLM3	8.1	8.2	8.2	8.0
Phenol Index	WSLM4	1.5	<0.5	<0.5	<0.5
SO4-- (acid sol)	ICPACS58	1090	1320	444	513
TPH FTIR (AR)	FTIRSW		1220	65	
Zinc	ICPSSS11	120	434	47	87

not UKAS accredited		CL/0002711	CL/0002708	CL/0002709	CL/0002712
Arsenic (Ms)		10.00	31	10.00	12.00
Boron.		2.5	<0.5	<0.5	1.5
Cadmium (Ms)		0.15	1.27	0.11	0.13
Mercury (Ms)		<0.01	0.46	0.04	<0.01
Selenium (MS)		<0.50	1.70	<0.50	<0.50

Results expressed as mg/kg Air Dried unless stated otherwise

SO4 Analysis not conducted in accordance with BS1377

AR denotes analysis conducted on As Received Sample

Date of Issue: 16/02/00

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Report 000368
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TEST REPORT

SOIL SAMPLE ANALYSIS

TES
Bretby



1252
1411

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
Mid Glamorgan
CF31 1JZ

TES Report No. 000368

Site: M4 Relief Road

Customer reference	CH	CH		
Depth (m)	23 0.16	23 0.46		
Date logged	02/02/00	02/02/00		
TES Bretby ID Number	CL/0002702	CL/0002705		

UKAS accredited	Test No.	CL/0002702	CL/0002705		
Chromium (total)	ICPSSS11	44	16		
CN- (total)	ICTSCN28	<1	<1		
Copper	ICPSSS11	743	46		
Lead	ICPSSS11	969	78		
Nickel	ICPSSS11	91	14		
Oil FTIR	FTIRSW	732			
PAH (screening)	PAHSCUV	832	35		
pH units	WSLM3	7.8	8.4		
Phenol Index	WSLM4	0.5	<0.5		
SO4-- (acid sol)	ICPACS58	1660	325		
TPH FTIR (AR)	FTIRSW	67			
Zinc	ICPSSS11	718	67		

not UKAS accredited		CL/0002702	CL/0002705		
Arsenic (Ms)		83	13.00		
Boron.		<0.5	<0.5		
Cadmium (Ms)		6.08	0.17		
Mercury (Ms)		1.32	0.08		
Selenium (MS)		1.21	0.53		

Results expressed as mg/kg Air Dried unless stated otherwise

SO4 Analysis not conducted in accordance with BS1377

AR denotes analysis conducted on As Received Sample

Date of Issue: 16/02/00

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TES Bretby
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TEST REPORT SOIL SAMPLE ANALYSIS



1252
1411

TES Report No. CL/000484

Site: M4 Relief Road

Exploration Associates
Unit 15
Crosby Yard
Bridgend
Mid Glamorgan
CF31 1JZ

The 24 samples described in this report were scheduled for analysis by TES Bretby on Friday, 11 February 2000. The analysis was completed by Thursday, 24 February 2000.

Tests marked as 'not UKAS accredited' and any opinions or interpretations expressed herein are outside the scope of any UKAS accreditation held by TES Bretby laboratories.

The following tables are contained in this report:

Table 1 Sample Descriptions
Table 2 Main Analysis Results
Tables of Semi-Volatile Organic Compounds (21 Pages)
Tables of Volatile Organic Compounds (5 Pages)

On behalf of
TES Bretby : J. Hannah
J Hannah Project Co-ordinator

Date of Issue: 24/02/00

Tests marked 'not UKAS accredited' in this report are not included in the UKAS Accreditation Schedule for our laboratory.

TES Bretby accepts no responsibility for the sampling related to the above results

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TES Bretby
Report 000484
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TEST REPORT

SOIL SAMPLE ANALYSIS

Exploration Associates
Unit 15
Crosby Yard
Bridgend
Mid Glamorgan
CF31 1JZ

TES Report No. 000484

Site: M4 Relief Road

Page No	Assessed Area	ID No. EFS/CL	Sample	Depth (m) from to	Description
1	M4 Relief Road	0003497	CH/ 5/	0.00	CH5 0.0
1	M4 Relief Road	0003498	CH/ 5/	0.80	CH5 0.8
1	M4 Relief Road	0003499	CH/ 6/	0.00	CH6 0.0
1	M4 Relief Road	0003500	CH/ 6/	1.50	CH6 1.5
2	M4 Relief Road	0003501	CH/ 7/	0.00	CH7 0.0
2	M4 Relief Road	0003502	CH/ 34/	0.15	CH34 0.15
2	M4 Relief Road	0003503	CH/ 34/	1.40	CH34 1.40
2	M4 Relief Road	0003504	CH/ 35/	0.00	CH35 0.00
3	M4 Relief Road	0003505	CH/ 35/	0.50	CH35 0.50
3	M4 Relief Road	0003506	CH/ 35/	0.80	CH35 0.80
3	M4 Relief Road	0003507	CH/ 36/	0.10	CH36 0.10
3	M4 Relief Road	0003508	CH/ 36/	0.70	CH36 0.70
4	M4 Relief Road	0003509	CH/ 37/	0.00	CH37 0.00
4	M4 Relief Road	0003510	CH/ 37/	1.20	CH37 1.20
4	M4 Relief Road	0003511	CH/ 38/	0.10	CH38 0.10
4	M4 Relief Road	0003512	CH/ 38/	1.30	CH38 1.30
5	M4 Relief Road	0003513	CH/ 39/	0.00	CH39 0.00
5	M4 Relief Road	0003514	CH/ 39/	1.00	CH39 1.00
5	M4 Relief Road	0003515	CH/ 40/	0.00	CH40 0.00
5	M4 Relief Road	0003516	CH/ 40/	0.50	CH40 0.50

Date of Issue: 24/02/00

Tests not marked 'not UKAS accredited' in this report are not included in the UKAS Accreditation Schedule for our laboratory.

TES Bretby accepts no responsibility for the sampling related to the above results

TES Bretby
Report 000484
Table 1
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TEST REPORT

SOIL SAMPLE ANALYSIS

1252
1411

Exploration Associates
Unit 15
Crosby Yard
Bridgend
Mid Glamorgan
CF31 1JZ

TES Report No. 000484

Site: M4 Relief Road

Page No	Assessed Area	ID No. EFS/CL	Sample	Depth (m) from to	Description
6	M4 Relief Road	0003517	CH/ 41/	0.05	CH41 0.05
6	M4 Relief Road	0003518	CH/ 41/	1.40	CH41 1.40
6	M4 Relief Road	0003519	CH/ 42/	0.20	CH42 0.20
6	M4 Relief Road	0003520	CH/ 42/	1.20	CH42 1.20

Date of Issue: 24/02/00

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TEST REPORT SOIL SAMPLE ANALYSIS



1252
1411

Exploration Associates
Unit 15
Crosby Yard
Bridgend
Mid Glamorgan
CF31 1JZ

TES Report No. 000484

Site: M4 Relief Road

Customer reference	CH 5	CH 5	CH 6	CH 6
Depth (m)	0.00	0.80	0.00	1.50
Date logged	11/02/00	11/02/00	11/02/00	11/02/00
TES Bretby ID Number	CL/0003497	CL/0003498	CL/0003499	CL/0003500

UKAS accredited	Test No.	CL/0003497	CL/0003498	CL/0003499	CL/0003500
Chromium (total)	ICPSSS11	25	29	16	16
CN- (total)	ICTSCN28	<1	<1	<1	<1
Copper	ICPSSS11	74	178	27	31
Lead	ICPSSS11	1610	1920	500	545
Nickel	ICPSSS11	23	27	16	17
Oil FTIR	FTIRSW	641	1946	338	183
PAH (screening)	PAHSCUV	75	201	79	72
pH units	WSLM3	8.3	8.0	8.4	8.9
Phenol Index	WSLM4	<0.5	0.6	<0.5	<0.5
SO4-- (acid sol)	ICPACS58	1110	1300	868	798
TPH FTIR (AR)	FTIRSW	1590	973	496	108
Zinc	ICPSSS11	186	280	119	130

not UKAS accredited		CL/0003497	CL/0003498	CL/0003499	CL/0003500
Arsenic (Ms)		9.90	13.00	7.70	6.60
Boron.		0.6	1.3	0.5	1.0
Cadmium (Ms)		1.10	1.32	0.51	0.27
Mercury (Ms)		0.42	0.41	0.15	0.12
Selenium (MS)		0.83	0.72	0.61	0.58

Results expressed as mg/kg Air Dried unless stated otherwise

SO4 Analysis not conducted in accordance with BS1377

AR denotes analysis conducted on As Received Sample

Date of Issue: 24/02/00

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TES Bretby accepts no responsibility for the sampling related to the above results

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TES Bretby
Report 000484
Table 2
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TEST REPORT

SOIL SAMPLE ANALYSIS

Exploration Associates
Unit 15
Crosby Yard
Bridgend
Mid Glamorgan
CF31 1JZ

TES Report No. 000484

Site: M4 Relief Road

Customer reference	CH 7	CH 34	CH 34	CH 35
Depth (m)	0.00	0.15	1.40	0.00
Date logged	11/02/00	11/02/00	11/02/00	11/02/00
TES Bretby ID Number	CL/0003501	CL/0003502	CL/0003503	CL/0003504

UKAS accredited	Test No.	CL/0003501	CL/0003502	CL/0003503	CL/0003504
Chromium (total)	ICPSSS11	38	45	38	56
CN- (total)	ICTSCN28	<1	<1	<1	<1
Copper	ICPSSS11	10	17	14	343
Cyanide (Free)	BGCN22		<1	<1	<1
Lead	ICPSSS11	43	59	45	95
Nickel	ICPSSS11	38	34	33	31
Oil FTIR	FTIRSW		178	104	280
PAH (screening)	PAHSCUV	<10	<10	<10	50
PCB - ARO	PCBAROEC		1.2	2.8	17
pH units	WSLM3	8.0	7.7	8.1	9.4
Phenol Index	WSLM4	<0.5	<0.5	<0.5	<0.5
SO4-- (acid sol)	ICPACS58	152	575	424	1630
Sulphide	ICTSCN28		5	<1	<1
SVOC	SVOCSW		*	*	*
Thiocyanate	WSLM5		8	<2	<2
TPH FTIR (AR)	FTIRSW		116	420	275
VOC	VOCSW				*
Zinc	ICPSSS11	99	113	100	239

not UKAS accredited		CL/0003501	CL/0003502	CL/0003503	CL/0003504
Arsenic (Ms)		18	16	13.00	6.90
Boron.		1.3	2.9	1.0	0.8
Cadmium (Ms)		0.08	0.13	0.13	1.37
Mercury (Ms)		0.08	0.18	0.43	15.2
Selenium (MS)			<0.50	<0.50	1.51

Results expressed as mg/kg Air Dried unless stated otherwise

SO4 Analysis not conducted in accordance with BS1377

AR denotes analysis conducted on As Received Sample

* denotes results included on attached sheet

Date of Issue: 24/02/00

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TES Bretby
Report 000484
Table 2
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TEST REPORT

SOIL SAMPLE ANALYSIS

Exploration Associates
 Unit 15
 Crosby Yard
 Bridgend
 Mid Glamorgan
 CF31 1JZ

1252
 1411

TES Report No. 000484

Site: M4 Relief Road

Customer reference	CH 35	CH 35	CH 36	CH 36
Depth (m)	0.50	0.80	0.10	0.70
Date logged	11/02/00	11/02/00	11/02/00	11/02/00
TES Bretby ID Number	CL/0003505	CL/0003506	CL/0003507	CL/0003508

UKAS accredited	Test No.	CL/0003505	CL/0003506	CL/0003507	CL/0003508
Chromium (total)	ICPSSS11	31	30	25	33
CN- (total)	ICTSCN28	<1	<1	<1	<1
Copper	ICPSSS11	22	12	22	10
Cyanide (Free)	BGCN22	<1	<1	<1	<1
Lead	ICPSSS11	71	45	128	37
Nickel	ICPSSS11	25	28	11	29
Oil FTIR	FTIRSW	365	238	1749	67
PAH (screening)	PAHSCUV	33	18	432	<10
PCB - ARO	PCBAROEC	18	7.8	27	1.5
pH units	WSLM3	7.8	7.8	9.5	8.2
Phenol Index	WSLM4	1.3	0.5	<0.5	<0.5
SO4-- (acid sol)	ICPACS58	944	485	2270	328
Sulphide	ICTSCN28	13	6	10	<1
SVOC	SVOCSW	*	*	*	*
Thiocyanate	WSLM5	13	4	<2	<2
TPH FTIR (AR)	FTIRSW	256	40	1530	195
VOC	VOCSW	*	*		
Zinc	ICPSSS11	112	90	303	80

not UKAS accredited		CL/0003505	CL/0003506	CL/0003507	CL/0003508
Arsenic (Ms)		15.00	13.00	6.00	8.90
Boron.		2.4	2.9	<0.5	1.5
Cadmium (Ms)		0.27	0.17	7.17	0.09
Mercury (Ms)		0.67	0.20	1.09	0.08
Selenium (MS)		0.89	<0.50	1.55	<0.50

Results expressed as mg/kg Air Dried unless stated otherwise

SO4 Analysis not conducted in accordance with BS1377

AR denotes analysis conducted on As Received Sample

* denotes results included on attached sheet

Date of Issue: 24/02/00

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 Report 000484
 Table 2
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TEST REPORT

SOIL SAMPLE ANALYSIS

 1252
 1411

 Exploration Associates
 Unit 15
 Crosby Yard
 Bridgend
 Mid Glamorgan
 CF31 1JZ

TES Report No. 000484

 Site: M4 Relief Road

Customer reference	CH 37	CH 37	CH 38	CH 38
Depth (m)	0.00	1.20	0.10	1.30
Date logged	11/02/00	11/02/00	11/02/00	11/02/00
TES Bretby ID Number	CL/0003509	CL/0003510	CL/0003511	CL/0003512

UKAS accredited	Test No.	CL/0003509	CL/0003510	CL/0003511	CL/0003512
Chromium (total)	ICPSSS11	18	37	221	37
CN- (total)	ICTSCN28	<1	<1	<1	<1
Copper	ICPSSS11	35	12	52	11
Cyanide (Free)	BGCN22	<1	<1	<1	<1
Lead	ICPSSS11	75	42	368	44
Nickel	ICPSSS11	14	32	29	34
Oil FTIR	FTIRSW	185	74	1120	77
PAH (screening)	PAHSCUV	23	953	<10	<10
PCB - ARO	PCBAROEC	117	41	117	2.1
pH units	WSLM3	9.7	8.5	9.8	8.6
Phenol Index	WSLM4	<0.5	<0.5	0.8	<0.5
SO4-- (acid sol)	ICPACS58	1450	304	1920	327
Sulphide	ICTSCN28	2	<1	25	<1
SVOC	SVOCSW	*	*	*	*
Thiocyanate	WSLM5	<2	<2	2	<2
TPH FTIR (AR)	FTIRSW	258	1870	843	31
Zinc	ICPSSS11	206	98	499	91

not UKAS accredited		CL/0003509	CL/0003510	CL/0003511	CL/0003512
Arsenic (Ms)		3.80	11.00	9.80	13.00
Boron.		<0.5	2.4	<0.5	2.5
Cadmium (Ms)		1.57	0.11	0.83	0.08
Mercury (Ms)		2.34	0.14	5.96	0.16
Selenium (MS)		1.39	<0.50	0.94	<0.50

Results expressed as mg/kg Air Dried unless stated otherwise

SO4 Analysis not conducted in accordance with BS1377

AR denotes analysis conducted on As Received Sample

* denotes results included on attached sheet

Date of Issue: 24/02/00

 Tests marked 'not UKAS accredited' in this report are not included
 in the UKAS Accreditation Schedule for our laboratory.

TES Bretby accepts no responsibility for the sampling related to the above results

 TES Bretby, P.O. Box 100, Burton-on-Trent, DE15 0XD Telephone: 01283 554400 Fax: 01283 554422
 TES Bretby is a division of Environmental Services Group Limited Registered in England Number 2880501

TES Bretby	
Report	000484
Table	2
Sheet	4/ 6

TEST REPORT

SOIL SAMPLE ANALYSIS

Exploration Associates
Unit 15
Crosby Yard
Bridgend
Mid Glamorgan
CF31 1JZ

1252
1411

TES Report No. 000484

Site: M4 Relief Road

Customer reference	CH 39 0.00	CH 39 1.00	CH 40 0.00	CH 40 0.50
Depth (m)				
Date logged	11/02/00	11/02/00	11/02/00	11/02/00
TES Bretby ID Number	CL/0003513	CL/0003514	CL/0003515	CL/0003516

UKAS accredited	Test No.	CL/0003513	CL/0003514	CL/0003515	CL/0003516
Chromium (total)	ICPSSS11	142	184	10	74
CN- (total)	ICTSCN28	<1	<1	<1	<1
Copper	ICPSSS11	40	34	5	20
Cyanide (Free)	BGCN22	<1	<1	<1	<1
Lead	ICPSSS11	497	98	31	53
Nickel	ICPSSS11	34	46	14	24
Oil FTIR	FTIRSW	285	2900	84	894
PAH (screening)	PAHSCUV	126	610	10	86
PCB - ARO	PCBAROEC	94	165	3.7	96
pH units	WSLM3	10.3	10.4	9.4	11.2
Phenol Index	WSLM4	1.3	1.6	<0.5	0.8
SO4-- (acid sol)	ICPACSS58	1600	2020	486	2440
Sulphide	ICTSCN28	363	122	<1	4
SVOC	SVOCSW	*	*	*	*
Thiocyanate	WSLM5	2	5	<2	<2
TPH FTIR (AR)	FTIRSW	245	609	48	524
Zinc	ICPSSS11	243	151	63	134

not UKAS accredited		CL/0003513	CL/0003514	CL/0003515	CL/0003516
Arsenic (Ms)		12.00	12.00	4.60	8.30
Boron.		<0.5	0.5	<0.5	0.7
Cadmium (Ms)		0.54	0.71	0.22	0.41
Mercury (Ms)		4.22	7.67	0.23	9.82
Selenium (MS)		1.49	1.38	1.28	0.89

Results expressed as mg/kg Air Dried unless stated otherwise

SO4 Analysis not conducted in accordance with BS1377

AR denotes analysis conducted on As Received Sample

* denotes results included on attached sheet

Date of Issue: 24/02/00

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TES Bretby is a division of Environmental Services Group Limited Registered in England Number 2880501

TES Bretby
Report 000484
Table 2
Sheet 5/ 6

TES**Bretby**

TEST REPORT

SOIL SAMPLE ANALYSIS



1252

1411

Exploration Associates
Unit 15
Crosby Yard
Bridgend
Mid Glamorgan
CF31 1JZ

TES Report No. 000484

Site: M4 Relief Road

Customer reference	CH 41	CH 41	CH 42	CH 42
Depth (m)	0.05	1.40	0.20	1.20
Date logged	11/02/00	11/02/00	11/02/00	11/02/00
TES Bretby ID Number	CL/0003517	CL/0003518	CL/0003519	CL/0003520

UKAS accredited	Test No.	CL/0003517	CL/0003518	CL/0003519	CL/0003520
Chromium (total)	ICPSSS11	120	272	37	43
CN- (total)	ICTSCN28	<1	<1	<1	<1
Copper	ICPSSS11	31	64	25	26
Cyanide (Free)	BGCN22	<1	<1	<1	<1
Lead	ICPSSS11	85	61	54	64
Nickel	ICPSSS11	28	32	26	33
Oil FTIR	FTIRSW	566	866	548	656
PAH (screening)	PAHSCUV	62	146	1470	1970
PCB - ARO	PCBAROEC	103	80	1010	1430
pH units	WSLM3	10.6	10.3	9.4	8.9
Phenol Index	WSLM4	<0.5	0.5	0.8	1.1
SO4-- (acid sol)	ICPACS58	1360	1760	1050	987
Sulphide	ICTSCN28	803	23	24	10
SVOC	SVOCSW	*	*	*	*
Thiocyanate	WSLM5	2	<2	<2	<2
TPH FTIR (AR)	FTIRSW	141	2400	127	244
VOC	VOCSW		*	*	
Zinc	ICPSSS11	386	488	128	133

not UKAS accredited		CL/0003517	CL/0003518	CL/0003519	CL/0003520
Arsenic (Ms)		9.80	8.10	7.40	7.70
Boron.		<0.5	<0.5	0.6	<0.5
Cadmium (Ms)		0.35	0.75	1.15	1.70
Mercury (Ms)		6.83	12.50	8.88	14.90
Selenium (MS)		1.37	0.91	1.08	1.03

Results expressed as mg/kg Air Dried unless stated otherwise

SO4 Analysis not conducted in accordance with BS1377

AR denotes analysis conducted on As Received Sample

* denotes results included on attached sheet

Date of Issue: 24/02/00

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TES Bretby is a division of Environmental Services Group Limited Registered in England Number 2880501

TES Bretby
Report 000484
Table 2
Sheet 6/ 6

Semi-Volatile Organic Compounds

Customer and Site Details:

Sample Details: CH/0034/ / 0.15- m
LIMS ID Number: CL03502
Job Number: 484

Exploration Associates: M4 Relief Road

Date Booked in: 11-Feb-00
Date Extracted: 16-Feb-00
Date Analysed: 17-Feb-00

Matrix:

Soil
Soxhlet
SW
Directory/Quant File: 0217ABN.MS5 0217CC2.D GPC (Y/N)

QC Batch Number: 219

Multiplier: 0.0667

1

N

1252

Target Compounds	CAS #	R.T. (min)	Concentration mg/kg	% Fit
Phenol	108-95-2	-	< 2.0	-
bis(2-Chloroethyl)ether	111-44-4	-	< 0.5	-
2-Chlorophenol	95-57-8	-	< 2.0	-
1,3-Dichlorobenzene	541-73-1	-	< 0.5	-
1,4-Dichlorobenzene	106-46-7	-	< 0.5	-
Benzyl alcohol	100-51-6	-	< 0.5	-
1,2-Dichlorobenzene	95-50-1	-	< 0.5	-
2-Methylphenol	95-48-7	-	< 0.5	-
bis(2-Chloroisopropyl)ether	108-60-1	-	< 0.5	-
Hexachloroethane	67-72-1	-	< 0.5	-
N-Nitroso-di-n-propylamine	621-64-7	-	< 0.5	-
3- & 4-Methylphenol	108-39-4/106-44-5	-	< 2.0	-
Nitrobenzene	98-95-3	-	< 0.5	-
Isophorone	78-59-1	-	< 0.5	-
2-Nitrophenol	88-75-5	-	< 2.0	-
2,4-Dimethylphenol	105-67-9	-	< 2.0	-
Benzoic Acid	65-85-0	-	< 10.0	-
bis(2-Chloroethoxy)methane	111-91-1	-	< 0.5	-
2,4-Dichlorophenol	120-83-2	-	< 2.0	-
1,2,4-Trichlorobenzene	120-82-1	-	< 0.5	-
Naphthalene	91-20-3	-	< 0.2	-
4-Chlorophenol *	106-48-9	-	< 2.0	-
4-Chloroaniline	106-47-8	-	< 0.5	-
Hexachlorobutadiene	87-68-3	-	< 0.5	-
4-Chloro-3-methylphenol	59-50-7	-	< 0.5	-
2-Methylnaphthalene	91-57-6	-	< 0.2	-
1-Methyl naphthalene *	90-12-0	-	< 0.2	-
Hexachlorocyclopentadiene	77-47-4	-	< 0.5	-
2,4,6-Trichlorophenol	88-06-2	-	< 2.0	-
2,4,5-Trichlorophenol	95-95-4	-	< 2.0	-
2-Chloronaphthalene	91-58-7	-	< 0.2	-
Biphenyl *	92-52-4	-	< 0.2	-
Diphenyl ether *	101-84-8	-	< 0.2	-
2-Nitroaniline	88-74-4	-	< 0.5	-
Acenaphthylene	208-96-8	-	< 0.2	-
Dimethylphthalate	131-11-3	-	< 0.5	-
2,6-Dinitrotoluene	606-20-2	-	< 0.5	-
Acenaphthene	83-32-9	-	< 0.2	-
3-Nitroaniline	99-09-2	-	< 0.5	-

* These compounds are not currently UKAS accredited.
Concentrations are reported on a wet weight basis.

Target Compounds	CAS #	R.T.	Concentration mg/kg	% Fit
2,4-Dinitrophenol	51-28-5	-	< 1.0	-
Dibenzofuran	132-64-9	-	< 0.5	-
4-Nitrophenol	100-02-7	-	< 5.0	-
2,4-Dinitrotoluene	121-14-2	-	< 0.5	-
Fluorene	86-73-7	-	< 0.2	-
Diethylphthalate	84-66-2	-	< 0.5	-
4-Chlorophenyl-phenylether	7005-72-4	-	< 0.5	-
4,6-Dinitro-2-methylphenol	534-52-1	-	< 5.0	-
4-Nitroaniline	100-01-6	-	< 0.5	-
N-Nitrosodiphenylamine	86-30-6	-	< 0.5	-
4-Bromophenyl-phenylether	101-55-3	-	< 0.5	-
Hexachlorobenzene	118-74-1	-	< 0.5	-
Pentachlorophenol	87-86-5	-	< 5.0	-
Phenanthrene	85-01-8	-	< 0.2	-
Anthracene	120-12-7	-	< 0.2	-
Di-n-butylphthalate	84-74-2	-	< 0.5	-
Fluoranthene	206-44-0	-	< 0.2	-
Pyrene	129-00-0	-	< 0.2	-
Butylbenzylphthalate	85-68-7	-	< 0.5	-
Benzofuran	56-55-3	-	< 0.2	-
Chrysene	218-01-9	-	< 0.2	-
3,3'-Dichlorobenzidine	91-94-1	-	< 2.0	-
bis(2-Ethylhexyl)phthalate	117-81-7	29.56	2.2	98
Di-n-octylphthalate	117-84-0	-	< 0.2	-
Benzofuran	205-99-2	-	< 0.2	-
Benzofuran	207-08-9	-	< 0.2	-
Benzofuran	50-32-8	-	< 0.2	-
Indeno[1,2,3-cd]pyrene	193-39-5	-	< 0.2	-
Dibenzofuran	53-70-3	-	< 0.2	-
Benzofuran	191-24-2	-	< 0.2	-

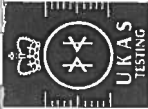
* compounds not currently UKAS accredited

Internal Standards	% Area
1,4-Dichlorobenzene-d4	113
Naphthalene-d8	104
Acenaphthene-d10	111
Phenanthrene-d10	105
Chrysene-d12	103
Perylene-d12	100

Surrogates	% Rec
2-Fluorophenol	80
Phenol-d5	85
Nitrobenzene-d5	83
2-Fluorobiphenyl	84
2,4,6-Tribromophenol	103
Terphenyl-d14	101



Semi-Volatile Organic Compounds



Customer and Site Details:

Sample Details:
LIMS ID Number:
Job Number:

Exploration Associates: M4 Relief Road

Date Booked in: 11-Feb-00
Date Extracted: 16-Feb-00
Date Analysed: 17-Feb-00

Matrix:

Soil
Soxhlet
SW
Directory/Quant File:

QC Batch Number:

Multiplier:

Dilution Factor:

20217ABN.MS2 0217CC03.D GPC (Y/N)

1252

219

0.0667

1

N

Target Compounds	CAS #	R.T. (min)	Concentration mg/kg	% Fit
Phenol	108-95-2	-	< 2.0	-
bis(2-Chloroethyl)ether	111-44-4	-	< 0.5	-
2-Chlorophenol	95-57-8	-	< 2.0	-
1,3-Dichlorobenzene	541-73-1	-	< 0.5	-
1,4-Dichlorobenzene	106-46-7	-	< 0.5	-
Benzyl alcohol	100-51-6	-	< 0.5	-
1,2-Dichlorobenzene	95-50-1	-	< 0.5	-
2-Methylphenol	95-48-7	-	< 0.5	-
bis(2-Chloroisopropyl)ether	108-60-1	-	< 0.5	-
Hexachloroethane	67-72-1	-	< 0.5	-
N-Nitroso-di-n-propylamine	621-64-7	-	< 0.5	-
3- & 4-Methylphenol	108-39-4/108-44-5	-	< 2.0	-
Nitrobenzene	98-95-3	-	< 0.5	-
Isophorone	78-59-1	-	< 0.5	-
2-Nitrophenol	88-75-5	-	< 2.0	-
2,4-Dimethylphenol	105-67-9	-	< 2.0	-
Benzoic Acid	65-85-0	-	< 10.0	-
bis(2-Chloroethoxy)methane	111-91-1	-	< 0.5	-
2,4-Dichlorophenol	120-83-2	-	< 2.0	-
1,2,4-Trichlorobenzene	120-82-1	-	< 0.5	-
Naphthalene	91-20-3	-	< 0.2	-
4-Chlorophenol *	106-48-9	-	< 2.0	-
4-Chloroaniline	106-47-8	-	< 0.5	-
Hexachlorobutadiene	87-68-3	-	< 0.5	-
4-Chloro-3-methylphenol	59-50-7	-	< 0.5	-
2-Methylnaphthalene	91-57-6	-	< 0.2	-
1-Methylnaphthalene *	90-12-0	-	< 0.2	-
Hexachlorocyclopentadiene	77-47-4	-	< 0.5	-
2,4,6-Trichlorophenol	88-06-2	-	< 2.0	-
2,4,5-Trichlorophenol	95-95-4	-	< 2.0	-
2-Chloronaphthalene	91-58-7	-	< 0.2	-
Biphenyl *	92-52-4	-	< 0.2	-
Diphenyl ether *	101-84-8	-	< 0.2	-
2-Nitroaniline	88-74-4	-	< 0.5	-
Acenaphthylene	208-96-8	-	< 0.2	-
Dimethylphthalate	131-11-3	-	< 0.5	-
2,6-Dinitrotoluene	606-20-2	-	< 0.5	-
Acenaphthene	83-32-9	-	< 0.2	-
3-Nitroaniline *	99-09-2	-	< 0.5	-

* These compounds are not currently UKAS accredited.

Concentrations are reported on a wet weight basis.

Target Compounds	CAS #	R.T.	Concentration mg/kg	% Fit
2,4-Dinitrophenol	51-28-5	-	< 1.0	-
Dibenzofuran	132-64-9	-	< 0.5	-
4-Nitrophenol	100-02-7	-	< 5.0	-
2,4-Dinitrotoluene	121-14-2	-	< 0.5	-
Fluorene	86-73-7	-	< 0.2	-
Diethylphthalate	84-66-2	-	< 0.5	-
4-Chlorophenyl-phenylether	7005-72-4	-	< 0.5	-
4,6-Dinitro-2-methylphenol	534-52-1	-	< 5.0	-
4-Nitroaniline	100-01-6	-	< 0.5	-
N-Nitrosodiphenylamine	86-30-6	-	< 0.5	-
4-Bromophenyl-phenylether	101-55-3	-	< 0.5	-
Hexachlorobenzene	118-74-1	-	< 0.5	-
Pentachlorophenol	87-86-5	-	< 5.0	-
Phenanthrene	85-01-8	-	< 0.2	-
Anthracene	120-12-7	-	< 0.2	-
Di-n-butylphthalate	84-74-2	-	< 0.5	-
Fluoranthene	206-44-0	-	< 0.2	-
Pyrene	129-00-0	-	< 0.2	-
Butylbenzylphthalate	85-68-7	-	< 0.5	-
Benzofuranthracene	56-55-3	-	< 0.2	-
Chrysene	218-01-9	-	< 0.2	-
3,3'-Dichlorobenzidine	91-94-1	-	< 2.0	-
bis(2-Ethylhexyl)phthalate	117-81-7	30.86	1.7	94
Di-n-octylphthalate	117-84-0	-	< 0.2	-
Benzobifluoranthene	205-99-2	-	< 0.2	-
Benzokifluoranthene	207-08-9	-	< 0.2	-
Benzofluoranthene	50-32-8	-	< 0.2	-
Indeno[1,2,3-cd]pyrene	193-39-5	-	< 0.2	-
Dibenzofluoranthene	53-70-3	-	< 0.2	-
Benzofluoranthene	191-24-2	-	< 0.2	-

* compounds not currently UKAS accredited

Internal Standards	% Area
1,4-Dichlorobenzene-d4	97
Naphthalene-d8	82
Acenaphthene-d10	99
Phenanthrene-d10	88
Chrysene-d12	90
Perylene-d12	92

Surrogates	% Rec
2-Fluorophenol	74
Phenol-d5	70
Nitrobenzene-d5	72
2-Fluorobiphenyl	79
2,4,6-Tribromophenol	97
Terphenyl-d14	97

Semi-Volatile Organic Compounds

Customer and Site Details:
Sample Details:
LIMS ID Number:
Job Number:

Exploration Associates : M4 Relief Road
CH/0035/0.00m
CL03504
484

Date Booked in: 11-Feb-00
Date Extracted: 16-Feb-00
Date Analysed: 18-Feb-00

Matrix: Soil
Ext Method: Soxhlet
Operator: cs
Directory/Quant File: 0218ABN MS2 0218CCC1.D GPC (Y/N)

QC Batch Number: 219
Multiplier: 0.133
Dilution Factor: 2
Y

Target Compounds	CAS #	R.T. (min)	Concentration mg/kg	% Fit
Phenol	108-95-2	-	< 4.0	-
bis(2-Chloroethyl)ether	111-44-4	-	< 1.0	-
2-Chlorophenol	95-57-8	-	< 4.0	-
1,3-Dichlorobenzene	541-73-1	-	< 1.0	-
1,4-Dichlorobenzene	106-46-7	-	< 1.0	-
Benzyl alcohol	100-51-6	-	< 1.0	-
1,2-Dichlorobenzene	95-50-1	-	< 1.0	-
2-Methylphenol	95-48-7	-	< 1.0	-
bis(2-Chloroisopropyl)ether	108-60-1	-	< 1.0	-
Hexachloroethane	67-72-1	-	< 1.0	-
N-Nitroso-di-n-propylamine	621-84-7	-	< 1.0	-
3- & 4-Methylphenol	108-39-4/106-44-5	-	< 4.0	-
Nitrobenzene	98-95-3	-	< 1.0	-
Isophorone	78-59-1	-	< 1.0	-
2-Nitrophenol	88-75-5	-	< 4.0	-
2,4-Dimethylphenol	105-67-9	-	< 4.0	-
Benzoic Acid	65-85-0	-	< 20.0	-
bis(2-Chloroethoxy)methane	111-91-1	-	< 1.0	-
2,4-Dichlorophenol	120-83-2	-	< 4.0	-
1,2,4-Trichlorobenzene	120-82-1	-	< 1.0	-
Naphthalene	91-20-3	-	< 0.4	-
4-Chlorophenol *	106-48-9	-	< 4.0	-
4-Chloroaniline	106-47-8	-	< 1.0	-
Hexachlorobutadiene	87-68-3	-	< 1.0	-
4-Chloro-3-methylphenol	59-50-7	-	< 1.0	-
2-Methylnaphthalene	91-57-6	-	< 0.4	-
1-Methylnaphthalene *	90-12-0	-	< 0.4	-
Hexachlorocyclopentadiene	77-47-4	-	< 1.0	-
2,4,6-Trichlorophenol	88-06-2	-	< 4.0	-
2,4,5-Trichlorophenol	95-95-4	-	< 4.0	-
2-Chloronaphthalene	91-58-7	-	< 0.4	-
Biphenyl *	92-52-4	16.40	0.5	96
Diphenyl ether *	101-84-8	-	< 0.4	-
2-Nitroaniline	88-74-4	-	< 1.0	-
Acenaphthylene	208-96-8	-	< 0.4	-
Dimethylphthalate	131-11-3	-	< 1.0	-
2,6-Dinitrotoluene	606-20-2	-	< 1.0	-
Acenaphthene	83-32-9	-	< 0.4	-
3-Nitroaniline *	99-09-2	-	< 1.0	-

* These compounds are not currently UKAS accredited.
Concentrations are reported on a wet weight basis.

Target Compounds	CAS #	R.T.	Concentration mg/kg	% Fit
2,4-Dinitrophenol	51-28-5	-	< 2.0	-
Dibenzofuran	132-64-9	-	< 1.0	-
4-Nitrophenol	100-02-7	-	< 10.0	-
2,4-Dinitrotoluene	121-14-2	-	< 1.0	-
Fluorene	86-73-7	-	< 0.4	-
Diethylphthalate	84-66-2	-	< 1.0	-
4-Chlorophenyl-phenylether	7005-72-4	-	< 1.0	-
4,6-Dinitro-2-methylphenol	534-52-1	-	< 10.0	-
4-Nitroaniline	100-01-6	-	< 1.0	-
N-Nitrosodiphenylamine	86-30-6	-	< 1.0	-
4-Bromophenyl-phenylether	101-55-3	-	< 1.0	-
Hexachlorobenzene	118-74-1	-	< 1.0	-
Pentachlorophenol	87-86-5	-	< 10.0	-
Phenanthrene	85-01-8	-	< 0.4	-
Anthracene	120-12-7	-	< 0.4	-
Di-n-butylphthalate	84-74-2	-	< 1.0	-
Fluoranthene	206-44-0	-	< 0.4	-
Pyrene	129-00-0	-	< 0.4	-
Butylbenzylphthalate	85-68-7	-	< 1.0	-
Benzol[a]anthracene	56-55-3	-	< 0.4	-
Chrysene	218-01-9	30.44	0.7	99
3,3'-Dichlorobenzidine	91-94-1	-	< 4.0	-
bis(2-Ethylhexyl)phthalate	117-81-7	30.83	1.9	93
Di-n-octylphthalate	117-84-0	-	< 0.4	-
Benzol[b]fluoranthene	205-99-2	-	< 0.4	-
Benzol[k]fluoranthene	207-08-9	-	< 0.4	-
Benzol[a]pyrene	50-32-8	-	< 0.4	-
Indeno[1,2,3-cd]pyrene	193-39-5	-	< 0.4	-
Dibenzo[a,h]anthracene	53-70-3	-	< 0.4	-
Benzol[g,h,i]perylene	191-24-2	-	< 0.4	-

* compounds not currently UKAS accredited

Internal Standards	% Area
1,4-Dichlorobenzene-d4	122
Naphthalene-d8	117
Acenaphthene-d10	130
Phenanthrene-d10	124
Chrysene-d12	132
Perylene-d12	130

Surrogates	% Rec
2-Fluorophenol	47
Phenol-d5	52
Nitrobenzene-d5	51
2-Fluorobiphenyl	55
2,4,6-Tribromophenol	78
Terphenyl-d14	77

Semi-Volatile Organic Compounds

Customer and Site Details:
Sample Details:
LIMS ID Number:
Job Number:

Exploration Associates: M4 Relief Road
CH/0035/ / 0.50- m
CL03505
484

11-Feb-00
16-Feb-00
22-Feb-00

Matrix:
Ext Method:
Operator:
Directory/Quant File:

Soil
Soxhlet
SW
0222ABN.MS5 0222CCC1.D GPC (Y/N)

QC Batch Number:
Multiplier:
Dilution Factor:

Target Compounds	CAS #	R.T. (min)	Concentration mg/kg	% Fit
Phenol	108-95-2	-	< 2.0	-
bis(2-Chloroethyl)ether	111-44-4	-	< 0.5	-
2-Chlorophenol	95-57-8	-	< 2.0	-
1,3-Dichlorobenzene	541-73-1	-	< 0.5	-
1,4-Dichlorobenzene	106-46-7	8.29	3.9	99
Benzyl alcohol	100-51-6	-	< 0.5	-
1,2-Dichlorobenzene	95-50-1	-	< 0.5	-
2-Methylphenol	95-48-7	-	< 0.5	-
bis(2-Chloroisopropyl)ether	108-60-1	-	< 0.5	-
Hexachloroethane	67-72-1	-	< 0.5	-
N-Nitroso-di-n-propylamine	621-64-7	-	< 0.5	-
3- & 4-Methylphenol	108-39-4/106-44-5	-	< 2.0	-
Nitrobenzene	98-95-3	-	< 0.5	-
Isophorone	78-59-1	-	< 0.5	-
2-Nitrophenol	88-75-5	-	< 2.0	-
2,4-Dimethylphenol	105-67-9	-	< 2.0	-
Benzoic Acid	65-85-0	-	< 10.0	-
bis(2-Chloroethoxy)methane	111-91-1	-	< 0.5	-
2,4-Dichlorophenol	120-83-2	-	< 2.0	-
1,2,4-Trichlorobenzene	120-82-1	-	< 0.5	-
Naphthalene	91-20-3	-	< 0.2	-
4-Chlorophenol *	106-48-9	11.90	6.0	90
4-Chloroaniline	106-47-8	12.00	19.2	100
Hexachlorobutadiene	87-68-3	-	< 0.5	-
4-Chloro-3-methylphenol	59-50-7	13.51	1.9	76
2-Methylnaphthalene	91-57-6	-	< 0.2	-
1-Methyl naphthalene *	90-12-0	-	< 0.2	-
Hexachlorocyclopentadiene	77-47-4	-	< 0.5	-
2,4,6-Trichlorophenol	88-06-2	-	< 2.0	-
2,4,5-Trichlorophenol	95-95-4	-	< 2.0	-
2-Chloronaphthalene	91-58-7	-	< 0.2	-
Biphenyl *	92-52-4	-	< 0.2	-
Diphenyl ether *	101-84-8	-	< 0.2	-
2-Nitroaniline	88-74-4	-	< 0.5	-
Acenaphthylene	208-96-8	-	< 0.2	-
Dimethylphthalate	131-11-3	-	< 0.5	-
2,6-Dinitrotoluene	606-20-2	-	< 0.5	-
Acenaphthene	83-32-9	-	< 0.2	-
3-Nitroaniline	99-09-2	-	< 0.5	-

* These compounds are not currently UKAS accredited.
Concentrations are reported on a wet weight basis.

Target Compounds	CAS #	R.T.	Concentration mg/kg	% Fit
2,4-Dinitrophenol	51-28-5	-	< 1.0	-
Dibenzofuran	132-64-9	-	< 0.5	-
4-Nitrophenol	100-02-7	-	< 5.0	-
2,4-Dinitrotoluene	121-14-2	-	< 0.5	-
Fluorene	86-73-7	-	< 0.2	-
Diethylphthalate	84-66-2	-	< 0.5	-
4-Chlorophenyl-phenylether	7005-72-4	-	< 0.5	-
4,6-Dinitro-2-methylphenol	534-52-1	-	< 5.0	-
4-Nitroaniline	100-01-6	-	< 0.5	-
N-Nitrosodiphenylamine	86-30-6	-	< 0.5	-
4-Bromophenyl-phenylether	101-55-3	-	< 0.5	-
Hexachlorobenzene	118-74-1	-	< 0.5	-
Pentachlorophenol	87-86-5	-	< 5.0	-
Phenanthrene	85-01-8	-	< 0.2	-
Anthracene	120-12-7	-	< 0.2	-
Di-n-butylphthalate	84-74-2	-	< 0.5	-
Fluoranthene	206-44-0	-	< 0.2	-
Pyrene	129-00-0	-	< 0.2	-
Butylbenzylphthalate	85-68-7	-	< 0.5	-
Benzo[a]anthracene	56-55-3	-	< 0.2	-
Chrysene	218-01-9	-	< 0.2	-
3,3'-Dichlorobenzidine	91-94-1	-	< 2.0	-
bis(2-Ethylhexyl)phthalate	117-81-7	29.51	2.6	98
Di-n-octylphthalate	117-84-0	-	< 0.2	-
Benzo[b]fluoranthene	205-99-2	-	< 0.2	-
Benzo[k]fluoranthene	207-08-9	-	< 0.2	-
Benzo[a]pyrene	50-32-8	-	< 0.2	-
Indeno[1,2,3-cd]pyrene	193-39-5	-	< 0.2	-
Dibenzo[a,h]anthracene	53-70-3	-	< 0.2	-
Benzo[g,h,i]perylene	191-24-2	-	< 0.2	-

* compounds not currently UKAS accredited

Internal Standards	% Area
1,4-Dichlorobenzene-d4	102
Naphthalene-d8	100
Acenaphthene-d10	110
Phenanthrene-d10	111
Chrysene-d12	124
Perylene-d12	137

Surrogates	% Rec
2-Fluorophenol	91
Phenol-d5	103
Nitrobenzene-d5	90
2-Fluorobiphenyl	84
2,4,6-Tribromophenol	111
Terphenyl-d14	93

Semi-Volatile Organic Compounds

Customer and Site Details:
Sample Details:
LIMS ID Number:
Job Number:

Exploration Associates: M4 Relief Road
CH/0035/ / 0.80- m
Date Booked in:
Date Extracted:
Date Analysed:

11-Feb-00
16-Feb-00
17-Feb-00

Matrix:
Ext Method:
Operator:
Directory/Quant File:

Soil
Soxhlet
SW
0217ABN.MS2

QC Batch Number:
Multiplier:
Dilution Factor:

219
0.0667
1
N

Target Compounds	CAS #	R.T. (min)	Concentration mg/kg	% Fit
Phenol	108-95-2	-	< 2.0	-
bis(2-Chloroethyl)ether	111-44-4	-	< 0.5	-
2-Chlorophenol	95-57-8	-	< 2.0	-
1,3-Dichlorobenzene	541-73-1	-	< 0.5	-
1,4-Dichlorobenzene	106-46-7	-	< 0.5	-
Benzyl alcohol	100-51-6	-	< 0.5	-
1,2-Dichlorobenzene	95-50-1	-	< 0.5	-
2-Methylphenol	95-48-7	-	< 0.5	-
bis(2-Chloroisopropyl)ether	108-60-1	-	< 0.5	-
Hexachloroethane	67-72-1	-	< 0.5	-
N-Nitroso-di-n-propylamine	621-84-7	-	< 0.5	-
3- & 4-Methylphenol	108-39-4/106-44-5	-	< 2.0	-
Nitrobenzene	98-95-3	-	< 0.5	-
Isophorone	78-59-1	-	< 0.5	-
2-Nitrophenol	88-75-5	-	< 2.0	-
2,4-Dimethylphenol	105-67-9	-	< 2.0	-
Benzoic Acid	65-85-0	-	< 10.0	-
bis(2-Chloroethoxy)methane	111-91-1	-	< 0.5	-
2,4-Dichlorophenol	120-83-2	-	< 2.0	-
1,2,4-Trichlorobenzene	120-82-1	-	< 0.5	-
Naphthalene	91-20-3	-	< 0.2	-
4-Chlorophenol *	106-48-9	-	< 2.0	-
4-Chloroaniline	106-47-8	-	< 0.5	-
Hexachlorobutadiene	87-68-3	-	< 0.5	-
4-Chloro-3-methylphenol	59-50-7	-	< 0.5	-
2-Methylnaphthalene	91-57-6	-	< 0.2	-
1-Methylnaphthalene *	90-12-0	-	< 0.2	-
Hexachlorocyclopentadiene	77-47-4	-	< 0.5	-
2,4,6-Trichlorophenol	88-06-2	-	< 2.0	-
2,4,5-Trichlorophenol	95-95-4	-	< 2.0	-
2-Chloronaphthalene	91-58-7	-	< 0.2	-
Biphenyl *	92-52-4	-	< 0.2	-
Diphenyl ether *	101-84-8	-	< 0.2	-
2-Nitroaniline	88-74-4	-	< 0.5	-
Acenaphthylene	208-96-8	-	< 0.2	-
Dimethylphthalate	131-11-3	-	< 0.5	-
2,6-Dinitrotoluene	606-20-2	-	< 0.5	-
Acenaphthene	83-32-9	-	< 0.2	-
3-Nitroaniline *	99-09-2	-	< 0.5	-

* These compounds are not currently UKAS accredited.
Concentrations are reported on a wet weight basis.

Target Compounds	CAS #	R.T.	Concentration mg/kg	% Fit
2,4-Dinitrophenol	51-28-5	-	< 1.0	-
Dibenzofuran	132-64-9	-	< 0.5	-
4-Nitrophenol	100-02-7	-	< 5.0	-
2,4-Dinitrotoluene	121-14-2	-	< 0.5	-
Fluorene	86-73-7	-	< 0.2	-
Diethylphthalate	84-66-2	-	< 0.5	-
4-Chlorophenyl-phenylether	7005-72-4	-	< 0.5	-
4,6-Dinitro-2-methylphenol	534-52-1	-	< 5.0	-
4-Nitroaniline	100-01-6	-	< 0.5	-
N-Nitrosodiphenylamine	86-30-6	-	< 0.5	-
4-Bromophenyl-phenylether	101-55-3	-	< 0.5	-
Hexachlorobenzene	118-74-1	-	< 0.5	-
Pentachlorophenol	87-86-5	-	< 5.0	-
Phenanthrene	85-01-8	-	< 0.2	-
Anthracene	120-12-7	-	< 0.2	-
Di-n-butylphthalate	84-74-2	-	< 0.5	-
Fluoranthene	206-44-0	-	< 0.2	-
Pyrene	129-00-0	-	< 0.2	-
Butylbenzylphthalate	85-68-7	-	< 0.5	-
Benzo[a]anthracene	56-55-3	-	< 0.2	-
Chrysene	218-01-9	-	< 0.2	-
3,3'-Dichlorobenzidine	91-94-1	-	< 2.0	-
bis(2-Ethylhexyl)phthalate	117-81-7	30.86	0.8	93
Di-n-octylphthalate	117-84-0	-	< 0.2	-
Benzo[b]fluoranthene	205-99-2	-	< 0.2	-
Benzo[k]fluoranthene	207-08-9	-	< 0.2	-
Benzo[a]pyrene	50-32-8	-	< 0.2	-
Indeno[1,2,3-cd]pyrene	193-39-5	-	< 0.2	-
Dibenzof[a,h]anthracene	53-70-3	-	< 0.2	-
Benzo[g,h,i]perylene	191-24-2	-	< 0.2	-

* compounds not currently UKAS accredited

Internal Standards	% Area
1,4-Dichlorobenzene-d4	103
Naphthalene-d8	87
Acenaphthene-d10	104
Phenanthrene-d10	95
Chrysene-d12	99
Perylene-d12	96

Surrogates	% Rec
2-Fluorophenol	80
Phenol-d5	75
Nitrobenzene-d5	80
2-Fluorobiphenyl	78
2,4,6-Tribromophenol	97
Terphenyl-d14	91

Semi-Volatile Organic Compounds

Customer and Site Details:
Sample Details: CH/0036/0.10m
LIMS ID Number: CL03507
Job Number: 484

Exploration Associates : M4 Relief Road
Date Booked in: 11-Feb-00
Date Extracted: 16-Feb-00
Date Analysed: 18-Feb-00

Matrix: Soil
Ext Method: Soxhlet
Operator: CS
Directory/Quant File: 0218ABN.MS2 0218CCC1.D GPC (Y/N)

QC Batch Number: 219
Multiplier: 0.333
Dilution Factor: 5
Y

Target Compounds	CAS #	R.T. (min)	Concentration mg/kg	% Fit
Phenol	108-95-2	-	< 10.0	-
bis(2-Chloroethyl)ether	111-44-4	-	< 2.5	-
2-Chlorophenol	95-57-8	-	< 10.0	-
1,3-Dichlorobenzene	541-73-1	-	< 2.5	-
1,4-Dichlorobenzene	106-46-7	-	< 2.5	-
Benzyl alcohol	100-51-6	-	< 2.5	-
1,2-Dichlorobenzene	95-50-1	-	< 2.5	-
2-Methylphenol	95-48-7	-	< 2.5	-
bis(2-Chloroisopropyl)ether	108-60-1	-	< 2.5	-
Hexachloroethane	67-72-1	-	< 2.5	-
N-Nitroso-di-n-propylamine	821-64-7	-	< 2.5	-
3- & 4-Methylphenol	108-39-4/108-44-5	-	< 10.0	-
Nitrobenzene	98-95-3	-	< 2.5	-
Isophorone	78-59-1	-	< 2.5	-
2-Nitrophenol	88-75-5	-	< 10.0	-
2,4-Dimethylphenol	105-67-9	-	< 10.0	-
Benzoic Acid	65-85-0	-	< 50.0	-
bis(2-Chloroethoxy)methane	111-91-1	-	< 2.5	-
2,4-Dichlorophenol	120-83-2	-	< 10.0	-
1,2,4-Trichlorobenzene	120-82-1	-	< 2.5	-
Naphthalene	91-20-3	-	< 1.0	-
4-Chlorophenol *	106-48-9	-	< 10.0	-
4-Chloroaniline	106-47-8	-	< 2.5	-
Hexachlorobutadiene	87-68-3	-	< 2.5	-
4-Chloro-3-methylphenol	59-50-7	-	< 2.5	-
2-Methylnaphthalene	91-57-6	-	< 1.0	-
1-Methylnaphthalene *	90-12-0	-	< 1.0	-
Hexachlorocyclopentadiene	77-47-4	-	< 2.5	-
2,4,6-Trichlorophenol	88-06-2	-	< 10.0	-
2,4,5-Trichlorophenol	95-95-4	-	< 10.0	-
2-Chloronaphthalene	91-58-7	-	< 1.0	-
Biphenyl *	92-52-4	-	< 1.0	-
Diphenyl ether *	101-84-8	-	< 1.0	-
2-Nitroaniline	88-74-4	-	< 2.5	-
Acenaphthylene	208-96-8	-	< 1.0	-
Dimethylphthalate	131-11-3	-	< 2.5	-
2,6-Dinitrotoluene	606-20-2	-	< 2.5	-
Acenaphthene	83-32-9	-	< 1.0	-
3-Nitroaniline *	99-09-2	-	< 2.5	-

* These compounds are not currently UKAS accredited.
Concentrations are reported on a wet weight basis.

Target Compounds	CAS #	R.T.	Concentration mg/kg	% Fit
2,4-Dinitrophenol	51-28-5	-	< 5.0	-
Dibenzofuran	132-64-9	-	< 2.5	-
4-Nitrophenol	100-02-7	-	< 25.0	-
2,4-Dinitrotoluene	121-14-2	-	< 2.5	-
Fluorene	86-73-7	-	< 1.0	-
Diethylphthalate	84-66-2	-	< 2.5	-
4-Chlorophenyl-phenylether	7005-72-4	-	< 2.5	-
4,6-Dinitro-2-methylphenol	534-52-1	-	< 25.0	-
4-Nitroaniline	100-01-6	-	< 2.5	-
N-Nitrosodiphenylamine	86-30-6	-	< 2.5	-
4-Bromophenyl-phenylether	101-55-3	-	< 2.5	-
Hexachlorobenzene	118-74-1	-	< 2.5	-
Pentachlorophenol	87-86-5	-	< 25.0	-
Phenanthrene	85-01-8	-	< 1.0	-
Anthracene	120-12-7	-	< 1.0	-
Di-n-butylphthalate	84-74-2	-	< 2.5	-
Fluoranthene	206-44-0	-	< 1.0	-
Pyrene	129-00-0	-	< 1.0	-
Butylbenzylphthalate	85-68-7	-	< 2.5	-
Benzo[a]anthracene	56-55-3	-	< 1.0	-
Chrysene	218-01-9	-	< 1.0	-
3,3'-Dichlorobenzidine	91-94-1	-	< 10.0	-
bis(2-Ethylhexyl)phthalate	117-81-7	30.82	4.7	95
Di-n-octylphthalate	117-84-0	-	< 1.0	-
Benzo[b]fluoranthene	205-99-2	-	< 1.0	-
Benzo[k]fluoranthene	207-08-9	-	< 1.0	-
Benzo[a]pyrene	50-32-8	-	< 1.0	-
Indeno[1,2,3-cd]pyrene	193-39-5	-	< 1.0	-
Dibenzo[a,h]anthracene	53-70-3	-	< 1.0	-
Benzo[g,h,i]perylene	191-24-2	-	< 1.0	-

* compounds not currently UKAS accredited

Internal Standards	% Area
1,4-Dichlorobenzene-d4	110
Naphthalene-d8	102
Acenaphthene-d10	117
Phenanthrene-d10	115
Chrysene-d12	116
Perylene-d12	109

Surrogates	% Rec
2-Fluorophenol	55
Phenol-d5	59
Nitrobenzene-d5	56
2-Fluorobiphenyl	62
2,4,6-Tribromophenol	71
Terphenyl-d14	89

Semi-Volatile Organic Compounds

Customer and Site Details:
Sample Details:
LIMS ID Number:
Job Number:

Exploration Associates: M4 Relief Road
CH/0036/ / 0.70- m
CL03508
484

11-Feb-00
16-Feb-00
17-Feb-00

Matrix:
Ext Method:
Operator:
Directory/Quant File:

Soil
Soxhlet
sw

QC Batch Number:
Multiplier:
Dilution Factor:

1252

219
0.0667
1
N

0217ABN:MS5 0217CCC2.D GPC (Y/N)

Target Compounds	CAS #	R.T. (min)	Concentration mg/kg	% Fit
Phenol	108-95-2	-	< 2.0	-
bis(2-Chloroethyl)ether	111-44-4	-	< 0.5	-
2-Chlorophenol	95-57-8	-	< 2.0	-
1,3-Dichlorobenzene	541-73-1	-	< 0.5	-
1,4-Dichlorobenzene	106-46-7	-	< 0.5	-
Benzyl alcohol	100-51-6	-	< 0.5	-
1,2-Dichlorobenzene	95-50-1	-	< 0.5	-
2-Methylphenol	95-48-7	-	< 0.5	-
bis(2-Chloroisopropyl)ether	108-60-1	-	< 0.5	-
Hexachloroethane	67-72-1	-	< 0.5	-
N-Nitroso-di-n-propylamine	621-64-7	-	< 0.5	-
3- & 4-Methylphenol	108-39-4/106-44-5	-	< 2.0	-
Nitrobenzene	98-95-3	-	< 0.5	-
Isophorone	78-59-1	-	< 0.5	-
2-Nitrophenol	88-75-5	-	< 2.0	-
2,4-Dimethylphenol	105-67-9	-	< 2.0	-
Benzoic Acid	65-85-0	-	< 10.0	-
bis(2-Chloroethoxy)methane	111-91-1	-	< 0.5	-
2,4-Dichlorophenol	120-83-2	-	< 2.0	-
1,2,4-Trichlorobenzene	120-82-1	-	< 0.5	-
Naphthalene	91-20-3	-	< 0.2	-
4-Chlorophenol *	106-48-9	-	< 2.0	-
4-Chloroaniline	106-47-8	-	< 0.5	-
Hexachlorobutadiene	87-68-3	-	< 0.5	-
4-Chloro-3-methylphenol	59-50-7	-	< 0.5	-
2-Methylnaphthalene	91-57-6	-	< 0.2	-
1-Methyl naphthalene *	90-12-0	-	< 0.2	-
Hexachlorocyclopentadiene	77-47-4	-	< 0.5	-
2,4,6-Trichlorophenol	88-06-2	-	< 2.0	-
2,4,5-Trichlorophenol	95-95-4	-	< 2.0	-
2-Chloronaphthalene	91-58-7	-	< 0.2	-
Biphenyl *	92-52-4	-	< 0.2	-
Diphenyl ether *	101-84-8	-	< 0.2	-
2-Nitroaniline	88-74-4	-	< 0.5	-
Acenaphthylene	208-96-8	-	< 0.2	-
Dimethylphthalate	131-11-3	-	< 0.5	-
2,6-Dinitrotoluene	606-20-2	-	< 0.5	-
Acenaphthene	83-32-9	-	< 0.2	-
3-Nitroaniline	99-09-2	-	< 0.5	-

* These compounds are not currently UKAS accredited.
Concentrations are reported on a wet weight basis.

Target Compounds	CAS #	R.T.	Concentration mg/kg	% Fit
2,4-Dinitrophenol	51-28-5	-	< 1.0	-
Dibenzofuran	132-64-9	-	< 0.5	-
4-Nitrophenol	100-02-7	-	< 5.0	-
2,4-Dinitrotoluene	121-14-2	-	< 0.5	-
Fluorene	86-73-7	-	< 0.2	-
Diethylphthalate	84-66-2	-	< 0.5	-
4-Chlorophenyl-phenylether	7005-72-4	-	< 0.5	-
4,6-Dinitro-2-methylphenol	534-52-1	-	< 5.0	-
4-Nitroaniline	100-01-6	-	< 0.5	-
N-Nitrosodiphenylamine	86-30-6	-	< 0.5	-
4-Bromophenyl-phenylether	101-55-3	-	< 0.5	-
Hexachlorobenzene	118-74-1	-	< 0.5	-
Pentachlorophenol	87-86-5	-	< 5.0	-
Phenanthrene	85-01-8	-	< 0.2	-
Anthracene	120-12-7	-	< 0.2	-
Di-n-butylphthalate	84-74-2	-	< 0.5	-
Fluoranthene	206-44-0	-	< 0.2	-
Pyrene	129-00-0	-	< 0.2	-
Butylbenzylphthalate	85-68-7	-	< 0.5	-
Benzofuran	56-55-3	-	< 0.2	-
Chrysene	218-01-9	-	< 0.2	-
3,3'-Dichlorobenzidine	91-94-1	-	< 2.0	-
bis(2-Ethylhexyl)phthalate	117-81-7	29.55	2.6	97
Di-n-octylphthalate	117-84-0	-	< 0.2	-
Benzofuran	205-99-2	-	< 0.2	-
Benzofuran	207-08-9	-	< 0.2	-
Benzofuran	50-32-8	-	< 0.2	-
Indeno[1,2,3-cd]pyrene	193-39-5	-	< 0.2	-
Dibenzofuran	53-70-3	-	< 0.2	-
Benzofuran	191-24-2	-	< 0.2	-

* compounds not currently UKAS accredited

Internal Standards	% Area
1,4-Dichlorobenzene-d4	112
Naphthalene-d8	105
Acenaphthene-d10	111
Phenanthrene-d10	104
Chrysene-d12	105
Perylene-d12	104

Surrogates	% Rec
2-Fluorophenol	72
Phenol-d5	79
Nitrobenzene-d5	80
2-Fluorobiphenyl	78
2,4,6-Tribromophenol	88
Terphenyl-d14	96

Semi-Volatile Organic Compounds

Customer and Site Details:
Sample Details:
LIMS ID Number:
Job Number:

Exploration Associates : M4 Relief Road
CH/0037/0.00m
CL03509
484

11-Feb-00
16-Feb-00
18-Feb-00

Matrix:
Ext Method:
Operator:
Directory/Quant File:

Soil
Soxhlet
CS
0218ABN MS4 0218CCCC1.D GPC (Y/N)

QC Batch Number: 219

Multiplier: 0.0667

Dilution Factor: 1

Y

Target Compounds	CAS #	R.T. (min)	Concentration mg/kg	% Fit
Phenol	108-95-2	-	< 2.0	-
bis(2-Chloroethyl)ether	111-44-4	-	< 0.5	-
2-Chlorophenol	95-57-8	-	< 2.0	-
1,3-Dichlorobenzene	541-73-1	-	< 0.5	-
1,4-Dichlorobenzene	106-46-7	-	< 0.5	-
Benzyl alcohol	100-51-6	-	< 0.5	-
1,2-Dichlorobenzene	95-50-1	-	< 0.5	-
2-Methylphenol	95-48-7	-	< 0.5	-
bis(2-Chloroisopropyl)ether	108-60-1	-	< 0.5	-
Hexachloroethane	67-72-1	-	< 0.5	-
N-Nitroso-di-n-propylamine	621-64-7	-	< 0.5	-
3- & 4-Methylphenol	108-39-4/108-44-5	-	< 2.0	-
Nitrobenzene	98-95-3	-	< 0.5	-
Isophorone	78-59-1	-	< 0.5	-
2-Nitrophenol	88-75-5	-	< 2.0	-
2,4-Dimethylphenol	105-67-9	-	< 2.0	-
Benzoic Acid	65-85-0	-	< 10.0	-
bis(2-Chloroethoxy)methane	111-91-1	-	< 0.5	-
2,4-Dichlorophenol	120-83-2	-	< 2.0	-
1,2,4-Trichlorobenzene	120-82-1	-	< 0.5	-
Naphthalene	91-20-3	11.75	0.3	91
4-Chlorophenol *	106-48-9	-	< 2.0	-
4-Chloroaniline *	106-47-8	-	< 0.5	-
Hexachlorobutadiene	87-68-3	-	< 0.5	-
4-Chloro-3-methylphenol	59-50-7	-	< 0.5	-
2-Methylnaphthalene	91-57-6	-	< 0.2	-
1-Methyl naphthalene *	90-12-0	-	< 0.2	-
Hexachlorocyclopentadiene	77-47-4	-	< 0.5	-
2,4,6-Trichlorophenol	88-06-2	-	< 2.0	-
2,4,5-Trichlorophenol	95-95-4	-	< 2.0	-
2-Chloronaphthalene	91-58-7	-	< 0.2	-
Biphenyl *	92-52-4	15.17	0.5	99
Diphenyl ether *	101-84-8	-	< 0.2	-
2-Nitroaniline	88-74-4	-	< 0.5	-
Acenaphthylene	208-96-8	-	< 0.2	-
Dimethylphthalate	131-11-3	-	< 0.5	-
2,6-Dinitrotoluene	606-20-2	-	< 0.5	-
Acenaphthene	83-32-9	-	< 0.2	-
3-Nitroaniline *	99-09-2	-	< 0.5	-

* These compounds are not currently UKAS accredited.
Concentrations are reported on a wet weight basis.

Target Compounds	CAS #	R.T.	Concentration mg/kg	% Fit
2,4-Dinitrophenol	51-28-5	-	< 1.0	-
Dibenzofuran	132-64-9	-	< 0.5	-
4-Nitrophenol	100-02-7	-	< 5.0	-
2,4-Dinitrotoluene	121-14-2	-	< 0.5	-
Fluorene	86-73-7	-	< 0.2	-
Diethylphthalate	84-66-2	-	< 0.5	-
4-Chlorophenyl-phenylether	7005-72-4	-	< 0.5	-
4,6-Dinitro-2-methylphenol	534-52-1	-	< 5.0	-
4-Nitroaniline	100-01-6	18.74	0.7	63
N-Nitrosodiphenylamine	86-30-6	-	< 0.5	-
4-Bromophenyl-phenylether	101-55-3	-	< 0.5	-
Hexachlorobenzene	118-74-1	-	< 0.5	-
Pentachlorophenol	87-86-5	-	< 5.0	-
Phenanthrene	85-01-8	-	< 0.2	-
Anthracene	120-12-7	-	< 0.2	-
Di-n-butylphthalate	84-74-2	-	< 0.5	-
Fluoranthene	206-44-0	-	< 0.2	-
Pyrene	129-00-0	-	< 0.2	-
Butylbenzylphthalate	85-68-7	-	< 0.5	-
Benzol(a)anthracene	56-55-3	-	< 0.2	-
Chrysene	218-01-9	29.05	0.3	98
3,3'-Dichlorobenzidine	91-94-1	-	< 2.0	-
bis(2-Ethylhexyl)phthalate	117-81-7	29.69	1.7	99
Di-n-octylphthalate	117-84-0	-	< 0.2	-
Benzol(b)fluoranthene	205-99-2	-	< 0.2	-
Benzol(k)fluoranthene	207-08-9	-	< 0.2	-
Benzol(a)pyrene	50-32-8	-	< 0.2	-
Indeno(1,2,3-cd)pyrene	193-39-5	37.35	0.3	99
Dibenzof(a,h)anthracene	53-70-3	37.55	0.2	95
Benzol(g,h,i)perylene	191-24-2	38.30	0.4	99

* compounds not currently UKAS accredited

Internal Standards	% Area
1,4-Dichlorobenzene-d4	108
Naphthalene-d8	98
Acenaphthene-d10	100
Phenanthrene-d10	101
Chrysene-d12	110
Perylene-d12	105

Surrogates	% Rec
2-Fluorophenol	64
Phenol-d5	66
Nitrobenzene-d5	47
2-Fluorobiphenyl	50
2,4,6-Tribromophenol	63
Triphenyl-d14	71

Semi-Volatile Organic Compounds

Customer and Site Details:
Sample Details:
LIMS ID Number:
Job Number:

Exploration Associates: M4 Relief Road
CH/0037/ / 1.20- m
Date Booked in:
Date Extracted:
Date Analysed:

Matrix: Soil
Ext Method: Soxhlet
Operator: SW
Directory/Quant File: 0217ABN.MS2 0217CCC3.D GPC (Y/N)

QC Batch Number: 219
Multiplier: 0.0667
Dilution Factor: 1
N

Target Compounds	CAS #	R.T. (min)	Concentration mg/kg	% Fit
Phenol	108-95-2	-	< 2.0	-
bis(2-Chloroethyl)ether	111-44-4	-	< 0.5	-
2-Chlorophenol	95-57-8	-	< 2.0	-
1,3-Dichlorobenzene	541-73-1	-	< 0.5	-
1,4-Dichlorobenzene	106-46-7	-	< 0.5	-
Benzyl alcohol	100-51-6	-	< 0.5	-
1,2-Dichlorobenzene	95-50-1	-	< 0.5	-
2-Methylphenol	95-48-7	-	< 0.5	-
bis(2-Chloroisopropyl)ether	108-60-1	-	< 0.5	-
Hexachloroethane	67-72-1	-	< 0.5	-
N-Nitroso-di-n-propylamine	621-64-7	-	< 0.5	-
3- & 4-Methylphenol	108-39-4/108-44-5	-	< 2.0	-
Nitrobenzene	98-95-3	-	< 0.5	-
Isophorone	78-59-1	-	< 0.5	-
2-Nitrophenol	88-75-5	-	< 2.0	-
2,4-Dimethylphenol	105-67-9	-	< 2.0	-
Benzoic Acid	65-85-0	-	< 10.0	-
bis(2-Chloroethoxy)methane	111-91-1	-	< 0.5	-
2,4-Dichlorophenol	120-83-2	-	< 2.0	-
1,2,4-Trichlorobenzene	120-82-1	-	< 0.5	-
Naphthalene	91-20-3	-	< 0.2	-
4-Chlorophenol *	106-48-9	-	< 2.0	-
4-Chloroaniline	106-47-8	-	< 0.5	-
Hexachlorobutadiene	87-68-3	-	< 0.5	-
4-Chloro-3-methylphenol	59-50-7	-	< 0.5	-
2-Methylnaphthalene	91-57-6	-	< 0.2	-
1-Methylnaphthalene *	90-12-0	-	< 0.2	-
Hexachlorocyclopentadiene	77-47-4	-	< 0.5	-
2,4,6-Trichlorophenol	88-06-2	-	< 2.0	-
2,4,5-Trichlorophenol	95-95-4	-	< 2.0	-
2-Chloronaphthalene	91-58-7	-	< 0.2	-
Biphenyl *	92-52-4	-	< 0.2	-
Diphenyl ether *	101-84-8	-	< 0.2	-
2-Nitroaniline	88-74-4	-	< 0.5	-
Acenaphthylene	208-96-8	-	< 0.2	-
Dimethylphthalate	131-11-3	-	< 0.5	-
2,6-Dinitrotoluene	606-20-2	-	< 0.5	-
Acenaphthene	83-32-9	-	< 0.2	-
3-Nitroaniline *	99-09-2	-	< 0.5	-

* These compounds are not currently UKAS accredited.
Concentrations are reported on a wet weight basis.

Target Compounds	CAS #	R.T.	Concentration mg/kg	% Fit
2,4-Dinitrophenol	51-28-5	-	< 1.0	-
Dibenzofuran	132-64-9	-	< 0.5	-
4-Nitrophenol	100-02-7	-	< 5.0	-
2,4-Dinitrotoluene	121-14-2	-	< 0.5	-
Fluorene	86-73-7	-	< 0.2	-
Diethylphthalate	84-66-2	-	< 0.5	-
4-Chlorophenyl-phenylether	7005-72-4	-	< 0.5	-
4,6-Dinitro-2-methylphenol	534-52-1	-	< 5.0	-
4-Nitroaniline	100-01-6	-	< 0.5	-
N-Nitrosodiphenylamine	86-30-6	-	< 0.5	-
4-Bromophenyl-phenylether	101-55-3	-	< 0.5	-
Hexachlorobenzene	118-74-1	-	< 0.5	-
Pentachlorophenol	87-86-5	-	< 5.0	-
Phenanthrene	85-01-8	-	< 0.2	-
Anthracene	120-12-7	-	< 0.2	-
Di-n-butylphthalate	84-74-2	-	< 0.5	-
Fluoranthene	206-44-0	-	< 0.2	-
Pyrene	129-00-0	-	< 0.2	-
Butylbenzylphthalate	85-68-7	-	< 0.5	-
Benzofuran	56-55-3	-	< 0.2	-
Chrysene	218-01-9	-	< 0.2	-
3,3'-Dichlorobenzidine	91-94-1	-	< 2.0	-
bis(2-Ethylhexyl)phthalate	117-81-7	30.86	2.1	94
Di-n-octylphthalate	117-84-0	-	< 0.2	-
Benzobiphenylene	205-99-2	-	< 0.2	-
Benzofuran	207-08-9	-	< 0.2	-
Benzofuran	50-32-8	-	< 0.2	-
Indeno[1,2,3-cd]pyrene	193-39-5	-	< 0.2	-
Dibenzofuran	59-70-3	-	< 0.2	-
Benzofuran	191-24-2	-	< 0.2	-

* compounds not currently UKAS accredited

Internal Standards	% Area
1,4-Dichlorobenzene-d4	95
Naphthalene-d8	81
Acenaphthene-d10	95
Phenanthrene-d10	86
Chrysene-d12	87
Perylene-d12	83

Surrogates	% Rec
2-Fluorophenol	76
Phenol-d5	66
Nitrobenzene-d5	71
2-Fluorobiphenyl	75
2,4,6-Tribromophenol	94
Terphenyl-d14	94



Semi-Volatile Organic Compounds

Customer and Site Details:
Sample Details:
LIMS ID Number:
Job Number:

Exploration Associates : M4 Relief Road
CH/0038/0.10m
CL03511
484

11-Feb-00
16-Feb-00
18-Feb-00

Matrix:
Ext Method:
Operator:
Directory/Quant File:

Soil
Soxhlet
CS
0218ABN.MS4 0218CCCC1.D GPC (Y/N)

QC Batch Number:

219

Multiplier:

0.333

Dilution Factor:

5

Y

Target Compounds	CAS #	R.T. (min)	Concentration mg/kg	% Fit
Phenol	108-95-2	-	< 10.0	-
bis(2-Chloroethyl)ether	111-44-4	-	< 2.5	-
2-Chlorophenol	95-57-8	-	< 10.0	-
1,3-Dichlorobenzene	541-73-1	-	< 2.5	-
1,4-Dichlorobenzene	106-46-7	-	< 2.5	-
Benzyl alcohol	100-51-6	-	< 2.5	-
1,2-Dichlorobenzene	95-50-1	-	< 2.5	-
2-Methylphenol	95-48-7	-	< 2.5	-
bis(2-Chloroisopropyl)ether	108-60-1	-	< 2.5	-
Hexachloroethane	67-72-1	-	< 2.5	-
N-Nitroso-di-n-propylamine	621-64-7	-	< 2.5	-
3- & 4-Methylphenol	108-39-4/106-44-5	-	< 10.0	-
Nitrobenzene	98-95-3	-	< 2.5	-
Isophorone	78-59-1	-	< 2.5	-
2-Nitrophenol	88-75-5	-	< 10.0	-
2,4-Dimethylphenol	105-67-9	-	< 10.0	-
Benzoic Acid	65-85-0	-	< 50.0	-
bis(2-Chloroethoxy)methane	111-91-1	-	< 2.5	-
2,4-Dichlorophenol	120-83-2	-	< 10.0	-
1,2,4-Trichlorobenzene	120-82-1	11.63	7.2	98
Naphthalene	91-20-3	-	< 1.0	-
4-Chlorophenol *	106-48-9	-	< 10.0	-
4-Chloroaniline *	106-47-8	-	< 2.5	-
Hexachlorobutadiene	87-68-3	-	< 2.5	-
4-Chloro-3-methylphenol	59-50-7	-	< 2.5	-
2-Methylnaphthalene	91-57-6	-	< 1.0	-
1-Methyl naphthalene *	90-12-0	-	< 1.0	-
Hexachlorocyclopentadiene	77-47-4	-	< 2.5	-
2,4,6-Trichlorophenol	88-06-2	-	< 10.0	-
2,4,5-Trichlorophenol	95-95-4	-	< 10.0	-
2-Chloronaphthalene	91-58-7	-	< 1.0	-
Biphenyl *	92-52-4	15.18	3.4	99
Diphenyl ether *	101-84-8	-	< 1.0	-
2-Nitroaniline	88-74-4	-	< 2.5	-
Acenaphthylene	208-96-8	-	< 1.0	-
Dimethylphthalate	131-11-3	-	< 2.5	-
2,6-Dinitrotoluene	606-20-2	-	< 2.5	-
Acenaphthene	83-32-9	-	< 1.0	-
3-Nitroaniline *	99-09-2	-	< 2.5	-

* These compounds are not currently UKAS accredited.
Concentrations are reported on a wet weight basis.

Target Compounds	CAS #	R.T.	Concentration mg/kg	% Fit
2,4-Dinitrophenol	51-28-5	-	< 5.0	-
Dibenzofuran	132-64-9	-	< 2.5	-
4-Nitrophenol	100-02-7	-	< 25.0	-
2,4-Dinitrotoluene	121-14-2	-	< 2.5	-
Fluorene	86-73-7	-	< 1.0	-
Diethylphthalate	84-66-2	-	< 2.5	-
4-Chlorophenyl-phenylether	7005-72-4	-	< 2.5	-
4,6-Dinitro-2-methylphenol	534-52-1	-	< 25.0	-
4-Nitroaniline	100-01-6	-	< 2.5	-
N-Nitrosodiphenylamine	86-30-6	-	< 2.5	-
4-Bromophenyl-phenylether	101-55-3	-	< 2.5	-
Hexachlorobenzene	118-74-1	-	< 2.5	-
Pentachlorophenol	87-86-5	-	< 25.0	-
Phenanthrene	85-01-8	-	< 1.0	-
Anthracene	120-12-7	-	< 1.0	-
Di-n-butylphthalate	84-74-2	-	< 2.5	-
Fluoranthene	206-44-0	24.70	1.3	99
Pyrene	129-00-0	-	< 1.0	-
Butylbenzylphthalate	85-68-7	-	< 2.5	-
Benzofuranthracene	56-55-3	-	< 1.0	-
Chrysene	218-01-9	29.05	2.0	100
3,3'-Dichlorobenzidine	91-94-1	-	< 10.0	-
bis(2-Ethylhexyl)phthalate	117-81-7	-	< 2.5	-
Di-n-octylphthalate	117-84-0	-	< 1.0	-
Benzofluoranthene	205-99-2	-	< 1.0	-
Benzokjfluoranthene	207-08-9	-	< 1.0	-
Benzofluorene	50-32-8	-	< 1.0	-
Indeno[1,2,3-cd]pyrene	193-39-5	-	< 1.0	-
Dibenzo[a,h]anthracene	53-70-3	-	< 1.0	-
Benzofluoranthene	191-24-2	-	< 1.0	-

* compounds not currently UKAS accredited

Internal Standards	% Area
1,4-Dichlorobenzene-d4	110
Naphthalene-d8	99
Acenaphthene-d10	99
Phenanthrene-d10	99
Chrysene-d12	128
Perylene-d12	113

Surrogates	% Rec
2-Fluorophenol	84
Phenol-d5	81
Nitrobenzene-d5	58
2-Fluorobiphenyl	70
2,4,6-Tribromophenol	75
Terphenyl-d14	85

Semi-Volatile Organic Compounds

Exploration Associates : M4 Relief Road

CH/0038/0.10m

CL03511

11-Feb-00

16-Feb-00

18-Feb-00

219

0218ABN.MS4

Job Number:

0.333

Dilution Factor:

GPC (Y/N):

Soil

Soxhlet

56

[illegible]

The compounds listed above have been tentatively identified by a computer based library search. Compounds identified in the sample are not reported if they also occur in the method blank. The % fit is an indication of the reliability of the compound assignment. Due to the similarity between mass spectra of some isomeric compounds assignments may not be correct. Other compounds may also be present but identification was not possible. Concentrations are semi-quantitative, assume a response factor of 1 and use the nearest internal standard. Concentrations are reported on a wet weight basis.

Semi-Volatile Organic Compounds

Customer and Site Details:
Sample Details:
LIMS ID Number:
Job Number:

Exploration Associates: M4 Relief Road
CH/0038/ / 1.30- m
Date Booked in:
Date Extracted:
Date Analysed:

Matrix: Soil
Ext Method: Soxhlet
Operator: SW
Directory/Quant File: 0217ABN.MS2 0217CCC3.D GPC (Y/N)

QC Batch Number: 219
Multiplier: 0.0667
Dilution Factor: 1
N

Target Compounds	CAS #	R.T. (min)	Concentration mg/kg	% Fit
Phenol	108-95-2	-	< 2.0	-
bis(2-Chloroethyl)ether	111-44-4	-	< 0.5	-
2-Chlorophenol	95-57-8	-	< 2.0	-
1,3-Dichlorobenzene	541-73-1	-	< 0.5	-
1,4-Dichlorobenzene	106-46-7	-	< 0.5	-
Benzyl alcohol	100-51-6	-	< 0.5	-
1,2-Dichlorobenzene	95-50-1	-	< 0.5	-
2-Methylphenol	95-48-7	-	< 0.5	-
bis(2-Chloroisopropyl)ether	108-60-1	-	< 0.5	-
Hexachloroethane	67-72-1	-	< 0.5	-
N-Nitroso-di-n-propylamine	621-64-7	-	< 0.5	-
3- & 4-Methylphenol	108-39-4/106-44-5	-	< 2.0	-
Nitrobenzene	98-95-3	-	< 0.5	-
Isophorone	78-59-1	-	< 0.5	-
2-Nitrophenol	88-75-5	-	< 2.0	-
2,4-Dimethylphenol	105-67-9	-	< 2.0	-
Benzoic Acid	65-85-0	-	< 10.0	-
bis(2-Chloroethoxy)methane	111-91-1	-	< 0.5	-
2,4-Dichlorophenol	120-83-2	-	< 2.0	-
1,2,4-Trichlorobenzene	120-82-1	-	< 0.5	-
Naphthalene	91-20-3	-	< 0.2	-
4-Chlorophenol *	106-48-9	-	< 2.0	-
4-Chloroaniline	106-47-8	-	< 0.5	-
Hexachlorobutadiene	87-68-3	-	< 0.5	-
4-Chloro-3-methylphenol	59-50-7	-	< 0.5	-
2-Methylnaphthalene	91-57-6	-	< 0.2	-
1-Methylnaphthalene *	90-12-0	-	< 0.2	-
Hexachlorocyclopentadiene	77-47-4	-	< 0.5	-
2,4,6-Trichlorophenol	88-06-2	-	< 2.0	-
2,4,5-Trichlorophenol	95-95-4	-	< 2.0	-
2-Chloronaphthalene	91-58-7	-	< 0.2	-
Biphenyl *	92-52-4	-	< 0.2	-
Diphenyl ether *	101-84-8	-	< 0.2	-
2-Nitroaniline	88-74-4	-	< 0.5	-
Acenaphthylene	208-96-8	-	< 0.2	-
Dimethylphthalate	131-11-3	-	< 0.5	-
2,6-Dinitrotoluene	606-20-2	-	< 0.5	-
Acenaphthene	83-32-9	-	< 0.2	-
3-Nitroaniline *	99-09-2	-	< 0.5	-

* These compounds are not currently UKAS accredited.
Concentrations are reported on a wet weight basis.

Target Compounds	CAS #	R.T.	Concentration mg/kg	% Fit
2,4-Dinitrophenol	51-28-5	-	< 1.0	-
Dibenzofuran	132-64-9	-	< 0.5	-
4-Nitrophenol	100-02-7	-	< 5.0	-
2,4-Dinitrotoluene	121-14-2	-	< 0.5	-
Fluorene	86-73-7	-	< 0.2	-
Diethylphthalate	84-66-2	-	< 0.5	-
4-Chlorophenyl-phenylether	7005-72-4	-	< 0.5	-
4,6-Dinitro-2-methylphenol	534-52-1	-	< 5.0	-
4-Nitroaniline	100-01-6	-	< 0.5	-
N-Nitrosodiphenylamine	86-30-6	-	< 0.5	-
4-Bromophenyl-phenylether	101-55-3	-	< 0.5	-
Hexachlorobenzene	118-74-1	-	< 0.5	-
Pentachlorophenol	87-86-5	-	< 5.0	-
Phenanthrene	85-01-8	-	< 0.2	-
Anthracene	120-12-7	-	< 0.2	-
Di-n-butylphthalate	84-74-2	-	< 0.5	-
Fluoranthene	206-44-0	-	< 0.2	-
Pyrene	129-00-0	-	< 0.2	-
Butylbenzylphthalate	85-68-7	-	< 0.5	-
Benzofuran	56-55-3	-	< 0.2	-
Chrysene	218-01-9	-	< 0.2	-
3,3'-Dichlorobenzidine	91-94-1	-	< 2.0	-
bis(2-Ethylhexyl)phthalate	117-81-7	30.86	2.2	95
Di-n-octylphthalate	117-84-0	-	< 0.2	-
Benzofluoranthene	205-99-2	-	< 0.2	-
Benzofluoranthene	207-08-9	-	< 0.2	-
Benzofluoranthene	50-32-8	-	< 0.2	-
Indeno[1,2,3-cd]pyrene	193-39-5	-	< 0.2	-
Dibenzofluoranthene	53-70-3	-	< 0.2	-
Benzofluoranthene	191-24-2	-	< 0.2	-

* compounds not currently UKAS accredited

Internal Standards	% Area
1,4-Dichlorobenzene-d4	97
Naphthalene-d8	81
Acenaphthene-d10	98
Phenanthrene-d10	91
Chrysene-d12	85
Perylene-d12	79

Surrogates	% Rec
2-Fluorophenol	77
Phenol-d5	70
Nitrobenzene-d5	73
2-Fluorobiphenyl	74
2,4,6-Tribromophenol	92
Terphenyl-d14	100

Semi-Volatile Organic Compounds

Customer and Site Details:
Sample Details:
LIMS ID Number:
Job Number:

Exploration Associates : M4 Relief Road
CH/0039/0.00m
CL03513
484

11-Feb-00
16-Feb-00
18-Feb-00

Matrix:
Ext Method:
Operator:
Directory/Quant File:

Soil
Soxhlet
CS
0218ABN.MS4 0218CCC1 D GPC (Y/N)

QC Batch Number:
Multiplier:
Dilution Factor:

219
0.0667
1
Y

Target Compounds	CAS #	R.T. (min)	Concentration mg/kg	% Fit
Phenol	108-95-2	-	< 2.0	-
bis(2-Chloroethyl)ether	111-44-4	-	< 0.5	-
2-Chlorophenol	95-57-8	-	< 2.0	-
1,3-Dichlorobenzene	541-73-1	-	< 0.5	-
1,4-Dichlorobenzene	106-46-7	-	< 0.5	-
Benzyl alcohol	100-51-6	-	< 0.5	-
1,2-Dichlorobenzene	95-50-1	-	< 0.5	-
2-Methylphenol	95-48-7	-	< 0.5	-
bis(2-Chloroisopropyl)ether	108-60-1	-	< 0.5	-
Hexachloroethane	67-72-1	-	< 0.5	-
N-Nitroso-di-n-propylamine	621-64-7	-	< 0.5	-
3- & 4-Methylphenol	108-39-4/108-44-5	-	< 2.0	-
Nitrobenzene	98-95-3	-	< 0.5	-
Isophorone	78-59-1	-	< 0.5	-
2-Nitrophenol	88-75-5	-	< 2.0	-
2,4-Dimethylphenol	105-67-9	-	< 2.0	-
Benzoic Acid	65-85-0	-	< 10.0	-
bis(2-Chloroethoxy)methane	111-91-1	-	< 0.5	-
2,4-Dichlorophenol	120-83-2	-	< 2.0	-
1,2,4-Trichlorobenzene	120-82-1	-	< 0.5	-
Naphthalene	91-20-3	-	< 0.2	-
4-Chlorophenol *	106-48-9	-	< 2.0	-
4-Chloroaniline *	106-47-8	-	< 0.5	-
Hexachlorobutadiene	87-68-3	-	< 0.5	-
4-Chloro-3-methylphenol	59-50-7	-	< 0.5	-
2-Methylnaphthalene	91-57-6	-	< 0.2	-
1-Methyl naphthalene *	90-12-0	-	< 0.2	-
Hexachlorocyclopentadiene	77-47-4	-	< 0.5	-
2,4,6-Trichlorophenol	88-06-2	-	< 2.0	-
2,4,5-Trichlorophenol	95-95-4	-	< 2.0	-
2-Chloronaphthalene	91-58-7	-	< 0.2	-
Biphenyl *	92-52-4	15.18	0.3	99
Diphenyl ether *	101-84-8	-	< 0.2	-
2-Nitroaniline	88-74-4	-	< 0.5	-
Acenaphthylene	208-96-8	-	< 0.2	-
Dimethylphthalate	131-11-3	-	< 0.5	-
2,6-Dinitrotoluene	606-20-2	-	< 0.5	-
Acenaphthene	83-32-9	-	< 0.2	-
3-Nitroaniline *	99-09-2	-	< 0.5	-

* These compounds are not currently UKAS accredited.
Concentrations are reported on a wet weight basis.

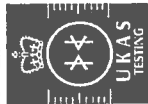
Target Compounds	CAS #	R.T.	Concentration mg/kg	% Fit
2,4-Dinitrophenol	51-28-5	-	< 1.0	-
Dibenzofuran	132-64-9	-	< 0.5	-
4-Nitrophenol	100-02-7	-	< 5.0	-
2,4-Dinitrotoluene	121-14-2	-	< 0.5	-
Fluorene	86-73-7	-	< 0.2	-
Diethylphthalate	84-66-2	-	< 0.5	-
4-Chlorophenyl-phenylether	7005-72-4	-	< 0.5	-
4,6-Dinitro-2-methylphenol	534-52-1	-	< 5.0	-
4-Nitroaniline	100-01-6	-	< 0.5	-
N-Nitrosodiphenylamine	86-30-6	-	< 0.5	-
4-Bromophenyl-phenylether	101-55-3	-	< 0.5	-
Hexachlorobenzene	118-74-1	-	< 0.5	-
Pentachlorophenol	87-86-5	-	< 5.0	-
Phenanthrene	85-01-8	21.18	1.7	99
Anthracene	120-12-7	21.31	0.4	93
Di-n-butylphthalate	84-74-2	-	< 0.5	-
Fluoranthene	206-44-0	24.71	2.2	98
Pyrene	129-00-0	25.34	1.5	100
Butylbenzylphthalate	85-68-7	-	< 0.5	-
Benzofluranthene	56-55-3	28.95	0.7	91
Chrysene	218-01-9	29.07	1.1	97
3,3'-Dichlorobenzidine	91-94-1	-	< 2.0	-
bis(2-Ethylhexyl)phthalate	117-81-7	29.71	4.1	99
Di-n-octylphthalate	117-84-0	-	< 0.2	-
Benzofluranthene	205-99-2	32.01	0.8	93
Benzofluranthene	207-08-9	32.07	0.6	93
Benzofluranthene	50-32-8	32.95	0.8	98
Indeno[1,2,3-cd]pyrene	193-39-5	37.41	0.3	99
Dibenzofluranthene	53-70-3	-	< 0.2	-
Benzofluranthene	191-24-2	38.35	0.6	99

* compounds not currently UKAS accredited

Internal Standards	% Area
1,4-Dichlorobenzene-d4	97
Naphthalene-d8	85
Acenaphthene-d10	91
Phenanthrene-d10	96
Chrysene-d12	107
Perylene-d12	96

Surrogates	% Rec
2-Fluorophenol	81
Phenol-d5	85
Nitrobenzene-d5	56
2-Fluorobiphenyl	65
2,4,6-Tribromophenol	78
Terphenyl-d14	88

Semi-Volatile Organic Compounds



1252

Customer and Site Details:

Sample Details:

LIMS ID Number:

Exploration Associates : M4 Relief Road

CH/0039/0.00m

CL03513

484

Job Number:

0.0667

Soil

Soxhlet

5

Date Booked in: 11-Feb-00

Date Booked in: 16-Feb-00

Date Analysed: 18-Feb-00

Date Analyzed.	OC Batch Number.
10/1	219

QC Batch Number: 0218ABN MSA
Director: Z13

[illegible]

The compounds listed above have been tentatively identified by a computer based library search.

Compounds identified in the sample are not reported if they also occur in the method blank.

The % fit is an indication of the reliability of the compound assignment. Compounds identified in the samples are not reported ninety days after testing.

Due to the similarity between mass spectra of some isomeric compounds assignments may not be correct.

Other compounds may also be present but identification was not possible. Due to the similarity between mass spectra of some unknown compounds

Other compounds may also be present but recognition was not possible. Concentrations are semi-quantitative, assume a response factor of 1 and use the nearest internal standard.

Concentrations are reported on a wet weight basis.

Semi-Volatile Organic Compounds



Customer and Site Details:
Sample Details:
LIMS ID Number:
Job Number:

Exploration Associates : M4 Relief Road
Date Booked in:
Date Extracted:
Date Analysed:

Matrix:
Ext Method:
Operator:
Directory/Quant File:

QC Batch Number:
Multiplier:
Dilution Factor:

219
0.133
2
Y

Target Compounds	CAS #	R.T. (min)	Concentration mg/kg	% Fit
Phenol	108-95-2	-	< 4.0	-
bis(2-Chloroethyl)ether	111-44-4	-	< 1.0	-
2-Chlorophenol	95-57-8	-	< 4.0	-
1,3-Dichlorobenzene	541-73-1	-	< 1.0	-
1,4-Dichlorobenzene	106-46-7	-	< 1.0	-
Benzyl alcohol	100-51-6	-	< 1.0	-
1,2-Dichlorobenzene	95-50-1	-	< 1.0	-
2-Methylphenol	95-48-7	-	< 1.0	-
bis(2-Chloroisopropyl)ether	108-60-1	-	< 1.0	-
Hexachloroethane	67-72-1	-	< 1.0	-
N-Nitroso-di-n-propylamine	621-64-7	-	< 1.0	-
3- & 4-Methylphenol	108-39-4/106-44-5	-	< 4.0	-
Nitrobenzene	98-95-3	-	< 1.0	-
Isophorone	78-59-1	-	< 1.0	-
2-Nitrophenol	88-75-5	-	< 4.0	-
2,4-Dimethylphenol	105-67-9	-	< 4.0	-
Benzoic Acid	65-85-0	-	< 20.0	-
bis(2-Chloroethoxy)methane	111-91-1	-	< 1.0	-
2,4-Dichlorophenol	120-83-2	-	< 4.0	-
1,2,4-Trichlorobenzene	120-82-1	-	< 1.0	-
Naphthalene	91-20-3	-	< 0.4	-
4-Chlorophenol *	106-48-9	-	< 4.0	-
4-Chloroaniline *	106-47-8	-	< 1.0	-
Hexachlorobutadiene	87-68-3	-	< 1.0	-
4-Chloro-3-methylphenol	59-50-7	-	< 1.0	-
2-Methylnaphthalene	91-57-6	-	< 0.4	-
1-Methyl naphthalene *	90-12-0	-	< 0.4	-
Hexachlorocyclopentadiene	77-47-4	-	< 1.0	-
2,4,6-Trichlorophenol	88-06-2	-	< 4.0	-
2,4,5-Trichlorophenol	95-95-4	-	< 4.0	-
2-Chloronaphthalene	91-58-7	-	< 0.4	-
Biphenyl *	92-52-4	15.18	0.6	98
Diphenyl ether *	101-84-8	-	< 0.4	-
2-Nitroaniline	88-74-4	-	< 1.0	-
Acenaphthylene	208-96-8	-	< 0.4	-
Dimethylphthalate	131-11-3	-	< 1.0	-
2,6-Dinitrotoluene	606-20-2	-	< 1.0	-
Acenaphthene	83-32-9	-	< 0.4	-
3-Nitroaniline *	99-09-2	-	< 1.0	-

* These compounds are not currently UKAS accredited.
Concentrations are reported on a wet weight basis.

Target Compounds	CAS #	R.T.	Concentration mg/kg	% Fit
2,4-Dinitrophenol	51-28-5	-	< 2.0	-
Dibenzofuran	132-64-9	-	< 1.0	-
4-Nitrophenol	100-02-7	-	< 10.0	-
2,4-Dinitrotoluene	121-14-2	-	< 1.0	-
Fluorene	86-73-7	-	< 0.4	-
Diethylphthalate	84-66-2	-	< 1.0	-
4-Chlorophenyl-phenylether	7005-72-4	-	< 1.0	-
4,6-Dinitro-2-methylphenol	534-52-1	-	< 10.0	-
4-Nitroaniline	100-01-6	-	< 1.0	-
N-Nitrosodiphenylamine	86-30-6	-	< 1.0	-
4-Bromophenyl-phenylether	101-55-3	-	< 1.0	-
Hexachlorobenzene	118-74-1	-	< 1.0	-
Pentachlorophenol	87-86-5	-	< 10.0	-
Phenanthrene	85-01-8	21.18	0.4	99
Anthracene	120-12-7	-	< 0.4	-
Di-n-butylphthalate	84-74-2	-	< 1.0	-
Fluoranthene	206-44-0	24.70	0.8	98
Pyrene	129-00-0	25.34	0.6	99
Butylbenzylphthalate	85-68-7	-	< 1.0	-
Benzolalanthracene	56-55-3	-	< 0.4	-
Chrysene	218-01-9	29.05	1.9	99
3,3'-Dichlorobenzidine	91-94-1	-	< 4.0	-
bis(2-Ethylhexyl)phthalate	117-81-7	29.70	6.9	99
Di-n-octylphthalate	117-84-0	-	< 0.4	-
Benzobiphenylene	205-99-2	32.00	0.5	97
Benzokjfluoranthene	207-08-9	32.06	0.4	97
Benzolalpyrene	50-32-8	32.94	0.6	94
Indenol1,2,3-cdpyrene	193-39-5	37.36	0.5	97
Dibenzofa,hlanthracene	53-70-3	-	< 0.4	-
Benzofg,h,i)perylene	191-24-2	38.31	0.4	98

* compounds not currently UKAS accredited

Internal Standards	% Area
1,4-Dichlorobenzene-d4	89
Naphthalene-d8	76
Acenaphthene-d10	83
Phenanthrene-d10	85
Chrysene-d12	109
Perylene-d12	95

Surrogates	% Rec
2-Fluorophenol	66
Phenol-d5	65
Nitrobenzene-d5	47
2-Fluorobiphenyl	60
2,4,6-Tribromophenol	78
Triphenyl-d14	82

Semi-Volatile Organic Compounds

Customer and Site Details:
Sample Details:
LIMS ID Number:
Job Number:

Exploration Associates: M4 Relief Road
CH/0040/ / 0.00- m
CL03515
484

11-Feb-00
16-Feb-00
17-Feb-00

Matrix:
Ext Method:
Operator:
Directory/Quant File:

Soil
Soxhlet
SW
0217ABN.MS2

QC Batch Number:
Multiplier:
Dilution Factor:

219
0.0667
1
N

Target Compounds	CAS #	R.T. (min)	Concentration mg/kg	% Fit
Phenol	108-95-2	-	< 2.0	-
bis(2-Chloroethyl)ether	111-44-4	-	< 0.5	-
2-Chlorophenol	95-57-8	-	< 2.0	-
1,3-Dichlorobenzene	541-73-1	-	< 0.5	-
1,4-Dichlorobenzene	106-46-7	-	< 0.5	-
Benzyl alcohol	100-51-6	-	< 0.5	-
1,2-Dichlorobenzene	95-50-1	-	< 0.5	-
2-Methylphenol	95-48-7	-	< 0.5	-
bis(2-Chloroisopropyl)ether	108-60-1	-	< 0.5	-
Hexachloroethane	67-72-1	-	< 0.5	-
N-Nitroso-di-n-propylamine	621-64-7	-	< 0.5	-
3- & 4-Methylphenol	108-39-4/106-44-5	-	< 2.0	-
Nitrobenzene	98-95-3	-	< 0.5	-
Isophorone	78-59-1	-	< 0.5	-
2-Nitrophenol	88-75-5	-	< 2.0	-
2,4-Dimethylphenol	105-67-9	-	< 2.0	-
Benzoic Acid	65-85-0	-	< 10.0	-
bis(2-Chloroethoxy)methane	111-91-1	-	< 0.5	-
2,4-Dichlorophenol	120-83-2	-	< 2.0	-
1,2,4-Trichlorobenzene	120-82-1	-	< 0.5	-
Naphthalene	91-20-3	-	< 0.2	-
4-Chlorophenol *	106-48-9	-	< 2.0	-
4-Chloroaniline	106-47-8	-	< 0.5	-
Hexachlorobutadiene	87-68-3	-	< 0.5	-
4-Chloro-3-methylphenol	59-50-7	-	< 0.5	-
2-Methylnaphthalene	91-57-6	-	< 0.2	-
1-Methylnaphthalene *	90-12-0	-	< 0.2	-
Hexachlorocyclopentadiene	77-47-4	-	< 0.5	-
2,4,6-Trichlorophenol	88-06-2	-	< 2.0	-
2,4,5-Trichlorophenol	95-95-4	-	< 2.0	-
2-Chloronaphthalene	91-58-7	-	< 0.2	-
Biphenyl *	92-52-4	-	< 0.2	-
Diphenyl ether *	101-84-8	-	< 0.2	-
2-Nitroaniline	88-74-4	-	< 0.5	-
Acenaphthylene	208-96-8	-	< 0.2	-
Dimethylphthalate	131-11-3	-	< 0.5	-
2,6-Dinitrotoluene	606-20-2	-	< 0.5	-
Acenaphthene	83-32-9	-	< 0.2	-
3-Nitroaniline *	99-09-2	-	< 0.5	-

* These compounds are not currently UKAS accredited.
Concentrations are reported on a wet weight basis.

Target Compounds	CAS #	R.T.	Concentration mg/kg	% Fit
2,4-Dinitrophenol	51-28-5	-	< 1.0	-
Dibenzofuran	132-64-9	-	< 0.5	-
4-Nitrophenol	100-02-7	-	< 5.0	-
2,4-Dinitrotoluene	121-14-2	-	< 0.5	-
Fluorene	86-73-7	-	< 0.2	-
Diethylphthalate	84-66-2	-	< 0.5	-
4-Chlorophenyl-phenylether	7005-72-4	-	< 0.5	-
4,6-Dinitro-2-methylphenol	534-52-1	-	< 5.0	-
4-Nitroaniline	100-01-6	-	< 0.5	-
N-Nitrosodiphenylamine	86-30-6	-	< 0.5	-
4-Bromophenyl-phenylether	101-55-3	-	< 0.5	-
Hexachlorobenzene	118-74-1	-	< 0.5	-
Pentachlorophenol	87-86-5	-	< 5.0	-
Phenanthrene	85-01-8	-	< 0.2	-
Anthracene	120-12-7	-	< 0.2	-
Di-n-butylphthalate	84-74-2	-	< 0.5	-
Fluoranthene	206-44-0	-	< 0.2	-
Pyrene	129-00-0	-	< 0.2	-
Butylbenzylphthalate	85-68-7	-	< 0.5	-
Benzofuranthracene	56-55-3	-	< 0.2	-
Chrysene	218-01-9	-	< 0.2	-
3,3'-Dichlorobenzidine	91-94-1	-	< 2.0	-
bis(2-Ethylhexyl)phthalate	117-81-7	-	< 0.5	-
Di-n-octylphthalate	117-84-0	-	< 0.2	-
Benzofluoranthene	205-99-2	-	< 0.2	-
Benzofluoranthene	207-08-9	-	< 0.2	-
Benzofluoranthene	50-32-8	-	< 0.2	-
Indenol 1,2,3-cdpyrene	193-39-5	-	< 0.2	-
Dibenzo[a,h]anthracene	53-70-3	-	< 0.2	-
Benzofluoranthene	191-24-2	-	< 0.2	-

* compounds not currently UKAS accredited

Internal Standards	% Area	Surrogates	% Rec
1,4-Dichlorobenzene-d4	103	2-Fluorophenol	85
Naphthalene-d8	86	Phenol-d5	78
Acenaphthene-d10	105	Nitrobenzene-d5	83
Phenanthrene-d10	95	2-Fluorobiphenyl	82
Chrysene-d12	102	2,4,6-Tribromophenol	98
Perylene-d12	99	Terphenyl-d14	93

Semi-Volatile Organic Compounds

Customer and Site Details:

Sample Details:
LIMS ID Number:
Job Number:

Exploration Associates : M4 Relief Road

Date Booked in:
Date Extracted:
Date Analysed:

11-Feb-00
16-Feb-00
18-Feb-00

Matrix:

Ext Method:
Operator:
Directory/Quant File:

Soil

Soxhlet
CS

QC Batch Number:

Multiplier:
Dilution Factor:

219

0.333
5

Target Compounds	CAS #	R.T. (min)	Concentration mg/kg	% Fit
Phenol	108-95-2	-	< 10.0	-
bis(2-Chloroethyl)ether	111-44-4	-	< 2.5	-
2-Chlorophenol	95-57-8	-	< 10.0	-
1,3-Dichlorobenzene	541-73-1	-	< 2.5	-
1,4-Dichlorobenzene	106-46-7	-	< 2.5	-
Benzyl alcohol	100-51-6	-	< 2.5	-
1,2-Dichlorobenzene	95-50-1	-	< 2.5	-
2-Methylphenol	95-48-7	-	< 2.5	-
bis(2-Chloroisopropyl)ether	108-60-1	-	< 2.5	-
Hexachloroethane	67-72-1	-	< 2.5	-
N-Nitroso-di-n-propylamine	621-64-7	-	< 2.5	-
3- & 4-Methylphenol	108-39-4/108-44-5	-	< 10.0	-
Nitrobenzene	98-95-3	-	< 2.5	-
Isophorone	78-59-1	-	< 2.5	-
2-Nitrophenol	88-75-5	-	< 10.0	-
2,4-Dimethylphenol	105-67-9	-	< 10.0	-
Benzoic Acid	65-85-0	-	< 50.0	-
bis(2-Chloroethoxy)methane	111-91-1	-	< 2.5	-
2,4-Dichlorophenol	120-83-2	-	< 10.0	-
1,2,4-Trichlorobenzene	120-82-1	11.63	2.6	99
Naphthalene	91-20-3	-	< 1.0	-
4-Chlorophenol *	106-48-9	-	< 10.0	-
4-Chloroaniline *	106-47-8	-	< 2.5	-
Hexachlorobutadiene	87-68-3	-	< 2.5	-
4-Chloro-3-methylphenol	59-50-7	-	< 2.5	-
2-Methylnaphthalene	91-57-6	-	< 1.0	-
1-Methyl naphthalene *	90-12-0	-	< 1.0	-
Hexachlorocyclopentadiene	77-47-4	-	< 2.5	-
2,4,6-Trichlorophenol	88-06-2	-	< 10.0	-
2,4,5-Trichlorophenol	95-95-4	-	< 10.0	-
2-Chloronaphthalene	91-58-7	-	< 1.0	-
Biphenyl *	92-52-4	-	< 1.0	-
Diphenyl ether *	101-84-8	-	< 1.0	-
2-Nitroaniline	88-74-4	-	< 2.5	-
Acenaphthylene	208-96-8	-	< 1.0	-
Dimethylphthalate	131-11-3	-	< 2.5	-
2,6-Dinitrotoluene	606-20-2	-	< 2.5	-
Acenaphthene	83-32-9	-	< 1.0	-
3-Nitroaniline *	99-09-2	-	< 2.5	-

* These compounds are not currently UKAS accredited.
Concentrations are reported on a wet weight basis.

Target Compounds	CAS #	R.T.	Concentration mg/kg	% Fit
2,4-Dinitrophenol	51-28-5	-	< 5.0	-
Dibenzofuran	132-64-9	-	< 2.5	-
4-Nitrophenol	100-02-7	-	< 25.0	-
2,4-Dinitrotoluene	121-14-2	-	< 2.5	-
Fluorene	86-73-7	-	< 1.0	-
Diethylphthalate	84-66-2	-	< 2.5	-
4-Chlorophenyl-phenylether	7005-72-4	-	< 2.5	-
4,6-Dinitro-2-methylphenol	534-52-1	-	< 25.0	-
4-Nitroaniline	100-01-6	-	< 2.5	-
N-Nitrosodiphenylamine	86-30-6	-	< 2.5	-
4-Bromophenyl-phenylether	101-55-3	-	< 2.5	-
Hexachlorobenzene	118-74-1	-	< 2.5	-
Pentachlorophenol	87-86-5	-	< 25.0	-
Phenanthrene	85-01-8	-	< 1.0	-
Anthracene	120-12-7	-	< 1.0	-
Di-n-butylphthalate	84-74-2	-	< 2.5	-
Fluoranthene	206-44-0	-	< 1.0	-
Pyrene	129-00-0	-	< 1.0	-
Butylbenzylphthalate	85-68-7	-	< 2.5	-
Benzofuran	56-55-3	-	< 1.0	-
Chrysene	218-01-9	-	< 1.0	-
3,3'-Dichlorobenzidine	91-94-1	-	< 10.0	-
bis(2-Ethylhexyl)phthalate	117-81-7	29.69	3.5	100
Di-n-octylphthalate	117-84-0	-	< 1.0	-
Benzobenzofuran	205-99-2	-	< 1.0	-
Benzofluoranthene	207-08-9	-	< 1.0	-
Benzofluoranthene	50-32-8	-	< 1.0	-
Indeno[1,2,3-cd]pyrene	193-39-5	-	< 1.0	-
Dibenzofluoranthene	53-70-3	-	< 1.0	-
Benzofluoranthene	191-24-2	-	< 1.0	-

* compounds not currently UKAS accredited

Internal Standards	% Area	Surrogates	% Rec
1,4-Dichlorobenzene-d4	119	2-Fluorophenol	89
Naphthalene-d8	107	Phenol-d5	88
Acenaphthene-d10	117	Nitrobenzene-d5	65
Phenanthrene-d10	117	2-Fluorobiphenyl	75
Chrysene-d12	141	2,4,6-Tribromophenol	74
Perylene-d12	115	Terphenyl-d14	93

Semi-Volatile Organic Compounds

Customer and Site Details:
Sample Details:
LIMS ID Number:
Job Number:

Exploration Associates: M4 Relief Road
CH/0041/ / 0.05- m
CL03517
484

11-Feb-00
16-Feb-00
22-Feb-00

Matrix:
Ext Method:
Operator:
Directory/Quant File:

Soil
Soxhlet
SW
0222ABN.MS5

QC Batch Number:
Multiplier:
Dilution Factor:

219
0.0667
1
N

Target Compounds	CAS #	R.T. (min)	Concentration mg/kg	% Fit
Phenol	108-95-2	-	< 2.0	-
bis(2-Chloroethyl)ether	111-44-4	-	< 0.5	-
2-Chlorophenol	95-57-8	-	< 2.0	-
1,3-Dichlorobenzene	541-73-1	-	< 0.5	-
1,4-Dichlorobenzene	106-46-7	-	< 0.5	-
Benzyl alcohol	100-51-6	-	< 0.5	-
1,2-Dichlorobenzene	95-50-1	-	< 0.5	-
2-Methylphenol	95-48-7	-	< 0.5	-
bis(2-Chloroisopropyl)ether	108-60-1	-	< 0.5	-
Hexachloroethane	67-72-1	-	< 0.5	-
N-Nitroso-di-n-propylamine	621-64-7	-	< 0.5	-
3- & 4-Methylphenol	108-39-4/106-44-5	-	< 2.0	-
Nitrobenzene	98-95-3	-	< 0.5	-
Isophorone	78-59-1	-	< 0.5	-
2-Nitrophenol	88-75-5	-	< 2.0	-
2,4-Dimethylphenol	105-67-9	-	< 2.0	-
Benzoic Acid	65-85-0	-	< 10.0	-
bis(2-Chloroethoxy)methane	111-91-1	-	< 0.5	-
2,4-Dichlorophenol	120-83-2	-	< 2.0	-
1,2,4-Trichlorobenzene	120-82-1	-	< 0.5	-
Naphthalene	91-20-3	-	< 0.2	-
4-Chlorophenol *	106-48-9	-	< 2.0	-
4-Chloroaniline	106-47-8	-	< 0.5	-
Hexachlorobutadiene	87-68-3	-	< 0.5	-
4-Chloro-3-methylphenol	59-50-7	-	< 0.5	-
2-Methylnaphthalene	91-57-6	-	< 0.2	-
1-Methyl naphthalene *	90-12-0	-	< 0.2	-
Hexachlorocyclopentadiene	77-47-4	-	< 0.5	-
2,4,6-Trichlorophenol	88-06-2	-	< 2.0	-
2,4,5-Trichlorophenol	95-95-4	-	< 2.0	-
2-Chloronaphthalene	91-58-7	-	< 0.2	-
Biphenyl *	92-52-4	-	< 0.2	-
Diphenyl ether *	101-84-8	-	< 0.2	-
2-Nitroaniline	88-74-4	-	< 0.5	-
Acenaphthylene	208-96-8	-	< 0.2	-
Dimethylphthalate	131-11-3	-	< 0.5	-
2,6-Dinitrotoluene	606-20-2	-	< 0.5	-
Acenaphthene	83-32-9	-	< 0.2	-
3-Nitroaniline	99-09-2	-	< 0.5	-

* These compounds are not currently UKAS accredited.

Concentrations are reported on a wet weight basis.

Target Compounds	CAS #	R.T.	Concentration mg/kg	% Fit
2,4-Dinitrophenol	51-28-5	-	< 1.0	-
Dibenzofuran	132-64-9	-	< 0.5	-
4-Nitrophenol	100-02-7	-	< 5.0	-
2,4-Dinitrotoluene	121-14-2	-	< 0.5	-
Fluorene	86-73-7	-	< 0.2	-
Diethylphthalate	84-66-2	-	< 0.5	-
4-Chlorophenyl-phenylether	7005-72-4	-	< 0.5	-
4,6-Dinitro-2-methylphenol	534-52-1	-	< 5.0	-
4-Nitroaniline	100-01-6	-	< 0.5	-
N-Nitrosodiphenylamine	86-30-6	-	< 0.5	-
4-Bromophenyl-phenylether	101-55-3	-	< 0.5	-
Hexachlorobenzene	118-74-1	-	< 0.5	-
Pentachlorophenol	87-86-5	-	< 5.0	-
Phenanthrene	85-01-8	-	< 0.2	-
Anthracene	120-12-7	-	< 0.2	-
Di-n-butylphthalate	84-74-2	-	< 0.5	-
Fluoranthene	206-44-0	-	< 0.2	-
Pyrene	129-00-0	-	< 0.2	-
Butylbenzylphthalate	85-68-7	-	< 0.5	-
Benzofuranthracene	56-55-3	-	< 0.2	-
Chrysene	218-01-9	-	< 0.2	-
3,3'-Dichlorobenzidine	91-94-1	-	< 2.0	-
bis(2-Ethylhexyl)phthalate	117-81-7	29.51	1.6	98
Di-n-octylphthalate	117-84-0	-	< 0.2	-
Benzobenzofluoranthene	205-99-2	-	< 0.2	-
Benzokjfluoranthene	207-08-9	-	< 0.2	-
Benzofluoranthene	50-32-8	-	< 0.2	-
Indeno[1,2,3-cd]pyrene	193-39-5	-	< 0.2	-
Dibenzofluoranthene	53-70-3	-	< 0.2	-
Benzofluoranthene	191-24-2	-	< 0.2	-

* compounds not currently UKAS accredited

Internal Standards	% Area	Surrogates	% Rec
1,4-Dichlorobenzene-d4	103	2-Fluorophenol	100
Naphthalene-d8	101	Phenol-d5	107
Acenaphthene-d10	113	Nitrobenzene-d5	94
Phenanthrene-d10	113	2-Fluorobiphenyl	86
Chrysene-d12	123	2,4,6-Tribromophenol	111
Perylene-d12	123	Terphenyl-d14	93

Semi-Volatile Organic Compounds

Customer and Site Details:
Sample Details:
LIMS ID Number:
Job Number:

Exploration Associates : M4 Relief Road
Date Booked in:
Date Extracted:
Date Analysed:

Matrix: Soil
Ext Method: Soxhlet
Operator: CS
Directory/Quant File: 0218ABN.MS4 0218CCCC1.D GPC (Y/N)

QC Batch Number: 219
Multiplier: 1.667
Dilution Factor: 25
Y

Target Compounds	CAS #	R.T. (min)	Concentration mg/kg	% Fit
Phenol	108-95-2	-	< 50.0	-
bis(2-Chloroethyl)ether	111-44-4	-	< 12.5	-
2-Chlorophenol	95-57-8	-	< 50.0	-
1,3-Dichlorobenzene	541-73-1	-	< 12.5	-
1,4-Dichlorobenzene	106-46-7	-	< 12.5	-
Benzyl alcohol	100-51-6	-	< 12.5	-
1,2-Dichlorobenzene	95-50-1	-	< 12.5	-
2-Methylphenol	95-48-7	-	< 12.5	-
bis(2-Chloroisopropyl)ether	108-60-1	-	< 12.5	-
Hexachloroethane	67-72-1	-	< 12.5	-
N-Nitroso-di-n-propylamine	621-64-7	-	< 12.5	-
3- & 4-Methylphenol	108-39-4/108-44-5	-	< 50.0	-
Nitrobenzene	98-95-3	-	< 12.5	-
Isophorone	78-59-1	-	< 12.5	-
2-Nitrophenol	88-75-5	-	< 50.0	-
2,4-Dimethylphenol	105-67-9	-	< 50.0	-
Benzoic Acid	65-85-0	-	< 250	-
bis(2-Chloroethoxy)methane	111-91-1	-	< 12.5	-
2,4-Dichlorophenol	120-83-2	-	< 50.0	-
1,2,4-Trichlorobenzene	120-82-1	11.63	14.9	98
Naphthalene	91-20-3	-	< 5.0	-
4-Chlorophenol *	106-48-9	-	< 50.0	-
4-Chloroaniline	106-47-8	-	< 12.5	-
Hexachlorobutadiene	87-68-3	-	< 12.5	-
4-Chloro-3-methylphenol	59-50-7	-	< 12.5	-
2-Methylnaphthalene	91-57-6	-	< 5.0	-
1-Methyl naphthalene *	90-12-0	-	< 5.0	-
Hexachlorocyclopentadiene	77-47-4	-	< 12.5	-
2,4,6-Trichlorophenol	88-06-2	-	< 50.0	-
2,4,5-Trichlorophenol	95-95-4	-	< 50.0	-
2-Chloronaphthalene	91-58-7	-	< 5.0	-
Biphenyl *	92-52-4	-	< 5.0	-
Diphenyl ether *	101-84-8	-	< 5.0	-
2-Nitroaniline	88-74-4	-	< 12.5	-
Acenaphthylene	208-96-8	-	< 5.0	-
Dimethylphthalate	131-11-3	-	< 12.5	-
2,6-Dinitrotoluene	606-20-2	-	< 12.5	-
Acenaphthene	83-32-9	-	< 5.0	-
3-Nitroaniline	99-09-2	-	< 12.5	-

* These compounds are not currently UKAS accredited.
Concentrations are reported on a wet weight basis.

Target Compounds	CAS #	R.T.	Concentration mg/kg	% Fit
2,4-Dinitrophenol	51-28-5	-	< 25.0	-
Dibenzofuran	132-64-9	-	< 12.5	-
4-Nitrophenol	100-02-7	-	< 12.5	-
2,4-Dinitrotoluene	121-14-2	-	< 12.5	-
Fluorene	86-73-7	-	< 5.0	-
Diethylphthalate	84-66-2	-	< 12.5	-
4-Chlorophenyl-phenylether	7005-72-4	-	< 12.5	-
4,6-Dinitro-2-methylphenol	534-52-1	-	< 12.5	-
4-Nitroaniline	100-01-6	-	< 12.5	-
N-Nitrosodiphenylamine	86-30-6	-	< 12.5	-
4-Bromophenyl-phenylether	101-55-3	-	< 12.5	-
Hexachlorobenzene	118-74-1	-	< 12.5	-
Pentachlorophenol	87-86-5	-	< 12.5	-
Phenanthrene	85-01-8	-	< 5.0	-
Anthracene	120-12-7	-	< 5.0	-
Di-n-butylphthalate	84-74-2	-	< 12.5	-
Fluoranthene	206-44-0	-	< 5.0	-
Pyrene	129-00-0	-	< 5.0	-
Butylbenzylphthalate	85-68-7	-	< 12.5	-
Benzol[a]anthracene	56-55-3	-	< 5.0	-
Chrysene	218-01-9	-	< 5.0	-
3,3'-Dichlorobenzidine	91-94-1	-	< 50.0	-
bis(2-Ethylhexyl)phthalate	117-81-7	-	< 12.5	-
Di-n-octylphthalate	117-84-0	-	< 5.0	-
Benzol[b]fluoranthene	205-99-2	-	< 5.0	-
Benzol[k]fluoranthene	207-08-9	-	< 5.0	-
Benzol[a]pyrene	50-32-8	-	< 5.0	-
Indeno[1,2,3-cd]pyrene	193-39-5	-	< 5.0	-
Dibenzol[a,h]anthracene	53-70-3	-	< 5.0	-
Benzol[g,h,i]perylene	191-24-2	-	< 5.0	-

* compounds not currently UKAS accredited

Internal Standards	% Area	Surrogates	% Rec
1,4-Dichlorobenzene-d4	115	2-Fluorophenol	93
Naphthalene-d8	102	Phenol-d5	83
Acenaphthene-d10	105	Nitrobenzene-d5	64
Phenanthrene-d10	111	2-Fluorobiphenyl	69
Chrysene-d12	125	2,4,6-Tribromophenol	54
Perylene-d12	103	Terphenyl-d14	79

Semi-Volatile Organic Compounds

Customer and Site Details:
Sample Details:
LIMS ID Number:
Job Number:

Exploration Associates : M4 Relief Road
CH/0042/0/20m
CL03519
484

11-Feb-00
16-Feb-00
18-Feb-00

Matrix:
Ext Method:
Operator:
Directory/Quant File:

Soil
Soxhlet
CS
0218ABN.MS4

QC Batch Number:
Multiplier:
Dilution Factor:
0218CCCC1.D GPC (Y/N)

219
0.0667
1
Y

Target Compounds	CAS #	R.T. (min)	Concentration mg/kg	% Fit
Phenol	108-95-2	-	< 2.0	-
bis(2-Chloroethyl)ether	111-44-4	-	< 0.5	-
2-Chlorophenol	95-57-8	-	< 2.0	-
1,3-Dichlorobenzene	541-73-1	-	< 0.5	-
1,4-Dichlorobenzene	106-46-7	-	< 0.5	-
Benzyl alcohol	100-51-6	-	< 0.5	-
1,2-Dichlorobenzene	95-50-1	-	< 0.5	-
2-Methylphenol	95-48-7	-	< 0.5	-
bis(2-Chloroisopropyl)ether	108-60-1	-	< 0.5	-
Hexachloroethane	67-72-1	-	< 0.5	-
N-Nitroso-di-n-propylamine	621-64-7	-	< 0.5	-
3- & 4-Methylphenol	108-39-4/106-44-5	-	< 2.0	-
Nitrobenzene	98-95-3	-	< 0.5	-
Isophorone	78-59-1	-	< 0.5	-
2-Nitrophenol	88-75-5	-	< 2.0	-
2,4-Dimethylphenol	105-67-9	-	< 2.0	-
Benzoic Acid	65-85-0	-	< 10.0	-
bis(2-Chloroethoxy)methane	111-91-1	-	< 0.5	-
2,4-Dichlorophenol	120-83-2	-	< 2.0	-
1,2,4-Trichlorobenzene	120-82-1	-	< 0.5	-
Naphthalene	91-20-3	-	< 0.2	-
4-Chlorophenol *	106-48-9	-	< 2.0	-
4-Chloroaniline *	106-47-8	-	< 0.5	-
Hexachlorobutadiene	87-68-3	-	< 0.5	-
4-Chloro-3-methylphenol	59-50-7	-	< 0.5	-
2-Methylnaphthalene	91-57-6	-	< 0.2	-
1-Methyl naphthalene *	90-12-0	-	< 0.2	-
Hexachlorocyclopentadiene	77-47-4	-	< 0.5	-
2,4,6-Trichlorophenol	88-06-2	-	< 2.0	-
2,4,5-Trichlorophenol	95-95-4	-	< 2.0	-
2-Chloronaphthalene	91-58-7	-	< 0.2	-
Biphenyl *	92-52-4	15.18	0.2	99
Diphenyl ether *	101-84-8	-	< 0.2	-
2-Nitroaniline	88-74-4	-	< 0.5	-
Acenaphthylene	208-96-8	-	< 0.2	-
Dimethylphthalate	131-11-3	-	< 0.5	-
2,6-Dinitrotoluene	606-20-2	-	< 0.5	-
Acenaphthene	83-32-9	-	< 0.2	-
3-Nitroaniline *	99-09-2	-	< 0.5	-

* These compounds are not currently UKAS accredited.
Concentrations are reported on a wet weight basis.

Target Compounds	CAS #	R.T.	Concentration mg/kg	% Fit
2,4-Dinitrophenol	51-28-5	-	< 1.0	-
Dibenzofuran	132-64-9	-	< 0.5	-
4-Nitrophenol	100-02-7	-	< 5.0	-
2,4-Dinitrotoluene	121-14-2	-	< 0.5	-
Fluorene	86-73-7	-	< 0.2	-
Diethylphthalate	84-66-2	-	< 0.5	-
4-Chlorophenyl-phenylether	7005-72-4	-	< 0.5	-
4,6-Dinitro-2-methylphenol	534-52-1	-	< 5.0	-
4-Nitroaniline	100-01-6	-	< 0.5	-
N-Nitrosodiphenylamine	86-30-6	-	< 0.5	-
4-Bromophenyl-phenylether	101-55-3	-	< 0.5	-
Hexachlorobenzene	118-74-1	-	< 0.5	-
Pentachlorophenol	87-86-5	-	< 5.0	-
Phenanthrene	85-01-8	21.18	0.3	100
Anthracene	120-12-7	-	< 0.2	-
Di-n-butylphthalate	84-74-2	-	< 0.5	-
Fluoranthene	206-44-0	24.71	0.8	97
Pyrene	129-00-0	25.33	0.7	99
Butylbenzylphthalate	85-68-7	-	< 0.5	-
Benzofluoranthene	56-55-3	28.95	0.3	96
Chrysene	218-01-9	29.06	0.8	97
3,3'-Dichlorobenzidine	91-94-1	-	< 2.0	-
bis(2-Ethylhexyl)phthalate	117-81-7	29.70	1.8	99
Di-n-octylphthalate	117-84-0	-	< 0.2	-
Benzofluoranthene	205-99-2	32.01	0.6	97
Benzofluoranthene	207-08-9	32.06	0.5	97
Benzofluoranthene	50-32-8	32.96	0.6	98
Indeno[1,2,3-cd]pyrene	193-39-5	37.38	0.5	97
Dibenzofluoranthene	53-70-3	-	< 0.2	-
Benzofluoranthene	191-24-2	38.32	0.4	100

* compounds not currently UKAS accredited

Internal Standards	% Area
1,4-Dichlorobenzene-d4	98
Naphthalene-d8	83
Acenaphthene-d10	94
Phenanthrene-d10	98
Chrysene-d12	116
Perylene-d12	88

Surrogates	% Rec
2-Fluorophenol	67
Phenol-d5	69
Nitrobenzene-d5	50
2-Fluorobiphenyl	58
2,4,6-Tribromophenol	72
Terphenyl-d14	80

Semi-Volatile Organic Compounds

Customer and Site Details:
Sample Details:
LIMS ID Number:
Job Number:

Exploration Associates : M4 Relief Road
CH/0042/1.20m
CL03520
484

11-Feb-00
16-Feb-00
18-Feb-00

Matrix:
Ext Method:
Operator:
Directory/Quant File:

Soil
Soxhlet
CS
0218ABN.MS4 0218CCCC2.D GPC (Y/N)

QC Batch Number:
Multiplier:
Dilution Factor:
219
0.0667
1
Y

Target Compounds	CAS #	R.T. (min)	Concentration mg/kg	% Fit
Phenol	108-95-2	-	< 2.0	-
bis(2-Chloroethyl)ether	111-44-4	-	< 0.5	-
2-Chlorophenol	95-57-8	-	< 2.0	-
1,3-Dichlorobenzene	541-73-1	-	< 0.5	-
1,4-Dichlorobenzene	106-46-7	-	< 0.5	-
Benzyl alcohol	100-51-6	-	< 0.5	-
1,2-Dichlorobenzene	95-50-1	-	< 0.5	-
2-Methylphenol	95-48-7	-	< 0.5	-
bis(2-Chloroisopropyl)ether	108-60-1	-	< 0.5	-
Hexachloroethane	67-72-1	-	< 0.5	-
N-Nitroso-di-n-propylamine	621-64-7	-	< 0.5	-
3- & 4-Methylphenol	108-39-4/106-44-5	-	< 2.0	-
Nitrobenzene	98-95-3	-	< 0.5	-
Isophorone	78-59-1	-	< 0.5	-
2-Nitrophenol	88-75-5	-	< 2.0	-
2,4-Dimethylphenol	105-67-9	-	< 2.0	-
Benzoic Acid	65-85-0	-	< 10.0	-
bis(2-Chloroethoxy)methane	111-91-1	-	< 0.5	-
2,4-Dichlorophenol	120-83-2	-	< 2.0	-
1,2,4-Trichlorobenzene	120-82-1	-	< 0.5	-
Naphthalene	91-20-3	-	< 2.0	-
4-Chlorophenol *	106-48-9	-	< 0.5	-
4-Chloroaniline *	106-47-8	-	< 0.5	-
Hexachlorobutadiene	87-68-3	-	< 0.5	-
4-Chloro-3-methylphenol	59-50-7	-	< 0.5	-
2-Methylnaphthalene	91-57-6	-	< 0.2	-
1-Methyl naphthalene *	90-12-0	-	< 0.2	-
Hexachlorocyclopentadiene	77-47-4	-	< 0.5	-
2,4,6-Trichlorophenol	88-06-2	-	< 2.0	-
95-95-4	95-95-4	-	< 2.0	-
2,4,5-Trichlorophenol	91-58-7	-	< 0.2	-
2-Chloronaphthalene	92-52-4	15.18	0.6	99
Biphenyl *	101-84-8	-	< 0.2	-
Diphenyl ether *	88-74-4	-	< 0.5	-
2-Nitroaniline	208-96-8	-	< 0.2	-
Acenaphthylene	131-11-3	-	< 0.5	-
Dimethylphthalate	606-20-2	-	< 0.5	-
2,6-Dinitrotoluene	83-32-9	-	< 0.2	-
Acenaphthene	99-09-2	-	< 0.5	-
3-Nitroaniline *				

* These compounds are not currently UKAS accredited.
Concentrations are reported on a wet weight basis.

Target Compounds	CAS #	R.T.	Concentration mg/kg	% Fit
2,4-Dinitrophenol	51-28-5	-	< 1.0	-
Dibenzofuran	132-64-9	-	< 0.5	-
4-Nitrophenol	100-02-7	-	< 0.5	-
2,4-Dinitrotoluene	121-14-2	-	< 0.5	-
Fluorene	86-73-7	-	< 0.2	-
Diethylphthalate	84-66-2	-	< 0.5	-
4-Chlorophenyl-phenylether	7005-72-4	-	< 0.5	-
4,6-Dinitro-2-methylphenol	534-52-1	-	< 5.0	-
4-Nitroaniline	100-01-6	-	< 0.5	-
N-Nitrosodiphenylamine	86-30-6	-	< 0.5	-
4-Bromophenyl-phenylether	101-55-3	-	< 0.5	-
Hexachlorobenzene	118-74-1	-	< 0.5	-
Pentachlorophenol	87-86-5	-	< 5.0	-
Phenanthrene	85-01-8	-	< 0.2	-
Anthracene	120-12-7	-	< 0.2	-
Di-n-butylphthalate	84-74-2	-	< 0.5	-
Fluoranthene	206-44-0	-	< 0.2	-
Pyrene	129-00-0	-	< 0.2	-
Butylbenzylphthalate	85-68-7	-	< 0.5	-
Benzol[a]anthracene	56-55-3	-	< 0.2	-
Chrysene	218-01-9	29.08	1.6	98
3,3-Dichlorobenzidine	91-94-1	-	< 2.0	-
bis(2-Ethylhexyl)phthalate	117-81-7	29.71	0.9	100
Di-n-octylphthalate	117-84-0	-	< 0.2	-
Benzol[b]fluoranthene	205-99-2	32.02	0.2	90
Benzol[k]fluoranthene	207-08-9	-	< 0.2	-
Benzol[a]pyrene	50-32-8	32.97	0.2	94
Indeno[1,2,3-cd]pyrene	193-39-5	-	< 0.2	-
Dibenzol[a,h]anthracene	53-70-3	-	< 0.2	-
Benzol[g,h,i]perylene	191-24-2	-	< 0.2	-

* compounds not currently UKAS accredited

Internal Standards	% Area	Surrogates	% Rec
1,4-Dichlorobenzene-d4	90	2-Fluorophenol	66
Naphthalene-d8	80	Phenol-d5	51
Acenaphthene-d10	86	Nitrobenzene-d5	49
Phenanthrene-d10	85	2-Fluorobiphenyl	59
Chrysene-d12	98	2,4,6-Tribromophenol	72
Perylene-d12	82	Terphenyl-d14	52

Volatile Organic Compounds by PTGCMS

Customer and Site Details: Exploration Associates : M4 Relief Road
Sample Details: CH/0035/ / 0.00-
LIMS ID Number: CL03504
Report Number: 484

Directory/Quant file: 0215VOC.MS\1 0215CCCC4.D
Date Booked in: 11-Feb-00
Date Analysed: 15-Feb-00
Operator: AB
Matrix: Soil
Method: Purge and Trap
Dilution: 5
Position: 9

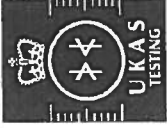
Target Compounds	CAS #	R.T. (min.)	Concentration µg/kg	% Fit
Dichlorodifluoromethane	75-71-8	-	< 5	-
Chloromethane	74-87-3	-	< 5	-
Vinyl Chloride	75-01-4	-	< 5	-
Bromomethane	74-83-9	-	< 5	-
Chloroethane	75-00-3	-	< 5	-
Trichlorofluoromethane	75-69-4	-	< 5	-
1,1-Dichloroethene	75-35-4	-	< 5	-
trans 1,2-Dichloroethene	156-60-5	-	< 5	-
1,1-Dichloroethane	75-34-3	-	< 5	-
2,2-Dichloropropane	594-20-7	-	< 5	-
cis 1,2-Dichloroethene	156-59-2	-	< 5	-
Bromochloromethane	74-97-5	-	< 5	-
Chloroform	67-66-3	-	< 5	-
1,1,1-Trichloroethane	71-55-6	-	< 5	-
Carbon Tetrachloride	56-23-5	-	< 5	-
1,1-Dichloropropene	563-58-6	-	< 5	-
Benzene	71-43-2	-	< 5	-
1,2-Dichloroethane	107-06-2	-	< 5	-
Trichloroethene	79-01-6	-	< 7	-
1,2-Dichloropropene	78-87-5	-	< 5	-
Dibromomethane	74-95-3	-	< 5	-
Bromodichloromethane	75-27-4	-	< 5	-
cis 1,3-Dichloropropene	10061-01-5	-	< 5	-
Toluene	108-88-3	-	< 5	-
trans 1,3-Dichloropropene	10061-02-6	-	< 5	-
1,1,2-Trichloroethane	79-00-5	-	< 5	-
Tetrachloroethene	127-18-4	-	< 5	-
1,3-Dichloropropene	142-28-29	-	< 5	-
Dibromochloromethane	124-48-1	-	< 5	-
1,2-Dibromoethane	106-93-4	-	< 5	-
Chlorobenzene	108-90-7	-	< 5	-
Ethylbenzene	100-41-4	-	< 5	-
1,1,1,2-Tetrachloroethane	630-20-6	-	< 5	-
m and p-Xylene	108-38-3/106-42-3	-	< 5	-
o-Xylene	95-47-6	-	< 5	-

Concentrations are reported on a wet weight basis

Target Compounds	CAS #	R.T. (min.)	Concentration µg/kg	% Fit
Styrene	100-42-5	-	< 5	-
Bromoform	75-25-2	-	< 5	-
iso-Propylbenzene	98-82-8	-	< 5	-
1,1,2,2-Tetrachloroethane	79-34-5	-	< 5	-
Propylbenzene	103-65-1	-	< 5	-
Bromobenzene	108-86-1	-	< 5	-
1,2,3-Trichloropropane	96-18-4	-	< 5	-
2-Chlorotoluene	95-49-8	-	< 5	-
1,3,5-Trimethylbenzene	108-67-8	-	< 5	-
4-Chlorotoluene	106-43-4	-	< 5	-
tert-Butylbenzene	98-06-6	-	< 5	-
1,2,4-Trimethylbenzene	95-63-6	-	< 5	-
sec-Butylbenzene	135-98-8	-	< 5	-
p-Isopropyltoluene	99-87-6	-	< 5	-
1,3-Dichlorobenzene	541-73-1	-	< 5	-
1,4-Dichlorobenzene	106-46-7	-	< 5	-
n-Butylbenzene	104-51-8	-	< 5	-
1,2-Dichlorobenzene	95-50-1	-	< 5	-
1,2-Dibromo-3-chloropropane	96-12-8	-	< 25	-
1,2,4-Trichlorobenzene	120-82-1	-	< 25	-
Hexachlorobutadiene	87-68-3	-	< 25	-
Naphthalene	91-20-3	-	< 25	-
1,2,3-Trichlorobenzene	87-61-6	-	< 25	-

Concentrations are reported on a wet weight basis

Internal standards	R.T.	Area %	Surrogates	% Rec
Pentafluorobenzene	12.24	111	Dibromofluoromethane	89
1,4-Difluorobenzene	13.39	102	Toluene-d8	102
Chlorobenzene-d5	17.69	97	Bromofluorobenzene	97
1,4-Dichlorobenzene-d4	21.15	93		



Volatile Organic Compounds by PTGCMS

Customer and Site Details: Exploration Associates : M4 Relief Road
Sample Details: CH/0035/ / 0.50-
LIMS ID Number: CL03505
Report Number: 484

Directory/Quant file: 0215VOC.MS1\
Date Booked in: 11-Feb-00
Date Analysed: 15-Feb-00
Operator: AB

Matrix: Soil
Method: Purge and Trap
Dilution: 5
Position: 10

1252

Target Compounds	CAS #	R.T. (min.)	Concentration µg/kg	% Fit
Dichlorodifluoromethane	75-71-8	-	< 5	-
Chloromethane	74-87-3	-	< 5	-
Vinyl Chloride	75-01-4	-	< 5	-
Bromomethane	74-83-9	-	< 5	-
Chloroethane	75-00-3	-	< 5	-
Trichlorofluoromethane	75-69-4	-	< 5	-
1,1-Dichloroethene	75-35-4	-	< 5	-
trans 1,2-Dichloroethene	156-60-5	-	< 5	-
1,1-Dichloroethane	75-34-3	-	< 5	-
2,2-Dichloropropane	594-20-7	-	< 5	-
cis 1,2-Dichloroethene	156-59-2	-	< 5	-
Bromochloromethane	74-97-5	-	< 5	-
Chloroform	67-66-3	-	< 5	-
1,1,1-Trichloroethane	71-55-6	-	< 5	-
Carbon Tetrachloride	56-23-5	-	< 5	-
1,1-Dichloropropene	563-58-6	-	< 5	-
Benzene	71-43-2	12.91	168	95
1,2-Dichloroethane	107-06-2	-	< 5	-
Trichloroethene	79-01-6	-	< 5	-
1,2-Dichloropropane	78-87-5	-	< 5	-
Dibromomethane	74-95-3	-	< 5	-
Bromodichloromethane	75-27-4	-	< 5	-
cis 1,3-Dichloropropene	10061-01-5	-	< 5	-
Toluene	108-88-3	15.66	23	98
trans 1,3-Dichloropropene	10061-02-6	-	< 5	-
1,1,2-Trichloroethane	79-00-5	-	< 5	-
Tetrachloroethene	127-18-4	-	< 5	-
1,3-Dichloropropane	142-28-29	-	< 5	-
Dibromochloromethane	124-48-1	-	< 5	-
1,2-Dibromoethane	106-93-4	-	< 5	-
Chlorobenzene	108-90-7	17.74	56	91
Ethylbenzene	100-41-4	17.82	571	97
1,1,1,2-Tetrachloroethane	630-20-6	-	< 5	-
m and p-Xylene	108-38-3/106-42-3	17.98	2,390	79
o-Xylene	95-47-6	18.6	496	92

Concentrations are reported on a wet weight basis

Target Compounds	CAS #	R.T. (min.)	Concentration µg/kg	% Fit
Styrene	100-42-5	-	< 5	-
Bromoform	75-25-2	-	< 5	-
iso-Propylbenzene	98-82-8	-	< 5	-
1,1,2,2-Tetrachloroethane	79-34-5	-	< 5	-
Propylbenzene	103-65-1	-	< 5	-
Bromobenzene	108-86-1	-	< 5	-
1,2,3-Trichloropropane	96-18-4	-	< 5	-
2-Chlorotoluene	95-49-8	-	< 5	-
1,3,5-Trimethylbenzene	108-67-8	-	< 5	-
4-Chlorotoluene	106-43-4	-	< 5	-
tert-Butylbenzene	98-06-6	-	< 5	-
1,2,4-Trimethylbenzene	95-63-6	-	< 5	-
sec-Butylbenzene	135-98-8	-	< 5	-
p-Isopropyltoluene	99-87-6	-	< 5	-
1,3-Dichlorobenzene	541-73-1	21.08	39	95
1,4-Dichlorobenzene	106-46-7	21.20	594	97
n-Butylbenzene	104-51-8	-	< 5	-
1,2-Dichlorobenzene	95-50-1	21.79	174	96
1,2-Dibromo-3-chloropropane	96-12-8	-	< 25	-
1,2,4-Trichlorobenzene	120-82-1	24.31	64	99
Hexachlorobutadiene	87-68-3	-	< 25	-
Naphthalene	91-20-3	-	< 25	-
1,2,3-Trichlorobenzene	87-61-6	-	< 25	-

Concentrations are reported on a wet weight basis

Internal standards	R.T.	Area %	Surrogates	% Rec
Pentafluorobenzene	12.25	101	Dibromofluoromethane	91
1,4-Difluorobenzene	13.39	99	Toluene-d8	100
Chlorobenzene-d5	17.70	89	Bromofluorobenzene	88
1,4-Dichlorobenzene-d4	21.16	67		

Volatile Organic Compounds by PTGCMS

Customer and Site Details: Exploration Associates : M4 Relief Road
Sample Details: CH/0035/ / 0.80-
LIMS ID Number: CL03506
Report Number: 484

Directory/Quant file: 0215VOC.MS1\ 0215CCCC4.D
Date Booked in: 11-Feb-00
Date Analysed: 15-Feb-00
Operator: AB

Matrix: Soil
Method: Purge and Trap
Dilution: 5
Position: 11

Target Compounds	CAS #	R.T. (min.)	Concentration µg/kg	% Fit
Dichlorodifluoromethane	75-71-8	-	< 5	-
Chloromethane	74-87-3	-	< 5	-
Vinyl Chloride	75-01-4	-	< 5	-
Bromomethane	74-83-9	-	< 5	-
Chloroethane	75-00-3	-	< 5	-
Trichlorofluoromethane	75-69-4	-	< 5	-
1,1-Dichloroethene	75-35-4	-	< 5	-
trans 1,2-Dichloroethene	156-60-5	-	< 5	-
1,1-Dichloroethane	75-34-3	-	< 5	-
2,2-Dichloropropane	594-20-7	-	< 5	-
cis 1,2-Dichloroethene	156-59-2	-	< 5	-
Bromochloromethane	74-97-5	-	< 5	-
Chloroform	67-66-3	-	< 5	-
1,1,1-Trichloroethane	71-55-6	-	< 5	-
Carbon Tetrachloride	56-23-5	-	< 5	-
1,1-Dichloropropene	563-58-6	-	< 5	-
Benzene	71-43-2	12.92	96	95
1,2-Dichloroethane	107-06-2	-	< 5	-
Trichloroethene	79-01-6	-	< 5	-
1,2-Dichloropropane	78-87-5	-	< 5	-
Dibromomethane	74-95-3	-	< 5	-
Bromodichloromethane	75-27-4	-	< 5	-
cis 1,3-Dichloropropene	10061-01-5	-	< 5	-
Toluene	108-88-3	15.66	10	96
trans 1,3-Dichloropropene	10061-02-6	-	< 5	-
1,1,2-Trichloroethane	79-00-5	-	< 5	-
Tetrachloroethene	127-18-4	-	< 5	-
1,3-Dichloropropane	142-28-29	-	< 5	-
Dibromochloromethane	124-48-1	-	< 5	-
1,2-Dibromoethane	106-93-4	-	< 5	-
Chlorobenzene	108-90-7	17.75	17	79
Ethylbenzene	100-41-4	17.83	244	96
1,1,1,2-Tetrachloroethane	630-20-6	-	< 5	-
m and p-Xylene	108-38-3/106-42-3	17.98	988	93
o-Xylene	95-47-6	18.6	213	90

Concentrations are reported on a wet weight basis

Target Compounds	CAS #	R.T. (min.)	Concentration µg/kg	% Fit
Styrene	100-42-5	-	< 5	-
Bromoform	75-25-2	-	< 5	-
iso-Propylbenzene	98-82-8	-	< 5	-
1,1,2,2-Tetrachloroethane	79-34-5	-	< 5	-
Propylbenzene	103-65-1	-	< 5	-
Bromobenzene	108-86-1	-	< 5	-
1,2,3-Trichloropropane	96-18-4	-	< 5	-
2-Chlorotoluene	95-49-8	-	< 5	-
1,3,5-Trimethylbenzene	108-67-8	-	< 5	-
4-Chlorotoluene	106-43-4	-	< 5	-
tert-Butylbenzene	98-06-6	-	< 5	-
1,2,4-Trimethylbenzene	95-63-6	-	< 5	-
sec-Butylbenzene	135-98-8	-	< 5	-
p-Isopropyltoluene	99-87-6	-	< 5	-
1,3-Dichlorobenzene	541-73-1	21.07	6	94
1,4-Dichlorobenzene	106-46-7	21.20	165	92
n-Butylbenzene	104-51-8	-	< 5	-
1,2-Dichlorobenzene	95-50-1	21.80	36	94
1,2-Dibromo-3-chloropropane	96-12-8	-	< 25	-
1,2,4-Trichlorobenzene	120-82-1	24.32	25	96
Hexachlorobutadiene	87-68-3	-	< 25	-
Naphthalene	91-20-3	-	< 25	-
1,2,3-Trichlorobenzene	87-61-6	-	< 25	-

Concentrations are reported on a wet weight basis

Internal standards	R.T.	Area %	Surrogates	% Rec
Pentafluorobenzene	12.24	79	Dibromofluoromethane	102
1,4-Difluorobenzene	13.40	79	Toluene-d8	101
Chlorobenzene-d5	17.69	76	Bromofluorobenzene	100
1,4-Dichlorobenzene-d4	21.17	67		

Volatile Organic Compounds by PTGCMS

Customer and Site Details: Exploration Associates : M4 Relief Road
Sample Details: CH/0041/ / 1.40-
LIMS ID Number: CL03518
Report Number: 484

Directory/Quant file:
Date Booked in:
Date Analysed:
Operator:

0215VOC.MS\1
11/2020
15-Feb-00
AB

Matrix:
Method:
Dilution:
Position:

Soil
Purge and Trap
5
12

Target Compounds	CAS #	R.T. (min.)	Concentration µg/kg	% Fit
Dichlorodifluoromethane	75-71-8	-	< 5	-
Chloromethane	74-87-3	-	< 5	-
Vinyl Chloride	75-01-4	-	< 5	-
Bromomethane	74-83-9	-	< 5	-
Chloroethane	75-00-3	-	< 5	-
Trichlorofluoromethane	75-69-4	-	< 5	-
1,1-Dichloroethene	75-35-4	-	< 5	-
trans 1,2-Dichloroethene	156-60-5	-	< 5	-
1,1-Dichloroethane	75-34-3	-	< 5	-
2,2-Dichloropropane	594-20-7	-	< 5	-
cis 1,2-Dichloroethene	156-59-2	-	< 5	-
Bromochloromethane	74-97-5	-	< 5	-
Chloroform	67-66-3	-	< 5	-
1,1,1-Trichloroethane	71-55-6	-	< 5	-
Carbon Tetrachloride	56-23-5	-	< 5	-
1,1-Dichloropropene	563-58-6	-	< 5	-
Benzene	71-43-2	-	< 5	-
1,2-Dichloroethane	107-06-2	-	< 5	-
Trichloroethene	79-01-6	-	< 5	-
1,2-Dichloropropane	78-87-5	-	< 5	-
Dibromomethane	74-95-3	-	< 5	-
Bromodichloromethane	75-27-4	-	< 5	-
cis 1,3-Dichloropropene	10061-01-5	-	< 5	-
Toluene	108-88-3	-	< 5	-
trans 1,3-Dichloropropene	10061-02-6	-	< 5	-
1,1,2-Trichloroethane	79-00-5	-	< 5	-
Tetrachloroethene	127-18-4	-	< 5	-
1,3-Dichloropropane	142-28-29	-	< 5	-
Dibromochloromethane	124-48-1	-	< 5	-
1,2-Dibromoethane	106-93-4	-	< 5	-
Chlorobenzene	108-90-7	-	< 5	-
Ethylbenzene	100-41-4	17.82	9	92
1,1,1,2-Tetrachloroethane	630-20-6	-	< 5	-
m and p-Xylene	108-38-3/106-42-3	17.98	12	86
o-Xylene	95-47-6	-	< 5	-

Concentrations are reported on a wet weight basis

Target Compounds	CAS #	R.T. (min.)	Concentration µg/kg	% Fit
Styrene	100-42-5	18.61	20	93
Bromoform	75-25-2	-	< 5	-
iso-Propylbenzene	98-82-8	-	< 5	-
1,1,2,2-Tetrachloroethane	79-34-5	-	< 5	-
Propylbenzene	103-65-1	-	< 5	-
Bromobenzene	108-86-1	-	< 5	-
1,2,3-Trichloropropane	96-18-4	-	< 5	-
2-Chlorotoluene	95-49-8	-	< 5	-
1,3,5-Trimethylbenzene	108-67-8	-	< 5	-
4-Chlorotoluene	106-43-4	-	< 5	-
tert-Butylbenzene	98-06-6	-	< 5	-
1,2,4-Trimethylbenzene	95-63-6	-	< 5	-
sec-Butylbenzene	135-98-8	-	< 5	-
p-Isopropyltoluene	99-87-6	-	< 5	-
1,3-Dichlorobenzene	541-73-1	21.07	6	76
1,4-Dichlorobenzene	106-46-7	21.20	46	76
n-Butylbenzene	104-51-8	-	< 5	-
1,2-Dichlorobenzene	95-50-1	21.79	5	82
1,2-Dibromo-3-chloropropane	96-12-8	-	< 25	-
1,2,4-Trichlorobenzene	120-82-1	24.32	72	96
Hexachlorobutadiene	87-68-3	-	< 25	-
Naphthalene	91-20-3	-	< 25	-
1,2,3-Trichlorobenzene	87-61-6	-	< 25	-

Concentrations are reported on a wet weight basis

Internal standards	R.T.	Area %	Surrogates	% Rec
Pentafluorobenzene	12.24	60	Dibromofluoromethane	114
1,4-Difluorobenzene	13.40	61	Toluene-d8	96
Chlorobenzene-d5	17.70	54	Bromofluorobenzene	101
1,4-Dichlorobenzene-d4	21.16	39		

Customer and Site Details: Exploration Associates: M4 Relief Road
Sample Details: CH/0042/ / 0.20-
LIMS ID Number: CL03519
Report Number: 484

Volatile Organic Compounds by PTGCMS

Directory/Quant file: 0217VOC.MS1\ 0217CCCC1.D
Date Booked in: 11-Feb-00
Date Analysed: 17-Feb-00
Operator: at

Matrix: Soil
Method: Purge and Trap
Dilution: 5
Position: 9

Target Compounds	CAS #	R.T. (min.)	Concentration µg/kg	% Fit
Dichlorodifluoromethane	75-71-8	-	< 5	-
Chloromethane	74-87-3	-	< 5	-
Vinyl Chloride	75-01-4	-	< 5	-
Bromomethane	74-83-9	-	< 5	-
Chloroethane	75-00-3	-	< 5	-
Trichlorofluoromethane	75-69-4	-	< 5	-
1,1-Dichloroethene	75-35-4	-	< 5	-
trans 1,2-Dichloroethene	156-60-5	-	< 5	-
1,1-Dichloroethane	75-34-3	-	< 5	-
2,2-Dichloropropane	594-20-7	-	< 5	-
cis 1,2-Dichloroethene	156-59-2	-	< 5	-
Bromochloromethane	74-97-5	-	< 5	-
Chloroform	67-66-3	-	< 5	-
1,1,1-Trichloroethane	71-55-6	-	< 5	-
Carbon Tetrachloride	56-23-5	-	< 5	-
1,1-Dichloropropene	563-58-6	-	< 5	-
Benzene	71-43-2	-	< 5	-
1,2-Dichloroethane	107-06-2	-	< 5	-
Trichloroethene	79-01-6	-	< 5	-
1,2-Dichloropropane	78-87-5	-	< 5	-
Dibromomethane	74-95-3	-	< 5	-
Bromodichloromethane	75-27-4	-	< 5	-
cis 1,3-Dichloropropene	10061-01-5	-	< 5	-
Toluene	108-88-3	-	< 5	-
trans 1,3-Dichloropropene	10061-02-6	-	< 5	-
1,1,2-Trichloroethane	79-00-5	-	< 5	-
Tetrachloroethene	127-18-4	-	< 5	-
1,3-Dichloropropane	142-28-29	-	< 5	-
Dibromochloromethane	124-48-1	-	< 5	-
1,2-Dibromoethane	106-93-4	-	< 5	-
Chlorobenzene	108-90-7	-	< 5	-
Ethylbenzene	100-41-4	-	< 5	-
1,1,2-Tetrachloroethane	630-20-6	-	< 5	-
m and p-Xylene	108-38-3/106-42-3	-	< 5	-
o-Xylene	95-47-6	-	< 5	-

Target Compounds	CAS #	R.T. (min.)	Concentration µg/kg	% Fit
Styrene	100-42-5	-	< 5	-
Bromoform	75-25-2	-	< 5	-
iso-Propylbenzene	98-82-8	-	< 5	-
1,1,2,2-Tetrachloroethane	79-34-5	-	< 5	-
Propylbenzene	103-65-1	-	< 5	-
Bromobenzene	108-86-1	-	< 5	-
1,2,3-Trichloropropane	96-18-4	-	< 5	-
2-Chlorotoluene	95-49-8	-	< 5	-
1,3,5-Trimethylbenzene	108-67-8	-	< 5	-
4-Chlorotoluene	106-43-4	-	< 5	-
tert-Butylbenzene	98-06-6	-	< 5	-
1,2,4-Trimethylbenzene	95-63-6	-	< 5	-
sec-Butylbenzene	135-98-8	-	< 5	-
p-Isopropyltoluene	99-87-6	-	< 5	-
1,3-Dichlorobenzene	541-73-1	-	< 5	-
1,4-Dichlorobenzene	106-46-7	-	< 5	-
n-Butylbenzene	104-51-8	-	< 5	-
1,2-Dichlorobenzene	95-50-1	-	< 5	-
1,2-Dibromo-3-chloropropane	96-12-8	-	< 25	-
1,2,4-Trichlorobenzene	120-82-1	-	< 25	-
Hexachlorobutadiene	87-68-3	-	< 25	-
Naphthalene	91-20-3	-	< 25	-
1,2,3-Trichlorobenzene	87-61-6	-	< 25	-

Concentrations are reported on a wet weight basis

Internal standards	R.T.	Area %	Surrogates	% Rec
Pentafluorobenzene	12.18	69	Dibromofluoromethane	129
1,4-Difluorobenzene	13.34	71	Toluene-d8	103
Chlorobenzene-d5	17.65	79	Bromofluorobenzene	113
1,4-Dichlorobenzene-d4	21.12	77		

Concentrations are reported on a wet weight basis



TEST REPORT SOIL SAMPLE ANALYSIS



1252
1411

TES Report No. CL/000542

Site: M4 Relief Road

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
Mid Glamorgan
CF31 1JZ

The 29 samples described in this report were scheduled for analysis by TES Bretby on Wednesday, 16 February 2000. The analysis was completed by Thursday, 2 March 2000.

Tests marked as 'not UKAS accredited' and any opinions or interpretations expressed herein are outside the scope of any UKAS accreditation held by TES Bretby laboratories.

The following tables are contained in this report:

Table 1 Sample Descriptions
Table 2 Main Analysis Results
Table of Volatile Organic Compounds (1 Page)

On behalf of
TES Bretby : J. Hannah
J Hannah Project Co-ordinator

Date of Issue: 02/03/00

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TEST REPORT

SOIL SAMPLE ANALYSIS



1252
1411

Exploration Associates
Unit 15
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Mid Glamorgan
CF31 1JZ

TES Report No. 000542

Site: M4 Relief Road

Page No	Assessed Area	ID No. EFS/CL	Sample	Depth (m) from to	Description
1	M4 Relief Road	0003912	CH/ 8/	1.00	CH8 1.0
1	M4 Relief Road	0003913	CH/ 9/	0.00	CH9 0.0
1	M4 Relief Road	0003914	CH/ 9/	1.50	CH9 1.5
1	M4 Relief Road	0003915	CH/ 10/	0.40	CH10 0.4
2	M4 Relief Road	0003916	CH/ 10/	1.40	CH10 1.4
2	M4 Relief Road	0003917	CH/ 11/	0.35	CH11 0.35
2	M4 Relief Road	0003918	CH/ 11/	1.65	CH11 1.65
2	M4 Relief Road	0003919	CH/ 12/	0.17	CH12 0.17
3	M4 Relief Road	0003920	CH/ 12/	0.67	CH12 0.67
3	M4 Relief Road	0003921	CH/ 13/	0.00	CH13 0.0
3	M4 Relief Road	0003922	CH/ 13/	0.77	CH13 0.77
3	M4 Relief Road	0003923	CH/ 14/	0.16	CH14 0.16
4	M4 Relief Road	0003924	CH/ 15/	0.30	CH15 0.3
4	M4 Relief Road	0003925	CH/ 15/A	0.17	CH15A 0.17
4	M4 Relief Road	0003926	CH/ 16/	0.20	CH16 0.2
4	M4 Relief Road	0003927	CH/ 16/	1.90	CH16 1.9
5	M4 Relief Road	0003928	CH/ 17/	0.60	CH17 0.6
5	M4 Relief Road	0003929	CH/ 17/	1.30	CH17 1.3
5	M4 Relief Road	0003930	CH/ 17/	1.90	CH17 1.9
5	M4 Relief Road	0003931	CH/ 18/	0.20	CH18 0.2

Date of Issue: 02/03/00

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TEST REPORT SOIL SAMPLE ANALYSIS



1252
1411

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
Mid Glamorgan
CF31 1JZ

TES Report No. 000542

Site: M4 Relief Road

Page No	Assessed Area	ID No. EFS/CL	Sample	Depth(m)		Description
				from	to	
6	M4 Relief Road	0003932	CH/ 18/	1.60		CH18 1.6
6	M4 Relief Road	0003933	CH/ 18/	4.00		CH18 4.0
6	M4 Relief Road	0003934	CH/ 19/	0.20		CH19 0.2
6	M4 Relief Road	0003935	CH/ 19/	2.00		CH19 2.0
7	M4 Relief Road	0003936	CH/ 19/A	0.30		CH19A 0.3
7	M4 Relief Road	0003937	CH/ 19/A	1.00		CH19A 1.0
7	M4 Relief Road	0003938	CH/ 24/	0.00		CH24 0.0
7	M4 Relief Road	0003939	CH/ 24/	0.75		CH24 0.75
8	M4 Relief Road	0003940	CH/ 25/	0.00		CH25 0.0

Date of Issue: 02/03/00

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TEST REPORT

SOIL SAMPLE ANALYSIS



1252
1411

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
Mid Glamorgan
CF31 1JZ

TES Report No. 000542

Site: M4 Relief Road

Customer reference	CH 8 1.00	CH 9 0.00	CH 9 1.50	CH 10 0.40
Depth (m)				
Date logged	16/02/00	16/02/00	16/02/00	16/02/00
TES Bretby ID Number	CL/0003912	CL/0003913	CL/0003914	CL/0003915

UKAS accredited	Test No.	CL/0003912	CL/0003913	CL/0003914	CL/0003915
Chromium (total)	ICPSSS11	45	40	41	27
CN- (total)	ICTSCN28	<1	<1	<1	1
Copper	ICPSSS11	12	20	19	181
Cyanide (Free)	BGCN22		<1	<1	<1
Lead	ICPSSS11	51	245	52	421
Nickel	ICPSSS11	38	19	34	31
Oil FTIR	FTIRSW		2230	184	547
PAH (screening)	PAHSCUV	<10	372	12	87
pH units	WSLM3	7.9	9.5	8.0	9.9
Phenol Index	WSLM4	<0.5	2.8	<0.5	<0.5
SO4-- (acid sol)	ICPACS58	283	6190	1070	2830
Sulphide	ICTSCN28		16	<1	502
TPH FTIR (AR)	FTIRSW		783	<20	136
Zinc	ICPSSS11	112	149	100	1040

not UKAS accredited		CL/0003912	CL/0003913	CL/0003914	CL/0003915
Arsenic (Ms)		13.00	8.70	13.00	21
Boron.		1.3	1.1	1.3	3.2
Cadmium (Ms)		0.11	0.38	0.13	11.20
Mercury (Ms)		0.04	0.14	0.02	3.77
Selenium (MS)			0.98	0.61	1.07

Results expressed as mg/kg Air Dried unless stated otherwise
SO4 Analysis not conducted in accordance with BS1377
AR denotes analysis conducted on As Received Sample

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TEST REPORT

SOIL SAMPLE ANALYSIS



1252
1411

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
Mid Glamorgan
CF31 1JZ

TES Report No. 000542

Site: M4 Relief Road

Customer reference	CH 10 1.40	CH 11 0.35	CH 11 1.65	CH 12 0.17
Depth (m)	16/02/00 CL/0003916	16/02/00 CL/0003917	16/02/00 CL/0003918	16/02/00 CL/0003919
Date logged				
TES Bretby ID Number				

UKAS accredited	Test No.	CL/0003916	CL/0003917	CL/0003918	CL/0003919
Chromium (total)	ICPSSS11	25	42	160	30
CN- (total)	ICTSCN28	<1	4	2	<1
Copper	ICPSSS11	34	64	402	169
Cyanide (Free)	BGCN22	<1	<1	<1	<1
Lead	ICPSSS11	90	120	390	304
Nickel	ICPSSS11	24	47	183	34
Oil FTIR	FTIRSW	1410	23800	116000	932
PAH (screening)	PAHSCUV	108	233	1620	170
pH units	WSLM3	8.4	10.1	9.7	8.8
Phenol Index	WSLM4	<0.5	<0.5	1.3	<0.5
SO4-- (acid sol)	ICPACS58	1310	2640	7650	1900
Sulphide	ICTSCN28	50	<1	58	114
TPH FTIR (AR)	FTIRSW	63	75	46300	405
Zinc	ICPSSS11	144	398	3230	170

not UKAS accredited	CL/0003916	CL/0003917	CL/0003918	CL/0003919
Arsenic (Ms)	10.00	11.00	86	40
Boron.	1.7	1.3	<0.5	1.8
Cadmium (Ms)	0.82	0.28	0.31	0.30
Mercury (Ms)	0.42	0.16	0.16	0.41
Selenium (MS)	0.64	0.61	0.76	1.07

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Wildmill
Bridgend
Mid Glamorgan
CF31 1JZ

TES Report No. 000542

Site: M4 Relief Road

Customer reference	CH 12 0.67	CH 13 0.00	CH 13 0.77	CH 14 0.16
Depth (m)				
Date logged	16/02/00	16/02/00	16/02/00	16/02/00
TES Bretby ID Number	CL/0003920	CL/0003921	CL/0003922	CL/0003923

UKAS accredited	Test No.	CL/0003920	CL/0003921	CL/0003922	CL/0003923
Chromium (total)	ICPSSS11	31	13	20	25
CN- (total)	ICTSCN28	<1	<1	<1	<1
Copper	ICPSSS11	100	42	18	141
Cyanide (Free)	BGCN22	<1	<1	<1	<1
Lead	ICPSSS11	118	154	69	173
Nickel	ICPSSS11	34	14	20	43
Oil FTIR	FTIRSW	809	247	1160	2260
PAH (screening)	PAHSCUV	177	143	332	111
pH units	WSLM3	7.9	8.5	8.7	8.8
Phenol Index	WSLM4	<0.5	<0.5	<0.5	<0.5
SO4-- (acid sol)	ICPACS58	1150	1410	2200	906
Sulphide	ICTSCN28	27	21	30	54
TPH FTIR (AR)	FTIRSW	303	220	113	169
Zinc	ICPSSS11	131	912	94	258

not UKAS accredited	CL/0003920	CL/0003921	CL/0003922	CL/0003923
Arsenic (Ms)	27	36	12.00	21
Boron.	0.8	<0.5	1.1	1.6
Cadmium (Ms)	0.53	9.74	0.25	0.41
Mercury (Ms)	0.19	0.27	0.07	1.27
Selenium (MS)	1.02	1.09	0.58	0.96

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TEST REPORT SOIL SAMPLE ANALYSIS



1252
1411

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
Mid Glamorgan
CF31 1JZ

TES Report No. 000542

Site: M4 Relief Road

Customer reference	CH 15 0.30	CH 15 A 0.17	CH 16 0.20	CH 16 1.90
Depth (m)				
Date logged	16/02/00	16/02/00	16/02/00	16/02/00
TES Bretby ID Number	CL/0003924	CL/0003925	CL/0003926	CL/0003927

UKAS accredited	Test No.	CL/0003924	CL/0003925	CL/0003926	CL/0003927
Chromium (total)	ICPSSS11	29	33	15	22
CN- (total)	ICTSCN28	<1	<1	<1	<1
Copper	ICPSSS11	80	128	78	61
Cyanide (Free)	BGCN22	<1	<1	<1	<1
Lead	ICPSSS11	126	98	98	60
Nickel	ICPSSS11	35	43	41	23
Oil FTIR	FTIRSW	242	229	157	1460
PAH (screening)	PAHSCUV	81	43	49	215
pH units	WSLM3	9.0	8.9	8.3	9.9
Phenol Index	WSLM4	<0.5	<0.5	<0.5	<0.5
SO4-- (acid sol)	ICPACS58	1120	974	680	1760
Sulphide	ICTSCN28	73	158	5	16
TPH FTIR (AR)	FTIRSW	106	140	335	711
Zinc	ICPSSS11	138	116	105	91

not UKAS accredited		CL/0003924	CL/0003925	CL/0003926	CL/0003927
Arsenic (Ms)		18	19	12.00	22
Boron.		2.5	2.1	1.7	1.8
Cadmium (Ms)		0.30	0.29	0.17	0.48
Mercury (Ms)		0.68	0.54	0.42	0.30
Selenium (MS)		0.88	0.87	1.25	1.01

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TEST REPORT

SOIL SAMPLE ANALYSIS



1252
1411

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
Mid Glamorgan
CF31 1JZ

TES Report No. 000542

Site: M4 Relief Road

Customer reference	CH 17 0.60	CH 17 1.30	CH 17 1.90	CH 18 0.20
Depth (m)				
Date logged	16/02/00	16/02/00	16/02/00	16/02/00
TES Bretby ID Number	CL/0003928	CL/0003929	CL/0003930	CL/0003931

UKAS accredited	Test No.	CL/0003928	CL/0003929	CL/0003930	CL/0003931
Chromium (total)	ICPSSS11	26	21	38	31
CN- (total)	ICTSCN28	<1	<1	<1	<1
Copper	ICPSSS11	27	104	22	16
Cyanide (Free)	BGCN22	<1	<1	<1	<1
Lead	ICPSSS11	86	92	61	63
Nickel	ICPSSS11	28	37	35	30
Oil FTIR	FTIRSW	126	251	128	95
PAH (screening)	PAHSCUV	515	564	45	17
pH units	WSLM3	8.3	10.3	7.7	8.1
Phenol Index	WSLM4	<0.5	0.5	0.5	0.5
SO4-- (acid sol)	ICPACS58	779	1740	796	1380
Sulphide	ICTSCN28	<1	14	7	1
TPH FTIR (AR)	FTIRSW	35	144	95	38
Zinc	ICPSSS11	119	142	112	96

not UKAS accredited	CL/0003928	CL/0003929	CL/0003930	CL/0003931
Arsenic (Ms)	9.20	11.00	15.00	11.00
Boron.	<0.5	1.0	2.9	1.1
Cadmium (Ms)	0.17	0.28	0.09	0.18
Mercury (Ms)	0.20	0.35	0.12	<0.01
Selenium (MS)	0.51	0.60	<0.50	0.61

Results expressed as mg/kg Air Dried unless stated otherwise
SO4 Analysis not conducted in accordance with BS1377
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TEST REPORT

SOIL SAMPLE ANALYSIS



1252
1411

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
Mid Glamorgan
CF31 1JZ

TES Report No. 000542

Site: M4 Relief Road

Customer reference	CH 18 1.60	CH 18 4.00	CH 19 0.20	CH 19 2.00
Depth (m)	16/02/00 CL/0003932	16/02/00 CL/0003933	16/02/00 CL/0003934	16/02/00 CL/0003935
Date logged				
TES Bretby ID Number				

UKAS accredited	Test No.	CL/0003932	CL/0003933	CL/0003934	CL/0003935
Chromium (total)	ICPSSS11	16	20	14	21
CN- (total)	ICTSCN28	2	<1	<1	<1
Copper	ICPSSS11	69	35	38	8
Cyanide (Free)	BGCN22	<1	<1	<1	<1
Lead	ICPSSS11	298	80	59	26
Nickel	ICPSSS11	21	24	37	21
Oil FTIR	FTIRSW	42	103	131	85
PAH (screening)	PAHSCUV	20	34	19	11
pH units	WSLM3	8.1	8.3	8.2	8.1
Phenol Index	WSLM4	<0.5	<0.5	<0.5	<0.5
SO4-- (acid sol)	ICPACS58	1960	891	498	1490
Sulphide	ICTSCN28	<1	21	2	8
TPH FTIR (AR)	FTIRSW	<20	30	91	<20
Zinc	ICPSSS11	163	149	61	51

not UKAS accredited	CL/0003932	CL/0003933	CL/0003934	CL/0003935
Arsenic (Ms)	22	15.00	13.00	6.20
Boron.	<0.5	1.1	1.1	1.3
Cadmium (Ms)	0.27	0.22	0.12	0.08
Mercury (Ms)	0.10	0.35	0.11	<0.01
Selenium (MS)	0.91	0.89	0.61	0.58

Results expressed as mg/kg Air Dried unless stated otherwise
SO4 Analysis not conducted in accordance with BS1377
AR denotes analysis conducted on As Received Sample

Date of Issue: 02/03/00

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TEST REPORT

SOIL SAMPLE ANALYSIS



1252
1411

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
Mid Glamorgan
CF31 1JZ

TES Report No. 000542

Site: M4 Relief Road

Customer reference	CH 19 A 0.30	CH 19 A 1.00	CH 24 0.00	CH 24 0.75
Depth (m)				
Date logged	16/02/00	16/02/00	16/02/00	16/02/00
TES Bretby ID Number	CL/0003936	CL/0003937	CL/0003938	CL/0003939

UKAS accredited	Test No.	CL/0003936	CL/0003937	CL/0003938	CL/0003939
Chromium (total)	ICPSSS11	17	19	99	44
CN- (total)	ICTSCN28	<1	<1	<1	<1
Copper	ICPSSS11	56	14	141	25
Cyanide (Free)	BGCN22	<1	<1	<1	<1
Lead	ICPSSS11	232	47	154	69
Nickel	ICPSSS11	35	18	51	39
Oil FTIR	FTIRSW	64	159	200	149
PAH (screening)	PAHSCUV	12	22	17	19
pH units	WSLM3	8.5	8.4	9.5	7.3
Phenol Index	WSLM4	<0.5	<0.5	<0.5	<0.5
SO4-- (acid sol)	ICPACS58	1550	1450	836	760
Sulphide	ICTSCN28	<1	4	40	6
TPH FTIR (AR)	FTIRSW	<20	99	<20	<20
Zinc	ICPSSS11	830	77	280	122

not UKAS accredited	CL/0003936	CL/0003937	CL/0003938	CL/0003939
Arsenic (Ms)	34	8.10	9.30	17
Boron.	2.3	2.1	<0.5	5.0
Cadmium (Ms)	1.72	0.21	0.87	0.20
Mercury (Ms)	0.08	0.03	0.14	0.14
Selenium (MS)	0.82	0.72	0.60	0.42

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TEST REPORT

SOIL SAMPLE ANALYSIS



1252
1411

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
Mid Glamorgan
CF31 1JZ

TES Report No. 000542

Site: M4 Relief Road

Customer reference	CH			
Depth (m)	25 0.00			
Date logged	16/02/00			
TES Bretby ID Number	CL/0003940			

UKAS accredited	Test No.	CL/0003940			
Chromium (total)	ICPSSS11	16			
CN- (total)	ICTSCN28	<1			
Copper	ICPSSS11	16			
Cyanide (Free)	BGCN22	<1			
Lead	ICPSSS11	56			
Nickel	ICPSSS11	9			
Oil FTIR	FTIRSW	1130			
PAH (screening)	PAHSCUV	191			
pH units	WSLM3	10.6			
Phenol Index	WSLM4	<0.5			
SO4-- (acid sol)	ICPACS58	2440			
Sulphide	ICTSCN28	12			
TPH FTIR (AR)	FTIRSW	302			
Zinc	ICPSSS11	69			

not UKAS accredited		CL/0003940			
Arsenic (Ms)		4.60			
Boron.		0.5			
Cadmium (Ms)		0.51			
Mercury (Ms)		0.14			
Selenium (MS)		0.64			

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Volatile Organic Compounds by PTGCMS

Customer and Site Details: Exploration Associates : M4 Relief Road
Sample Details: CH 11 / 1.65m
LIMS ID Number: CL03918
Report Number: 000542

Directory/Quant file:
Date Booked in:
Date Analysed:
Operator:

0229VOC.MS1\ 0229CCC1.D
16-Feb-00
29-Feb-00
D.WALTON

Matrix: Soil
Method: Purge and Trap
Dilution: 5
Position: 2

Target Compounds	CAS #	R.T. (min.)	Concentration µg/kg	% Fit
Dichlorodifluoromethane	75-71-8	-	< 5	-
Chloromethane	74-87-3	-	< 5	-
Vinyl Chloride	75-01-4	-	< 5	-
Bromomethane	74-83-9	-	< 5	-
Chloroethane	75-00-3	-	< 5	-
Trichlorofluoromethane	75-69-4	-	< 5	-
1,1-Dichloroethene	75-35-4	-	< 5	-
trans 1,2-Dichloroethene	156-60-5	-	< 5	-
1,1-Dichloroethane	75-34-3	-	< 5	-
2,2-Dichloropropane	594-20-7	-	< 5	-
cis 1,2-Dichloroethene	156-59-2	-	< 5	-
Bromochloromethane	74-97-5	-	< 5	-
Chloroform	67-66-3	-	< 5	-
1,1,1-Trichloroethane	71-55-6	-	< 5	-
Carbon Tetrachloride	56-23-5	-	< 5	-
1,1-Dichloropropene	563-58-6	-	< 5	-
Benzene	71-43-2	-	< 5	-
1,2-Dichloroethane	107-06-2	-	< 5	-
Trichloroethene	79-01-6	-	< 5	-
1,2-Dichloropropene	78-87-5	-	< 5	-
Dibromomethane	74-95-3	-	< 5	-
Bromodichloromethane	75-27-4	-	< 5	-
cis 1,3-Dichloropropene	10061-01-5	-	< 5	-
Toluene	108-88-3	15.67	120	99
trans 1,3-Dichloropropene	10061-02-6	-	< 5	-
1,1,2-Trichloroethane	79-00-5	-	< 5	-
Tetrachloroethene	127-18-4	-	< 5	-
1,3-Dichloropropane	142-28-29	-	< 5	-
Dibromochloromethane	124-48-1	-	< 5	-
1,2-Dibromoethane	106-93-4	-	< 5	-
Chlorobenzene	108-90-7	-	< 5	-
Ethylbenzene	100-41-4	17.85	9	92
1,1,1,2-Tetrachloroethane	630-20-6	-	< 5	-
m and p-Xylene	108-38-3/106-42-3	18.01	37	97
o-Xylene	95-47-6	18.61	15	96

Concentrations are reported on a wet weight basis

Target Compounds	CAS #	R.T. (min.)	Concentration µg/kg	% Fit
Styrene	100-42-5	-	< 5	-
Bromoform	75-25-2	-	< 5	-
iso-Propylbenzene	98-82-8	-	< 5	-
1,1,2,2-Tetrachloroethane	79-34-5	-	< 5	-
Propylbenzene	103-65-1	-	< 5	-
Bromobenzene	108-86-1	-	< 5	-
1,2,3-Trichloropropane	96-18-4	-	< 5	-
2-Chlorotoluene	95-49-8	-	< 5	-
1,3,5-Trimethylbenzene	108-67-8	19.96	26	99
4-Chlorotoluene	106-43-4	-	< 5	-
tert-Butylbenzene	98-06-6	-	< 5	-
1,2,4-Trimethylbenzene	95-63-6	20.56	81	99
sec-Butylbenzene	135-98-8	20.81	7	87
p-Isopropyltoluene	99-87-6	21.00	12	97
1,3-Dichlorobenzene	541-73-1	-	< 5	-
1,4-Dichlorobenzene	106-46-7	-	< 5	-
n-Butylbenzene	104-51-8	21.63	12	80
1,2-Dichlorobenzene	95-50-1	21.81	23	90
1,2-Dibromo-3-chloropropane	96-12-8	-	< 25	-
1,2,4-Trichlorobenzene	120-82-1	-	< 25	-
Hexachlorobutadiene	87-68-3	-	< 25	-
Naphthalene	91-20-3	24.81	51	100
1,2,3-Trichlorobenzene	87-61-6	-	< 25	-

Concentrations are reported on a wet weight basis

Internal standards	R.T.	Area %	Surrogates	% Rec
Pentafluorobenzene	12.27	114	Dibromofluoromethane	96
1,4-Difluorobenzene	13.42	108	Toluene-d8	95
Chlorobenzene-d5	17.72	95	Bromofluorobenzene	91
1,4-Dichlorobenzene-d4	21.17	74		



TEST REPORT SOIL SAMPLE ANALYSIS



1252
1411

Amended Report TES Report No. CL/000543

Site: M4 Relief Road

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
Mid Glamorgan
CF31 1JZ

The 34 samples described in this report were scheduled for analysis by TES Bretby on Wednesday, 16 February 2000. The analysis was completed by Tuesday, 4 April 2000.

Tests marked as 'not UKAS accredited' and any opinions or interpretations expressed herein are outside the scope of any UKAS accreditation held by TES Bretby laboratories.

The following tables are contained in this report:

Table 1 Sample Descriptions
Table 2 Main Analysis Results

On behalf of
TES Bretby : J. Hannah
J Hannah Project Co-ordinator

Date of Issue: 04/04/00

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TEST REPORT

SOIL SAMPLE ANALYSIS



1252
1411

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
Mid Glamorgan
CF31 1JZ

Amended Report
TES Report No. 000543

Site: M4 Relief Road

Page No	Assessed Area	ID No. EFS/CL	Sample	Depth(m) from - to	Description
1	M4 Relief Road	0003941	CH/ 25/	0.80	CH25 0.8
1	M4 Relief Road	0003942	CH/ 26/	0.30	CH26 0.3
1	M4 Relief Road	0003943	CH/ 27/	0.55	CH27 0.55
1	M4 Relief Road	0003944	CH/ 27/	1.10	CH27 1.1
2	M4 Relief Road	0003945	CH/ 28/	0.22	CH28 0.22
2	M4 Relief Road	0003946	CH/ 28/	1.90	CH28 1.9
2	M4 Relief Road	0003947	CH/ 29/	0.40	CH29 0.4
2	M4 Relief Road	0003948	CH/ 30/	0.00	CH30 0.0
3	M4 Relief Road	0003949	CH/ 30/	1.00	CH30 1.0
3	M4 Relief Road	0003950	CH/ 31/	0.00	CH31 0.0
3	M4 Relief Road	0003951	CH/ 32/	0.00	CH32 0.0
3	M4 Relief Road	0003952	CH/ 32/	0.50	CH32 0.5
4	M4 Relief Road	0003953	CH/ 33/	0.20	CH33 0.2
4	M4 Relief Road	0003954	CH/ 33/	1.50	CH33 1.5
4	M4 Relief Road	0003955	CH/ 43/	0.20	CH43 0.2
4	M4 Relief Road	0003956	CH/ 43/	1.30	CH43 1.3
5	M4 Relief Road	0003957	CH/ 44/	0.30	CH44 0.3
5	M4 Relief Road	0003958	CH/ 44/	1.00	CH44 1.0
5	M4 Relief Road	0003959	CH/ 45/	0.20	CH45 0.2
5	M4 Relief Road	0003961	CH/ 46/	0.50	CH46 0.5

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Report 000543
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TEST REPORT

SOIL SAMPLE ANALYSIS



1252
1411

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
Mid Glamorgan
CF31 1JZ

Amended Report
TES Report No. 000543

Site: M4 Relief Road

Page No	Assessed Area	ID No. EFS/CL	Sample	Depth (m) from to	Description
6	M4 Relief Road	0003960	CH/ 46/	1.20	CH46 1.2
6	M4 Relief Road	0003962	CH/ 47/	0.10	CH47 0.1
6	M4 Relief Road	0003963	CH/ 47/	1.10	CH47 1.1
6	M4 Relief Road	0003964	CH/ 47/A	0.20	CH47A 0.2
7	M4 Relief Road	0003965	CH/ 48/	0.05	CH48 0.5
7	M4 Relief Road	0003966	CH/ 48/	1.00	CH48 1.0
7	M4 Relief Road	0003967	CH/ 49/	0.05	CH49 0.05
7	M4 Relief Road	0003968	CH/ 49/	0.60	CH49 0.6
8	M4 Relief Road	0003970	CH/ 50/	0.05	CH50 0.05
8	M4 Relief Road	0003969	CH/ 50/	0.50	CH50 0.5
8	M4 Relief Road	0003971	CH/ 51/	0.00	CH51 0.0
8	M4 Relief Road	0003972	CH/ 55/	0.00	CH55 0.0
9	M4 Relief Road	0003973	CH/ 55/	0.25	CH55 0.25
9	M4 Relief Road	0003974	CH/ 56/	0.00	CH56 0.0

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TEST REPORT

SOIL SAMPLE ANALYSIS

1252
1411

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
Mid Glamorgan
CF31 1JZ

Amended Report
TES Report No. 000543

Site: M4 Relief Road

Customer reference	CH 25 0.80	CH 26 0.30	CH 27 0.55	CH 27 1.10
Depth (m)				
Date logged	16/02/00	16/02/00	16/02/00	16/02/00
TES Bretby ID Number	CL/0003941	CL/0003942	CL/0003943	CL/0003944

UKAS accredited	Test No.	CL/0003941	CL/0003942	CL/0003943	CL/0003944
Chromium (total)	ICPSSS11	35	13	120	42
CN- (total)	ICTSCN28	<1	<1	<1	<1
Copper	ICPSSS11	23	19	147	15
Cyanide (Free)	BGCN22		<1	<1	<1
Lead	ICPSSS11	68	45	157	49
Nickel	ICPSSS11	32	10	53	33
Oil FTIR	FTIRSW		137	485	105
PAH (screening)	PAHSCUV	71	40	56	<10
pH units	WSLM3	7.6	10.3	8.5	8.6
Phenol Index	WSLM4	<0.5	<0.5	<0.5	<0.5
SO4-- (acid sol)	ICPACS58	843	1110	626	533
Sulphide	ICTSCN28		16	11	4
TPH FTIR (AR)	FTIRSW		<20	346	<20
Zinc	ICPSSS11	129	53	188	92

not UKAS accredited		CL/0003941	CL/0003942	CL/0003943	CL/0003944
Arsenic (Ms)		15.00	6.43	14.40	11.30
Boron.		1.6	1.0	1.8	2.6
Cadmium (Ms)		0.21	0.10	0.35	0.12
Mercury (Ms)		0.14	0.09	0.14	0.04
Selenium (MS)			0.94	0.60	<0.50

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TEST REPORT

SOIL SAMPLE ANALYSIS



1252
1411

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
Mid Glamorgan
CF31 1JZ

Amended Report
TES Report No. 000543

Site: M4 Relief Road

Customer reference	CH 28 0.22	CH 28 1.90	CH 29 0.40	CH 30 0.00
Depth (m)				
Date logged	16/02/00	16/02/00	16/02/00	16/02/00
TES Bretby ID Number	CL/0003945	CL/0003946	CL/0003947	CL/0003948

UKAS accredited	Test No.	CL/0003945	CL/0003946	CL/0003947	CL/0003948
Chromium (total)	ICPSSS11	16	14	33	32
CN- (total)	ICTSCN28	<1	<1	<1	<1
Copper	ICPSSS11	43	10	52	60
Cyanide (Free)	BGCN22	<1	<1	<1	<1
Lead	ICPSSS11	62	26	60	125
Nickel	ICPSSS11	13	10	29	23
Oil FTIR	FTIRSW	175	137	443	152
PAH (screening)	PAHSCUV	58	20	21	10
pH units	WSLM3	11.2	11.1	10.0	9.2
Phenol Index	WSLM4	<0.5	<0.5	<0.5	<0.5
SO4-- (acid sol)	ICPACS58	2240	1370	674	760
Sulphide	ICTSCN28	<1	1	19	19
TPH FTIR (AR)	FTIRSW	66	53	88	<20
Zinc	ICPSSS11	89	44	63	139

not UKAS accredited		CL/0003945	CL/0003946	CL/0003947	CL/0003948
Arsenic (Ms)		16.7	5.52	8.18	8.85
Boron.		0.6	0.6	0.7	1.5
Cadmium (Ms)		0.22	0.17	0.18	0.36
Mercury (Ms)		2.91	0.13	0.80	0.38
Selenium (MS)		0.99	1.01	0.67	0.83

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SOIL SAMPLE ANALYSIS



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Bridgend
Mid Glamorgan
CF31 1JZ

Amended Report
TES Report No. 000543

Site: M4 Relief Road

Customer reference	CH 30	CH 31	CH 32	CH 32
Depth (m)	1.00	0.00	0.00	0.50
Date logged	16/02/00	16/02/00	16/02/00	16/02/00
TES Bretby ID Number	CL/0003949	CL/0003950	CL/0003951	CL/0003952

UKAS accredited	Test No.	CL/0003949	CL/0003950	CL/0003951	CL/0003952
Chromium (total)	ICPSSS11	45	2170	1790	41
CN- (total)	ICTSCN28	<1	<1	<1	<1
Copper	ICPSSS11	77	536	221	172
Cyanide (Free)	BGCN22	<1	<1	<1	<1
Lead	ICPSSS11	111	182	166	417
Nickel	ICPSSS11	29	141	49	32
Oil FTIR	FTIRSW	477	160	514	131
PAH (screening)	PAHSCUV	66	10	26	44
PCB - ARO	PCBAROEC		0.2	0.2	0.5
pH units	WSLM3	10.4	11.7	11.2	8.8
Phenol Index	WSLM4	<0.5	<0.5	<0.5	<0.5
SO4-- (acid sol)	ICPACS58	1190	2440	1150	1420
Sulphide	ICTSCN28	20	14	38	<1
TPH FTIR (AR)	FTIRSW	132	105	1190	94
Zinc	ICPSSS11	168	475	390	389

not UKAS accredited		CL/0003949	CL/0003950	CL/0003951	CL/0003952
Arsenic (Ms)		12.00	21.3	12.80	29.5
Boron.		3.2	23	2.1	1.3
Cadmium (Ms)		0.68	1.09	0.57	0.29
Mercury (Ms)		1.21	0.38	3.21	16.9
Selenium (MS)		0.97	2.52	1.79	0.97

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TEST REPORT

SOIL SAMPLE ANALYSIS



1252
1411

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
Mid Glamorgan
CF31 1JZ

Amended Report
TES Report No. 000543

Site: M4 Relief Road

Customer reference	CH 33	CH 33	CH 43	CH 43
Depth (m)	0.20	1.50	0.20	1.30
Date logged	16/02/00	16/02/00	16/02/00	16/02/00
TES Bretby ID Number	CL/0003953	CL/0003954	CL/0003955	CL/0003956

UKAS accredited	Test No.	CL/0003953	CL/0003954	CL/0003955	CL/0003956
Chromium (total)	ICPSSS11	40	30	41	36
CN- (total)	ICTSCN28	<1	<1	<1	<1
Copper	ICPSSS11	16	19	15	11
Cyanide (Free)	BGCN22	<1	<1	<1	<1
Lead	ICPSSS11	56	52	44	41
Nickel	ICPSSS11	33	21	31	33
Oil FTIR	FTIRSW	141	234	906	72
PAH (screening)	PAHSCUV	16	38	2370	60
PCB - ARO	PCBAROEC	1.1	1.9	136	1.8
pH units	WSLM3	8.0	8.0	9.3	8.0
Phenol Index	WSLM4	<0.5	0.5	11.2	<0.5
SO4-- (acid sol)	ICPACS58	694	620	1340	435
Sulphide	ICTSCN28	<1	17	7	<1
TPH FTIR (AR)	FTIRSW	28	22	330	60
Zinc	ICPSSS11	103	121	91	82

not UKAS accredited		CL/0003953	CL/0003954	CL/0003955	CL/0003956
Arsenic (Ms)		13.30	7.68	10.20	12.30
Boron.		1.1	2.0	0.7	2.2
Cadmium (Ms)		0.12	0.54	0.10	0.07
Mercury (Ms)		0.64	2.14	0.07	0.24
Selenium (MS)		0.47	0.94	0.36	<0.50

Results expressed as mg/kg Air Dried unless stated otherwise
SO4 Analysis not conducted in accordance with BS1377
AR denotes analysis conducted on As Received Sample

Date of Issue: 04/04/00

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TES Bretby accepts no responsibility for the sampling related to the above results

TES Bretby, P.O. Box 100, Burton-on-Trent, DE15 0XD Telephone: 01283 554400 Fax: 01283 554422
TES Bretby is a division of Environmental Services Group Limited Registered in England Number 2880501

TES Bretby
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TEST REPORT

SOIL SAMPLE ANALYSIS



1252
1411

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
Mid Glamorgan
CF31 1JZ

Amended Report
TES Report No. 000543

Site: M4 Relief Road

Customer reference	CH 44	CH 44	CH 45	CH 46
Depth (m)	0.30	1.00	0.20	0.50
Date logged	16/02/00	16/02/00	16/02/00	16/02/00
TES Bretby ID Number	CL/0003957	CL/0003958	CL/0003959	CL/0003961

UKAS accredited	Test No.	CL/0003957	CL/0003958	CL/0003959	CL/0003961
Chromium (total)	ICPSSS11	49	50	76	87
CN- (total)	ICTSCN28	<1	<1	<1	<1
Copper	ICPSSS11	41	16	19	31
Cyanide (Free)	BGCN22	<1	<1	<1	<1
Lead	ICPSSS11	62	46	45	124
Nickel	ICPSSS11	28	15	11	37
Oil FTIR	FTIRSW	805	657	251	487
PAH (screening)	PAHSCUV	2310	1430	55	1940
PCB - ARO	PCBAROEC	58	19.9	2.0	35
pH units	WSLM3	9.8	9.6	11.8	9.7
Phenol Index	WSLM4	0.6	0.7	<0.5	1.5
SO4-- (acid sol)	ICPACS58	1690	2550	2950	1560
Sulphide	ICTSCN28	6	17	11	25
TPH FTIR (AR)	FTIRSW	879	162	72	88
Zinc	ICPSSS11	123	70	37	145

not UKAS accredited		CL/0003957	CL/0003958	CL/0003959	CL/0003961
Arsenic (Ms)		7.91	8.15	8.74	7.34
Boron.		0.7	1.0	0.6	0.6
Cadmium (Ms)		0.47	0.40	0.14	0.50
Mercury (Ms)		34.0	7.66	0.19	2.50
Selenium (MS)		1.03	1.02	0.87	1.14

Results expressed as mg/kg Air Dried unless stated otherwise
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AR denotes analysis conducted on As Received Sample

Date of Issue: 04/04/00

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TES Bretby
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TEST REPORT

SOIL SAMPLE ANALYSIS



1252
1411

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
Mid Glamorgan
CF31 1JZ

Amended Report
TES Report No. 000543

Site: M4 Relief Road

Customer reference	CH 46	CH 47	CH 47	CH 47 A
Depth (m)	1.20	0.10	1.10	0.20
Date logged	16/02/00	16/02/00	16/02/00	16/02/00
TES Bretby ID Number	CL/0003960	CL/0003962	CL/0003963	CL/0003964

UKAS accredited	Test No.	CL/0003960	CL/0003962	CL/0003963	CL/0003964
Chromium (total)	ICPSSS11	101	54	51	67
CN- (total)	ICTSCN28	<1	<1	<1	<1
Copper	ICPSSS11	30	22	18	20
Cyanide (Free)	BGCN22	<1	<1	<1	<1
Lead	ICPSSS11	131	54	84	55
Nickel	ICPSSS11	44	24	36	22
Oil FTIR	FTIRSW	780	401	93	294
PAH (screening)	PAHSCUV	2400	47	26	97
PCB - ARO	PCBAROEC	44	238	7.0	72
pH units	WSLM3	8.8	10.0	8.1	9.8
Phenol Index	WSLM4	0.6	1.1	<0.5	<0.5
SO4-- (acid sol)	ICPACS58	1450	1210	1050	1100
Sulphide	ICTSCN28	7	44	9	29
TPH FTIR (AR)	FTIRSW	58	273	24	114
Zinc	ICPSSS11	143	125	118	115

not UKAS accredited		CL/0003960	CL/0003962	CL/0003963	CL/0003964
Arsenic (Ms)		8.18	9.03	13.60	7.52
Boron.		<0.5	1.0	1.3	1.2
Cadmium (Ms)		0.72	0.46	0.14	0.43
Mercury (Ms)		2.84	1.02	0.14	2.27
Selenium (MS)		1.24	0.98	<0.50	0.91

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TEST REPORT

SOIL SAMPLE ANALYSIS



1252
1411

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
Mid Glamorgan
CF31 1JZ

Amended Report
TES Report No. 000543

Site: M4 Relief Road

Customer reference	CH 48	CH 48	CH 49	CH 49
Depth (m)	0.05	1.00	0.05	0.60
Date logged	16/02/00	16/02/00	16/02/00	16/02/00
TES Bretby ID Number	CL/0003965	CL/0003966	CL/0003967	CL/0003968

UKAS accredited	Test No.	CL/0003965	CL/0003966	CL/0003967	CL/0003968
Chromium (total)	ICPSSS11	349	42	214	35
CN- (total)	ICTSCN28	1	<1	4	<1
Copper	ICPSSS11	30	11	20	14
Iron	ICPSSS11			32500	
Lead	ICPSSS11	56	44	134	53
Nickel	ICPSSS11	7	32	8	29
Oil FTIR	FTIRSW	74		42	
PAH (screening)	PAHSCUV	72	<10	85	<10
pH units	WSLM3	10.6	8.6	10.5	8.2
Phenol Index	WSLM4	<0.5	<0.5	<0.5	<0.5
SO4-- (acid sol)	ICPACS58	4810	668	5090	672
Sulphide	ICTSCN28	762		583	
TPH FTIR (AR)	FTIRSW	25		63	
Zinc	ICPSSS11	41	116	589	102

not UKAS accredited		CL/0003965	CL/0003966	CL/0003967	CL/0003968
Arsenic (Ms)		3.41	10.60	7.16	10.70
Boron.		1.8	1.6	1.5	1.4
Cadmium (Ms)		0.13	0.08	0.64	0.12
Mercury (Ms)		0.15	0.05	0.04	0.02
Selenium (MS)		1.85	<0.50	1.69	<0.50

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TES Bretby
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TEST REPORT

SOIL SAMPLE ANALYSIS



1252
1411

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
Mid Glamorgan
CF31 1JZ

Amended Report
TES Report No. 000543

Site: M4 Relief Road

Customer reference	CH 50 0.05	CH 50 0.50	CH 51 0.00	CH 55 0.00
Depth (m)				
Date logged	16/02/00	16/02/00	16/02/00	16/02/00
TES Bretby ID Number	CL/0003970	CL/0003969	CL/0003971	CL/0003972

UKAS accredited	Test No.	CL/0003970	CL/0003969	CL/0003971	CL/0003972
Chromium (total)	ICPSSS11	60	50	706	11
CN- (total)	ICTSCN28	4	<1	3	<1
Copper	ICPSSS11	19	12	24	10
Iron	ICPSSS11	16700		14300	
Lead	ICPSSS11	180	45	62	36
Nickel	ICPSSS11	9	32	5	8
Oil FTIR	FTIRSW	74		585	15900
PAH (screening)	PAHSCUV	102	15	82	4980
pH units	WSLM3	9.9	9.6	12.2	9.9
Phenol Index	WSLM4	<0.5	<0.5	<0.5	<0.5
SO4-- (acid sol)	ICPACS58	7330	636	4020	2100
Sulphide	ICTSCN28	30		6	144
TPH FTIR (AR)	FTIRSW	181		198	6100
Zinc	ICPSSS11	520	103	69	58

not UKAS accredited		CL/0003970	CL/0003969	CL/0003971	CL/0003972
Arsenic (Ms)		11.10	9.63	2.82	5.22
Boron.		0.7	1.0	0.9	0.9
Cadmium (Ms)		1.50	0.11	0.13	0.23
Mercury (Ms)		0.02	0.04	0.07	0.01
Selenium (MS)		1.33	0.40	1.63	0.84

Results expressed as mg/kg Air Dried unless stated otherwise
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Date of Issue: 04/04/00

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TES Bretby
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TEST REPORT

SOIL SAMPLE ANALYSIS



1252
1411

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
Mid Glamorgan
CF31 1JZ

Amended Report
TES Report No. 000543

Site: M4 Relief Road

Customer reference	CH	CH		
Depth (m)	55 0.25	56 0.00		
Date logged	16/02/00	16/02/00		
TES Bretby ID Number	CL/0003973	CL/0003974		

UKAS accredited	Test No.	CL/0003973	CL/0003974		
Chromium (total)	ICPSSS11	20	23		
CN- (total)	ICTSCN28	<1	<1		
Copper	ICPSSS11	3	18		
Lead	ICPSSS11	19	38		
Nickel	ICPSSS11	24	18		
Oil FTIR	FTIRSW	292	182		
PAH (screening)	PAHSCUV	12	<10		
pH units	WSLM3	8.5	8.1		
Phenol Index	WSLM4	<0.5	<0.5		
SO4-- (acid sol)	ICPACS58	644	337		
Sulphide	ICTSCN28	4	<1		
TPH FTIR (AR)	FTIRSW	468	185		
Zinc	ICPSSS11	24	52		

not UKAS accredited		CL/0003973	CL/0003974		
Arsenic (Ms)		8.69	9.74		
Boron.		0.5	0.6		
Cadmium (Ms)		0.04	0.24		
Mercury (Ms)		<0.01	0.03		
Selenium (MS)		0.85	<0.50		

Results expressed as mg/kg Air Dried unless stated otherwise
SO4 Analysis not conducted in accordance with BS1377
AR denotes analysis conducted on As Received Sample

Date of Issue: 04/04/00

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TES Bretby
Report 000543
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Semi-Volatile Organic Compounds



Customer and Site Details:
Sample Details:
LIMS ID Number:
Job Number:

Exploration Associates: M4 Relief Road
CH/0033/ / 0.20- m
CL03953
534

Date Booked in: 16-Feb-00
Date Extracted: 18-Feb-00
Date Analysed: 24-Feb-00

Matrix: Soil
Ext Method: Soxhlet
Operator: SW
Directory/Quant File: 0224ABN.MS5 0224CCC1.D GPC (Y/N)

QC Batch Number: 226
Multiplier: 0.0667
Dilution Factor: 1
N

Target Compounds	CAS #	R.T. (min)	Concentration mg/kg	% Fit
Phenol	108-95-2	-	< 2.0	-
bis(2-Chloroethyl)ether	111-44-4	-	< 0.5	-
2-Chlorophenol	95-57-8	-	< 2.0	-
1,3-Dichlorobenzene	541-73-1	-	< 0.5	-
1,4-Dichlorobenzene	106-46-7	-	< 0.5	-
Benzyl alcohol	100-51-6	-	< 0.5	-
1,2-Dichlorobenzene	95-50-1	-	< 0.5	-
2-Methylphenol	95-48-7	-	< 0.5	-
bis(2-Chloroisopropyl)ether	108-60-1	-	< 0.5	-
Hexachloroethane	67-72-1	-	< 0.5	-
N-Nitroso-di-n-propylamine	621-64-7	-	< 0.5	-
3- & 4-Methylphenol	108-39-4/108-44-5	-	< 2.0	-
Nitrobenzene	98-95-3	-	< 0.5	-
Isophorone	78-59-1	-	< 0.5	-
2-Nitrophenol	88-75-5	-	< 2.0	-
2,4-Dimethylphenol	105-67-9	-	< 2.0	-
Benzoic Acid	65-85-0	-	< 10.0	-
bis(2-Chloroethoxy)methane	111-91-1	-	< 0.5	-
2,4-Dichlorophenol	120-83-2	-	< 2.0	-
1,2,4-Trichlorobenzene	120-82-1	-	< 0.5	-
Naphthalene	91-20-3	-	< 0.2	-
4-Chlorophenol *	106-48-9	-	< 2.0	-
4-Chloroaniline	106-47-8	-	< 0.5	-
Hexachlorobutadiene	87-68-3	-	< 0.5	-
4-Chloro-3-methylphenol	59-50-7	-	< 0.5	-
2-Methylnaphthalene	91-57-6	-	< 0.2	-
1-Methyl naphthalene *	90-12-0	-	< 0.2	-
Hexachlorocyclopentadiene	77-47-4	-	< 0.5	-
2,4,6-Trichlorophenol	88-06-2	-	< 2.0	-
2,4,5-Trichlorophenol	95-95-4	-	< 2.0	-
2-Chloronaphthalene	91-58-7	-	< 0.2	-
Biphenyl *	92-52-4	-	< 0.2	-
Diphenyl ether *	101-84-8	-	< 0.2	-
2-Nitroaniline	88-74-4	-	< 0.5	-
Acenaphthylene	208-96-8	-	< 0.2	-
Dimethylphthalate	131-11-3	-	< 0.5	-
2,6-Dinitrotoluene	606-20-2	-	< 0.5	-
Acenaphthene	83-32-9	-	< 0.2	-
3-Nitroaniline	99-09-2	-	< 0.5	-

* These compounds are not currently UKAS accredited.
Concentrations are reported on a wet weight basis.

Target Compounds	CAS #	R.T.	Concentration mg/kg	% Fit
2,4-Dinitrophenol	51-28-5	-	< 1.0	-
Dibenzofuran	132-64-9	-	< 0.5	-
4-Nitrophenol	100-02-7	-	< 5.0	-
2,4-Dinitrotoluene	121-14-2	-	< 0.5	-
Fluorene	86-73-7	-	< 0.2	-
Diethylphthalate	84-66-2	-	< 0.5	-
4-Chlorophenyl-phenylether	7005-72-4	-	< 0.5	-
4,6-Dinitro-2-methylphenol	534-52-1	-	< 5.0	-
4-Nitroaniline	100-01-6	-	< 0.5	-
N-Nitrosodiphenylamine	86-30-6	-	< 0.5	-
4-Bromophenyl-phenylether	101-55-3	-	< 0.5	-
Hexachlorobenzene	118-74-1	-	< 0.5	-
Pentachlorophenol	87-86-5	-	< 5.0	-
Phenanthrene	85-01-8	-	< 0.2	-
Anthracene	120-12-7	-	< 0.2	-
Di-n-butylphthalate	84-74-2	-	< 0.5	-
Fluoranthene	208-44-0	-	< 0.2	-
Pyrene	129-00-0	-	< 0.5	-
Butylbenzylphthalate	85-68-7	-	< 0.2	-
Benzolajanthracene	56-55-3	-	< 0.2	-
Chrysene	218-01-9	-	< 0.2	-
3,3'-Dichlorobenzidine	91-94-1	-	< 2.0	-
bis(2-Ethylhexyl)phthalate	117-81-7	-	< 0.5	-
Di-n-octylphthalate	117-84-0	-	< 0.2	-
Benzobifluoranthene	205-99-2	-	< 0.2	-
Benzokifluoranthene	207-08-9	-	< 0.2	-
Benzofalapyrene	50-32-8	-	< 0.2	-
Indenol1,2,3-cdpyrene	193-39-5	-	< 0.2	-
Dibenzo[a,h]anthracene	53-70-3	-	< 0.2	-
Benzofg,h,iperylene	191-24-2	-	< 0.2	-

* compounds not currently UKAS accredited

Internal Standards	% Area
1,4-Dichlorobenzene-d4	95
Naphthalene-d8	90
Acenaphthene-d10	94
Phenanthrene-d10	91
Chrysene-d12	102
Perylene-d12	113

Surrogates	% Rec
2-Fluorophenol	50
Phenol-d5	71
Nitrobenzene-d5	74
2-Fluorobiphenyl	78
2,4,6-Tribromophenol	40
Terphenyl-d14	90

Semi-Volatile Organic Compounds

Customer and Site Details:
Sample Details:
LIMS ID Number:
Job Number:

Exploration Associates: M4 Relief Road
CH/0033/ / 1.50- m
CL03954
543

16-Feb-00
18-Feb-00
24-Feb-00

Date Booked in:
Date Extracted:
Date Analysed:

Matrix:
Ext Method:
Operator:
Directory/Quant File:

Soil
Soxhlet
SW
0224ABN.MS4

QC Batch Number:
Multiplier:
Dilution Factor:

226
0.0667
1
N

0224CCCC2.D GPC (Y/N)

Target Compounds	CAS #	R.T. (min)	Concentration mg/kg	% Fit
Phenol	108-95-2	-	< 2.0	-
bis(2-Chloroethyl)ether	111-44-4	-	< 0.5	-
2-Chlorophenol	95-57-8	-	< 2.0	-
1,3-Dichlorobenzene	541-73-1	-	< 0.5	-
1,4-Dichlorobenzene	106-46-7	-	< 0.5	-
Benzyl alcohol	100-51-6	-	< 0.5	-
1,2-Dichlorobenzene	95-50-1	-	< 0.5	-
2-Methylphenol	95-48-7	-	< 0.5	-
bis(2-Chloroisopropyl)ether	108-60-1	-	< 0.5	-
Hexachloroethane	67-72-1	-	< 0.5	-
N-Nitroso-di-n-propylamine	621-64-7	-	< 0.5	-
3- & 4-Methylphenol	108-39-4/106-44-5	-	< 2.0	-
Nitrobenzene	98-95-3	-	< 0.5	-
Isophorone	78-59-1	-	< 0.5	-
2-Nitrophenol	88-75-5	-	< 2.0	-
2,4-Dimethylphenol	105-67-9	-	< 2.0	-
Benzoic Acid	65-85-0	-	< 0.5	-
bis(2-Chloroethoxy)methane	111-91-1	-	< 10.0	-
2,4-Dichlorophenol	120-83-2	-	< 2.0	-
1,2,4-Trichlorobenzene	120-82-1	-	< 0.5	-
Naphthalene	91-20-3	-	< 0.2	-
4-Chlorophenol *	106-48-9	-	< 2.0	-
4-Chloroaniline *	106-47-8	-	< 0.5	-
Hexachlorobutadiene	87-68-3	-	< 0.5	-
4-Chloro-3-methylphenol	59-50-7	-	< 0.5	-
2-Methylnaphthalene *	91-57-6	-	< 0.2	-
1-Methyl naphthalene	90-12-0	-	< 0.2	-
Hexachlorocyclopentadiene	77-47-4	-	< 0.5	-
2,4,6-Trichlorophenol	88-06-2	-	< 2.0	-
2,4,5-Trichlorophenol	95-95-4	-	< 2.0	-
2-Chloronaphthalene	91-58-7	-	< 0.2	-
Biphenyl *	92-52-4	-	< 0.2	-
Diphenyl ether *	101-84-8	-	< 0.2	-
2-Nitroaniline	88-74-4	-	< 0.5	-
Acenaphthylene	208-96-8	-	< 0.2	-
Dimethylphthalate	131-11-3	-	< 0.5	-
2,6-Dinitrotoluene	606-20-2	-	< 0.5	-
Acenaphthene *	83-32-9	-	< 0.2	-
3-Nitroaniline *	99-09-2	-	< 0.5	-

* These compounds are not currently UKAS accredited.
Concentrations are reported on a wet weight basis.

Target Compounds	CAS #	R.T.	Concentration mg/kg	% Fit
2,4-Dinitrophenol	51-28-5	-	< 1.0	-
Dibenzofuran	132-64-9	-	< 0.5	-
4-Nitrophenol	100-02-7	-	< 5.0	-
2,4-Dinitrotoluene	121-14-2	-	< 0.5	-
Fluorene	86-73-7	-	< 0.2	-
Diethylphthalate	84-66-2	-	< 0.5	-
4-Chlorophenyl-phenylether	7005-72-4	-	< 0.5	-
4,6-Dinitro-2-methylphenol	534-52-1	-	< 5.0	-
4-Nitroaniline	100-01-6	-	< 0.5	-
N-Nitrosodiphenylamine	86-30-6	-	< 0.5	-
4-Bromophenyl-phenylether	101-55-3	-	< 0.5	-
Hexachlorobenzene	118-74-1	-	< 0.5	-
Pentachlorophenol	87-86-5	-	< 0.2	-
Phenanthrene	85-01-8	-	< 0.2	-
Anthracene	120-12-7	-	< 0.2	-
Di-n-butylphthalate	84-74-2	-	< 0.5	-
Fluoranthene	206-44-0	-	< 0.2	-
Pyrene	129-00-0	-	< 0.2	-
Butylbenzylphthalate	85-68-7	-	< 0.5	-
Benzo[a]anthracene	56-55-3	-	< 0.2	-
Chrysene	218-01-9	-	< 0.2	-
3,3'-Dichlorobenzidine	91-94-1	-	< 0.5	-
bis(2-Ethylhexyl)phthalate	117-81-7	-	< 0.2	-
Di-n-octylphthalate	117-84-0	-	< 0.2	-
Benzo[b]fluoranthene	205-99-2	-	< 0.2	-
Benzo[k]fluoranthene	207-08-9	-	< 0.2	-
Benzo[a]pyrene	50-32-8	-	< 0.2	-
Indeno[1,2,3-cd]pyrene	193-39-5	-	< 0.2	-
Dibenzof[a,h]anthracene	53-70-3	-	< 0.2	-
Benzo[g,h,i]perylene	191-24-2	-	< 0.2	-

* compounds not currently UKAS accredited

Internal Standards	% Area
1,4-Dichlorobenzene-d4	90
Naphthalene-d8	80
Acenaphthene-d10	80
Phenanthrene-d10	79
Chrysene-d12	85
Perylene-d12	81

Surrogates	% Rec
2-Fluorophenol	76
Phenol-d5	86
Nitrobenzene-d5	80
2-Fluorobiphenyl	82
2,4,6-Tribromophenol	77
Terphenyl-d14	93



Semi-Volatile Organic Compounds

Customer and Site Details:
Sample Details:
LIMS ID Number:
Job Number:

Exploration Associates: M4 Relief Road
CH/00431 / 0.20- m
Date Booked in:
Date Extracted:
Date Analysed:

16-Feb-00
18-Feb-00
28-Feb-00

Matrix:
Soil
Soxhlet
SW
Operator:
Directory/Quant File:

QC Batch Number:
Multiplier:
Dilution Factor:
0.6667
10
Y

Target Compounds	CAS #	R.T. (min)	Concentration mg/kg	% Fit
Phenol	108-95-2	-	< 20.0	-
bis(2-Chloroethyl)ether	111-44-4	-	< 5.0	-
2-Chlorophenol	95-57-8	-	< 20.0	-
1,3-Dichlorobenzene	541-73-1	-	< 5.0	-
1,4-Dichlorobenzene	106-46-7	-	< 5.0	-
Benzyl alcohol	100-51-6	-	< 5.0	-
1,2-Dichlorobenzene	95-50-1	-	< 5.0	-
2-Methylphenol	95-48-7	-	< 5.0	-
bis(2-Chloroisopropyl)ether	108-60-1	-	< 5.0	-
Hexachloroethane	67-72-1	-	< 5.0	-
N-Nitroso-di-n-propylamine	621-64-7	-	< 20.0	-
3- & 4-Methylphenol	108-39-4/106-44-5	-	< 5.0	-
Nitrobenzene	98-95-3	-	< 5.0	-
Isophorone	78-59-1	-	< 20.0	-
2-Nitrophenol	88-75-5	-	< 20.0	-
2,4-Dimethylphenol	105-67-9	-	< 100	-
Benzoic Acid	65-85-0	-	< 5.0	-
bis(2-Chloroethoxy)methane	111-91-1	-	< 20.0	-
2,4-Dichlorophenol	120-83-2	-	< 5.0	-
1,2,4-Trichlorobenzene	120-82-1	-	< 2.0	-
Naphthalene	91-20-3	-	< 20.0	-
4-Chlorophenol *	106-48-9	-	< 5.0	-
4-Chloroaniline	106-47-8	-	< 5.0	-
Hexachlorobutadiene	87-68-3	-	< 5.0	-
4-Chloro-3-methylphenol	59-50-7	-	< 2.0	-
2-Methylnaphthalene	91-57-6	-	< 2.0	-
1-Methylnaphthalene *	90-12-0	-	< 5.0	-
Hexachlorocyclopentadiene	77-47-4	-	< 20.0	-
2,4,6-Trichlorophenol	88-06-2	-	< 20.0	-
2,4,5-Trichlorophenol	95-95-4	-	< 2.0	-
2-Chloronaphthalene	91-58-7	-	9.6	96
Biphenyl *	92-52-4	16.01	< 2.0	-
Diphenyl ether *	101-84-8	-	< 5.0	-
2-Nitroaniline	88-74-4	-	< 2.0	-
Acenaphthylene	208-96-8	-	< 5.0	-
Dimethylphthalate	131-11-3	-	< 5.0	-
2,6-Dinitrotoluene	606-20-2	-	< 2.0	-
Acenaphthene	83-32-9	-	< 5.0	-
3-Nitroaniline	99-09-2	-	< 5.0	-

* These compounds are not currently UKAS accredited.
Concentrations are reported on a wet weight basis.

Target Compounds	CAS #	R.T.	Concentration mg/kg	% Fit
2,4-Dinitrophenol	51-28-5	-	< 10.0	-
Dibenzofuran	132-64-9	-	< 5.0	-
4-Nitrophenol	100-02-7	-	< 50.0	-
2,4-Dinitrotoluene	121-14-2	-	< 5.0	-
Fluorene	86-73-7	19.28	2.2	95
Diethylphthalate	84-66-2	-	< 5.0	-
4-Chlorophenyl-phenylether	7005-72-4	-	< 50.0	-
4,6-Dinitro-2-methylphenol	534-52-1	-	< 5.0	-
4-Nitroaniline	100-01-6	-	< 5.0	-
N-Nitrosodiphenylamine	86-30-6	-	< 5.0	-
4-Bromophenyl-phenylether	101-55-3	-	< 5.0	-
Hexachlorobenzene	118-74-1	-	< 5.0	-
Pentachlorophenol	87-86-5	-	< 50.0	-
Phenanthrene	85-01-8	22.11	2.9	98
Anthracene	120-12-7	-	< 2.0	-
Di-n-butylphthalate	84-74-2	-	< 2.0	-
Fluoranthene	206-44-0	-	< 2.0	-
Pyrene	129-00-0	-	< 5.0	-
Butylbenzylphthalate	85-68-7	-	< 2.0	-
Benzofuranthracene	56-55-3	-	44.0	99
Chrysene	218-01-9	30.03	< 20.0	-
3,3'-Dichlorobenzidine	91-94-1	-	6.9	93
bis(2-Ethylhexyl)phthalate	117-81-7	30.43	< 2.0	-
Di-n-octylphthalate	117-84-0	-	< 2.0	-
Benzofluoranthene	205-99-2	-	< 2.0	-
Benzofluoranthene	207-08-9	-	< 2.0	-
Benzofluoranthene	50-32-8	-	< 2.0	-
Indeno[1,2,3-cd]pyrene	193-39-5	-	< 2.0	-
Dibenzofluoranthene	53-70-3	-	< 2.0	-
Benzofluoranthene	191-24-2	-	< 2.0	-

* compounds not currently UKAS accredited

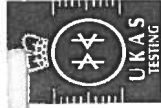
Internal Standards	% Area
1,4-Dichlorobenzene-d4	102
Naphthalene-d8	98
Acenaphthene-d10	108
Phenanthrene-d10	103
Chrysene-d12	107
Perylene-d12	111

Surrogates	% Rec
2-Fluorophenol	43 D
Phenol-d5	61
Nitrobenzene-d5	58
2-Fluorobiphenyl	81
2,4,6-Tribromophenol	D
Terphenyl-d14	18 D



The compounds listed above have been tentatively identified by a computer based library search. Compounds identified in the sample are not reported if they also occur in the method blank. The % fit is an indication of the reliability of the compound assignment. Due to the similarity between mass spectra of some isomeric compounds assignments may not be correct. Other compounds may also be present but identification was not possible. Concentrations are semi-quantitative, assume a response factor of 1 and use the nearest internal standard. Concentrations are reported on a wet weight basis.

Semi-Volatile Organic Compounds



1252

Customer and Site Details:
Sample Details:
LIMS ID Number:
Job Number:

Exploration Associates: M4 Relief Road
CH/0043/ / 1.30- m
Date Booked in:
Date Extracted:
Date Analysed:

Matrix:
Ext Method:
Operator:
Directory/Quant File:

Soil
Soxhlet
SW
0224ABN.MS5
0224CCC1.D GPC (Y/N)

QC Batch Number:
Multiplier:
Dilution Factor:

Target Compounds	CAS #	R.T. (min)	Concentration mg/kg	% Fit
Phenol	108-95-2	-	< 2.0	-
bis(2-Chloroethyl)ether	111-44-4	-	< 0.5	-
2-Chlorophenol	95-57-8	-	< 2.0	-
1,3-Dichlorobenzene	541-73-1	-	< 0.5	-
1,4-Dichlorobenzene	106-46-7	-	< 0.5	-
Benzyl alcohol	100-51-6	-	< 0.5	-
1,2-Dichlorobenzene	95-50-1	-	< 0.5	-
2-Methylphenol	95-48-7	-	< 0.5	-
bis(2-Chloroisopropyl)ether	108-60-1	-	< 0.5	-
Hexachloroethane	67-72-1	-	< 0.5	-
N-Nitroso-di-n-propylamine	621-64-7	-	< 0.5	-
3- & 4-Methylphenol	108-39-4/106-44-5	-	< 2.0	-
Nitrobenzene	98-95-3	-	< 0.5	-
Isophorone	78-59-1	-	< 0.5	-
2-Nitrophenol	88-75-5	-	< 2.0	-
2,4-Dimethylphenol	105-67-9	-	< 2.0	-
Benzoic Acid	65-85-0	-	< 10.0	-
bis(2-Chloroethoxy)methane	111-91-1	-	< 0.5	-
2,4-Dichlorophenol	120-83-2	-	< 2.0	-
1,2,4-Trichlorobenzene	120-82-1	-	< 0.5	-
Naphthalene	91-20-3	-	< 0.2	-
4-Chlorophenol *	106-48-9	-	< 2.0	-
4-Chloroaniline	106-47-8	-	< 0.5	-
Hexachlorobutadiene	87-68-3	-	< 0.5	-
4-Chloro-3-methylphenol	59-50-7	-	< 0.5	-
2-Methylnaphthalene	91-57-6	-	< 0.2	-
1-Methyl naphthalene *	90-12-0	-	< 0.2	-
Hexachlorocyclopentadiene	77-47-4	-	< 0.5	-
2,4,6-Trichlorophenol	88-06-2	-	< 2.0	-
2,4,5-Trichlorophenol	95-95-4	-	< 2.0	-
2-Chloronaphthalene	91-58-7	-	< 0.2	-
Biphenyl *	92-52-4	-	< 0.2	-
Diphenyl ether *	101-84-8	-	< 0.2	-
2-Nitroaniline	88-74-4	-	< 0.5	-
Acenaphthylene	208-96-8	-	< 0.2	-
Dimethylphthalate	131-11-3	-	< 0.5	-
2,6-Dinitrotoluene	606-20-2	-	< 0.5	-
Acenaphthene	83-32-9	-	< 0.2	-
3-Nitroaniline	99-09-2	-	< 0.5	-

* These compounds are not currently UKAS accredited.
Concentrations are reported on a wet weight basis.

Target Compounds	CAS #	R.T.	Concentration mg/kg	% Fit
2,4-Dinitrophenol	51-28-5	-	< 1.0	-
Dibenzofuran	132-64-9	-	< 0.5	-
4-Nitrophenol	100-02-7	-	< 5.0	-
2,4-Dinitrotoluene	121-14-2	-	< 0.5	-
Fluorene	86-73-7	-	< 0.2	-
Diethylphthalate	84-66-2	-	< 0.5	-
4-Chlorophenyl-phenylether	7005-72-4	-	< 0.5	-
4,6-Dinitro-2-methylphenol	534-52-1	-	< 5.0	-
4-Nitroaniline	100-01-6	-	< 0.5	-
N-Nitrosodiphenylamine	86-30-6	-	< 0.5	-
4-Bromophenyl-phenylether	101-55-3	-	< 0.5	-
Hexachlorobenzene	118-74-1	-	< 0.5	-
Pentachlorophenol	87-86-5	-	< 5.0	-
Phenanthrene	85-01-8	-	< 0.2	-
Anthracene	120-12-7	-	< 0.2	-
Di-n-butylphthalate	84-74-2	-	< 0.5	-
Fluoranthene	206-44-0	-	< 0.2	-
Pyrene	129-00-0	-	< 0.2	-
Butylbenzylphthalate	85-68-7	-	< 0.5	-
Benzofluoranthene	56-55-3	-	< 0.2	-
Chrysene	218-01-9	-	< 0.2	-
3,3'-Dichlorobenzidine	91-94-1	-	< 2.0	-
bis(2-Ethylhexyl)phthalate	117-81-7	29.36	5.6	96
Di-n-octylphthalate	117-84-0	-	< 0.2	-
Benzofluoranthene	205-99-2	-	< 0.2	-
Benzofluoranthene	207-08-9	-	< 0.2	-
Benzofluoranthene	50-32-8	-	< 0.2	-
Indeno[1,2,3-cd]pyrene	193-39-5	-	< 0.2	-
Dibenzofluoranthene	53-70-3	-	< 0.2	-
Benzofluoranthene	191-24-2	-	< 0.2	-

* compounds not currently UKAS accredited

Internal Standards	% Area
1,4-Dichlorobenzene-d4	96
Naphthalene-d8	92
Acenaphthene-d10	96
Phenanthrene-d10	92
Chrysene-d12	101
Perylene-d12	108

Surrogates	% Rec
2-Fluorophenol	67
Phenol-d5	79
Nitrobenzene-d5	74
2-Fluorobiphenyl	78
2,4,6-Tribromophenol	70
Terphenyl-d14	94

Semi-Volatile Organic Compounds

Customer and Site Details:

Sample Details:
LIMS ID Number:
Job Number:

Exploration Associates: M4 Relief Road

CH/0044/1/0.30- m
Date Booked in:
Date Extracted:
Date Analysed:

16-Feb-00
18-Feb-00
24-Feb-00

Matrix:

Ext Method:
Operator:
Directory/Quant File:

Soil
Soxhlet
SW
0224ABN.MS4

QC Batch Number:

Multiplier:

Dilution Factor:

226

0.1333

2

Y

Target Compounds	CAS #	R.T. (min)	Concentration mg/kg	% Fit
Phenol	108-95-2	-	< 4.0	-
bis(2-Chloroethyl)ether	111-44-4	-	< 1.0	-
2-Chlorophenol	95-57-8	-	< 4.0	-
1,3-Dichlorobenzene	541-73-1	-	< 1.0	-
1,4-Dichlorobenzene	106-46-7	-	< 1.0	-
Benzyl alcohol	100-51-6	-	< 1.0	-
1,2-Dichlorobenzene	95-50-1	-	< 1.0	-
2-Methylphenol	95-48-7	-	< 1.0	-
bis(2-Chloroisopropyl)ether	108-60-1	-	< 1.0	-
Hexachloroethane	67-72-1	-	< 1.0	-
N-Nitroso-di-n-propylamine	621-64-7	-	< 4.0	-
3- & 4-Methylphenol	108-39-4/106-44-5	-	< 1.0	-
Nitrobenzene	98-95-3	-	< 1.0	-
Isophorone	78-59-1	-	< 1.0	-
2-Nitrophenol	88-75-5	-	< 4.0	-
2,4-Dimethylphenol	105-67-9	-	< 4.0	-
Benzoic Acid	65-85-0	-	< 20.0	-
bis(2-Chloroethoxy)methane	111-91-1	-	< 1.0	-
2,4-Dichlorophenol	120-83-2	-	< 4.0	-
1,2,4-Trichlorobenzene	120-82-1	-	< 1.0	-
Naphthalene	91-20-3	-	< 0.4	-
4-Chlorophenol *	106-48-9	-	< 4.0	-
4-Chloroaniline *	106-47-8	-	< 1.0	-
Hexachlorobutadiene	87-68-3	-	< 1.0	-
4-Chloro-3-methylphenol	59-50-7	-	< 1.0	-
2-Methylnaphthalene	91-57-6	-	< 0.4	-
1-Methyl naphthalene *	90-12-0	-	< 0.4	-
Hexachlorocyclopentadiene	77-47-4	-	< 1.0	-
2,4,6-Trichlorophenol	88-06-2	-	< 4.0	-
2,4,5-Trichlorophenol	95-95-4	-	< 4.0	-
2-Chloronaphthalene	91-58-7	-	< 0.4	-
Biphenyl *	92-52-4	15.37	4.8	100
Diphenyl ether *	101-84-8	-	< 0.4	-
2-Nitroaniline	88-74-4	-	< 1.0	-
Acenaphthylene	208-96-8	-	< 0.4	-
Dimethylphthalate	131-11-3	-	< 1.0	-
2,6-Dinitrotoluene	606-20-2	-	< 1.0	-
Acenaphthene	83-32-9	-	< 0.4	-
3-Nitroaniline *	99-09-2	-	< 1.0	-

* These compounds are not currently UKAS accredited.
Concentrations are reported on a wet weight basis.

Target Compounds	CAS #	R.T.	Concentration mg/kg	% Fit
2,4-Dinitrophenol	51-28-5	-	< 2.0	-
Dibenzofuran	132-64-9	-	< 1.0	-
4-Nitrophenol	100-02-7	-	< 10.0	-
2,4-Dinitrotoluene	121-14-2	-	< 1.0	-
Fluorene	86-73-7	18.57	0.8	100
Diethylphthalate	84-66-2	-	< 1.0	-
4-Chlorophenyl-phenylether	7005-72-4	-	< 1.0	-
4,6-Dinitro-2-methylphenol	534-52-1	-	< 10.0	-
4-Nitroaniline	100-01-6	-	< 1.0	-
N-Nitrosodiphenylamine	86-30-6	-	< 1.0	-
4-Bromophenyl-phenylether	101-55-3	-	< 1.0	-
Hexachlorobenzene	118-74-1	-	< 1.0	-
Pentachlorophenol	87-86-5	-	< 10.0	-
Phenanthrene	85-01-8	21.37	1.1	99
Anthracene	120-12-7	-	< 0.4	-
Di-n-butylphthalate	84-74-2	-	< 1.0	-
Fluoranthene	206-44-0	24.91	0.6	99
Pyrene	129-00-0	25.54	0.5	92
Butylbenzylphthalate	85-68-7	-	< 0.4	-
Benzofuran	56-55-3	-	< 1.0	-
Chrysene	218-01-9	29.34	20.4	98
3,3'-Dichlorobenzidine	91-94-1	-	< 4.0	-
bis(2-Ethylhexyl)phthalate	117-81-7	29.89	4.4	98
Di-n-octylphthalate	117-84-0	-	< 0.4	-
Benzobifluoranthene	205-99-2	-	< 0.4	-
Benzokjfluoranthene	207-08-9	32.29	0.7	M
Benzofluoranthene	50-32-8	33.23	0.6	87
Indeno[1,2,3-cd]pyrene	193-39-5	37.85	0.4	98
Dibenzofluoranthene	53-70-3	-	< 0.4	-
Benzofluoranthene	191-24-2	-	< 0.4	-

* compounds not currently UKAS accredited

Internal Standards	% Area
1,4-Dichlorobenzene-d4	87
Naphthalene-d8	79
Acenaphthene-d10	82
Phenanthrene-d10	86
Chrysene-d12	87
Perylene-d12	75

Surrogates	% Rec
2-Fluorophenol	53
Phenol-d5	69
Nitrobenzene-d5	65
2-Fluorobiphenyl	80
2,4,6-Tribromophenol	78
Terphenyl-d14	14



Semi-Volatile Organic Compounds

Customer and Site Details:
Sample Details:
LIMS ID Number:
Job Number:

Exploration Associates: M4 Relief Road
CH/0044/ / 1.00- m
Date Booked in:
Date Extracted:
Date Analysed:

16-Feb-00
18-Feb-00
24-Feb-00

Matrix: Soil
Ext Method: Soxhlet
Operator: SW
Directory/Quant File: 0224ABN.MS4
QC Batch Number: 226
Multiplier: 0.1333
Dilution Factor: 2
Y

Target Compounds	CAS #	R.T. (min)	Concentration mg/kg	% Fit
Phenol	108-95-2	-	< 4.0	-
bis(2-Chloroethyl)ether	111-44-4	-	< 1.0	-
2-Chlorophenol	95-57-8	-	< 4.0	-
1,3-Dichlorobenzene	541-73-1	-	< 1.0	-
1,4-Dichlorobenzene	106-46-7	-	< 1.0	-
Benzyl alcohol	100-51-6	-	< 1.0	-
1,2-Dichlorobenzene	95-50-1	-	< 1.0	-
2-Methylphenol	95-48-7	-	< 1.0	-
bis(2-Chloroisopropyl)ether	108-60-1	-	< 1.0	-
Hexachloroethane	67-72-1	-	< 1.0	-
N-Nitroso-di-n-propylamine	621-64-7	-	< 1.0	-
3- & 4-Methylphenol	108-39-4/106-44-5	-	< 4.0	-
Nitrobenzene	98-95-3	-	< 1.0	-
Isophorone	78-59-1	-	< 1.0	-
2-Nitrophenol	88-75-5	-	< 4.0	-
2,4-Dimethylphenol	105-67-9	-	< 4.0	-
Benzoic Acid	65-85-0	-	< 20.0	-
bis(2-Chloroethoxy)methane	111-91-1	-	< 1.0	-
2,4-Dichlorophenol	120-83-2	-	< 4.0	-
1,2,4-Trichlorobenzene	120-82-1	-	< 1.0	-
Naphthalene	91-20-3	-	< 0.4	-
4-Chlorophenol *	106-48-9	-	< 4.0	-
4-Chloroaniline *	106-47-8	-	< 1.0	-
Hexachlorobutadiene	87-68-3	-	< 1.0	-
4-Chloro-3-methylphenol	59-50-7	-	< 1.0	-
2-Methylnaphthalene	91-57-6	-	< 0.4	-
1-Methyl naphthalene *	90-12-0	-	< 0.4	-
Hexachlorocyclopentadiene	77-47-4	-	< 1.0	-
2,4,6-Trichlorophenol	88-06-2	-	< 4.0	-
2,4,5-Trichlorophenol	95-95-4	-	< 4.0	-
2-Chloronaphthalene	91-58-7	-	< 0.4	-
Biphenyl *	92-52-4	15.37	7.8	99
Diphenyl ether *	101-84-8	-	< 0.4	-
2-Nitroaniline	88-74-4	-	< 1.0	-
Acenaphthylene	208-96-8	-	< 0.4	-
Dimethylphthalate	131-11-3	-	< 1.0	-
2,6-Dinitrotoluene	606-20-2	-	< 1.0	-
Acenaphthene	83-32-9	-	< 0.4	-
3-Nitroaniline *	99-09-2	-	< 1.0	-

* These compounds are not currently UKAS accredited.
Concentrations are reported on a wet weight basis.

Target Compounds	CAS #	R.T.	Concentration mg/kg	% Fit
2,4-Dinitrophenol	51-28-5	-	< 2.0	-
Dibenzofuran	132-64-9	-	< 1.0	-
4-Nitrophenol	100-02-7	-	< 10.0	-
2,4-Dinitrotoluene	121-14-2	-	< 1.0	-
Fluorene	86-73-7	18.58	1.3	99
Diethylphthalate	84-66-2	-	< 1.0	-
4-Chlorophenyl-phenylether	7005-72-4	-	< 1.0	-
4,6-Dinitro-2-methylphenol	534-52-1	-	< 10.0	-
4-Nitroaniline	100-01-6	-	< 1.0	-
N-Nitrosodiphenylamine	86-30-6	-	< 1.0	-
4-Bromophenyl-phenylether	101-55-3	-	< 1.0	-
Hexachlorobenzene	118-74-1	-	< 1.0	-
Pentachlorophenol	87-86-5	-	< 10.0	-
Phenanthrene	85-01-8	21.37	1.7	99
Anthracene	120-12-7	-	< 0.4	-
Di-n-butylphthalate	84-74-2	-	< 1.0	-
Fluoranthene	208-44-0	24.91	1.2	92
Pyrene	129-00-0	25.54	0.6	93
Butylbenzylphthalate	85-68-7	-	< 1.0	-
Benzofuranthracene	56-55-3	29.22	0.5	63
Chrysene	218-01-9	29.36	27.8	96
3,3'-Dichlorobenzidine	91-94-1	-	< 4.0	-
bis(2-Ethylhexyl)phthalate	117-81-7	29.89	7.7	99
Di-n-octylphthalate	117-84-0	-	< 0.4	-
Benzofluoranthene	205-99-2	32.26	0.6	97
Benzokjfluoranthene	207-08-9	32.31	1.2	M
Benzofluorene	50-32-8	-	< 0.4	-
Indenol1,2,3-copolyrene	193-39-5	37.82	0.7	96
Dibenzofluoranthracene	53-70-3	-	< 0.4	-
Benzofluoranthene	191-24-2	38.66	0.6	98

* compounds not currently UKAS accredited

Internal Standards	% Area
1,4-Dichlorobenzene-d4	82
Naphthalene-d8	74
Acenaphthene-d10	79
Phenanthrene-d10	84
Chrysene-d12	84
Perylene-d12	73

Surrogates	% Rec
2-Fluorophenol	57
Phenol-d5	73
Nitrobenzene-d5	67
2-Fluorobiphenyl	82
2,4,6-Tribromophenol	89
Terphenyl-d14	4

Semi-Volatile Organic Compounds

Customer and Site Details:
Sample Details:
LIMS ID Number:
Job Number:

Exploration Associates: M4 Relief Road
CH/0045/ / 0.20- m
CL03959
543

Date Booked in: 16-Feb-00
Date Extracted: 18-Feb-00
Date Analysed: 24-Feb-00

Matrix: Soil
Ext Method: Soxhlet
Operator: SW
Directory/Quant File: 0224ABN.MS5

QC Batch Number: 226
Multiplier: 0.1667
Dilution Factor: 2.5
N

Target Compounds	CAS #	R.T. (min)	Concentration mg/kg	% Fit
Phenol	108-95-2	-	< 5.0	-
bis(2-Chloroethyl)ether	111-44-4	-	< 1.3	-
2-Chlorophenol	95-57-8	-	< 5.0	-
1,3-Dichlorobenzene	541-73-1	-	< 1.3	-
1,4-Dichlorobenzene	106-46-7	-	< 1.3	-
Benzyl alcohol	100-51-6	-	< 1.3	-
1,2-Dichlorobenzene	95-50-1	-	< 1.3	-
2-Methylphenol	95-48-7	-	< 1.3	-
bis(2-Chloroisopropyl)ether	108-60-1	-	< 1.3	-
Hexachloroethane	67-72-1	-	< 1.3	-
N-Nitroso-di-n-propylamine	621-64-7	-	< 1.3	-
3- & 4-Methylphenol	108-39-4/106-44-5	-	< 5.0	-
Nitrobenzene	98-95-3	-	< 1.3	-
Isophorone	78-59-1	-	< 1.3	-
2-Nitrophenol	88-75-5	-	< 5.0	-
2,4-Dimethylphenol	105-67-9	-	< 5.0	-
Benzoic Acid	65-85-0	-	< 25.0	-
bis(2-Chloroethoxy)methane	111-91-1	-	< 1.3	-
2,4-Dichlorophenol	120-83-2	-	< 5.0	-
1,2,4-Trichlorobenzene	120-82-1	-	< 1.3	-
Naphthalene	91-20-3	-	< 0.5	-
4-Chlorophenol *	106-48-9	-	< 5.0	-
4-Chloroaniline	106-47-8	-	< 1.3	-
Hexachlorobutadiene	87-68-3	-	< 1.3	-
4-Chloro-3-methylphenol	59-50-7	-	< 1.3	-
2-Methylnaphthalene	91-57-6	-	< 0.5	-
1-Methyl naphthalene *	90-12-0	-	< 0.5	-
Hexachlorocyclopentadiene	77-47-4	-	< 1.3	-
2,4,6-Trichlorophenol	88-06-2	-	< 5.0	-
2,4,5-Trichlorophenol	95-95-4	-	< 5.0	-
2-Chloronaphthalene	91-58-7	-	< 0.5	-
Biphenyl *	92-52-4	-	< 0.5	-
Diphenyl ether *	101-84-8	-	< 0.5	-
2-Nitroaniline	88-74-4	-	< 1.3	-
Acenaphthylene	208-96-8	-	< 0.5	-
Dimethylphthalate	131-11-3	-	< 1.3	-
2,6-Dinitrotoluene	606-20-2	-	< 1.3	-
Acenaphthene	83-32-9	-	< 0.5	-
3-Nitroaniline	99-09-2	-	< 1.3	-

* These compounds are not currently UKAS accredited.
Concentrations are reported on a wet weight basis.

Target Compounds	CAS #	R.T.	Concentration mg/kg	% Fit
2,4-Dinitrophenol	51-28-5	-	< 2.5	-
Dibenzofuran	132-64-9	-	< 1.3	-
4-Nitrophenol	100-02-7	-	< 12.5	-
2,4-Dinitrotoluene	121-14-2	-	< 1.3	-
Fluorene	86-73-7	-	< 0.5	-
Diethylphthalate	84-66-2	-	< 1.3	-
4-Chlorophenyl-phenylether	7005-72-4	-	< 1.3	-
4,6-Dinitro-2-methylphenol	534-52-1	-	< 12.5	-
4-Nitroaniline	100-01-6	-	< 1.3	-
N-Nitrosodiphenylamine	86-30-6	-	< 1.3	-
4-Bromophenyl-phenylether	101-55-3	-	< 1.3	-
Hexachlorobenzene	118-74-1	-	< 1.3	-
Pentachlorophenol	87-86-5	-	< 12.5	-
Phenanthrene	85-01-8	-	< 0.5	-
Anthracene	120-12-7	-	< 0.5	-
Di-n-butylphthalate	84-74-2	-	< 1.3	-
Fluoranthene	206-44-0	-	< 0.5	-
Pyrene	129-00-0	-	< 0.5	-
Butylbenzylphthalate	85-68-7	-	< 1.3	-
Benzofuranthracene	56-55-3	-	< 0.5	-
Chrysene	218-01-9	-	< 0.5	-
3,3'-Dichlorobenzidine	91-94-1	-	< 5.0	-
bis(2-Ethylhexyl)phthalate	117-81-7	-	< 1.3	-
Di-n-octylphthalate	117-84-0	-	< 0.5	-
Benzofluoranthene	205-99-2	-	< 0.5	-
Benzokjfluoranthene	207-08-9	-	< 0.5	-
Benzofluoranthene	50-32-8	-	< 0.5	-
Indenof1,2,3-cdipyrene	193-39-5	-	< 0.5	-
Dibenzofluoranthene	53-70-3	-	< 0.5	-
Benzofluoranthene	191-24-2	-	< 0.5	-

* compounds not currently UKAS accredited

Internal Standards	% Area
1,4-Dichlorobenzene-d4	95
Naphthalene-d8	90
Acenaphthene-d10	93
Phenanthrene-d10	93
Chrysene-d12	108
Perylene-d12	128

Surrogates	% Rec
2-Fluorophenol	54
Phenol-d5	69
Nitrobenzene-d5	64
2-Fluorobiphenyl	90
2,4,6-Tribromophenol	34
Terphenyl-d14	92



Semi-Volatile Organic Compounds

Customer and Site Details:
Sample Details:
LIMS ID Number:
Job Number:

Exploration Associates: M4 Relief Road
CH/0046/ / 0.50- m
CL03961
543

16-Feb-00
18-Feb-00
24-Feb-00

Matrix:
Ext Method:
Operator:
Directory/Quant File:

Soil
Soxhlet
SW
0224ABN.MS4 0224CCC2.D GPC (Y/N)

QC Batch Number: 226
Multiplier: 0.6667
Dilution Factor: 10
Y

Target Compounds	CAS #	R.T. (min)	Concentration mg/kg	% Fit
Phenol	108-95-2	-	< 20.0	-
bis(2-Chloroethyl)ether	111-44-4	-	< 5.0	-
2-Chlorophenol	95-57-8	-	< 20.0	-
1,3-Dichlorobenzene	541-73-1	-	< 5.0	-
1,4-Dichlorobenzene	106-46-7	-	< 5.0	-
Benzyl alcohol	100-51-6	-	< 5.0	-
1,2-Dichlorobenzene	95-50-1	-	< 5.0	-
2-Methylphenol	95-48-7	-	< 5.0	-
bis(2-Chloroisopropyl)ether	108-60-1	-	< 5.0	-
Hexachloroethane	67-72-1	-	< 5.0	-
N-Nitroso-di-n-propylamine	621-64-7	-	< 5.0	-
3- & 4-Methylphenol	108-39-4/106-44-5	-	< 20.0	-
Nitrobenzene	98-95-3	-	< 5.0	-
Isophorone	78-59-1	-	< 5.0	-
2-Nitrophenol	88-75-5	-	< 20.0	-
2,4-Dimethylphenol	105-67-9	-	< 20.0	-
Benzoic Acid	65-85-0	-	< 100	-
bis(2-Chloroethoxy)methane	111-91-1	-	< 5.0	-
2,4-Dichlorophenol	120-83-2	-	< 20.0	-
1,2,4-Trichlorobenzene	120-82-1	-	< 5.0	-
Naphthalene	91-20-3	-	< 2.0	-
4-Chlorophenol *	106-48-9	-	< 20.0	-
4-Chloroaniline *	106-47-8	-	< 5.0	-
Hexachlorobutadiene	87-68-3	-	< 5.0	-
4-Chloro-3-methylphenol	59-50-7	-	< 5.0	-
2-Methylnaphthalene	91-57-6	-	< 2.0	-
1-Methyl naphthalene *	90-12-0	-	< 2.0	-
Hexachlorocyclopentadiene	77-47-4	-	< 5.0	-
2,4,6-Trichlorophenol	88-06-2	-	< 20.0	-
2,4,5-Trichlorophenol	95-95-4	-	< 20.0	-
2-Chloronaphthalene	91-58-7	-	< 2.0	-
Biphenyl *	92-52-4	15.37	50.7	100
Diphenyl ether *	101-84-8	-	< 2.0	-
2-Nitroaniline	88-74-4	-	< 5.0	-
Acenaphthylene	208-96-8	-	< 2.0	-
Dimethylphthalate	131-11-3	-	< 5.0	-
2,6-Dinitrotoluene	606-20-2	-	< 5.0	-
Acenaphthene	83-32-9	-	< 2.0	-
3-Nitroaniline *	99-09-2	-	< 5.0	-

* These compounds are not currently UKAS accredited.
Concentrations are reported on a wet weight basis.

Target Compounds	CAS #	R.T.	Concentration mg/kg	% Fit
2,4-Dinitrophenol	51-28-5	-	< 10.0	-
Dibenzofuran	132-64-9	-	< 5.0	-
4-Nitrophenol	100-02-7	-	< 50.0	-
2,4-Dinitrotoluene	121-14-2	-	< 5.0	-
Fluorene	86-73-7	18.57	22.5	99
Diethylphthalate	84-66-2	-	< 5.0	-
4-Chlorophenyl-phenylether	7005-72-4	-	< 5.0	-
4,6-Dinitro-2-methylphenol	534-52-1	-	< 50.0	-
4-Nitroaniline	100-01-6	-	< 5.0	-
N-Nitrosodiphenylamine	86-30-6	-	< 5.0	-
4-Bromophenyl-phenylether	101-55-3	-	< 5.0	-
Hexachlorobenzene	118-74-1	-	< 5.0	-
Pentachlorophenol	87-86-5	-	< 5.0	-
Phenanthrene	85-01-8	21.38	37.3	100
Anthracene	120-12-7	21.49	4.1	99
Di-n-butylphthalate	84-74-2	-	< 5.0	-
Fluoranthene	206-44-0	24.91	2.8	97
Pyrene	129-00-0	-	< 2.0	-
Butylbenzylphthalate	65-68-7	-	< 5.0	-
Benzol[a]anthracene	56-55-3	-	< 2.0	-
Chrysene	218-01-9	29.39	158	98
3,3'-Dichlorobenzidine	91-94-1	-	< 20.0	-
bis(2-Ethylhexyl)phthalate	117-81-7	-	< 5.0	-
Di-n-octylphthalate	117-84-0	-	< 2.0	-
Benzol[b]fluoranthene	205-99-2	-	< 2.0	-
Benzol[k]fluoranthene	207-08-9	-	< 2.0	-
Benzol[a]pyrene	50-32-8	-	< 2.0	-
Indeno[1,2,3-cd]pyrene	193-39-5	-	< 2.0	-
Dibenzol[a,h]anthracene	53-70-3	-	< 2.0	-
Benzol[g,h,i]perylene	191-24-2	-	< 2.0	-

* compounds not currently UKAS accredited

Internal Standards	% Area
1,4-Dichlorobenzene-d4	83
Naphthalene-d8	73
Acenaphthene-d10	76
Phenanthrene-d10	74
Chrysene-d12	88
Perylene-d12	92

Surrogates	% Rec
2-Fluorophenol	46 D
Phenol-d5	60
Nitrobenzene-d5	69
2-Fluorobiphenyl	91
2,4,6-Tribromophenol	76
Terphenyl-d14	D



Semi-Volatile Organic Compounds

Customer and Site Details:
Sample Details:
LIMS ID Number:
Job Number:

Exploration Associates: M4 Relief Road
CH/0046/ / 1,20- m
CL03960
543

16-Feb-00
18-Feb-00
24-Feb-00

Matrix: Soil
Ext Method: Soxhlet
Operator: SW
Directory/Quant File: 0224ABN.MS4 0224CCC3.D GPC (Y/N)

QC Batch Number: 226
Multiplier: 0.0667
Dilution Factor: 1
Y

Target Compounds	CAS #	R.T. (min)	Concentration mg/kg	% Fit
Phenol	108-95-2	-	< 2.0	-
bis(2-Chloroethyl)ether	111-44-4	-	< 0.5	-
2-Chlorophenol	95-57-8	-	< 2.0	-
1,3-Dichlorobenzene	541-73-1	-	< 0.5	-
1,4-Dichlorobenzene	106-46-7	-	< 0.5	-
Benzyl alcohol	100-51-6	-	< 0.5	-
1,2-Dichlorobenzene	95-50-1	-	< 0.5	-
2-Methylphenol	95-48-7	-	< 0.5	-
bis(2-Chloroisopropyl)ether	108-60-1	-	< 0.5	-
Hexachloroethane	67-72-1	-	< 0.5	-
N-Nitroso-di-n-propylamine	621-64-7	-	< 2.0	-
3- & 4-Methylphenol	108-39-4/106-44-5	-	< 0.5	-
Nitrobenzene	98-95-3	-	< 0.5	-
Isophorone	78-59-1	-	< 0.5	-
2-Nitrophenol	88-75-5	-	< 2.0	-
2,4-Dimethylphenol	105-67-9	-	< 2.0	-
Benzoic Acid	65-85-0	-	< 10.0	-
bis(2-Chloroethoxy)methane	111-91-1	-	< 0.5	-
2,4-Dichlorophenol	120-83-2	-	< 2.0	-
1,2,4-Trichlorobenzene	120-82-1	-	< 0.5	-
Naphthalene	91-20-3	-	< 0.2	-
4-Chlorophenol *	106-48-9	-	< 2.0	-
4-Chloroaniline *	106-47-8	-	< 0.5	-
Hexachlorobutadiene	87-68-3	-	< 0.5	-
4-Chloro-3-methylphenol	59-50-7	-	< 0.5	-
2-Methylnaphthalene	91-57-6	-	< 0.2	-
1-Methyl naphthalene *	90-12-0	-	< 0.2	-
Hexachlorocyclopentadiene	77-47-4	-	< 0.5	-
2,4,6-Trichlorophenol	88-06-2	-	< 2.0	-
2,4,5-Trichlorophenol	95-95-4	-	< 0.2	-
2-Chloronaphthalene	91-58-7	-	< 0.2	-
Biphenyl *	92-52-4	15.36	1.4	99
Diphenyl ether *	101-84-8	-	< 0.2	-
2-Nitroaniline	88-74-4	-	< 0.5	-
Acenaphthylene	208-96-8	-	< 0.2	-
Dimethylphthalate	131-11-3	-	< 0.5	-
2,6-Dinitrotoluene	606-20-2	-	< 0.5	-
Acenaphthene	83-32-9	-	< 0.2	-
3-Nitroaniline *	99-09-2	-	< 0.5	-

* These compounds are not currently UKAS accredited.
Concentrations are reported on a wet weight-basis.

Target Compounds	CAS #	R.T.	Concentration mg/kg	% Fit
2,4-Dinitrophenol	51-28-5	-	< 1.0	-
Dibenzofuran	132-64-9	-	< 0.5	-
4-Nitrophenol	100-02-7	-	< 5.0	-
2,4-Dinitrotoluene	121-14-2	-	< 0.5	-
Fluorene	86-73-7	18.57	0.3	98
Diethylphthalate	84-66-2	-	< 0.5	-
4-Chlorophenyl-phenylether	7005-72-4	-	< 0.5	-
4,6-Dinitro-2-methylphenol	534-52-1	-	< 5.0	-
4-Nitroaniline	100-01-6	-	< 0.5	-
N-Nitrosodiphenylamine	86-30-6	-	< 0.5	-
4-Bromophenyl-phenylether	101-55-3	-	< 0.5	-
Hexachlorobenzene	118-74-1	-	< 0.5	-
Pentachlorophenol	87-86-5	-	< 5.0	-
Phenanthrene	85-01-8	21.37	0.5	99
Anthracene	120-12-7	-	< 0.2	-
Di-n-butylphthalate	84-74-2	-	< 0.5	-
Fluoranthene	206-44-0	24.91	0.6	97
Pyrene	129-00-0	25.54	0.5	85
Butylbenzylphthalate	85-66-7	-	< 0.5	-
Benzo[a]anthracene	56-55-3	29.16	0.3	93
Chrysene	218-01-9	29.29	2.9	99
3,3'-Dichlorobenzidine	91-94-1	-	< 2.0	-
bis(2-Ethylhexyl)phthalate	117-81-7	-	< 0.5	-
Di-n-octylphthalate	117-84-0	-	< 0.2	-
Benzo[b]fluoranthene	205-99-2	32.22	0.3	98
Benzo[k]fluoranthene	207-08-9	32.29	0.4	99
Benzo[a]pyrene	50-32-8	33.21	0.3	92
Indeno[1,2,3-cd]pyrene	193-39-5	37.86	0.2	94
Dibenzo[a,h]anthracene	53-70-3	-	< 0.2	-
Benzo[g,h,i]perylene	191-24-2	38.64	0.2	99

* compounds not currently UKAS accredited

Internal Standards	% Area
1,4-Dichlorobenzene-d4	91
Naphthalene-d8	82
Acenaphthene-d10	85
Phenanthrene-d10	84
Chrysene-d12	93
Perylene-d12	81

Surrogates	% Rec
2-Fluorophenol	39
Phenol-d5	59
Nitrobenzene-d5	68
2-Fluorobiphenyl	76
2,4,6-Tribromophenol	44
Terphenyl-d14	6

[illegible]

The compounds listed above have been tentatively identified by a computer based library search. Compounds identified in the sample are not reported if they also occur in the method blank. The % fit is an indication of the reliability of the compound assignment. Due to the similarity between mass spectra of some isomeric compounds assignments may not be correct. Other compounds may also be present but identification was not possible. Concentrations are semi-quantitative, assume a response factor of 1 and use the nearest internal standard. Concentrations are reported on a wet weight basis.

Semi-Volatile Organic Compounds

Customer and Site Details:

Sample Details:
LIMS ID Number:
Job Number:

Exploration Associates: M4 Relief Road
CH/00471/ / 0.10- m
Date Booked in:
Date Extracted:
Date Analysed:

Matrix:
Ext Method:
Operator:
Directory/Quant File:

Soil
Soxhlet
SW
0224ABN.MS4

QC Batch Number:
Multiplier:
Dilution Factor:

226
0.0667
1
Y

Target Compounds	CAS #	R.T. (min)	Concentration mg/kg	% Fit
Phenol	108-95-2	-	< 2.0	-
bis(2-Chloroethyl)ether	111-44-4	-	< 0.5	-
2-Chlorophenol	95-57-8	-	< 2.0	-
1,3-Dichlorobenzene	541-73-1	-	< 0.5	-
1,4-Dichlorobenzene	106-46-7	-	< 0.5	-
Benzyl alcohol	100-51-6	-	< 0.5	-
1,2-Dichlorobenzene	95-50-1	-	< 0.5	-
2-Methylphenol	95-48-7	-	< 0.5	-
bis(2-Chloroisopropyl)ether	108-60-1	-	< 0.5	-
Hexachloroethane	67-72-1	-	< 0.5	-
N-Nitroso-di-n-propylamine	621-64-7	-	< 0.5	-
3- & 4-Methylphenol	108-39-4/106-44-5	-	< 2.0	-
Nitrobenzene	98-95-3	-	< 0.5	-
Isophorone	78-59-1	-	< 0.5	-
2-Nitrophenol	88-75-5	-	< 2.0	-
2,4-Dimethylphenol	105-67-9	-	< 2.0	-
Benzoic Acid	65-85-0	-	< 10.0	-
bis(2-Chloroethoxy)methane	111-91-1	-	< 0.5	-
2,4-Dichlorophenol	120-83-2	-	< 2.0	-
1,2,4-Trichlorobenzene	120-82-1	-	< 0.5	-
Naphthalene	91-20-3	-	< 0.2	-
4-Chlorophenol *	106-48-9	-	< 2.0	-
4-Chloroaniline	106-47-8	-	< 0.5	-
Hexachlorobutadiene	87-68-3	-	< 0.5	-
4-Chloro-3-methylphenol	59-50-7	-	< 0.5	-
2-Methylnaphthalene	91-57-6	-	< 0.2	-
1-Methyl naphthalene *	90-12-0	-	< 0.2	-
Hexachlorocyclopentadiene	77-47-4	-	< 0.5	-
2,4,6-Trichlorophenol	88-06-2	-	< 2.0	-
2,4,5-Trichlorophenol	95-95-4	-	< 2.0	-
2-Chloronaphthalene	91-58-7	-	< 0.2	-
Biphenyl *	92-52-4	15.36	0.5	99
Diphenyl ether *	101-84-8	-	< 0.2	-
2-Nitroaniline	88-74-4	-	< 0.5	-
Acenaphthylene	208-96-8	-	< 0.2	-
Dimethylphthalate	131-11-3	-	< 0.5	-
2,6-Dinitrotoluene	606-20-2	-	< 0.5	-
Acenaphthene	83-32-9	-	< 0.2	-
3-Nitroaniline	99-09-2	-	< 0.5	-

* These compounds are not currently UKAS accredited.
Concentrations are reported on a wet weight basis.

Target Compounds	CAS #	R.T.	Concentration mg/kg	% Fit
2,4-Dinitrophenol	51-28-5	-	< 1.0	-
Dibenzofuran	132-64-9	-	< 0.5	-
4-Nitrophenol	100-02-7	-	< 5.0	-
2,4-Dinitrotoluene	121-14-2	-	< 0.5	-
Fluorene	86-73-7	-	< 0.2	-
Diethylphthalate	84-66-2	-	< 0.5	-
4-Chlorophenyl-phenylether	7005-72-4	-	< 0.5	-
4,6-Dinitro-2-methylphenol	534-52-1	-	< 5.0	-
4-Nitroaniline	100-01-6	-	< 0.5	-
N-Nitrosodiphenylamine	86-30-6	-	< 0.5	-
4-Bromophenyl-phenylether	101-55-3	-	< 0.5	-
Hexachlorobenzene	118-74-1	-	< 0.5	-
Pentachlorophenol	87-86-5	-	< 5.0	-
Phenanthrene	85-01-8	-	< 0.2	-
Anthracene	120-12-7	-	< 0.2	-
Di-n-butylphthalate	84-74-2	-	< 0.5	-
Fluoranthene	206-44-0	-	< 0.2	-
Pyrene	129-00-0	-	< 0.2	-
Butylbenzylphthalate	85-68-7	-	< 0.5	-
Benzofluoranthene	56-55-3	-	< 0.2	-
Chrysene	218-01-9	29.26	0.3	97
3,3'-Dichlorobenzidine	91-94-1	-	< 2.0	-
bis(2-Ethylhexyl)phthalate	117-81-7	29.88	3.8	99
Di-n-octylphthalate	117-84-0	-	< 0.2	-
Benzofluoranthene	205-99-2	-	< 0.2	-
Benzofluoranthene	207-08-9	-	< 0.2	-
Benzofluoranthene	50-32-8	-	< 0.2	-
Indenol(1,2,3-cd)pyrene	193-39-5	-	< 0.2	-
Dibenzofluoranthene	53-70-3	-	< 0.2	-
Benzofluoranthene	191-24-2	-	< 0.2	-

* compounds not currently UKAS accredited

Internal Standards	% Area
1,4-Dichlorobenzene-d4	91
Naphthalene-d8	85
Acenaphthene-d10	87
Phenanthrene-d10	89
Chrysene-d12	86
Perylene-d12	81

Surrogates	% Rec
2-Fluorophenol	54
Phenol-d5	65
Nitrobenzene-d5	63
2-Fluorobiphenyl	78
2,4,6-Tribromophenol	89
Terphenyl-d14	99

 1252

Directory:

The compounds listed above have been tentatively identified by a computer based library search. Compounds identified in the sample are not reported if they also occur in the method blank. The % fit is an indication of the reliability of the compound assignment. Due to the similarity between mass spectra of some isomeric compounds assignments may not be correct. Other compounds may also be present but identification was not possible. Concentrations are semi-quantitative, assume a response factor of 1 and use the nearest internal standard. Concentrations are reported on a wet weight basis.

Semi-Volatile Organic Compounds

Customer and Site Details:
Sample Details:
LIMS ID Number:
Job Number:

Exploration Associates: M4 Relief Road
CH/00477/1/1.10- m
Date Booked in:
Date Extracted:
Date Analysed:

16-Feb-00
18-Feb-00
25-Feb-00

Matrix:
Ext Method:
Operator:
Directory/Quant File:

Soil
Soxhlet
SW
0225ABN.MS4 0225CCC1.D GPC (Y/N)

QC Batch Number:
Multiplier:
Dilution Factor:

226
0.0667
1
N

Target Compounds	CAS #	R.T. (min)	Concentration mg/kg	% Fit
Phenol	108-95-2	-	< 2.0	-
bis(2-Chloroethyl)ether	111-44-4	-	< 0.5	-
2-Chlorophenol	95-57-8	-	< 2.0	-
1,3-Dichlorobenzene	541-73-1	-	< 0.5	-
1,4-Dichlorobenzene	106-46-7	-	< 0.5	-
Benzyl alcohol	100-51-6	-	< 0.5	-
1,2-Dichlorobenzene	95-50-1	-	< 0.5	-
2-Methylphenol	95-48-7	-	< 0.5	-
bis(2-Chloroisopropyl)ether	108-60-1	-	< 0.5	-
Hexachloroethane	67-72-1	-	< 0.5	-
N-Nitroso-di-n-propylamine	621-64-7	-	< 0.5	-
3- & 4-Methylphenol	108-39-4/106-44-5	-	< 2.0	-
Nitrobenzene	98-95-3	-	< 0.5	-
Isophorone	78-59-1	-	< 0.5	-
2-Nitrophenol	88-75-5	-	< 2.0	-
2,4-Dimethylphenol	105-67-9	-	< 2.0	-
Benzoic Acid	65-85-0	-	< 10.0	-
bis(2-Chloroethoxy)methane	111-91-1	-	< 0.5	-
2,4-Dichlorophenol	120-83-2	-	< 2.0	-
1,2,4-Trichlorobenzene	120-82-1	-	< 0.5	-
Naphthalene	91-20-3	-	< 0.2	-
4-Chlorophenol *	106-48-9	-	< 2.0	-
4-Chloroaniline *	106-47-8	-	< 0.5	-
Hexachlorobutadiene	87-68-3	-	< 0.5	-
4-Chloro-3-methylphenol	59-50-7	-	< 0.5	-
2-Methylnaphthalene	91-57-6	-	< 0.2	-
1-Methyl naphthalene *	90-12-0	-	< 0.2	-
Hexachlorocyclopentadiene	77-47-4	-	< 0.5	-
2,4,6-Trichlorophenol	88-06-2	-	< 2.0	-
2,4,5-Trichlorophenol	95-95-4	-	< 2.0	-
2-Chloronaphthalene	91-58-7	-	< 0.2	-
Biphenyl *	92-52-4	-	< 0.2	-
Diphenyl ether *	101-84-8	-	< 0.2	-
2-Nitroaniline	88-74-4	-	< 0.5	-
Acenaphthylene	208-96-8	-	< 0.2	-
Dimethylphthalate	131-11-3	-	< 0.5	-
2,6-Dinitrotoluene	606-20-2	-	< 0.5	-
Acenaphthene	83-32-9	-	< 0.2	-
3-Nitroaniline *	99-09-2	-	< 0.5	-

* These compounds are not currently UKAS accredited.
Concentrations are reported on a wet weight basis.

Target Compounds	CAS #	R.T.	Concentration mg/kg	% Fit
2,4-Dinitrophenol	51-28-5	-	< 1.0	-
Dibenzofuran	132-64-9	-	< 0.5	-
4-Nitrophenol	100-02-7	-	< 5.0	-
2,4-Dinitrotoluene	121-14-2	-	< 0.5	-
Fluorene	86-73-7	-	< 0.2	-
Diethylphthalate	84-66-2	-	< 0.5	-
4-Chlorophenyl-phenylether	7005-72-4	-	< 0.5	-
4,6-Dinitro-2-methylphenol	534-52-1	-	< 5.0	-
4-Nitroaniline	100-01-6	-	< 0.5	-
N-Nitrosodiphenylamine	86-30-6	-	< 0.5	-
4-Bromophenyl-phenylether	101-55-3	-	< 0.5	-
Hexachlorobenzene	118-74-1	-	< 0.5	-
Pentachlorophenol	87-86-5	-	< 5.0	-
Phenanthrene	85-01-8	-	< 0.2	-
Anthracene	120-12-7	-	< 0.2	-
Di-n-butylphthalate	84-74-2	-	< 0.5	-
Fluoranthene	206-44-0	-	< 0.2	-
Pyrene	129-00-0	-	< 0.2	-
Butylbenzylphthalate	85-68-7	-	< 0.5	-
Benzofuranthracene	56-55-3	-	< 0.2	-
Chrysene	218-01-9	-	< 0.2	-
3,3'-Dichlorobenzidine	91-94-1	-	< 2.0	-
bis(2-Ethylhexyl)phthalate	117-81-7	29.83	8.1	99
Di-n-octylphthalate	117-84-0	-	< 0.2	-
Benzobifluoranthene	205-99-2	-	< 0.2	-
Benzokjfluoranthene	207-08-9	-	< 0.2	-
Benzofluorene	50-32-8	-	< 0.2	-
Indeno[1,2,3-cd]pyrene	193-39-5	-	< 0.2	-
Dibenzofluoranthene	53-70-3	-	< 0.2	-
Benzofluoranthene	191-24-2	-	< 0.2	-

* compounds not currently UKAS accredited

Internal Standards	% Area	Surrogates	% Rec
1,4-Dichlorobenzene-d4	84	2-Fluorophenol	67
Naphthalene-d8	76	Phenol-d5	79
Acenaphthene-d10	76	Nitrobenzene-d5	76
Phenanthrene-d10	71	2-Fluorobiphenyl	78
Chrysene-d12	68	2,4,6-Tribromophenol	71
Perylene-d12	64	Terphenyl-d14	101

Semi-Volatile Organic Compounds



1252

Customer and Site Details:
Sample Details:
LIMS ID Number:
Job Number:

Exploration Associates: M4 Relief Road
CH/0047/A/10.20- m
Date Booked in:
Date Extracted:
Date Analysed:

16-Feb-00
18-Feb-00
25-Feb-00

Matrix:
Ext Method:
Operator:
Directory/Quant File:

Soil
Soxhlet
SW
0225ABN.MS4

QC Batch Number:
Multiplier:
Dilution Factor:

226
0.0667
1
Y

Target Compounds	CAS #	R.T. (min)	Concentration mg/kg	% Fit
Phenol	108-95-2	-	< 2.0	-
bis(2-Chloroethyl)ether	111-44-4	-	< 0.5	-
2-Chlorophenol	95-57-8	-	< 2.0	-
1,3-Dichlorobenzene	541-73-1	-	< 0.5	-
1,4-Dichlorobenzene	106-46-7	-	< 0.5	-
Benzyl alcohol	100-51-6	-	< 0.5	-
1,2-Dichlorobenzene	95-50-1	-	< 0.5	-
2-Methylphenol	95-48-7	-	< 0.5	-
bis(2-Chloroisopropyl)ether	108-60-1	-	< 0.5	-
Hexachloroethane	67-72-1	-	< 0.5	-
N-Nitroso-di-n-propylamine	621-64-7	-	< 0.5	-
3- & 4-Methylphenol	108-39-4/106-44-5	-	< 2.0	-
Nitrobenzene	98-95-3	-	< 0.5	-
Isophorone	78-59-1	-	< 0.5	-
2-Nitrophenol	88-75-5	-	< 2.0	-
2,4-Dimethylphenol	105-67-9	-	< 2.0	-
Benzoic Acid	65-85-0	-	< 10.0	-
bis(2-Chloroethoxy)methane	111-91-1	-	< 0.5	-
2,4-Dichlorophenol	120-83-2	-	< 2.0	-
1,2,4-Trichlorobenzene	120-82-1	-	< 0.5	-
Naphthalene	91-20-3	-	< 0.2	-
4-Chlorophenol *	106-48-9	-	< 2.0	-
4-Chloroaniline *	106-47-8	-	< 0.5	-
Hexachlorobutadiene	87-68-3	-	< 0.5	-
4-Chloro-3-methylphenol	59-50-7	-	< 0.5	-
2-Methylnaphthalene	91-57-6	-	< 0.2	-
1-Methyl naphthalene *	90-12-0	-	< 0.2	-
Hexachlorocyclopentadiene	77-47-4	-	< 0.5	-
2,4,6-Trichlorophenol	88-06-2	-	< 2.0	-
2,4,5-Trichlorophenol	95-95-4	-	< 2.0	-
2-Chloronaphthalene	91-58-7	-	< 0.2	-
Biphenyl *	92-52-4	15.31	0.4	99
Diphenyl ether *	101-84-8	-	< 0.2	-
2-Nitroaniline	88-74-4	-	< 0.5	-
Acenaphthylene	208-96-8	-	< 0.2	-
Dimethylphthalate	131-11-3	-	< 0.5	-
2,6-Dinitrotoluene	606-20-2	-	< 0.5	-
Acenaphthene	83-32-9	-	< 0.2	-
3-Nitroaniline *	99-09-2	-	< 0.5	-

* These compounds are not currently UKAS accredited.
Concentrations are reported on a wet weight basis.

Target Compounds	CAS #	R.T.	Concentration mg/kg	% Fit
2,4-Dinitrophenol	51-28-5	-	< 1.0	-
Dibenzofuran	132-64-9	-	< 0.5	-
4-Nitrophenol	100-02-7	-	< 5.0	-
2,4-Dinitrotoluene	121-14-2	-	< 0.5	-
Fluorene	86-73-7	-	< 0.2	-
Diethylphthalate	84-66-2	-	< 0.5	-
4-Chlorophenyl-phenylether	7005-72-4	-	< 0.5	-
4,6-Dinitro-2-methylphenol	534-52-1	-	< 5.0	-
4-Nitroaniline	100-01-6	-	< 0.5	-
N-Nitrosodiphenylamine	86-30-6	-	< 0.5	-
4-Bromophenyl-phenylether	101-55-3	-	< 0.5	-
Hexachlorobenzene	118-74-1	-	< 0.5	-
Pentachlorophenol	87-86-5	-	< 5.0	-
Phenanthrene	85-01-8	-	< 0.2	-
Anthracene	120-12-7	-	< 0.2	-
Di-n-butylphthalate	84-74-2	-	< 0.5	-
Fluoranthene	206-44-0	24.86	0.3	99
Pyrene	129-00-0	25.49	0.2	93
Butylbenzylphthalate	85-68-7	-	< 0.5	-
Benzofuranthracene	56-55-3	-	< 0.2	-
Chrysene	218-01-9	29.20	0.4	98
3,3'-Dichlorobenzidine	91-94-1	-	< 2.0	-
bis(2-Ethylhexyl)phthalate	117-81-7	29.82	0.7	98
Di-n-octylphthalate	117-84-0	-	< 0.2	-
Benzofluoranthene	205-99-2	32.16	0.2	99
Benzokifluoranthene	207-08-9	-	< 0.2	-
Benzofluoranthene	50-32-8	33.13	0.2	96
Indeno[1,2,3-cd]pyrene	193-39-5	37.67	0.2	97
Dibenz[a,h]anthracene	53-70-3	-	< 0.2	-
Benzofg,h,ijperylene	191-24-2	-	< 0.2	-

* compounds not currently UKAS accredited

Internal Standards	% Area
1,4-Dichlorobenzene-d4	82
Naphthalene-d8	75
Acenaphthene-d10	81
Phenanthrene-d10	80
Chrysene-d12	81
Perylene-d12	74

Surrogates	% Rec
2-Fluorophenol	32
Phenol-d5	46
Nitrobenzene-d5	46
2-Fluorobiphenyl	60
2,4,6-Tribromophenol	63
Terphenyl-d14	98



Volatile Organic Compounds by PTGCMs

Customer and Site Details: Exploration Associates: M4 Relief Road
Sample Details: CH 46 / 0.50m
LIMS ID Number: CL03961
Report Number: 000543

Directory/Quant file: 0225VOC.MS1\ 0225CCCC1.D
Date Booked in: 16-Feb-00
Date Analysed: 26-Feb-00
Operator: D.WALTON

Matrix: Soil
Method: Purge and Trap
Dilution: 5
Position: 1

1252

Target Compounds	CAS #	R.T. (min.)	Concentration µg/kg	% Fit
Dichlorodifluoromethane	75-71-8	-	< 5	-
Chloromethane	74-87-3	-	< 5	-
Vinyl Chloride	75-01-4	-	< 5	-
Bromomethane	74-83-9	-	< 5	-
Chloroethane	75-00-3	-	< 5	-
Trichlorofluoromethane	75-69-4	-	< 5	-
1,1-Dichloroethene	75-35-4	-	< 5	-
trans 1,2-Dichloroethene	156-60-5	-	< 5	-
1,1-Dichloroethane	75-34-3	-	< 5	-
2,2-Dichloropropane	594-20-7	-	< 5	-
cis 1,2-Dichloroethene	156-59-2	-	< 5	-
Bromochloromethane	74-97-5	-	< 5	-
Chloroform	67-66-3	-	< 5	-
1,1,1-Trichloroethane	71-55-6	-	< 5	-
Carbon Tetrachloride	56-23-5	-	< 5	-
1,1-Dichloropropane	563-58-6	-	< 5	-
Benzene	71-43-2	-	< 5	-
1,2-Dichloroethane	107-06-2	-	< 5	-
Trichloroethene	79-01-6	-	< 5	-
1,2-Dichloropropane	78-87-5	-	< 5	-
Dibromomethane	74-95-3	-	< 5	-
Bromodichloromethane	75-27-4	-	< 5	-
cis 1,3-Dichloropropene	10061-01-5	-	< 5	-
Toluene	108-88-3	-	< 5	-
trans 1,3-Dichloropropene	10061-02-6	-	< 5	-
1,1,2-Trichloroethane	79-00-5	-	< 5	-
Tetrachloroethene	127-18-4	-	< 5	-
1,3-Dichloropropane	142-28-29	-	< 5	-
Dibromochloromethane	124-48-1	-	< 5	-
1,2-Dibromoethane	106-93-4	-	< 5	-
Chlorobenzene	108-90-7	-	< 5	-
Ethylbenzene	100-41-4	-	< 5	-
1,1,1,2-Tetrachloroethane	630-20-6	-	< 5	-
m and p-Xylene	108-38-3/106-42-3	-	< 5	-
o-Xylene	95-47-6	-	< 5	-

Concentrations are reported on a wet weight basis

Target Compounds	CAS #	R.T. (min.)	Concentration µg/kg	% Fit
Styrene	100-42-5	-	< 5	-
Bromoform	75-25-2	-	< 5	-
iso-Propylbenzene	98-82-8	-	< 5	-
1,1,2,2-Tetrachloroethane	79-34-5	-	< 5	-
Propylbenzene	103-65-1	-	< 5	-
Bromobenzene	108-86-1	-	< 5	-
1,2,3-Trichloropropane	96-18-4	-	< 5	-
2-Chlorotoluene	95-49-8	-	< 5	-
1,3,5-Trimethylbenzene	108-67-8	-	< 5	-
4-Chlorotoluene	106-43-4	-	< 5	-
tert-Butylbenzene	98-06-6	-	< 5	-
1,2,4-Trimethylbenzene	95-63-6	-	< 5	-
sec-Butylbenzene	135-98-8	-	< 5	-
p-Isopropyltoluene	99-87-6	-	< 5	-
1,3-Dichlorobenzene	541-73-1	-	< 5	-
1,4-Dichlorobenzene	106-46-7	-	< 5	-
n-Butylbenzene	104-51-8	-	< 5	-
1,2-Dichlorobenzene	95-50-1	-	< 5	-
1,2-Dibromo-3-chloropropane	96-12-8	-	< 25	-
1,2,4-Trichlorobenzene	120-82-1	-	< 25	-
Hexachlorobutadiene	87-68-3	-	< 25	-
Naphthalene	91-20-3	-	< 25	-
1,2,3-Trichlorobenzene	87-61-6	-	< 25	-

Concentrations are reported on a wet weight basis

Internal standards	R.T.	Area %	Surrogates	% Rec
Pentafluorobenzene	12.14	88	Dibromofluoromethane	100
1,4-Difluorobenzene	13.28	85	Toluene-d8	99
Chlorobenzene-d5	17.60	81	Bromofluorobenzene	95
1,4-Dichlorobenzene-d4	21.06	72		



TEST REPORT SOIL SAMPLE ANALYSIS



1252
1411

Amended Report TES Report No. CL/000581

Site: M4 Relief Road

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
Mid Glamorgan
CF31 1JZ

The 5 samples described in this report were scheduled for analysis by TES Bretby on Thursday, 17 February 2000. The analysis was completed by Tuesday, 4 April 2000.

Tests marked as 'not UKAS accredited' and any opinions or interpretations expressed herein are outside the scope of any UKAS accreditation held by TES Bretby laboratories.

The following tables are contained in this report:

Table 1 Sample Descriptions
Table 2 Main Analysis Results

On behalf of
TES Bretby : J. Hannah
J Hannah Project Co-ordinator

Date of Issue: 04/04/00

Tests marked 'not UKAS accredited' in this report are not included in the UKAS Accreditation Schedule for our laboratory.

TES Bretby accepts no responsibility for the sampling related to the above results

TES Bretby, P.O. Box 100, Burton-on-Trent, DE15 0XD Telephone: 01283 554400 Fax: 01283 554422
TES Bretby is a division of Environmental Services Group Limited Registered in England Number 2880501

TES Bretby
Report 000581
Control Page
Sheet 1/ 1



TEST REPORT

SOIL SAMPLE ANALYSIS



1252

1411

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
Mid Glamorgan
CF31 1JZ

Amended Report
TES Report No. 000581

Site: M4 Relief Road

Page No	Assessed Area	ID No. EFS/CL	Sample	Depth (m) from to	Description
1	M4 Relief Road	0004281	CH/ 52/A	0.30	CH52A 0.3
1	M4 Relief Road	0004282	CH/ 52/A	1.20	CH52A 1.2
1	M4 Relief Road	0004283	CH/ 53/A	0.40	CH53A 0.4
1	M4 Relief Road	0004284	CH/ 53/A	1.10	CH53A 1.1
2	M4 Relief Road	0004285	CH/ 54/A	0.40	CH54A 0.4

Date of Issue: 04/04/00

Tests not marked 'not UKAS accredited' in this report are not included in the UKAS Accreditation Schedule for our laboratory.

TES Bretby accepts no responsibility for the sampling related to the above results

TES Bretby
Report 000581
Table 1
Sheet 1/ 1



TEST REPORT

SOIL SAMPLE ANALYSIS



1252
1411

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
Mid Glamorgan
CF31 1JZ

Amended Report
TES Report No. 000581

Site: M4 Relief Road

Customer reference	CH 52 A 0.30	CH 52 A 1.20	CH 53 A 0.40	CH 53 A 1.10
Depth (m)				
Date logged	17/02/00	17/02/00	17/02/00	17/02/00
TES Bretby ID Number	CL/0004281	CL/0004282	CL/0004283	CL/0004284

UKAS accredited	Test No.	CL/0004281	CL/0004282	CL/0004283	CL/0004284
Chromium (total)	ICPSSS11	53	647	670	48
CN- (total)	ICTSCN28	<1	<1	<1	<1
Copper	ICPSSS11	15	23	19	13
Cyanide (Free)	BGCN22	<1	<1	<1	<1
Elemental Sulphur	ELESULP	47	<20	<20	<20
Iron	ICPSSS11			131000	
Lead	ICPSSS11	59	65	71	53
Nickel	ICPSSS11	42	4	7	38
Oil FTIR	FTIRSW	95	133	164	43
PAH (screening)	PAHSCUV	11	12	26	<10
pH units	WSLM3	8.0	12.3	12.2	9.0
Phenol Index	WSLM4	<0.5	<0.5	<0.5	<0.5
SO4-- (acid sol)	ICPACS58	678	1750	2660	644
Sulphide	ICTSCN28	<1	48	66	<1
TPH FTIR (AR)	FTIRSW	37	190	225	55
Zinc	ICPSSS11	118	44	61	130

not UKAS accredited		CL/0004281	CL/0004282	CL/0004283	CL/0004284
Arsenic (Ms)		15.2	2.23	1.15	13.20
Boron.		1.6	2.0	6.0	1.0
Cadmium (Ms)		0.15	0.14	1.60	0.17
Mercury (Ms)		<0.01	<0.01	0.02	<0.01
Selenium (MS)		<0.50	1.46	1.60	<0.50

Results expressed as mg/kg Air Dried unless stated otherwise
SO4 Analysis not conducted in accordance with BS1377
AR denotes analysis conducted on As Received Sample

Date of Issue: 04/04/00

Tests marked 'not UKAS accredited' in this report are not included
in the UKAS Accreditation Schedule for our laboratory.

TES Bretby accepts no responsibility for the sampling related to the above results

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TES Bretby
Report 000581
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TEST REPORT SOIL SAMPLE ANALYSIS



1252
1411

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
Mid Glamorgan
CF31 1JZ

Amended Report
TES Report No. 000581

Site: M4 Relief Road

Customer reference	CH			
Depth (m)	54 A 0.40			
Date logged	17/02/00			
TES Bretby ID Number	CL/0004285			

UKAS accredited	Test No.	CL/0004285			
Chromium (total)	ICPSSS11	36			
CN- (total)	ICTSCN28	<1			
Copper	ICPSSS11	10			
Lead	ICPSSS11	35			
Nickel	ICPSSS11	27			
Oil FTIR	FTIRSW	77			
PAH (screening)	PAHSCUV	<10			
pH units	WSLM3	10.0			
Phenol Index	WSLM4	<0.5			
SO4-- (acid sol)	ICPACS58	866			
TPH FTIR (AR)	FTIRSW	<20			
Zinc	ICPSSS11	69			

not UKAS accredited		CL/0004285			
Arsenic (Ms)		8.30			
Boron.		0.8			
Cadmium (Ms)		0.13			
Mercury (Ms)		<0.01			
Selenium (MS)		<0.50			

Results expressed as mg/kg Air Dried unless stated otherwise
SO4 Analysis not conducted in accordance with BS1377
AR denotes analysis conducted on As Received Sample

Date of Issue: 04/04/00

Tests marked 'not UKAS accredited' in this report are not included
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TES Bretby accepts no responsibility for the sampling related to the above results

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TES Bretby
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TEST REPORT LEACHATE SAMPLE ANALYSIS

TES Report No. CL/000368

Site: M4 RELIEF RD

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
CF31 1JZ

The Sample described in this report was scheduled for analysis by TES Bretby on Thursday, 30 March 2000. The analysis was completed by Wednesday, 12 April 2000.

Tests marked as 'not UKAS accredited' and any opinions or interpretations expressed herein are outside the scope of any UKAS accreditation held by TES Bretby laboratories.

The following tables are contained in this report:

Table 1 Sample Descriptions
Table 2 Main Analysis Results

On behalf of
TES Bretby :
J Elstub


Project Co-ordinator

Date of Issue: 12/04/00

Tests marked 'not UKAS accredited' in this report are not included in the UKAS Accreditation Schedule for our laboratory.
TES Bretby accepts no responsibility for the sampling related to the above results.



TEST REPORT

LEACHATE SAMPLE ANALYSIS



1252

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
CF31 1JZ

TES Report No. CL/000368

Client: EXPLORATION ASS.

Site: M4 RELIEF RD

Page No	ID No. W/EX/		Sample Date
1	0007513	JOB 000368 CL/0002702	
1	0007514	JOB 000368 Blank	

Date of Issue: 12/04/00

Tests marked 'not UKAS accredited' in this report are not included
in the UKAS Accreditation Schedule for our laboratory.

TES Bretby accepts no responsibility for the sampling related to the above results.

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TES Bretby	
Report Number	
CL/000368	
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TEST REPORT

LEACHATE SAMPLE ANALYSIS



1252

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
CF31 1JZ

TES Report No. W/EXR/001544

Client: EXPLORATION ASS.

Site: M4 RELIEF RD

NDW

Customer Reference	CH			
Depth (m)	23 0.16	blank		
TES Bretby ID Number	Leachate W/EX/.. Soil Sample EFS/CL/..	0007513 0002702	0007514	

UKAS accredited	Test No.	0007513	0007514		
pH units	WSLM3	7.4	7.6		
TPH FTIR	FTIRSW	5.9	<0.2		
TPH GC	TPHFID	I.S	I.S		

not UKAS accredited		0007513	0007514		
Copper as Cu (Dissolved)		0.007	<0.001		
Zinc as Zn (Dissolved)		0.021	<0.001		
Arsenic as As (Dissolved)		0.004	<0.001		

Results expressed as mg/l unless stated, for the sample as received
.S Insufficient Sample for analysis

Date of Issue: 12/04/00

Tests marked 'not UKAS accredited' in this report are not included
in the UKAS Accreditation Schedule for our laboratory.

TES Bretby accepts no responsibility for the sampling related to the above results.

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TES Bretby
EFS/CL/000368
W/EXR/001544
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TEST REPORT LEACHATE SAMPLE ANALYSIS



1252

TES Report No. CL/000484

Site: M4 RELIEF RD

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
CF31 1JZ

The 4 Samples described in this report were scheduled for analysis by TES Bretby on Thursday, 30 March 2000. The analysis was completed by Thursday, 13 April 2000.

Tests marked as 'not UKAS accredited' and any opinions or interpretations expressed herein are outside the scope of any UKAS accreditation held by TES Bretby laboratories.

The following tables are contained in this report:

Table 1 Sample Descriptions
Table 2 Main Analysis Results

On behalf of
TES Bretby :
J Elstub


Project Co-ordinator

Date of Issue: 13/04/00

Tests marked 'not UKAS accredited' in this report are not included in the UKAS Accreditation Schedule for our laboratory.
TES Bretby accepts no responsibility for the sampling related to the above results.

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TES Bretby
Report Number
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TEST REPORT

LEACHATE SAMPLE ANALYSIS



1252

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
CF31 1JZ

TES Report No. CL/000484

Client: EXPLORATION ASS.

Site: M4 RELIEF RD

Page No	ID No. W/EX/		Sample Date
1	0007515	JOB 000484 CL/0003507	
1	0007516	JOB 000484 CL/0003511	
1	0007517	JOB 000484 CL/0003513	
1	0007518	JOB 000484 CL/0003520	
2	0007519	JOB 000484 Blank	

Date of Issue: 13/04/00

Tests marked 'not UKAS accredited' in this report are not included

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TES Bretby accepts no responsibility for the sampling related to the above results.

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TES Bretby
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Table 1
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TEST REPORT

LEACHATE SAMPLE ANALYSIS

1252

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
CF31 1JZ

TES Report No. W/EXR/001545

Client: EXPLORATION ASS.
Site: M4 RELIEF RD

SOLUTION

Customer Reference	CH 36	CH 38	CH 39	CH 42
Depth (m)	0.10	0.10	0.00	1.20
TES Bretby ID Number	0007515	0007516	0007517	0007518
Leachate W/EX/.. Soil Sample EFS/CL/..	0003507	0003511	0003513	0003520

UKAS accredited	Test No.	0007515	0007516	0007517	0007518
pH units	WSLM3	7.9	7.8	7.9	8.1
TPH FTIR	FTIRSW	0.5	<0.3	<0.3	<0.3
TPH GC	TPHFID	I.S	I.S	I.S	I.S
Pentachlorophenol	SVOC SW	<0.050	<0.050	<0.050	<0.050

not UKAS accredited		0007515	0007516	0007517	0007518
Copper as Cu (Dissolved)		0.003	0.003	0.007	0.004
Zinc as Zn (Dissolved)		0.008	0.016	0.023	0.011
PCB - ARO		<0.0008	<0.0008	<0.001	<0.001

Results expressed as mg/l unless stated, for the sample as received
I.S Insufficient Sample for analysis

Date of Issue: 13/04/00

Tests marked 'not UKAS accredited' in this report are not included
in the UKAS Accreditation Schedule for our laboratory.
TES Bretby accepts no responsibility for the sampling related to the above results.

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TES Bretby
EFS/CL/000484
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Table 2
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TEST REPORT

LEACHATE SAMPLE ANALYSIS



1252

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
CF31 1JZ

TES Report No. W/EXR/001545

Client: EXPLORATION ASS.
Site: M4 RELIEF RD

Customer Reference				
Depth (m)	blank			
TES Bretby ID Number Leachate W/EX/.. Soil Sample EFS/CL/..	0007519			

UKAS accredited	Test No.	0007519			
pH units	WSLM3	8.2			
TPH FTIR	FTIRSW	<0.2			
TPH GC	TPHFID	I.S			
Pentachlorophenol	SVOC SW	<0.050			

not UKAS accredited		0007519			
Copper as Cu (Dissolved)		≤0.001			
Zinc as Zn (Dissolved)		<0.001			
PCB - ARO		<0.0006			

Results expressed as mg/l unless stated, for the sample as received
I.S Insufficient Sample for analysis

Date of Issue: 13/04/00

Tests marked 'not UKAS accredited' in this report are not included
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TES Bretby accepts no responsibility for the sampling related to the above results.

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TES Bretby
EFS/CL/000484
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TEST REPORT

LEACHATE SAMPLE ANALYSIS

TES Report No. CL/000542

Site: M4 RELIEF RD

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
CF31 1JZ

The 4 Samples described in this report were scheduled for analysis by TES Bretby on Thursday, 30 March 2000. The analysis was completed by Friday, 14 April 2000.

Tests marked as 'not UKAS accredited' and any opinions or interpretations expressed herein are outside the scope of any UKAS accreditation held by TES Bretby laboratories.

The following tables are contained in this report:

Table 1 Sample Descriptions
Table 2 Main Analysis Results

On behalf of
TES Bretby :
J Elstub


Project Co-ordinator

Date of Issue: 14/04/00

TEST REPORT

LEACHATE SAMPLE ANALYSIS

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
CF31 1JZ

TES Report No. CL/000542

Client: EXPLORATION ASSOCIATES
Site: M4 RELIEF RD

Page No	ID No. W/EX/		Sample Date
1	0007497	JOB 000542 CL/0003913	
1	0007498	JOB 000542 CL/0003915	
1	0007499	JOB 000542 CL/0003917	
1	0007500	JOB 000542 CL/0003918	
2	0007501	JOB 000542 Blank	

Date of Issue: 14/04/00

TES Brethby accepts no responsibility for the sampling related to the above results.
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TES Brethby
Report Number
CL/000542
Table 1
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TEST REPORT

LEACHATE SAMPLE ANALYSIS

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
CF31 1JZ

TES Report No. W/EXR/001541

Client: EXPLORATION ASSOCIATES
Site: M4 RELIEF RD

Customer Reference		CH 9	CH 10	CH 11	CH 11
Depth (m)		0.00	0.40	0.35	1.65
TES Bretby ID Number		0007497	0007498	0007499	0007500
Leachate W/EX/.. Soil Sample EFS/CL/..		0003913	0003915	0003917	0003918
UKAS accredited		Test No.	0007497	0007498	0007499
TPH FTIR		FTIRSW	0.4	0.6	3.7
not UKAS accredited			0007497	0007498	0007499
pH units			7.6	8.0	7.9
Copper as Cu (Dissolved)			0.002	0.006	0.002
Zinc as Zn (Dissolved)			0.034	0.011	0.020
Arsenic as As (Dissolved)			0.002	0.005	0.002
TPH GC			<0.2	<0.1	<0.1

		0007500	0003918
		7.7	0.011
		0.025	0.009
		3.7	

Results expressed as mg/l unless stated, for the sample as received

Note - pH Units not UKAS accredited due to QC failure

Date of Issue: 14/04/00

TES Bretby accepts no responsibility for the sampling related to the above results.
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TES Bretby
EFS/CL/000542
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TEST REPORT

LEACHATE SAMPLE ANALYSIS

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
CF31 1JZ

TES Report No. W/EXR/001541

Client: EXPLORATION ASSOCIATES
Site: M4 RELIEF RD

Customer Reference					
Depth (m)		blank			
TES Bretby ID Number	Leachate W/EX/.. Soil Sample EFS/CL/..	0007501			
UKAS accredited	Test No.	0007501			
TPH FTIR	FTIRSW	<0.2			
not UKAS accredited		0007501			
pH units		8.0			
Copper as Cu (Dissolved)		<0.001			
Zinc as Zn (Dissolved)		<0.001			
Arsenic as As (Dissolved)		<0.001			
TPH GC		<0.1			

Results expressed as mg/l unless stated, for the sample as received

Note - pH Units not UKAS accredited due to QC failure

Date of Issue: 14/04/00

TES Bretby accepts no responsibility for the sampling related to the above results.
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TES Bretby	
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TEST REPORT

LEACHATE SAMPLE ANALYSIS

TES Report No. CL/000543

Site: M4 RELIEF RD

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
CF31 1JZ

The 8 Samples described in this report were scheduled for analysis by TES Bretby on Thursday, 30 March 2000. The analysis was completed by Friday, 14 April 2000.

Tests marked as 'not UKAS accredited' and any opinions or interpretations expressed herein are outside the scope of any UKAS accreditation held by TES Bretby laboratories.

The following tables are contained in this report:

Table 1 Sample Descriptions
Table 2 Main Analysis Results

On behalf of
TES Bretby : 
P Thompson Project Co-ordinator

Date of Issue: 14/04/00

TEST REPORT

LEACHATE SAMPLE ANALYSIS

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
CF31 1JZ

TES Report No. CL/000543

Client: EXPLORATION ASS.
Site: M4 RELIEF RD

Page No	ID No. W/EX/		Sample Date
1	0007504	JOB 000543 CL/0003950	
1	0007505	JOB 000543 CL/0003951	
1	0007506	JOB 000543 CL/0003955	
1	0007507	JOB 000543 CL/0003962	
2	0007508	JOB 000543 CL/0003967	
2	0007509	JOB 000543 CL/0003970	
2	0007510	JOB 000543 CL/0003971	
2	0007511	JOB 000543 CL/0003972	
3	0007512	JOB 000543 Blank	

Date of Issue: 14/04/00

TES Bretby accepts no responsibility for the sampling related to the above results.
TES Bretby, P.O. Box 100, Burton-on-Trent, DE15 0XD
Telephone: 01283 554400 Fax: 01283 554422

TES Bretby
Report Number
CL/000543
Table 1
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TEST REPORT

LEACHATE SAMPLE ANALYSIS

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
CF31 1JZ

TES Report No. W/EXR/001543

Client: EXPLORATION ASS.
Site: M4 RELIEF RD

ND-E

SOL

Customer Reference	CH 31 0.00	CH 32 0.00	CH 43 0.20	CH 47 0.10
Depth (m)				
TES Bretby ID Number	0007504	0007505	0007506	0007507
Leachate W/EX/.. Soil Sample EFS/CL/..	0003950	0003951	0003955	0003962

UKAS accredited	Test No.	0007504	0007505	0007506	0007507
pH units	WSLM3	10.8	9.4	8.8	8.6
TPH FTIR	FTIRSW			4.0	<0.2
Pentachlorophenol	SVOC SW			<0.083	<0.050

not UKAS accredited		0007504	0007505	0007506	0007507
Chromium as Cr (Dissolved)		0.035	0.005		
Copper as Cu (Dissolved)		0.014	0.008	0.004	0.003
Zinc as Zn (Dissolved)	-			0.006	0.010
Boron as B (Dissolved)		0.09	0.14		
PCB - ARO				<0.002	<0.0009
TPH GC		<0.1	<0.1	<0.2	<0.2

Results expressed as mg/l unless stated, for the sample as received

Date of Issue: 14/04/00

TES Bretby accepts no responsibility for the sampling related to the above results.
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TES Bretby
EFS/CL/000543
W/EXR/001543
Table 2
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TEST REPORT

LEACHATE SAMPLE ANALYSIS

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
CF31 1JZ

TES Report No. W/EXR/001543

Client: EXPLORATION ASS.

Site: M4 RELIEF RD

*LLANWERN
EPR*

EPR

MD

Customer Reference	CH 49	CH 50	CH 51	CH 55
Depth (m)	0.05	0.05	0.00	0.00
TES Bretby ID Number	0007508	0007509	0007510	0007511
Leachate W/EX/.. Soil Sample EFS/CL/..	0003967	0003970	0003971	0003972

UKAS accredited	Test No.	0007508	0007509	0007510	0007511
pH units	WSLM3	8.4	8.5	10.9	9.8
Total Sulphur as SO4 (Dissolved)	ICPPWW13	13.6	12.7	28.0	
Iron as Fe (Dissolved)	ICPPWW13	0.03	0.17	<0.01	
TPH FTIR	FTIRSW				0.8

not UKAS accredited		0007508	0007509	0007510	0007511
Zinc as Zn (Dissolved)		0.005	0.026		
TPH GC					I.S

Results expressed as mg/l unless stated, for the sample as received

.S Insufficient Sample for analysis

Date of Issue: 14/04/00

TES Bretby accepts no responsibility for the sampling related to the above results.
TES Bretby, P.O. Box 100, Burton-on-Trent, DE15 0XD
Telephone: 01283 554400 Fax: 01283 554422

TES Bretby
EFS/CL/000543
W/EXR/001543
Table 2
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TEST REPORT

LEACHATE SAMPLE ANALYSIS

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
CF31 1JZ

TES Report No. W/EXR/001543

Client: EXPLORATION ASS.
Site: M4 RELIEF RD

Customer Reference				
Depth (m)	blank			
TES Bretby ID Number Leachate W/EX/.. Soil Sample EFS/CL/..	0007512			

UKAS accredited	Test No.	0007512			
pH units	WSLM3	8.8			
Total Sulphur as SO4 (Dissolved)	ICPPWW13	<0.06			
Iron as Fe (Dissolved)	ICPPWW13	<0.01			
TPH FTIR	FTIRSW	<0.2			
Pentachlorophenol	SVOCSW	<0.050			

not UKAS accredited		0007512			
Chromium as Cr (Dissolved)		<0.001			
Copper as Cu (Dissolved)		<0.001			
Zinc as Zn (Dissolved)		<0.001			
Boron as B (Dissolved)		<0.05			
PCB - ARO		<0.0005			
TPH GC		<0.1			

Results expressed as mg/l unless stated, for the sample as received

Date of Issue: 14/04/00

TES Bretby accepts no responsibility for the sampling related to the above results.
TES Bretby, P.O. Box 100, Burton-on-Trent, DE15 0XD
Telephone: 01283 554400 Fax: 01283 554422

TES Bretby
EFS/CL/000543
W/EXR/001543
Table 2
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TEST REPORT

LEACHATE SAMPLE ANALYSIS

TES Report No. CL/000581

Site: M4 RELIEF RD

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
CF31 1JZ

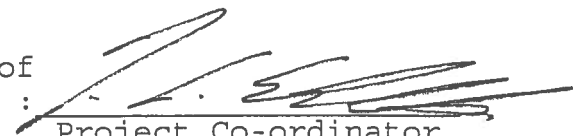
The Sample described in this report was scheduled for analysis by TES Bretby on Thursday, 30 March 2000. The analysis was completed by Monday, 10 April 2000.

Tests marked as 'not UKAS accredited' and any opinions or interpretations expressed herein are outside the scope of any UKAS accreditation held by TES Bretby laboratories.

The following tables are contained in this report:

Table 1 Sample Descriptions
Table 2 Main Analysis Results

On behalf of
TES Bretby :
J Elstub



Project Co-ordinator

Date of Issue: 10/04/00

Tests marked 'not UKAS accredited' in this report are not included in the UKAS Accreditation Schedule for our laboratory.

TES Bretby accepts no responsibility for the sampling related to the above results.

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TES Bretby
Report Number
CL/000581
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TEST REPORT

LEACHATE SAMPLE ANALYSIS



1252

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
CF31 1JZ

TES Report No. CL/000581

Client: EXPLORATION ASS.
Site: M4 RELIEF RD

Page No	ID No. W/EX/	Sample Date
1	0007502	JOB 000581 CL/0004283
1	0007503	JOB 000581 Blank

Date of Issue: 10/04/00

Tests marked 'not UKAS accredited' in this report are not included

in the UKAS Accreditation Schedule for our laboratory.

TES Bretby accepts no responsibility for the sampling related to the above results.

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TES Bretby
Report Number
CL/000581
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TEST REPORT

LEACHATE SAMPLE ANALYSIS



1252

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
CF31 1JZ

TES Report No. W/EXR/001542

Client: EXPLORATION ASS.
Site: M4 RELIEF RD

E.P.L.

Customer Reference	CH			
Depth (m)	53 A 0.40	blank		
TES Bretby ID Number	Leachate W/EX/.. Soil Sample EFS/CL/..	0007502 0004283	0007503	
UKAS accredited	Test No.	0007502	0007503	
pH units	WSLM3	11.8	10.2	
Total Sulphur as SO4 (Dissolved)	ICPPWW13	13.9	<0.06	
Iron as Fe (Dissolved)	ICPPWW13	<0.01	<0.01	

Results expressed as mg/l unless stated, for the sample as received

Date of Issue: 10/04/00

Tests marked 'not UKAS accredited' in this report are not included
in the UKAS Accreditation Schedule for our laboratory.

TES Bretby accepts no responsibility for the sampling related to the above results.

TES Bretby, P.O. Box 100, Burton-on-Trent, DE15 0XD Telephone: 01283 554400 Fax: 01283 554422
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TES Bretby
EFS/CL/000581
W/EXR/001542
Table 2
Sheet 1/ 1

TEST REPORT

WATER SAMPLE ANALYSIS

TES Report No. W/EXR/000694

Site: M4 RELIEF RD

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
CF31 1JZ

The 6 Samples described in this report were scheduled for analysis by TES Bretby on Thursday, 10 February 2000. The analysis was completed by Monday, 28 February 2000.

Tests marked as 'not UKAS accredited' and any opinions or interpretations expressed herein are outside the scope of any UKAS accreditation held by TES Bretby laboratories.

The following tables are contained in this report:

Table 1 Sample Descriptions
Table 2 Main Analysis Results
Table 3 Phenol Results
Tables of Poly Aromatic Hydrocarbons (6 Pages)
Tables of Volatile Organic Compounds (3 Pages)

On behalf of
TES Bretby :
J Elstub


Project Co-ordinator

Date of Issue: 28/02/00

TEST REPORT

WATER SAMPLE ANALYSIS

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
CF31 1JZ

TES Report No. W/EXR/000694

Client: EXPLORATION ASS.
Site: M4 RELIEF RD

Page No	ID No. W/EX/		Sample Date
1	0003196	Site 02 Reen Sample	31/01/00
1	0003195	BH CH 05 @ 2.70m	31/01/00
1	0003197	BH CH 08 @ 1.50m	31/01/00
1	0003198	BH CH 35 @ 0.80m	31/01/00
2	0003199	BH CH 40 @ 1.10m	31/01/00
2	0003200	BH CH 41 @ 1.10m	31/01/00

Date of Issue: 28/02/00

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TES Bretby
Report Number
W/EXR/000694
Table 1
Sheet 1/ 1

TEST REPORT

WATER SAMPLE ANALYSIS

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
CF31 1JZ

TES Report No. W/EXR/000694

Client: EXPLORATION ASS.
Site: M4 RELIEF RD

Sample reference	Sample 01	Sample 02	Sample 03	Sample 04
Sample Date	31/01/00	31/01/00	31/01/00	31/01/00
TES Bretby ID Number	0003196	0003195	0003197	0003198

UKAS accredited	Test No.	0003196	0003195	0003197	0003198
pH units	WSLM3	7.1	7.0	7.0	7.8
Chloride as Cl	WSLM1				75
Total Sulphur as SO ₄ (Dissolved)	ICPPWW13	26.0	18.6	24.0	91.4
Ammoniacal Nitrogen as N	AMMKA40				17
Thiocyanate as SCN	WSLM5				0.6
Sulphide (Free) as S	ISESW24				<0.2
Biochemical Oxygen Demand	WSLM20				17
TPH FTIR	FTIRSW	0.09	0.33	0.06	0.17
Phenol Index as C ₆ H ₅ OH	WSLM4	<0.05	<0.05	<0.05	1.06
TPH GC	TPHFID	0.12	0.6	<0.1	10.0
Volatile Organic Compounds	VOCSW				*
Pentachlorophenol	SVOC SW				4.275

not UKAS accredited		0003196	0003195	0003197	0003198
Nickel as Ni (Dissolved)		0.004	0.087	0.007	0.007
Chromium as Cr (Dissolved)		<0.001	0.001	<0.001	0.002
Cadmium as Cd (Dissolved)		<0.0001	<0.0001	<0.0001	<0.0001
Copper as Cu (Dissolved)		0.004	0.006	0.003	0.004
Lead as Pb (Dissolved)		<0.001	<0.001	<0.001	<0.001
Zinc as Zn (Dissolved)		0.032	0.075	0.043	0.030
Arsenic as As (Dissolved)		0.002	0.022	0.002	0.009
Boron as B (Dissolved)		0.06	0.24	0.22	0.33
Mercury as Hg (Dissolved)		<0.0001	<0.0001	<0.0001	<0.0001
Selenium as Se (Dissolved)		<0.001	0.007	<0.001	0.002
Cyanide (Free) as CN					<0.025
Cyanide (Total) as CN		1.3	<0.025	<0.025	0.05
Complex Cyanide as CN					0.05
PCB - ARO					0.066
PAH HPLC		*	*	*	*

Results expressed as mg/l unless stated, for the sample as received

* See attached sheets for results

Sample 01: Site 02 Reen Sample

Sample 02: BH CH 05 @ 2.70m

Sample 03: BH CH 08 @ 1.50m

Sample 04: BH CH 35 @ 0.80m

Date of Issue: 28/02/00

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TES Bretby
Report Number
W/EXR/000694
Table 2
Sheet 1/2

TEST REPORT

WATER SAMPLE ANALYSIS

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
CF31 1JZ

TES Report No. W/EXR/000694

Client: EXPLORATION ASS.

Site: M4 RELIEF RD

Sample reference		Sample 05	Sample 06		
Sample Date		31/01/00	31/01/00		
TES Bretby ID Number	W/EX/..	0003199	0003200		

UKAS accredited	Test No.	0003199	0003200		
pH units	WSLM3	8.9	7.4		
Chloride as Cl	WSLM1	21	17		
Total Sulphur as SO ₄ (Dissolved)	ICPPWW13	41.1	23.3		
Ammoniacal Nitrogen as N	AMMKA40	<0.2	<0.2		
Thiocyanate as SCN	WSLM5	<0.2	<0.2		
Sulphide (Free) as S	ISESW24	<0.2	0.8		
Biochemical Oxygen Demand	WSLM20	4.0	<2.0		
TPH FTIR	FTIRSW	0.18	0.66		
Phenol Index as C ₆ H ₅ OH	WSLM4	<0.05	<0.05		
TPH GC	TPHFID	0.8	8.0		
Volatile Organic Compounds	VOCSW	*	*		
Pentachlorophenol	SVOCSW	<0.050	<0.063		

* Min.
0.1

- All
6-comp

not UKAS accredited		0003199	0003200		
Nickel as Ni (Dissolved)		0.006	0.010		
Chromium as Cr (Dissolved)		0.003	0.004		
Cadmium as Cd (Dissolved)		<0.0001	<0.0001		
Copper as Cu (Dissolved)		0.010	0.003		
Lead as Pb (Dissolved)		<0.001	<0.001		
Zinc as Zn (Dissolved)		0.016	0.034		
Arsenic as As (Dissolved)		0.007	0.012		
Boron as B (Dissolved)		0.07	0.06		
Mercury as Hg (Dissolved)		<0.0001	<0.0001		
Selenium as Se (Dissolved)		<0.001	<0.001		
Cyanide (Free) as CN		<0.025	<0.025		
Cyanide (Total) as CN		0.05	<0.025		
Complex Cyanide as CN		0.05	<0.025		
PCB - ARO		0.048	0.022		
PAH HPLC		*	*		

Results expressed as mg/l unless stated, for the sample as received

* See attached sheets for results

Sample 05: BH CH 40 @ 1.10m

Sample 06: BH CH 41 @ 1.10m

Date of Issue: 28/02/00

TES Bretby accepts no responsibility for the sampling related to the above results.
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TES Bretby	
Report Number	W/EXR/000694
Table	2
Sheet	2 / 2

TEST REPORT

WATER SAMPLE ANALYSIS

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
CF31 1JZ

TES Report No. W/EXR/000694

Client: EXPLORATION ASS.
Site: M4 RELIEF RD

Phenol Analysis

Sample reference	Sample 01	Sample 02	Sample 03	
Sample Date	31/01/00	31/01/00	31/01/00	
TES Bretby ID Number	0003198	0003199	0003200	
W/EX/..				
not UKAS accredited	0003198	0003199	0003200	
Phenol	1.083	<0.0005	0.0006	
Cresols	4.14	<0.0005	0.0024	
Dimethylphenols	<0.05	<0.0005	<0.0005	
Trimethylphenols	0.9780	<0.0005	<0.0005	

Results expressed as mg/l unless stated, for the sample as received

Sample 01: BH CH 35 @ 0.80m

Sample 02: BH CH 40 @ 1.10m

Sample 03: BH CH 41 @ 1.10m

Date of Issue: 28/02/00

TES Bretby accepts no responsibility for the sampling related to the above results.
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Telephone: 01283 554400 Fax: 01283 554422

TES Bretby
Report Number
W/EXR/000694
Table 3
Sheet 1/ 1



Polycyclic Aromatic Hydrocarbons by HPLC

Customer and Site Details:	Exploration Ass. : M4 Relief Road	Report Number:	694
Sample Details:	BH CH 05 @ 2.70m	Date Booked in:	10-Feb-00
LIMS ID Number:	EX3195	Date Extracted:	15-Feb-00
QC Batch Number:	282	Date Analysed:	16-Feb-00
Directory:	0215PAH.LC2	Matrix:	Water
Dilution (fluorescence):	1.0	Ext Method:	Sep Funnel
Dilution (UV):	1.0		

Target Compounds	CAS #	Concentration µg/l	Fluorescence / UV
Naphthalene	91-20-3	1.19	Fluorescence
Acenaphthylene	208-96-8	< 1.00	UV
Acenaphthene	83-32-9	0.16	Fluorescence
Fluorene	86-73-7	0.11	Fluorescence
Phenanthrene	85-01-8	0.15	Fluorescence
Anthracene	120-12-7	< 0.05	Fluorescence
Fluoranthene	206-44-0	0.08	Fluorescence
Pyrene	129-00-0	< 0.05	Fluorescence
Benzo[a]anthracene	56-55-3	< 0.05	Fluorescence
Chrysene	218-01-9	< 0.05	Fluorescence
Benzo[b]fluoranthene	205-99-2	< 0.05	Fluorescence
Benzo[k]fluoranthene	207-08-9	< 0.05	Fluorescence
Benzo[a]pyrene	50-32-8	< 0.05	Fluorescence
Indeno[1,2,3-cd]pyrene	193-39-5	< 0.05	Fluorescence
Dibenzo[a,h]anthracene	53-70-3	< 0.05	Fluorescence
Benzo[g,h,i]perylene	191-24-2	< 0.05	Fluorescence
Total PAHs (USEPA 16)	-	< 3.18	-

Data may be obtained from either the fluorescence detector or when the response is very high a less sensitive UV detector (this is indicated for each compound).
Any dilutions necessary during analysis have been included in the calculations.
Reported results do not require any corrections.



Polycyclic Aromatic Hydrocarbons by HPLC

Customer and Site Details:	Exploration Ass. : M4 Relief Road	Report Number:	694
Sample Details:	Site 02 Reen Sample	Date Booked in:	10-Feb-00
LIMS ID Number:	EX3196	Date Extracted:	15-Feb-00
QC Batch Number:	282	Date Analysed:	16-Feb-00
Directory:	0215PAH.LC2	Matrix:	Water
Dilution (fluorescence):	1.0	Ext Method:	Sep Funnel
Dilution (UV):	1.0		

Target Compounds	CAS #	Concentration µg/l	Fluorescence / UV
Naphthalene	91-20-3	< 0.50	Fluorescence
Acenaphthylene	208-96-8	< 1.00	UV
Acenaphthene	83-32-9	< 0.05	Fluorescence
Fluorene	86-73-7	< 0.05	Fluorescence
Phenanthrene	85-01-8	< 0.05	Fluorescence
Anthracene	120-12-7	< 0.05	Fluorescence
Fluoranthene	206-44-0	< 0.05	Fluorescence
Pyrene	129-00-0	< 0.05	Fluorescence
Benzo[a]anthracene	56-55-3	< 0.05	Fluorescence
Chrysene	218-01-9	< 0.05	Fluorescence
Benzo[b]fluoranthene	205-99-2	< 0.05	Fluorescence
Benzo[k]fluoranthene	207-08-9	< 0.05	Fluorescence
Benzo[a]pyrene	50-32-8	< 0.05	Fluorescence
Indeno[1,2,3-cd]pyrene	193-39-5	< 0.05	Fluorescence
Dibenzo[a,h]anthracene	53-70-3	< 0.05	Fluorescence
Benzo[g,h,i]perylene	191-24-2	< 0.05	Fluorescence
Total PAHs (USEPA 16)	-	< 2.20	-

Data may be obtained from either the fluorescence detector or when the response is very high a less sensitive UV detector (this is indicated for each compound).
Any dilutions necessary during analysis have been included in the calculations.
Reported results do not require any corrections.



Polycyclic Aromatic Hydrocarbons by HPLC

Customer and Site Details:	Exploration Ass. : M4 Relief Road	Report Number:	694
Sample Details:	BH CH 08 @ 1.50m	Date Booked in:	10-Feb-00
LIMS ID Number:	EX3197	Date Extracted:	15-Feb-00
QC Batch Number:	282	Date Analysed:	16-Feb-00
Directory:	0215PAH.LC2	Matrix:	Water
Dilution (fluorescence):	1.0	Ext Method:	Sep Funnel
Dilution (UV):	1.0		

Target Compounds	CAS #	Concentration µg/l	Fluorescence / UV
Naphthalene	91-20-3	< 0.50	Fluorescence
Acenaphthylene	208-96-8	< 1.00	UV
Acenaphthene	83-32-9	< 0.05	Fluorescence
Fluorene	86-73-7	< 0.05	Fluorescence
Phenanthrene	85-01-8	< 0.05	Fluorescence
Anthracene	120-12-7	< 0.05	Fluorescence
Fluoranthene	206-44-0	< 0.05	Fluorescence
Pyrene	129-00-0	< 0.05	Fluorescence
Benzo[a]anthracene	56-55-3	< 0.05	Fluorescence
Chrysene	218-01-9	< 0.05	Fluorescence
Benzo[b]fluoranthene	205-99-2	< 0.05	Fluorescence
Benzo[k]fluoranthene	207-08-9	< 0.05	Fluorescence
Benzo[a]pyrene	50-32-8	< 0.05	Fluorescence
Indeno[1,2,3-cd]pyrene	193-39-5	< 0.05	Fluorescence
Dibenzo[a,h]anthracene	53-70-3	< 0.05	Fluorescence
Benzo[g,h,i]perylene	191-24-2	< 0.05	Fluorescence
Total PAHs (USEPA 16)	-	< 2.20	-

Data may be obtained from either the fluorescence detector or when the response is very high a less sensitive UV detector (this is indicated for each compound). Any dilutions necessary during analysis have been included in the calculations. Reported results do not require any corrections.

Polycyclic Aromatic Hydrocarbons by HPLC

Customer and Site Details:	Exploration Ass. : M4 Relief Road	Report Number:	694
Sample Details:	BH CH 35 @ 0.80m	Date Booked in:	10-Feb-00
LIMS ID Number:	EX3198	Date Extracted:	15-Feb-00
QC Batch Number:	282	Date Analysed:	16-Feb-00
Directory:	0215PAH.LC2	Matrix:	Water
Dilution (fluorescence):	10.0	Ext Method:	Sep Funnel
Dilution (UV):	1.0		

Target Compounds	CAS #	Concentration µg/l	Fluorescence / UV
Naphthalene	91-20-3	< 5.00	Fluorescence
Acenaphthylene	208-96-8	< 1.00	UV
Acenaphthene	83-32-9	0.54	Fluorescence
Fluorene	86-73-7	1.07	Fluorescence
Phenanthrene	85-01-8	< 0.50	Fluorescence
Anthracene —	120-12-7	0.64	Fluorescence
Fluoranthene	206-44-0	< 0.50	Fluorescence
Pyrene	129-00-0	< 0.50	Fluorescence
Benzo[a]anthracene	56-55-3	< 0.50	Fluorescence
Chrysene	218-01-9	< 0.50	Fluorescence
Benzo[b]fluoranthene	205-99-2	< 0.50	Fluorescence
Benzo[k]fluoranthene	207-08-9	< 0.50	Fluorescence
Benzo[a]pyrene	50-32-8	< 0.50	Fluorescence
Indeno[1,2,3-cd]pyrene	193-39-5	< 0.50	Fluorescence
Dibenzo[a,h]anthracene	53-70-3	< 0.50	Fluorescence
Benzo[g,h,i]perylene	191-24-2	< 0.50	Fluorescence
Total PAHs (USEPA 16)	-	< 13.75	-

Data may be obtained from either the fluorescence detector or when the response is very high a less sensitive UV detector (this is indicated for each compound).
Any dilutions necessary during analysis have been included in the calculations.
Reported results do not require any corrections.



Polycyclic Aromatic Hydrocarbons by HPLC

Customer and Site Details:	Exploration Ass. : M4 Relief Road	Report Number:	694
Sample Details:	BH CH 40 @ 1.10m	Date Booked in:	10-Feb-00
LIMS ID Number:	EX3199	Date Extracted:	15-Feb-00
QC Batch Number:	282	Date Analysed:	16-Feb-00
Directory:	0215PAH.LC2	Matrix:	Water
Dilution (fluorescence):	1.0	Ext Method:	Sep Funnel
Dilution (UV):	1.0		

Target Compounds	CAS #	Concentration µg/l	Fluorescence / UV
Naphthalene	91-20-3	< 0.50	Fluorescence
Acenaphthylene	208-96-8	< 1.00	UV
Acenaphthene	83-32-9	< 0.05	Fluorescence
Fluorene	86-73-7	< 0.05	Fluorescence
Phenanthrene	85-01-8	< 0.05	Fluorescence
Anthracene	120-12-7	< 0.05	Fluorescence
Fluoranthene	206-44-0	< 0.05	Fluorescence
Pyrene	129-00-0	< 0.05	Fluorescence
Benzo[a]anthracene	56-55-3	< 0.05	Fluorescence
Chrysene	218-01-9	< 0.05	Fluorescence
Benzo[b]fluoranthene	205-99-2	< 0.05	Fluorescence
Benzo[k]fluoranthene	207-08-9	< 0.05	Fluorescence
Benzo[a]pyrene	50-32-8	< 0.05	Fluorescence
Indeno[1,2,3-cd]pyrene	193-39-5	< 0.05	Fluorescence
Dibenzo[a,h]anthracene	53-70-3	< 0.05	Fluorescence
Benzo[g,h,i]perylene	191-24-2	< 0.05	Fluorescence
Total PAHs (USEPA 16)	-	< 2.20	-

Data may be obtained from either the fluorescence detector or when the response is very high a less sensitive UV detector (this is indicated for each compound).
Any dilutions necessary during analysis have been included in the calculations.
Reported results do not require any corrections.

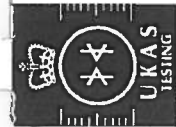


Polycyclic Aromatic Hydrocarbons by HPLC

Customer and Site Details:	Exploration Ass. : M4 Relief Road	Report Number:	694
Sample Details:	BH CH 41 @ 1.10m	Date Booked in:	10-Feb-00
LIMS ID Number:	EX3200	Date Extracted:	15-Feb-00
QC Batch Number:	282	Date Analysed:	16-Feb-00
Directory:	0215PAH.LC2	Matrix:	Water
Dilution (fluorescence):	1.0	Ext Method:	Sep Funnel
Dilution (UV):	1.0		

Target Compounds	CAS #	Concentration µg/l	Fluorescence / UV
Naphthalene	91-20-3	< 0.50	Fluorescence
Acenaphthylene	208-96-8	< 1.00	UV
Acenaphthene	83-32-9	< 0.05	Fluorescence
Fluorene	86-73-7	0.07	Fluorescence
Phenanthrene	85-01-8	0.17	Fluorescence
Anthracene	120-12-7	< 0.05	Fluorescence
Fluoranthene	206-44-0	0.21	Fluorescence
Pyrene	129-00-0	0.19	Fluorescence
Benzo[a]anthracene	56-55-3	0.50	Fluorescence
Chrysene	218-01-9	0.56	Fluorescence
Benzo[b]fluoranthene	205-99-2	0.16	Fluorescence
Benzo[k]fluoranthene	207-08-9	0.06	Fluorescence
Benzo[a]pyrene	50-32-8	0.16	Fluorescence
Indeno[1,2,3-cd]pyrene	193-39-5	0.07	Fluorescence
Dibenzo[a,h]anthracene	53-70-3	< 0.05	Fluorescence
Benzo[g,h,i]perylene	191-24-2	0.07	Fluorescence
Total PAHs (USEPA 16)	-	< 3.87	-

Data may be obtained from either the fluorescence detector or when the response is very high a less sensitive UV detector (this is indicated for each compound). Any dilutions necessary during analysis have been included in the calculations. Reported results do not require any corrections.



Volatile Organic Compounds by PTGCMS

Customer and Site Details:
Sample Details:
LIMS ID Number:
Report Number:

Exploration Associates : M4 Relief Road
BH CH 35 / 0.80m
EX03198
000694

Directory/Quant file:
Date Booked in:
Date Analysed:
Operator:

0214VOC.MS3\ 0214CCC1.D
10-Feb-00
14-Feb-00
D.WALTON

Matrix:
Method:
Dilution:
Position:

1252

Water

Purge and Trap

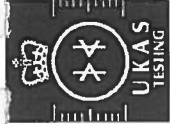
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6

Target Compounds	CAS #	R.T. (min.)	Concentration µg/l	% Fit
Dichlorodifluoromethane	75-71-8	-	< 5	-
Chloromethane	74-87-3	-	< 5	-
Vinyl Chloride	75-01-4	-	< 5	-
Bromomethane	74-83-9	-	< 5	-
Chloroethane	75-00-3	-	< 5	-
Trichlorofluoromethane	75-69-4	-	< 5	-
1,1-Dichloroethene	75-35-4	-	< 5	-
trans 1,2-Dichloroethene	156-60-5	-	< 5	-
1,1-Dichloroethane	75-34-3	-	< 5	-
2,2-Dichloropropane	594-20-7	-	< 5	-
cis 1,2-Dichloroethene	156-59-2	-	< 5	-
Bromochloromethane	74-97-5	-	< 5	-
Chloroform	67-66-3	-	< 5	-
1,1,1-Trichloroethane	71-55-6	-	< 5	-
Carbon Tetrachloride	56-23-5	-	< 5	-
1,1-Dichloropropene	563-58-6	-	< 5	-
Benzene	71-43-2	12.55	130	94
1,2-Dichloroethane	107-06-2	-	< 5	-
Trichloroethene	79-01-6	-	< 5	-
1,2-Dichloropropane	78-87-5	-	< 5	-
Dibromomethane	74-95-3	-	< 5	-
Bromodichloromethane	75-27-4	-	< 5	-
cis 1,3-Dichloropropene	10061-01-5	-	< 5	-
Toluene	108-88-3	15.33	18	94
trans 1,3-Dichloropropene	10061-02-6	-	< 5	-
1,1,2-Trichloroethane	79-00-5	-	< 5	-
Tetrachloroethane	127-18-4	-	< 5	-
1,3-Dichloropropane	142-28-29	-	< 5	-
Dibromochloromethane	124-48-1	-	< 5	-
1,2-Dibromomethane	106-93-4	-	< 5	-
Chlorobenzene	108-90-7	17.42	6	61
Ethylbenzene	100-41-4	17.52	255	97
1,1,1,2-Tetrachloroethane	630-20-6	-	< 5	-
m and p-Xylene	108-38-3/106-42-3	17.67	872	98
o-Xylene	95-47-6	18.29	245	93

Target Compounds	CAS #	R.T. (min.)	Concentration µg/l	% Fit
Styrene	100-42-5	-	< 5	-
Bromoform	75-25-2	-	< 5	-
iso-Propylbenzene	98-82-8	-	< 5	-
1,1,2,2-Tetrachloroethane	79-34-5	-	< 5	-
Propylbenzene	103-65-1	-	< 5	-
Bromobenzene	108-86-1	-	< 5	-
1,2,3-Trichloropropane	96-18-4	-	< 5	-
2-Chlorotoluene	95-49-8	-	< 5	-
1,3,5-Trimethylbenzene	108-67-8	-	< 5	-
4-Chlorotoluene	106-43-4	-	< 5	-
tert-Butylbenzene	98-06-6	-	< 5	-
1,2,4-Trimethylbenzene	95-63-6	-	< 5	-
sec-Butylbenzene	135-98-8	-	< 5	-
p-Isopropyltoluene	99-87-6	-	< 5	-
1,3-Dichlorobenzene	541-73-1	20.79	6	79
1,4-Dichlorobenzene	106-46-7	20.91	45	83
n-Butylbenzene	104-51-8	-	< 5	-
1,2-Dichlorobenzene	95-50-1	21.51	12	96
1,2-Dibromo-3-chloropropane	96-12-8	-	< 25	-
1,2,4-Trichlorobenzene	120-82-1	24.01	204	97
Hexachlorobutadiene	87-68-3	-	< 25	-
Naphthalene	91-20-3	-	< 25	-
1,2,3-Trichlorobenzene	87-61-6	-	< 25	-

Internal standards	R.T.	Area %	Surrogates	% Rec
Pentafluorobenzene	11.89	82	Dibromofluoromethane	109
1,4-Difluorobenzene	13.04	86	Toluene-d8	101
Chlorobenzene-d5	17.39	87	Bromofluorobenzene	97
1,4-Dichlorobenzene-d4	20.87	86		



Volatile Organic Compounds by PTGCMS

Customer and Site Details:

Exploration Associates : M4 Relief Road

Sample Details:

BH CH 40 / 1.10m

LIMS ID Number:

EX03199

Report Number:

000694

Directory/Quant file:

0214VOC.MS3\

0214CCC1.D

Date Booked in:

10-Feb-00

Date Analysed:

14-Feb-00

Operator:

D.WALTON

Matrix:

Water

Method:

Purge and Trap

Dilution:

1

Position:

7

1252

Target Compounds	CAS #	R.T. (min.)	Concentration µg/l	% Fit
Dichlorodifluoromethane	75-71-8	-	< 1	-
Chloromethane	74-87-3	-	< 1	-
Vinyl Chloride	75-01-4	-	< 1	-
Bromomethane	74-83-9	-	< 1	-
Chloroethane	75-00-3	-	< 1	-
Trichlorofluoromethane	75-69-4	-	< 1	-
1,1-Dichloroethene	75-35-4	-	< 1	-
trans 1,2-Dichloroethene	156-60-5	-	< 1	-
1,1-Dichloroethane	75-34-3	-	< 1	-
2,2-Dichloropropane	594-20-7	-	< 1	-
cis 1,2-Dichloroethene	156-59-2	-	< 1	-
Bromochloromethane	74-97-5	-	< 1	-
Chloroform	67-66-3	-	< 1	-
1,1,1-Trichloroethane	71-55-6	-	< 1	-
Carbon Tetrachloride	56-23-5	-	< 1	-
1,1-Dichloropropene	563-58-6	-	< 1	-
Benzene	71-43-2	-	< 1	-
1,2-Dichloroethane	107-06-2	-	< 1	-
Trichloroethene	79-01-6	-	< 1	-
1,2-Dichloropropane	78-87-5	-	< 1	-
Dibromomethane	74-95-3	-	< 1	-
Bromodichloromethane	75-27-4	-	< 1	-
cis 1,3-Dichloropropene	10061-01-5	-	< 1	-
Toluene	108-88-3	-	< 1	-
trans 1,3-Dichloropropene	10061-02-6	-	< 1	-
1,1,2-Trichloroethane	79-00-5	-	< 1	-
Tetrachloroethene	127-18-4	-	< 1	-
1,3-Dichloropropane	142-28-29	-	< 1	-
Dibromochloromethane	124-48-1	-	< 1	-
1,2-Dibromoethane	106-93-4	-	< 1	-
Chlorobenzene	108-90-7	-	< 1	-
Ethylbenzene	100-41-4	-	< 1	-
1,1,1,2-Tetrachloroethane	630-20-6	-	< 1	-
m and p-Xylene	108-38-3/106-42-3	-	< 1	-
o-Xylene	95-47-6	-	< 1	-

Target Compounds	CAS #	R.T. (min.)	Concentration µg/l	% Fit
Styrene	100-42-5	-	< 1	-
Bromoform	75-25-2	-	< 1	-
iso-Propylbenzene	98-82-8	-	< 1	-
1,1,2,2-Tetrachloroethane	79-34-5	-	< 1	-
Propylbenzene	103-65-1	-	< 1	-
Bromobenzene	108-86-1	-	< 1	-
1,2,3-Trichloropropane	96-18-4	-	< 1	-
2-Chlorotoluene	95-49-8	-	< 1	-
1,3,5-Trimethylbenzene	108-67-8	-	< 1	-
4-Chlorotoluene	106-43-4	-	< 1	-
tert-Butylbenzene	98-06-6	-	< 1	-
1,2,4-Trimethylbenzene	95-63-6	-	< 1	-
sec-Butylbenzene	135-98-8	-	< 1	-
p-Isopropyltoluene	99-87-6	-	< 1	-
1,3-Dichlorobenzene	541-73-1	-	< 1	-
1,4-Dichlorobenzene	106-46-7	-	< 1	-
n-Butylbenzene	104-51-8	-	< 1	-
1,2-Dichlorobenzene	95-50-1	-	< 1	-
1,2-Dibromo-3-chloropropane	96-12-8	-	< 5	-
1,2,4-Trichlorobenzene	120-82-1	-	< 5	-
Hexachlorobutadiene	87-68-3	-	< 5	-
Naphthalene	91-20-3	-	< 5	-
1,2,3-Trichlorobenzene	87-61-6	-	< 5	-

Internal standards	R.T.	Area %	Surrogates	% Rec
Pentafluorobenzene	11.89	86	Dibromofluoromethane	107
1,4-Difluorobenzene	13.05	90	Toluene-d8	99
Chlorobenzene-d5	17.38	91	Bromofluorobenzene	97
1,4-Dichlorobenzene-d4	20.87	88		



Volatile Organic Compounds by PTGCMS

Customer and Site Details:
Sample Details:
LIMS ID Number:
Report Number:

Exploration Associates : M4 Relief Road
BH CH 41 / 1.10m
EX03200
000694

Directory/Quant file:
Date Booked in:
Date Analysed:
Operator:

0215VOC.MS3\ 0214CCCC1.D
10-Feb-00
15-Feb-00
D.WALTON

Matrix: Water
Method: Purge and Trap
Dilution: 1
Position: 10

Target Compounds	CAS #	R.T. (min.)	Concentration µg/l	% Fit
Dichlorodifluoromethane	75-71-8	-	< 1	-
Chloromethane	74-87-3	-	< 1	-
Vinyl Chloride	75-01-4	-	< 1	-
Bromomethane	74-83-9	-	< 1	-
Chloroethane	75-00-3	-	< 1	-
Trichlorofluoromethane	75-69-4	-	< 1	-
1,1-Dichloroethane	75-35-4	-	< 1	-
trans 1,2-Dichloroethene	156-60-5	-	< 1	-
1,1-Dichloroethane	75-34-3	-	< 1	-
2,2-Dichloropropane	594-20-7	-	< 1	-
cis 1,2-Dichloroethene	156-59-2	-	< 1	-
Bromochloromethane	74-97-5	-	< 1	-
Chloroform	67-66-3	-	< 1	-
1,1,1-Trichloroethane	71-55-6	-	< 1	-
Carbon Tetrachloride	56-23-5	-	< 1	-
1,1-Dichloropropene	563-58-6	-	< 1	-
Benzene	71-43-2	-	< 1	-
1,2-Dichloroethane	107-06-2	-	< 1	-
Trichloroethane	79-01-6	-	< 1	-
1,2-Dichloropropene	78-87-5	-	< 1	-
Dibromomethane	74-95-3	-	< 1	-
Bromodichloromethane	75-27-4	-	< 1	-
cis 1,3-Dichloropropene	10061-01-5	-	< 1	-
Toluene	108-88-3	-	< 1	-
trans 1,3-Dichloropropene	10061-02-6	-	< 1	-
1,1,2-Trichloroethane	79-00-5	-	< 1	-
Tetrachloroethane	127-18-4	-	< 1	-
1,3-Dichloropropene	142-28-29	-	< 1	-
Dibromochloromethane	124-48-1	-	< 1	-
1,2-Dibromoethane	106-93-4	-	< 1	-
Chlorobenzene	108-90-7	-	< 1	-
Ethylbenzene	100-41-4	-	< 1	-
1,1,1,2-Tetrachloroethane	630-20-6	-	< 1	-
m and p-Xylene	108-38-3/106-42-3	-	< 1	-
o-Xylene	95-47-6	-	< 1	-

Target Compounds	CAS #	R.T. (min.)	Concentration µg/l	% Fit
Styrene	100-42-5	-	< 1	-
Bromoform	75-25-2	-	< 1	-
iso-Propylbenzene	98-82-8	-	< 1	-
1,1,2,2-Tetrachloroethane	79-34-5	-	< 1	-
Propylbenzene	103-65-1	-	< 1	-
Bromobenzene	108-86-1	-	< 1	-
1,2,3-Trichloropropane	96-18-4	-	< 1	-
2-Chlorotoluene	95-49-8	-	< 1	-
1,3,5-Trimethylbenzene	108-67-8	-	< 1	-
4-Chlorotoluene	106-43-4	-	< 1	-
tert-Butylbenzene	98-06-6	-	< 1	-
1,2,4-Trimethylbenzene	95-53-6	-	< 1	-
sec-Butylbenzene	135-98-8	-	< 1	-
p-Isopropyltoluene	99-87-6	-	< 1	-
1,3-Dichlorobenzene	541-73-1	-	< 1	-
1,4-Dichlorobenzene	106-46-7	-	< 1	-
n-Butylbenzene	104-51-8	-	< 1	-
1,2-Dichlorobenzene	95-50-1	-	< 1	-
1,2-Dibromo-3-chloropropane	96-12-8	-	< 5	-
1,2,4-Trichlorobenzene	120-82-1	-	< 5	-
Hexachlorobutadiene	87-68-3	-	< 5	-
Naphthalene	91-20-3	-	< 5	-
1,2,3-Trichlorobenzene	87-61-6	-	< 5	-

Internal standards	R.T.	Area %	Surrogates	% Rec
Pentafluorobenzene	11.89	82	Dibromofluoromethane	109
1,4-Difluorobenzene	13.04	85	Toluene-d8	100
Chlorobenzene-d5	17.38	87	Bromofluorobenzene	100
1,4-Dichlorobenzene-d4	20.86	89		

TEST REPORT

WATER SAMPLE ANALYSIS

TES Report No. W/EXR/000771

Site: M4 RELIEF RD

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
CF31 1JZ

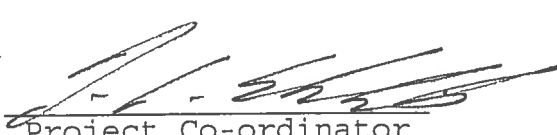
The 5 Samples described in this report were scheduled for analysis by TES Bretby on Wednesday, 16 February 2000. The analysis was completed by Monday, 6 March 2000.

Tests marked as 'not UKAS accredited' and any opinions or interpretations expressed herein are outside the scope of any UKAS accreditation held by TES Bretby laboratories.

The following tables are contained in this report:

Table 1 Sample Descriptions
Table 2 Main Analysis Results
Table 3 Phenol Results
Tables of Poly Aromatic Hydrocarbons (5 Pages)
Tables of Volatile Organic Compounds (1 Page)

On behalf of
TES Bretby :
J Elstub


Project Co-ordinator

Date of Issue: 06/03/00

TEST REPORT

WATER SAMPLE ANALYSIS

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
CF31 1JZ

TES Report No. W/EXR/000771

Client: EXPLORATION ASS.
Site: M4 RELIEF RD

Page No	ID No. W/EX/		Sample Date
1	0003674	BH CH 12 @ 1.35m	07/02/00
1	0003675	BH CH 17 @ 2.70m	
1	0003676	BH CH 32 @ 0.10m	
1	0003677	BH CH 45 @ 1.30m	
2	0003678	BH CH 51 @ 0.80m	

Date of Issue: 06/03/00

TES Bretby accepts no responsibility for the sampling related to the above results.
TES Bretby, P.O. Box 100, Burton-on-Trent, DE15 0XD
Telephone: 01283 554400 Fax: 01283 554422

TES Bretby
Report Number
W/EXR/000771
Table 1
Sheet 1/ 1

TEST REPORT

WATER SAMPLE ANALYSIS

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
CF31 1JZ

TES Report No. W/EXR/000771

Client: EXPLORATION ASS.
Site: M4 RELIEF RD

Sample reference	Sample 01	Sample 02	Sample 03	Sample 04
Sample Date	07/02/00			
TES Bretby ID Number	0003674	0003675	0003676	0003677

UKAS accredited	Test No.	0003674	0003675	0003676	0003677
pH units	WSLM3	7.4	7.7	10.8	8.6
Chloride as Cl	WSLM1				24
Total Sulphur as SO ₄ (Dissolved)	ICPPWW13	79.8	253	65.5	4.07
Ammoniacal Nitrogen as N	AMMKA40				<0.2
Cyanide (Total) as CN	ICTSCN28		<0.025	<0.025	<0.025
Thiocyanate as SCN	WSLM5				0.2
Sulphide (Free) as S	ISESW24				<0.2
Chemical Oxygen Demand	WSLM11				33
Biochemical Oxygen Demand	WSLM20				3.0
TPH FTIR	FTIRSW		0.08	4.15	0.41
Phenol Index as C ₆ H ₅ OH	WSLM4	<0.05	<0.05	0.06	<0.05
TPH GC	TPHFID	0.22	0.12	14.0	5.7
Volatile Organic Compounds	VOCSW				*
Pentachlorophenol	SVOC SW				<0.050

not UKAS accredited		0003674	0003675	0003676	0003677
Nickel as Ni (Dissolved)		0.007	0.009	0.009	0.004
Chromium as Cr (Dissolved)		<0.001	<0.001	0.001	0.018
Cadmium as Cd (Dissolved)		<0.0001	<0.0001	<0.0001	<0.0001
Copper as Cu (Dissolved)		0.013	0.002	0.019	0.009
Lead as Pb (Dissolved)		<0.001	<0.001	0.010	<0.001
Zinc as Zn (Dissolved)		0.007	0.007	0.003	0.009
Arsenic as As (Dissolved)		0.002	0.003	0.007	0.009
Boron as B (Dissolved)		0.07	0.36	0.76	<0.05
Mercury as Hg (Dissolved)		<0.0001	<0.0001	0.0003	<0.0001
Selenium as Se (Dissolved)		<0.001	<0.001	0.004	<0.001
Cyanide (Free) as CN					<0.025
Complex Cyanide as CN					<0.025
PCB - ARO				0.0008	0.0023
PAH HPLC		*	*	*	*

Results expressed as mg/l unless stated, for the sample as received

* See attached sheets for results

Sample 01: BH CH 12 @ 1.35m

Sample 02: BH CH 17 @ 2.70m

Sample 03: BH CH 32 @ 0.10m

Sample 04: BH CH 45 @ 1.30m

Date of Issue: 06/03/00

TES Bretby accepts no responsibility for the sampling related to the above results.
TES Bretby, P.O. Box 100, Burton-on-Trent, DE15 0XD
Telephone: 01283 554400 Fax: 01283 554422

TES Bretby	
Report Number	W/EXR/000771
Table	2
Sheet	1/ 2

TEST REPORT

WATER SAMPLE ANALYSIS

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
CF31 1JZ

TES Report No. W/EXR/000771

Client: EXPLORATION ASS.
Site: M4 RELIEF RD

Sample reference	Sample 05			
Sample Date				
TES Bretby ID Number	W/EX/..	0003678		

UKAS accredited	Test No.	0003678			
pH units	WSLM3	10.1			
Total Sulphur as SO4 (Dissolved)	ICPPWW13	62.0			
Cyanide (Total) as CN	ICTSCN28	0.09			
Sulphide (Free) as S	ISESW24	<0.2			
TPH FTIR	FTIRSW	0.09			
Phenol Index as C6H5OH	WSLM4	<0.05			
TPH GC	TPHFID	0.57			

not UKAS accredited		0003678			
Nickel as Ni (Dissolved)		0.004			
Chromium as Cr (Dissolved)		0.005			
Cadmium as Cd (Dissolved)		<0.0001			
Copper as Cu (Dissolved)		0.009			
Lead as Pb (Dissolved)		<0.001			
Zinc as Zn (Dissolved)		0.026			
Arsenic as As (Dissolved)		0.006			
Boron as B (Dissolved)		0.19			
Mercury as Hg (Dissolved)		<0.0001			
Selenium as Se (Dissolved)		<0.001			
PAH HPLC		*			

Results expressed as mg/l unless stated, for the sample as received

* See attached sheets for results

Sample 05: BH CH 51 @ 0.80m

Date of Issue: 06/03/00

TES Bretby accepts no responsibility for the sampling related to the above results.
TES Bretby, P.O. Box 100, Burton-on-Trent, DE15 0XD
Telephone: 01283 554400 Fax: 01283 554422

TES Bretby	
Report Number	W/EXR/000771
Table	2
Sheet	2 / 2

TEST REPORT

WATER SAMPLE ANALYSIS

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
CF31 1JZ

TES Report No. W/EXR/000771

Client: EXPLORATION ASS.
Site: M4 RELIEF RD

Phenol Analysis

Sample reference	Sample 01			
Sample Date				
TES Bretby ID Number	W/EX/..	0003677		
not UKAS accredited	0003677			
Phenol	<0.0005			
Cresols	<0.0005			
Dimethylphenols	<0.0005			
Trimethylphenols	<0.0005			

Results expressed as mg/l unless stated, for the sample as received
Sample 01: BH CH 45 @ 1.30m

Date of Issue: 06/03/00

TES Bretby accepts no responsibility for the sampling related to the above results.
TES Bretby, P.O. Box 100, Burton-on-Trent, DE15 0XD
Telephone: 01283 554400 Fax: 01283 554422

TES Bretby
Report Number
W/EXR/000771
Table 3
Sheet 1/ 1



Polycyclic Aromatic Hydrocarbons by HPLC

Customer and Site Details:	Exploration Ass : M4 Relief Road	Report Number:	771
Sample Details:	BH CH 12 @ 1.35m	Date Booked in:	16-Feb-00
LIMS ID Number:	EX3674	Date Extracted:	23-Feb-00
QC Batch Number:	286	Date Analysed:	24-Feb-00
Directory:	0223PAH.LC1	Matrix:	Water
Dilution (fluorescence):	1.0	Ext Method:	Sep Funnel
Dilution (UV):	1.0		

Target Compounds	CAS #	Concentration µg/l	Fluorescence / UV
Naphthalene	91-20-3	< 0.50	Fluorescence
Acenaphthylene	208-96-8	< 1.00	UV
Acenaphthene	83-32-9	< 0.05	Fluorescence
Fluorene	86-73-7	< 0.05	Fluorescence
Phenanthrene	85-01-8	< 0.05	Fluorescence
Anthracene	120-12-7	< 0.05	Fluorescence
Fluoranthene	206-44-0	< 0.05	Fluorescence
Pyrene	129-00-0	< 0.05	Fluorescence
Benzo[a]anthracene	56-55-3	< 0.05	Fluorescence
Chrysene	218-01-9	< 0.05	Fluorescence
Benzo[b]fluoranthene	205-99-2	< 0.05	Fluorescence
Benzo[k]fluoranthene	207-08-9	< 0.05	Fluorescence
Benzo[a]pyrene	50-32-8	< 0.05	Fluorescence
Indeno[1,2,3-cd]pyrene	193-39-5	< 0.05	Fluorescence
Dibenzo[a,h]anthracene	53-70-3	< 0.05	Fluorescence
Benzo[g,h,i]perylene	191-24-2	< 0.05	Fluorescence
Total PAHs (USEPA 16)	-	< 2.20	-

Data may be obtained from either the fluorescence detector or when the response is very high a less sensitive UV detector (this is indicated for each compound).
Any dilutions necessary during analysis have been included in the calculations.
Reported results do not require any corrections.



Polycyclic Aromatic Hydrocarbons by HPLC

Customer and Site Details:	Exploration Ass : M4 Relief Road	Report Number:	771
Sample Details:	BH CH 17 @ 2.70m	Date Booked in:	16-Feb-00
LIMS ID Number:	EX3675	Date Extracted:	23-Feb-00
QC Batch Number:	286	Date Analysed:	24-Feb-00
Directory:	0223PAH.LC1	Matrix:	Water
Dilution (fluorescence):	1.0	Ext Method:	Sep Funnel
Dilution (UV):	1.0		

Target Compounds	CAS #	Concentration µg/l	Fluorescence / UV
Naphthalene	91-20-3	< 0.50	Fluorescence
Acenaphthylene	208-96-8	< 1.00	UV
Acenaphthene	83-32-9	< 0.05	Fluorescence
Fluorene	86-73-7	< 0.05	Fluorescence
Phenanthrene	85-01-8	< 0.05	Fluorescence
Anthracene	120-12-7	< 0.05	Fluorescence
Fluoranthene	206-44-0	< 0.05	Fluorescence
Pyrene	129-00-0	< 0.05	Fluorescence
Benzo[a]anthracene	56-55-3	< 0.05	Fluorescence
Chrysene	218-01-9	< 0.05	Fluorescence
Benzo[b]fluoranthene	205-99-2	< 0.05	Fluorescence
Benzo[k]fluoranthene	207-08-9	< 0.05	Fluorescence
Benzo[a]pyrene	50-32-8	< 0.05	Fluorescence
Indeno[1,2,3-cd]pyrene	193-39-5	< 0.05	Fluorescence
Dibenzo[a,h]anthracene	53-70-3	< 0.05	Fluorescence
Benzo[g,h,i]perylene	191-24-2	< 0.05	Fluorescence
Total PAHs (USEPA 16)	-	< 2.20	-

Data may be obtained from either the fluorescence detector or when the response is very high a less sensitive UV detector (this is indicated for each compound).
Any dilutions necessary during analysis have been included in the calculations.
Reported results do not require any corrections.



Polycyclic Aromatic Hydrocarbons by HPLC

Customer and Site Details:	Exploration Ass : M4 Relief Road		
Sample Details:	BH CH 32 @ 0.10m	Report Number:	771
LIMS ID Number:	EX3676	Date Booked in:	16-Feb-00
QC Batch Number:	286	Date Extracted:	23-Feb-00
Directory:	0223PAH.LC1	Date Analysed:	24-Feb-00
Dilution (fluorescence):	1.0	Matrix:	Water
Dilution (UV):	1.0	Ext Method:	Sep Funnel

Target Compounds	CAS #	Concentration µg/l	Fluorescence / UV
Naphthalene	91-20-3	< 1.00	UV
Acenaphthylene	208-96-8	< 1.00	UV
Acenaphthene	83-32-9	< 0.05	Fluorescence
Fluorene	86-73-7	0.63	Fluorescence
Phenanthrene	85-01-8	0.38	Fluorescence
Anthracene	120-12-7	0.06	Fluorescence
Fluoranthene	206-44-0	0.17	Fluorescence
Pyrene	129-00-0	0.21	Fluorescence
Benzo[a]anthracene	56-55-3	0.12	Fluorescence
Chrysene	218-01-9	< 0.05	Fluorescence
Benzo[b]fluoranthene	205-99-2	< 0.05	Fluorescence
Benzo[k]fluoranthene	207-08-9	< 0.05	Fluorescence
Benzo[a]pyrene	50-32-8	< 0.05	Fluorescence
Indeno[1,2,3-cd]pyrene	193-39-5	< 0.05	Fluorescence
Dibenzo[a,h]anthracene	53-70-3	< 0.05	Fluorescence
Benzo[g,h,i]perylene	191-24-2	< 0.05	Fluorescence
Total PAHs (USEPA 16)	-	< 3.97	-

Data may be obtained from either the fluorescence detector or when the response is very high a less sensitive UV detector (this is indicated for each compound).
Any dilutions necessary during analysis have been included in the calculations.
Reported results do not require any corrections.



Polycyclic Aromatic Hydrocarbons by HPLC

Customer and Site Details:	Exploration Ass : M4 Relief Road	Report Number:	771
Sample Details:	BH CH 45 @ 1.30m	Date Booked in:	16-Feb-00
LIMS ID Number:	EX3677	Date Extracted:	23-Feb-00
QC Batch Number:	286	Date Analysed:	24-Feb-00
Directory:	0223PAH.LC1	Matrix:	Water
Dilution (fluorescence):	1.0	Ext Method:	Sep Funnel
Dilution (UV):	1.0		

Target Compounds	CAS #	Concentration µg/l	Fluorescence / UV
Naphthalene	91-20-3	< 0.50	Fluorescence
Acenaphthylene	208-96-8	< 1.00	UV
Acenaphthene	83-32-9	< 0.05	Fluorescence
Fluorene	86-73-7	< 0.05	Fluorescence
Phenanthrene	85-01-8	< 0.05	Fluorescence
Anthracene	120-12-7	< 0.05	Fluorescence
Fluoranthene	206-44-0	< 0.05	Fluorescence
Pyrene	129-00-0	< 0.05	Fluorescence
Benzo[a]anthracene	56-55-3	< 0.05	Fluorescence
Chrysene	218-01-9	< 0.05	Fluorescence
Benzo[b]fluoranthene	205-99-2	< 0.05	Fluorescence
Benzo[k]fluoranthene	207-08-9	< 0.05	Fluorescence
Benzo[a]pyrene	50-32-8	< 0.05	Fluorescence
Indeno[1,2,3-cd]pyrene	193-39-5	< 0.05	Fluorescence
Dibenzo[a,h]anthracene	53-70-3	< 0.05	Fluorescence
Benzo[g,h,i]perylene	191-24-2	< 0.05	Fluorescence
Total PAHs (USEPA 16)	-	< 2.20	-

Data may be obtained from either the fluorescence detector or when the response is very high a less sensitive UV detector (this is indicated for each compound).
Any dilutions necessary during analysis have been included in the calculations.
Reported results do not require any corrections.



Polycyclic Aromatic Hydrocarbons by HPLC

Customer and Site Details:	Exploration Ass : M4 Relief Road	Report Number:	771
Sample Details:	BH CH 51 @ 0.80m	Date Booked in:	16-Feb-00
LIMS ID Number:	EX3678	Date Extracted:	23-Feb-00
QC Batch Number:	286	Date Analysed:	24-Feb-00
Directory:	0223PAH.LC1	Matrix:	Water
Dilution (fluorescence):	1.0	Ext Method:	Sep Funnel
Dilution (UV):	1.0		

Target Compounds	CAS #	Concentration µg/l	Fluorescence / UV
Naphthalene	91-20-3	< 0.50	Fluorescence
Acenaphthylene	208-96-8	< 1.00	UV
Acenaphthene	83-32-9	< 0.05	Fluorescence
Fluorene	86-73-7	< 0.05	Fluorescence
Phenanthrene	85-01-8	< 0.05	Fluorescence
Anthracene	120-12-7	< 0.05	Fluorescence
Fluoranthene	206-44-0	< 0.05	Fluorescence
Pyrene	129-00-0	< 0.05	Fluorescence
Benzo[a]anthracene	56-55-3	< 0.05	Fluorescence
Chrysene	218-01-9	< 0.05	Fluorescence
Benzo[b]fluoranthene	205-99-2	< 0.05	Fluorescence
Benzo[k]fluoranthene	207-08-9	< 0.05	Fluorescence
Benzo[a]pyrene	50-32-8	< 0.05	Fluorescence
Indeno[1,2,3-cd]pyrene	193-39-5	< 0.05	Fluorescence
Dibenzo[a,h]anthracene	53-70-3	< 0.05	Fluorescence
Benzo[g,h,i]perylene	191-24-2	< 0.05	Fluorescence
Total PAHs (USEPA 16)	-	< 2.20	-

Data may be obtained from either the fluorescence detector or when the response is very high a less sensitive UV detector (this is indicated for each compound).
Any dilutions necessary during analysis have been included in the calculations.
Reported results do not require any corrections.

Volatile Organic Compounds by PTGCMS

Customer and Site Details: Exploration Associates : M4 Relief Road
Sample Details: BH CH 45 / 1.30m
LIMS ID Number: EX03677
Report Number: 000771

Directory/Quant file: 0221VOC.MS3\ 0221CCC1.D
Date Booked in: 16-Feb-00
Date Analysed: 22-Feb-00
Operator: D.WALTON

Matrix: Water
Method: Purge and Trap
Dilution: 1
Position: 10

Target Compounds	CAS #	R.T. (min.)	Concentration µg/l	% Fit
Dichlorodifluoromethane	75-71-8	-	< 1	-
Chloromethane	74-87-3	-	< 1	-
Vinyl Chloride	75-01-4	-	< 1	-
Bromomethane	74-83-9	-	< 1	-
Chloroethane	75-00-3	-	< 1	-
Trichlorofluoromethane	75-69-4	-	< 1	-
1,1-Dichloroethene	75-35-4	-	< 1	-
trans 1,2-Dichloroethene	156-60-5	-	< 1	-
1,1-Dichloroethane	75-34-3	-	< 1	-
2,2-Dichloropropane	594-20-7	-	< 1	-
cis 1,2-Dichloroethene	156-59-2	-	< 1	-
Bromochloromethane	74-97-5	-	< 1	-
Chloroform	67-66-3	-	< 1	-
1,1,1-Trichloroethane	71-55-6	-	< 1	-
Carbon Tetrachloride	56-23-5	-	< 1	-
1,1-Dichloropropene	563-58-6	-	< 1	-
Benzene	71-43-2	-	< 1	-
1,2-Dichloroethane	107-06-2	-	< 1	-
Trichloroethene	79-01-6	-	< 1	-
1,2-Dichloropropane	78-87-5	-	< 1	-
Dibromomethane	74-95-3	-	< 1	-
Bromodichloromethane	75-27-4	-	< 1	-
cis 1,3-Dichloropropene	10061-01-5	-	< 1	-
Toluene	108-88-3	-	< 1	-
trans 1,3-Dichloropropene	10061-02-6	-	< 1	-
1,1,2-Trichloroethane	79-00-5	-	< 1	-
Tetrachloroethene	127-18-4	-	< 1	-
1,3-Dichloropropane	142-28-29	-	< 1	-
Dibromochloromethane	124-48-1	-	< 1	-
1,2-Dibromoethane	106-93-4	-	< 1	-
Chlorobenzene	108-90-7	-	< 1	-
Ethylbenzene	100-41-4	-	< 1	-
1,1,1,2-Tetrachloroethane	630-20-6	-	< 1	-
m and p-Xylene	108-38-3/106-42-3	-	< 1	-
o-Xylene	95-47-6	-	< 1	-

Target Compounds	CAS #	R.T. (min.)	Concentration µg/l	% Fit
Styrene	100-42-5	-	< 1	-
Bromoform	75-25-2	-	< 1	-
iso-Propylbenzene	98-82-8	-	< 1	-
1,1,2,2-Tetrachloroethane	79-34-5	-	< 1	-
Propylbenzene	103-65-1	-	< 1	-
Bromobenzene	108-86-1	-	< 1	-
1,2,3-Trichloropropane	96-18-4	-	< 1	-
2-Chlorotoluene	95-49-8	-	< 1	-
1,3,5-Trimethylbenzene	108-67-8	-	< 1	-
4-Chlorotoluene	106-43-4	-	< 1	-
tert-Butylbenzene	98-06-6	-	< 1	-
1,2,4-Trimethylbenzene	95-63-6	-	< 1	-
sec-Butylbenzene	135-98-8	-	< 1	-
p-Isopropyltoluene	99-87-6	-	< 1	-
1,3-Dichlorobenzene	541-73-1	-	< 1	-
1,4-Dichlorobenzene	106-46-7	-	< 1	-
n-Butylbenzene	104-51-8	-	< 1	-
1,2-Dichlorobenzene	95-50-1	-	< 1	-
1,2-Dibromo-3-chloropropane	96-12-8	-	< 5	-
1,2,4-Trichlorobenzene	120-82-1	-	< 5	-
Hexachlorobutadiene	87-68-3	-	< 5	-
Naphthalene	91-20-3	-	< 5	-
1,2,3-Trichlorobenzene	87-61-6	-	< 5	-

Internal standards	R.T.	Area %	Surrogates	% Rec
Pentafluorobenzene	11.89	94	Dibromofluoromethane	109
1,4-Difluorobenzene	13.04	97	Toluene-d8	97
Chlorobenzene-d5	17.38	93	Bromofluorobenzene	89
1,4-Dichlorobenzene-d4	20.87	82		

TEST REPORT

WATER SAMPLE ANALYSIS

TES Report No. W/EXR/000800

Site: M4 RELIEF RD

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
CF31 1JZ

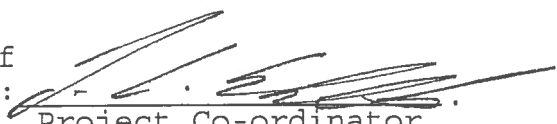
The 4 Samples described in this report were scheduled for analysis by TES Bretby on Thursday, 17 February 2000. The analysis was completed by Friday, 3 March 2000.

Tests marked as 'not UKAS accredited' and any opinions or interpretations expressed herein are outside the scope of any UKAS accreditation held by TES Bretby laboratories.

The following tables are contained in this report:

Table 1 Sample Descriptions
Table 2 Main Analysis Results
Table 3 Phenol Results
Tables of Poly Aromatic Hydrocarbons (4 Pages)
Table of Volatile Organic Compounds (1 Page)

On behalf of
TES Bretby :
J Elstub


Project Co-ordinator

Date of Issue: 03/03/00

TEST REPORT

WATER SAMPLE ANALYSIS

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
CF31 1JZ

TES Report No. W/EXR/000800

Client: EXPLORATION ASS.

Site: M4 RELIEF RD

Page No	ID No. W/EX/		Sample Date
1	0003829	BH CH 52 @ 0.00m	
1	0003830	BH CH 52A @ 0.80m	
1	0003831	Site No 7 Pond Sample	
1	0003832	Site No 9 Reen Sample	

Date of Issue: 03/03/00

TES Bretby accepts no responsibility for the sampling related to the above results.
TES Bretby, P.O. Box 100, Burton-on-Trent, DE15 0XD
Telephone: 01283 554400 Fax: 01283 554422

TES Bretby
Report Number
W/EXR/000800
Table 1
Sheet 1/ 1

TEST REPORT

WATER SAMPLE ANALYSIS

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
CF31 1JZ

TES Report No. W/EXR/000800

Client: EXPLORATION ASS.
Site: M4 RELIEF RD

Sample reference Sample Date TES Bretby ID Number	W/EX/..	Sample 01	Sample 02	Sample 03	Sample 04
		0003829	0003830	0003831	0003832
UKAS accredited	Test No.	0003829	0003830	0003831	0003832
pH units	WSLM3	7.5	11.9	7.9	7.2
Total Sulphur as SO ₄ (Dissolved)	ICPPWW13	38.5	26.4	60.7	87.3
Thiocyanate as SCN	WSLM5		0.2	0.2	
Sulphide (Free) as S	ISESW24	<0.2	<0.2	<0.2	<0.2
Chemical Oxygen Demand	WSLM11			16	
Biochemical Oxygen Demand	WSLM20			3.5	
TPH FTIR	FTIRSW	0.08	0.08	<0.05	<0.05
Phenol Index as C ₆ H ₅ OH	WSLM4	<0.05	<0.05	<0.05	<0.05
TPH GC	TPHFID	0.38	0.27	<0.1	<0.1
Volatile Organic Compounds	VOCSW			*	

not UKAS accredited		0003829	0003830	0003831	0003832
Nickel as Ni (Dissolved)		0.005	0.002	0.006	0.005
Chromium as Cr (Dissolved)		<0.001	<0.001	<0.001	<0.001
Cadmium as Cd (Dissolved)		<0.0001	<0.0001	<0.0001	<0.0001
Copper as Cu (Dissolved)		0.003	0.002	0.003	0.009
Lead as Pb (Dissolved)		<0.001	<0.001	<0.001	0.005
Zinc as Zn (Dissolved)		0.050	<0.001	0.034	0.029
Arsenic as As (Dissolved)		<0.001	<0.001	0.003	<0.001
Boron as B (Dissolved)		0.06	<0.05	0.18	0.09
Mercury as Hg (Dissolved)		<0.0001	<0.0001	<0.0001	<0.0001
Selenium as Se (Dissolved)		<0.001	0.003	<0.001	<0.001
Cyanide (Free) as CN			<0.025	<0.025	
Cyanide (Total) as CN		<0.025	<0.025	<0.025	<0.025
Complex Cyanide as CN			<0.025	<0.025	
PCB - ARO				<0.001	
PAH HPLC		*	*	*	*

Results expressed as mg/l unless stated, for the sample as received

* See attached sheets for results

Sample 01: BH CH 52 @ 0.00m

Sample 02: BH CH 52A @ 0.80m

Sample 03: Site No 7 Pond Sample

Sample 04: Site No 9 Reen Sample

Date of Issue: 03/03/00

TES Bretby accepts no responsibility for the sampling related to the above results.
TES Bretby, P.O. Box 100, Burton-on-Trent, DE15 0XD
Telephone: 01283 554400 Fax: 01283 554422

TES Bretby
Report Number
W/EXR/000800
Table 2
Sheet 1/1

TEST REPORT

WATER SAMPLE ANALYSIS

Exploration Associates
Unit 15
Crosby Yard
Wildmill
Bridgend
CF31 1JZ

TES Report No. W/EXR/000800

Client: EXPLORATION ASS.
Site: M4 RELIEF RD

Phenol Analysis

Sample reference	Sample 01	Sample 02		
Sample Date				
TES Bretby ID Number W/EX/..	0003830	0003831		
not UKAS accredited	0003830	0003831		
Phenol	0.0204	<0.0005		
Cresols	0.0401	<0.0005		
Dimethylphenols	0.0186	<0.0005		
Trimethylphenols	<0.0005	<0.0005		

Results expressed as mg/l unless stated, for the sample as received

Sample 01: BH CH 52A @ 0.80m

Sample 02: Site No 7 Pond Sample

Date of Issue: 03/03/00

TES Bretby accepts no responsibility for the sampling related to the above results.
TES Bretby, P.O. Box 100, Burton-on-Trent, DE15 0XD
Telephone: 01283 554400 Fax: 01283 554422

TES Bretby
Report Number
W/EXR/000800
Table 3
Sheet 1/ 1



Polycyclic Aromatic Hydrocarbons by HPLC

Customer and Site Details:	Exploration Ass. : M4 Relief Road	Report Number:	800
Sample Details:	BH CH 52 @ 0.00m	Date Booked in:	17-Feb-00
LIMS ID Number:	EX3829	Date Extracted:	28-Feb-00
QC Batch Number:	288	Date Analysed:	29-Feb-00
Directory:	0228PAH.LC2	Matrix:	Water
Dilution (fluorescence):	1.0	Ext Method:	Sep Funnel
Dilution (UV):	1.0		

Target Compounds	CAS #	Concentration µg/l	Fluorescence / UV
Naphthalene	91-20-3	< 0.50	Fluorescence
Acenaphthylene	208-96-8	< 1.00	UV
Acenaphthene	83-32-9	< 0.05	Fluorescence
Fluorene	86-73-7	< 0.05	Fluorescence
Phenanthrene	85-01-8	< 0.05	Fluorescence
Anthracene	120-12-7	< 0.05	Fluorescence
Fluoranthene	206-44-0	< 0.05	Fluorescence
Pyrene	129-00-0	< 0.05	Fluorescence
Benzo[a]anthracene	56-55-3	< 0.05	Fluorescence
Chrysene	218-01-9	< 0.05	Fluorescence
Benzo[b]fluoranthene	205-99-2	< 0.05	Fluorescence
Benzo[k]fluoranthene	207-08-9	< 0.05	Fluorescence
Benzo[a]pyrene	50-32-8	< 0.05	Fluorescence
Indeno[1,2,3-cd]pyrene	193-39-5	< 0.05	Fluorescence
Dibenzo[a,h]anthracene	53-70-3	< 0.05	Fluorescence
Benzo[g,h,i]perylene	191-24-2	< 0.05	Fluorescence
Total PAHs (USEPA 16)	-	< 2.20	-

Data may be obtained from either the fluorescence detector or when the response is very high a less sensitive UV detector (this is indicated for each compound). Any dilutions necessary during analysis have been included in the calculations. Reported results do not require any corrections.



Polycyclic Aromatic Hydrocarbons by HPLC

Customer and Site Details:	Exploration Ass. : M4 Relief Road	Report Number:	800
Sample Details:	BH CH 52A @ 0.80m	Date Booked in:	17-Feb-00
LIMS ID Number:	EX3830	Date Extracted:	28-Feb-00
QC Batch Number:	288	Date Analysed:	29-Feb-00
Directory:	0228PAH.LC2	Matrix:	Water
Dilution (fluorescence):	1.0	Ext Method:	Sep Funnel
Dilution (UV):	1.0		

Target Compounds	CAS #	Concentration µg/l	Fluorescence / UV
Naphthalene	91-20-3	0.74	Fluorescence
Acenaphthylene	208-96-8	< 1.00	UV
Acenaphthene	83-32-9	< 0.05	Fluorescence
Fluorene	86-73-7	0.47	Fluorescence
Phenanthrene	85-01-8	0.69	Fluorescence
Anthracene	120-12-7	0.07	Fluorescence
Fluoranthene	206-44-0	0.16	Fluorescence
Pyrene	129-00-0	0.13	Fluorescence
Benzo[a]anthracene	56-55-3	0.10	Fluorescence
Chrysene	218-01-9	0.06	Fluorescence
Benzo[b]fluoranthene	205-99-2	0.07	Fluorescence
Benzo[k]fluoranthene	207-08-9	< 0.05	Fluorescence
Benzo[a]pyrene	50-32-8	0.06	Fluorescence
Indeno[1,2,3-cd]pyrene	193-39-5	< 0.05	Fluorescence
Dibenzo[a,h]anthracene	53-70-3	< 0.05	Fluorescence
Benzo[g,h,i]perylene	191-24-2	< 0.05	Fluorescence
Total PAHs (USEPA 16)	-	< 3.80	-

Data may be obtained from either the fluorescence detector or when the response is very high a less sensitive UV detector (this is indicated for each compound). Any dilutions necessary during analysis have been included in the calculations. Reported results do not require any corrections.



Polycyclic Aromatic Hydrocarbons by HPLC

Customer and Site Details:	Exploration Ass. : M4 Relief Road		
Sample Details:	Site No 7 Pond Sample	Report Number:	800
LIMS ID Number:	EX3831	Date Booked in:	17-Feb-00
QC Batch Number:	288	Date Extracted:	28-Feb-00
Directory:	0228PAH.LC2	Date Analysed:	29-Feb-00
Dilution (fluorescence):	1.0	Matrix:	Water
Dilution (UV):	1.0	Ext Method:	Sep Funnel

Target Compounds	CAS #	Concentration µg/l	Fluorescence / UV
Naphthalene	91-20-3	< 0.50	Fluorescence
Acenaphthylene	208-96-8	< 1.00	UV
Acenaphthene	83-32-9	< 0.05	Fluorescence
Fluorene	86-73-7	< 0.05	Fluorescence
Phenanthrene	85-01-8	< 0.05	Fluorescence
Anthracene	120-12-7	< 0.05	Fluorescence
Fluoranthene	206-44-0	< 0.05	Fluorescence
Pyrene	129-00-0	< 0.05	Fluorescence
Benzo[a]anthracene	56-55-3	< 0.05	Fluorescence
Chrysene	218-01-9	< 0.05	Fluorescence
Benzo[b]fluoranthene	205-99-2	< 0.05	Fluorescence
Benzo[k]fluoranthene	207-08-9	< 0.05	Fluorescence
Benzo[a]pyrene	50-32-8	< 0.05	Fluorescence
Indeno[1,2,3-cd]pyrene	193-39-5	< 0.05	Fluorescence
Dibenzo[a,h]anthracene	53-70-3	< 0.05	Fluorescence
Benzo[g,h,i]perylene	191-24-2	< 0.05	Fluorescence
Total PAHs (USEPA 16)	-	< 2.20	-

Data may be obtained from either the fluorescence detector or when the response is very high a less sensitive UV detector (this is indicated for each compound). Any dilutions necessary during analysis have been included in the calculations. Reported results do not require any corrections.

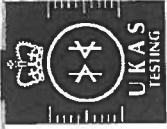


Polycyclic Aromatic Hydrocarbons by HPLC

Customer and Site Details:	Exploration Ass. : M4 Relief Road	Report Number:	800
Sample Details:	Site No 9 Reen Sample	Date Booked in:	17-Feb-00
LIMS ID Number:	EX3832	Date Extracted:	28-Feb-00
QC Batch Number:	288	Date Analysed:	29-Feb-00
Directory:	0228PAH.LC2	Matrix:	Water
Dilution (fluorescence):	1.0	Ext Method:	Sep Funnel
Dilution (UV):	1.0		

Target Compounds	CAS #	Concentration µg/l	Fluorescence / UV
Naphthalene	91-20-3	< 0.50	Fluorescence
Acenaphthylene	208-96-8	< 1.00	UV
Acenaphthene	83-32-9	< 0.05	Fluorescence
Fluorene	86-73-7	< 0.05	Fluorescence
Phenanthrene	85-01-8	< 0.05	Fluorescence
Anthracene	120-12-7	< 0.05	Fluorescence
Fluoranthene	206-44-0	< 0.05	Fluorescence
Pyrene	129-00-0	< 0.05	Fluorescence
Benzo[a]anthracene	56-55-3	< 0.05	Fluorescence
Chrysene	218-01-9	< 0.05	Fluorescence
Benzo[b]fluoranthene	205-99-2	< 0.05	Fluorescence
Benzo[k]fluoranthene	207-08-9	< 0.05	Fluorescence
Benzo[a]pyrene	50-32-8	< 0.05	Fluorescence
Indeno[1,2,3-cd]pyrene	193-39-5	< 0.05	Fluorescence
Dibenzo[a,h]anthracene	53-70-3	< 0.05	Fluorescence
Benzo[g,h,i]perylene	191-24-2	< 0.05	Fluorescence
Total PAHs (USEPA 16)	-	< 2.20	-

Data may be obtained from either the fluorescence detector or when the response is very high a less sensitive UV detector (this is indicated for each compound). Any dilutions necessary during analysis have been included in the calculations. Reported results do not require any corrections.



Volatile Organic Compounds by PTGCMS

Customer and Site Details:
Sample Details:
LIMS ID Number:
Report Number:

Exploration Associates : M4 Relief Road
Site No7 Pond Sample
EX03831
000800

Directory/Quant file:
Date Booked in:
Date Analysed:
Operator:

0222VOC.MS3\ 0221CCC3.D
17-Feb-00
23-Feb-00
D.WALTON

Matrix:
Method:
Dilution:
Position:

Water
Purge and Trap
1
4

1252

Target Compounds	CAS #	R.T. (min.)	Concentration µg/l	% Fit
Dichlorodifluoromethane	75-71-8	-	< 1	-
Chloromethane	74-87-3	-	< 1	-
Vinyl Chloride	75-01-4	-	< 1	-
Bromomethane	74-83-9	-	< 1	-
Chloroethane	75-00-3	-	< 1	-
Trichlorofluoromethane	75-69-4	-	< 1	-
1,1-Dichloroethane	75-35-4	-	< 1	-
trans 1,2-Dichloroethane	156-60-5	-	< 1	-
1,1-Dichloroethane	75-34-3	-	< 1	-
2,2-Dichloropropane	594-20-7	-	< 1	-
cis 1,2-Dichloroethane	156-59-2	-	< 1	-
Bromochloromethane	74-97-5	-	< 1	-
Chloroform	67-66-3	11.68	7	98
1,1,1-Trichloroethane	71-55-6	-	< 1	-
Carbon Tetrachloride	56-23-5	-	< 1	-
1,1-Dichloropropene	563-58-6	-	< 1	-
Benzene	71-43-2	-	< 1	-
1,2-Dichloroethane	107-06-2	-	< 1	-
Trichloroethene	79-01-6	-	< 1	-
1,2-Dichloropropane	78-87-5	-	< 1	-
Dibromomethane	74-95-3	-	< 1	-
Bromodichloromethane	75-27-4	14.20	2	96
cis 1,3-Dichloropropene	10061-01-5	-	< 1	-
Toluene	108-88-3	-	< 1	-
trans 1,3-Dichloropropene	10061-02-6	-	< 1	-
1,1,2-Trichloroethane	79-00-5	-	< 1	-
Tetrachloroethene	127-18-4	-	< 1	-
1,3-Dichloropropane	142-28-29	-	< 1	-
Dibromochloromethane	124-48-1	-	< 1	-
1,2-Dibromoethane	106-93-4	-	< 1	-
Chlorobenzene	108-90-7	-	< 1	-
Ethylbenzene	100-41-4	-	< 1	-
1,1,1,2-Tetrachloroethane	630-20-6	-	< 1	-
m and p-Xylene	108-38-3/106-42-3	-	< 1	-
o-Xylene	95-47-6	-	< 1	-

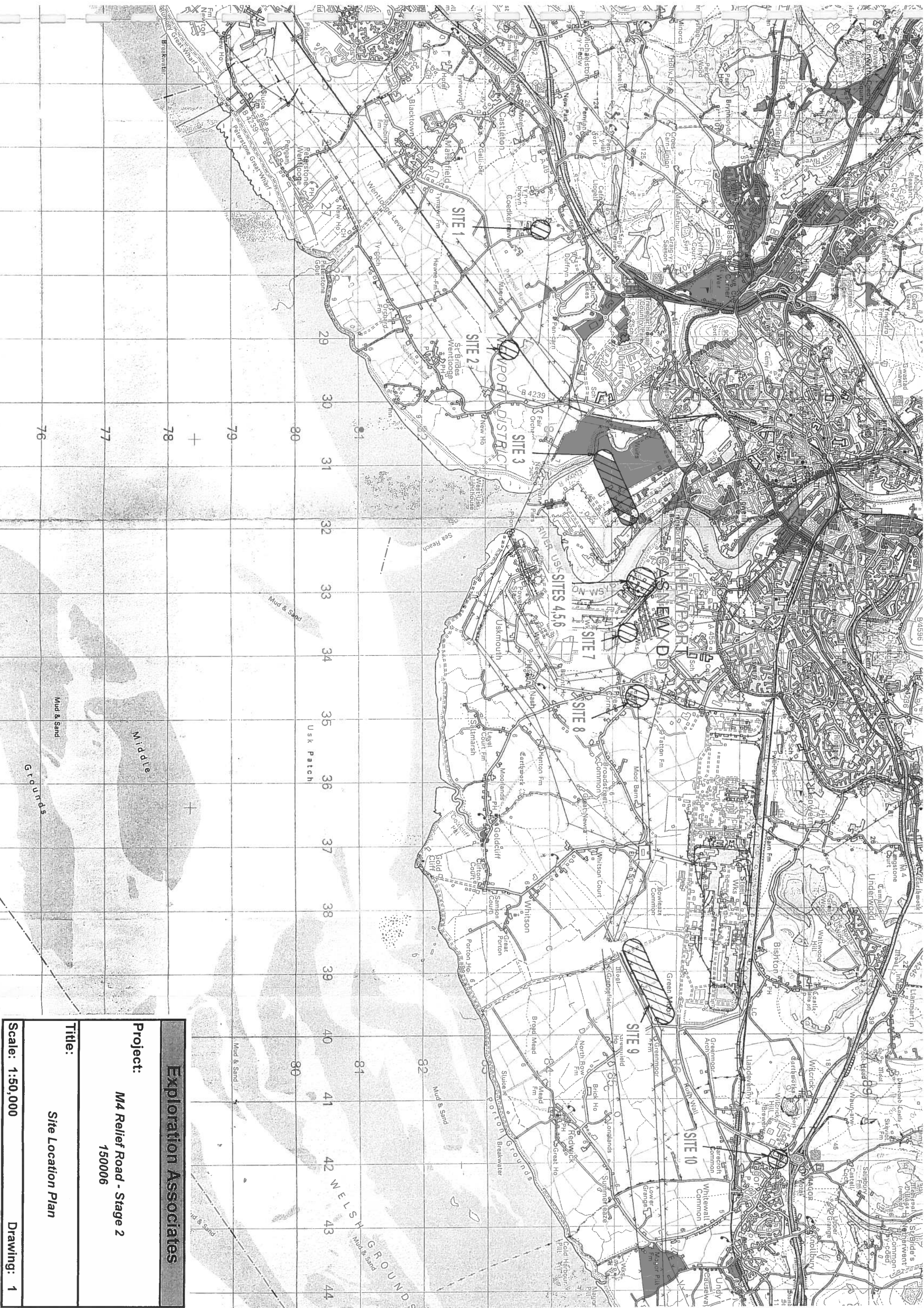
Target Compounds	CAS #	R.T. (min.)	Concentration µg/l	% Fit
Styrene	100-42-5	-	< 1	-
Bromoform	75-25-2	-	< 1	-
iso-Propylbenzene	98-82-8	-	< 1	-
1,1,2,2-Tetrachloroethane	79-34-5	-	< 1	-
Propylbenzene	103-65-1	-	< 1	-
Bromobenzene	108-86-1	-	< 1	-
1,2,3-Trichloropropane	96-18-4	-	< 1	-
2-Chlorotoluene	95-49-8	-	< 1	-
1,3,5-Trimethylbenzene	108-67-8	-	< 1	-
4-Chlorotoluene	106-43-4	-	< 1	-
tert-Butylbenzene	98-06-6	-	< 1	-
1,2,4-Trimethylbenzene	95-63-6	-	< 1	-
sec-Butylbenzene	135-98-8	-	< 1	-
p-Isopropyltoluene	99-87-6	-	< 1	-
1,3-Dichlorobenzene	541-73-1	-	< 1	-
1,4-Dichlorobenzene	106-46-7	-	< 1	-
n-Butylbenzene	104-51-8	-	< 1	-
1,2-Dichlorobenzene	95-50-1	-	< 1	-
1,2-Dibromo-3-chloropropane	96-12-8	-	< 5	-
1,2,4-Trichlorobenzene	120-82-1	-	< 5	-
Hexachlorobutadiene	87-68-3	-	< 5	-
Naphthalene	91-20-3	-	< 5	-
1,2,3-Trichlorobenzene	87-61-6	-	< 5	-

Internal standards	R.T.	Area %	Surrogates	% Rec
Pentafluorobenzene	11.89	124	Dibromofluoromethane	92
1,4-Difluorobenzene	13.04	126	Toluene-d8	95
Chlorobenzene-d5	17.37	113	Bromofluorobenzene	92
1,4-Dichlorobenzene-d4	20.87	95		

ENCLOSURE D

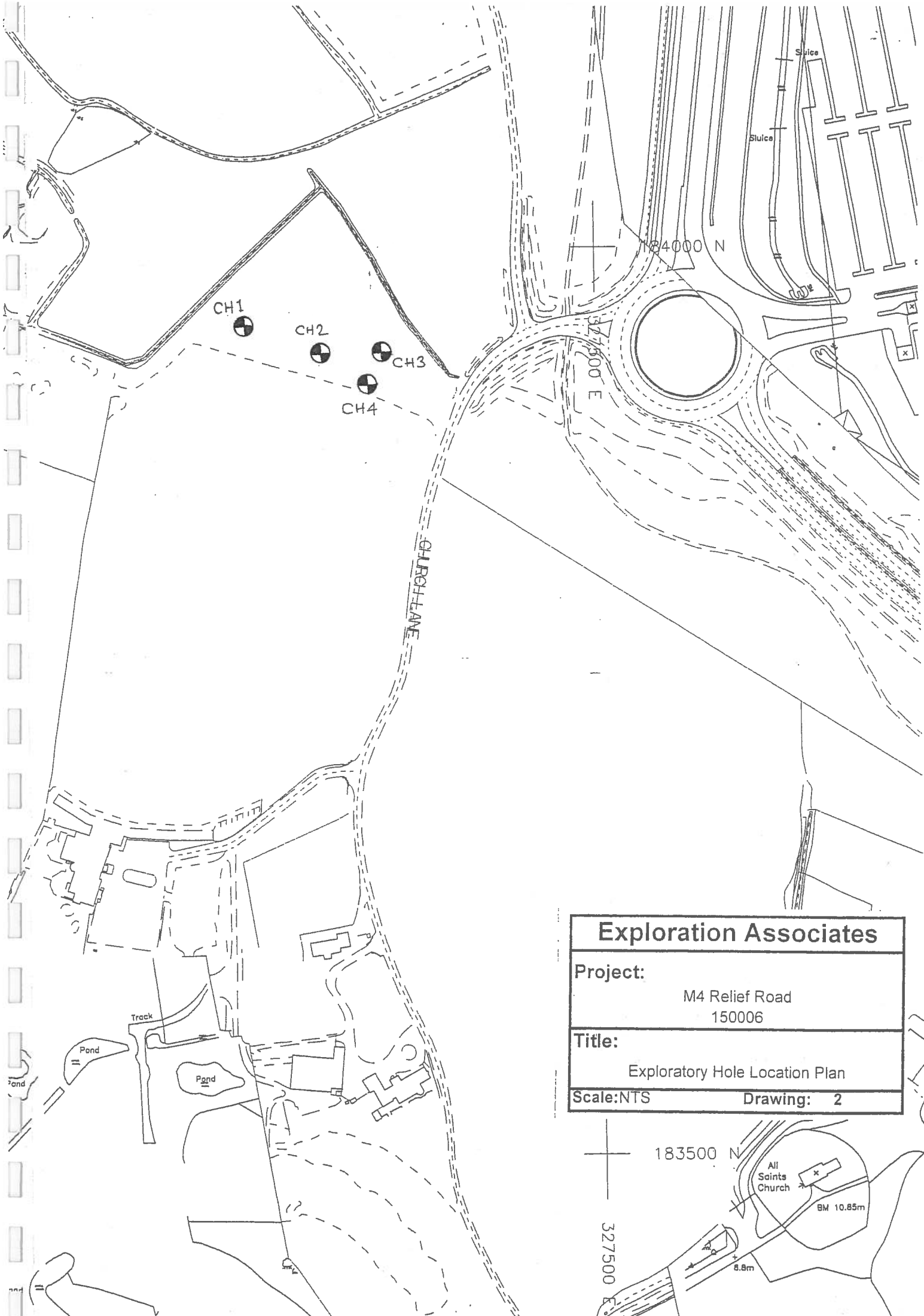
Drawings

Site Location Plan	1
Exploratory Hole Location Plan	2 to 13



Exploration Associates	
Project:	M4 Relief Road - Stage 2
	150006
Title:	Site Location Plan
Scale: 1:50,000	Drawing: 1

APPENDIX
General Notes



Exploration Associates

Project:

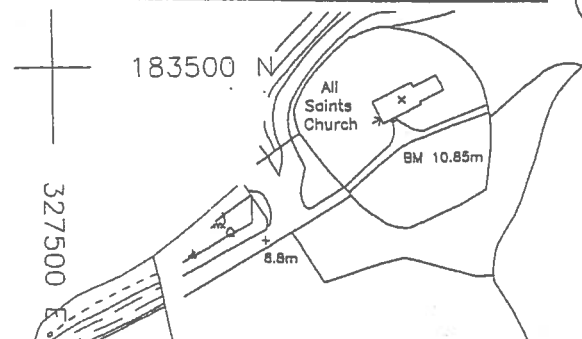
M4 Relief Road
150006

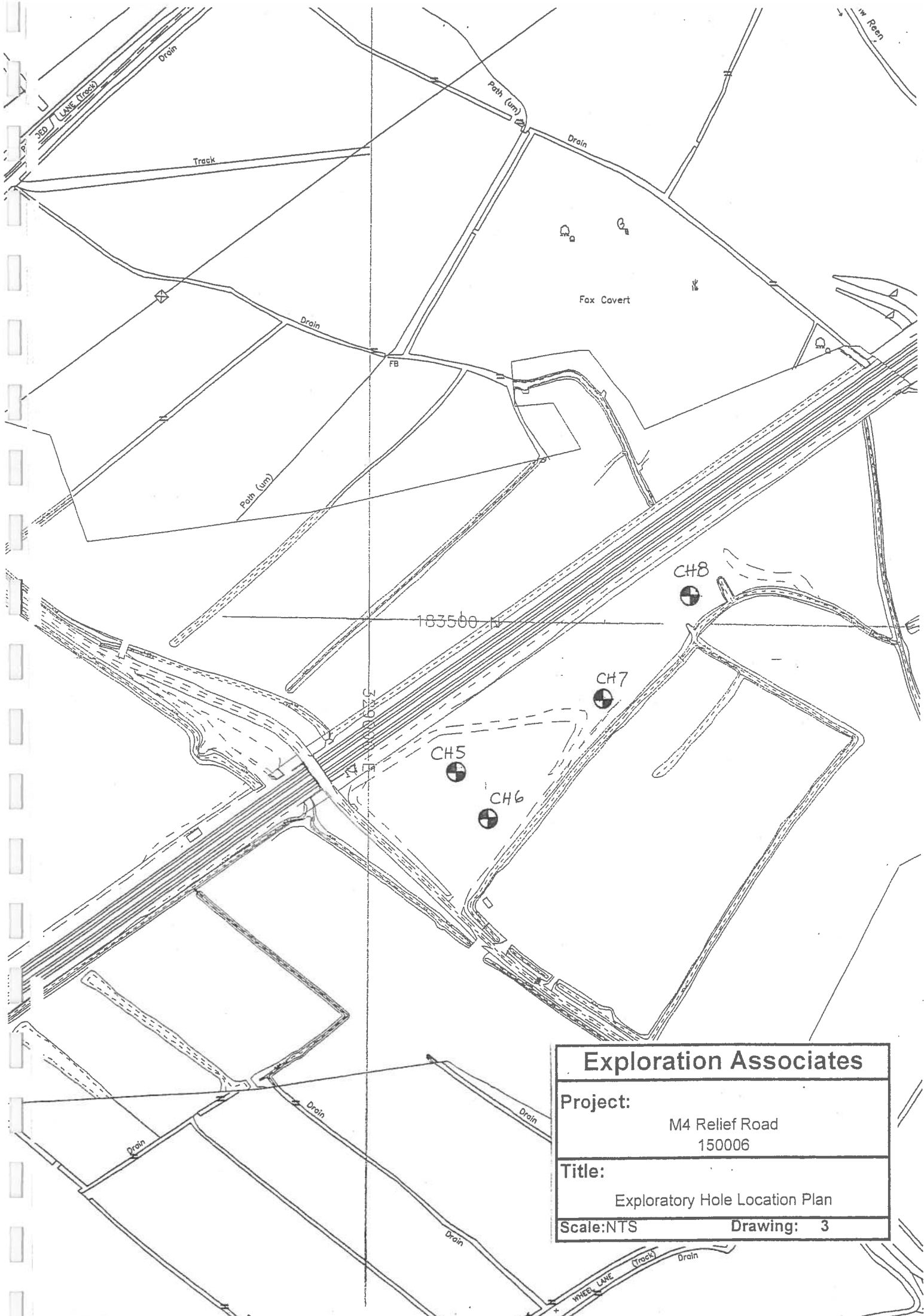
Title:

Exploratory Hole Location Plan

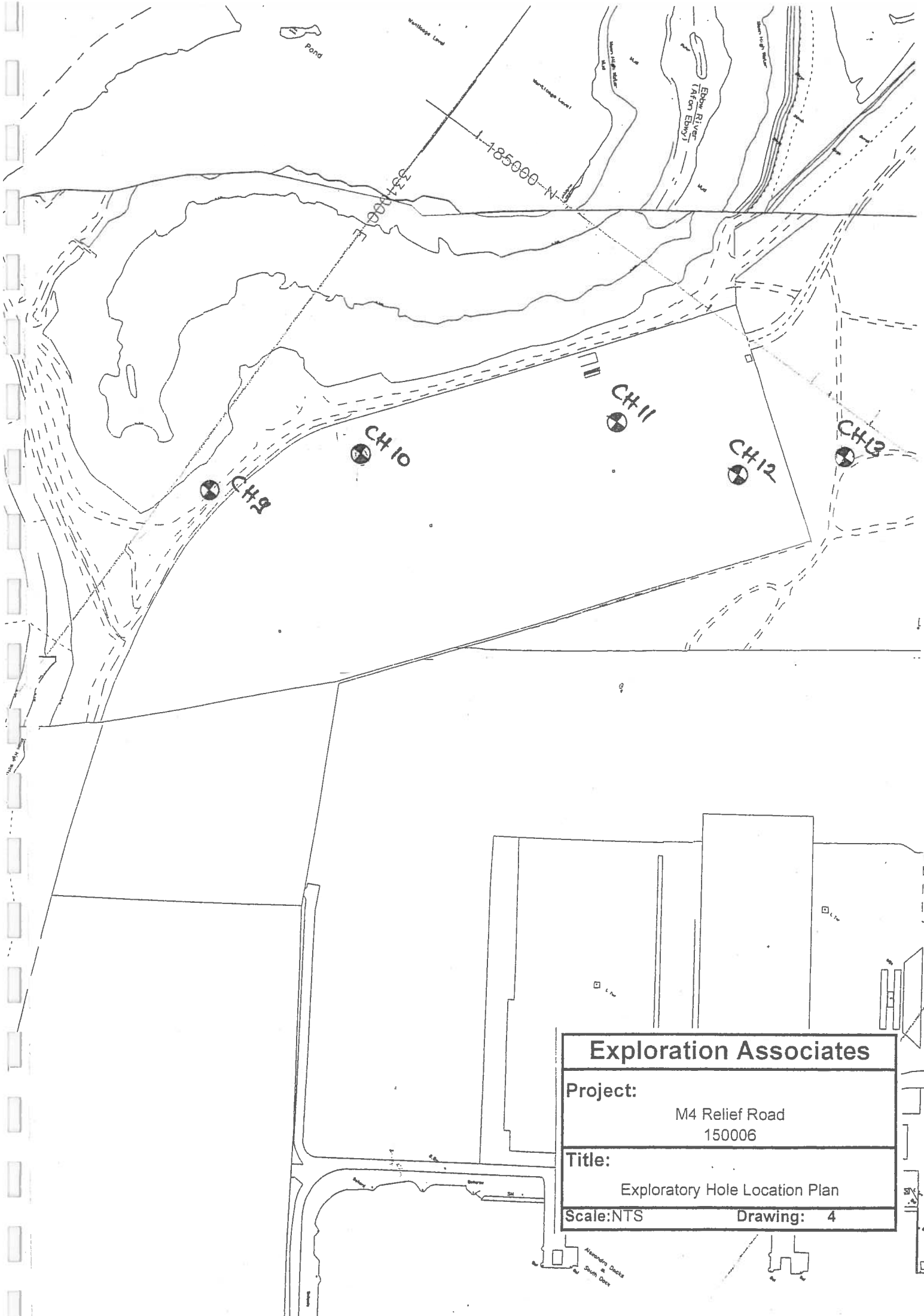
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Drawing: 2





Exploration Associates	
Project:	M4 Relief Road 150006
Title:	Exploratory Hole Location Plan
Scale: NTS	Drawing: 3



Exploration Associates

Project:

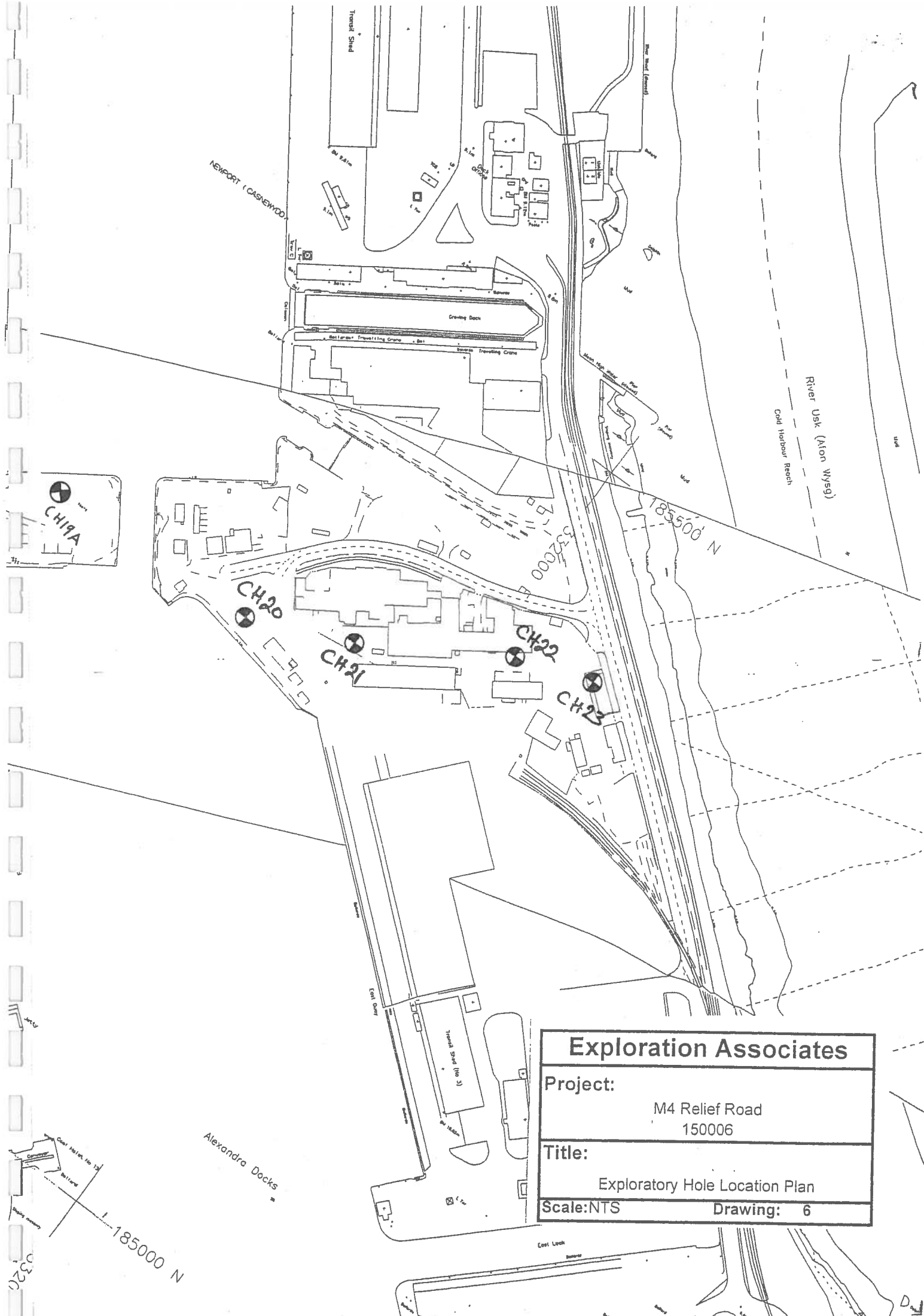
M4 Relief Road
150006

Title:

Exploratory Hole Location Plan

Scale: NTS

Drawing: 4



Exploration Associates	
Project:	M4 Relief Road 150006
Title:	Exploratory Hole Location Plan
Scale:NTS	Drawing: 6

Stevenson Street
Industrial Estate

CH24

CH25

CH26

CH27

CH28

CH29

155500 N

372500 F

River Usk (Afon Wysg)
Cold Harbour Reach

Exploration Associates

Project:

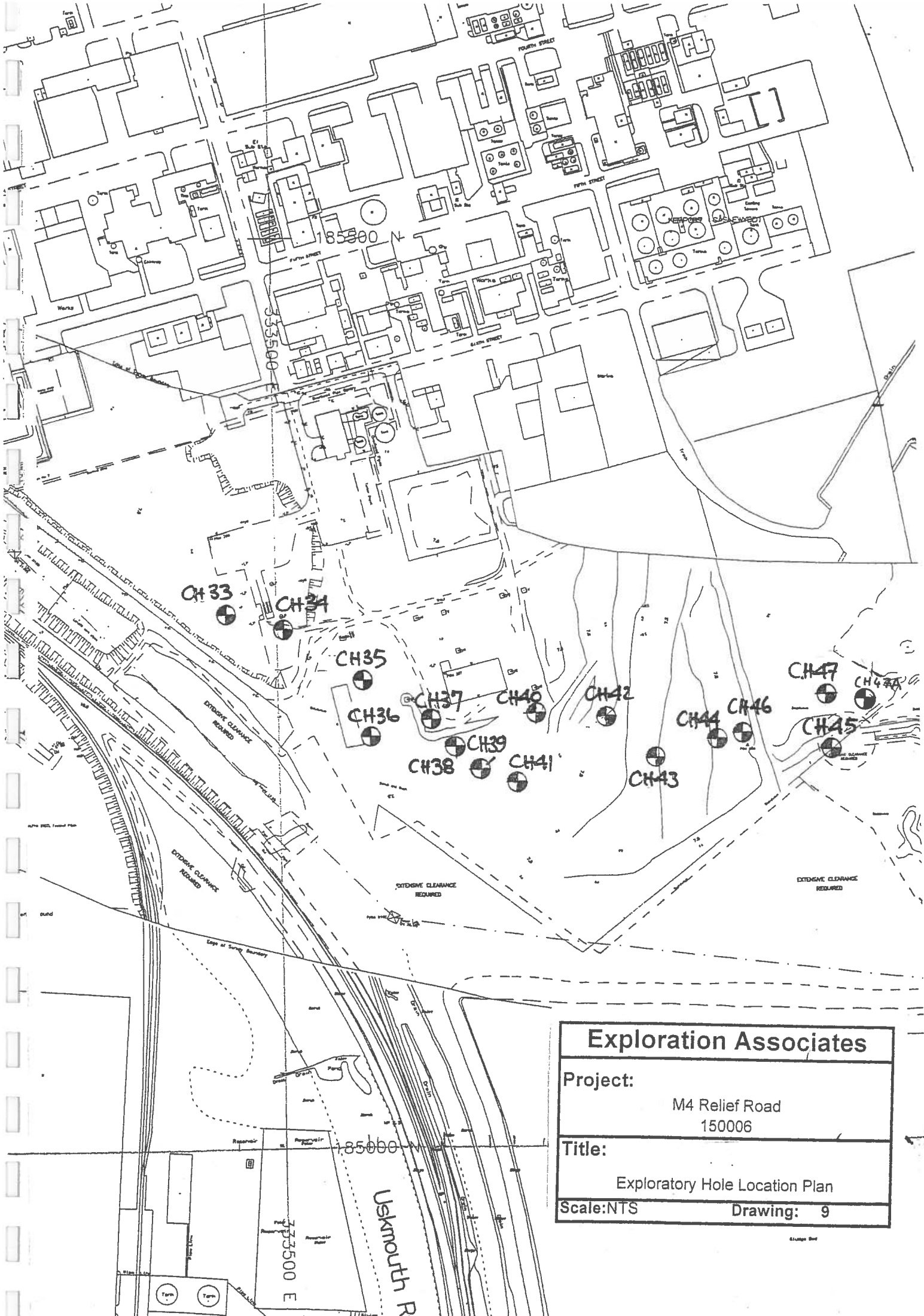
M4 Relief Road
150006

Title:

Exploratory Hole Location Plan

Scale: NTS

Drawing: 7



Exploration Associates

Project:

M4 Relief Road
150006

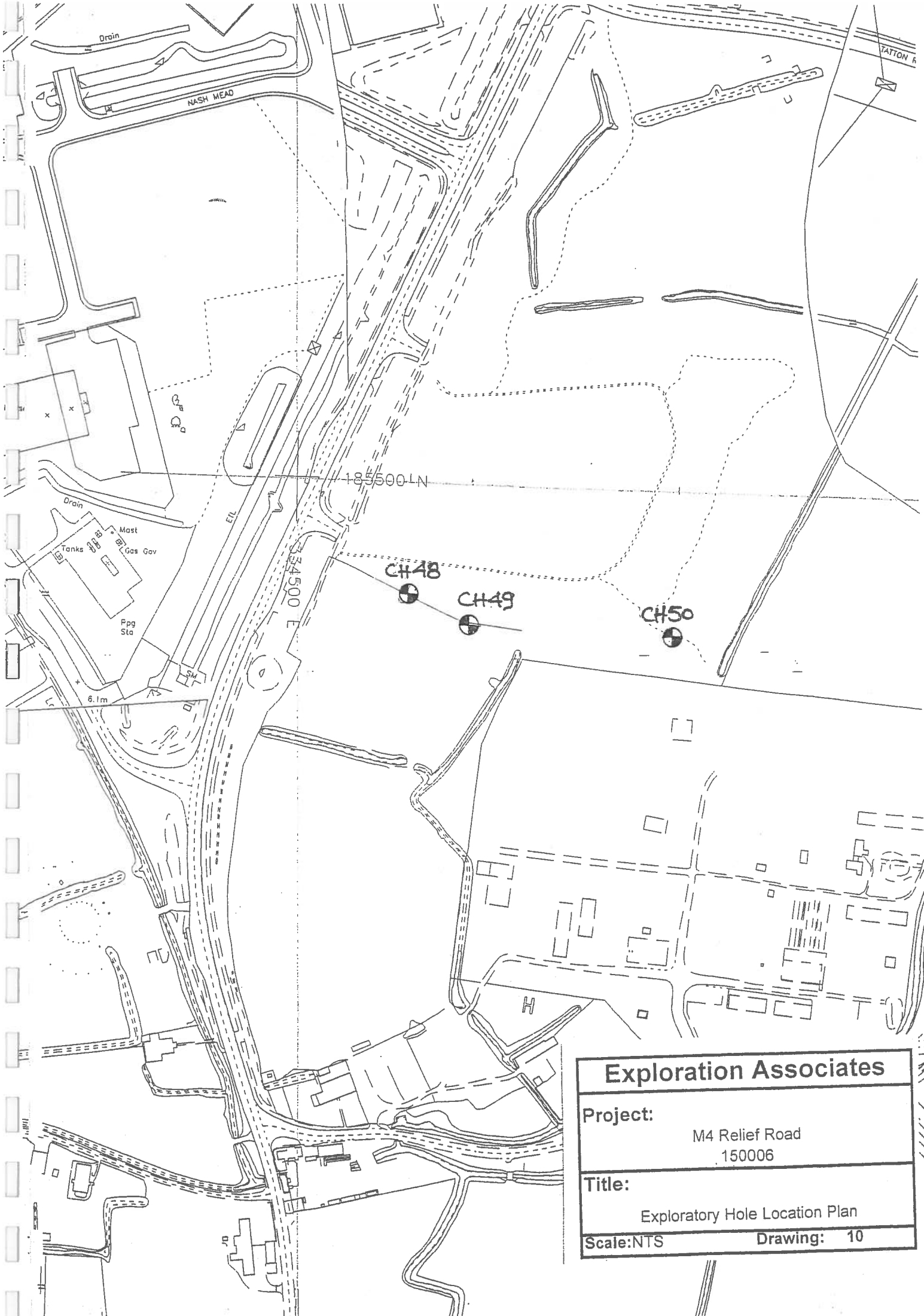
Title:

Exploratory Hole Location Plan

Scale: NTS

Drawing: 9

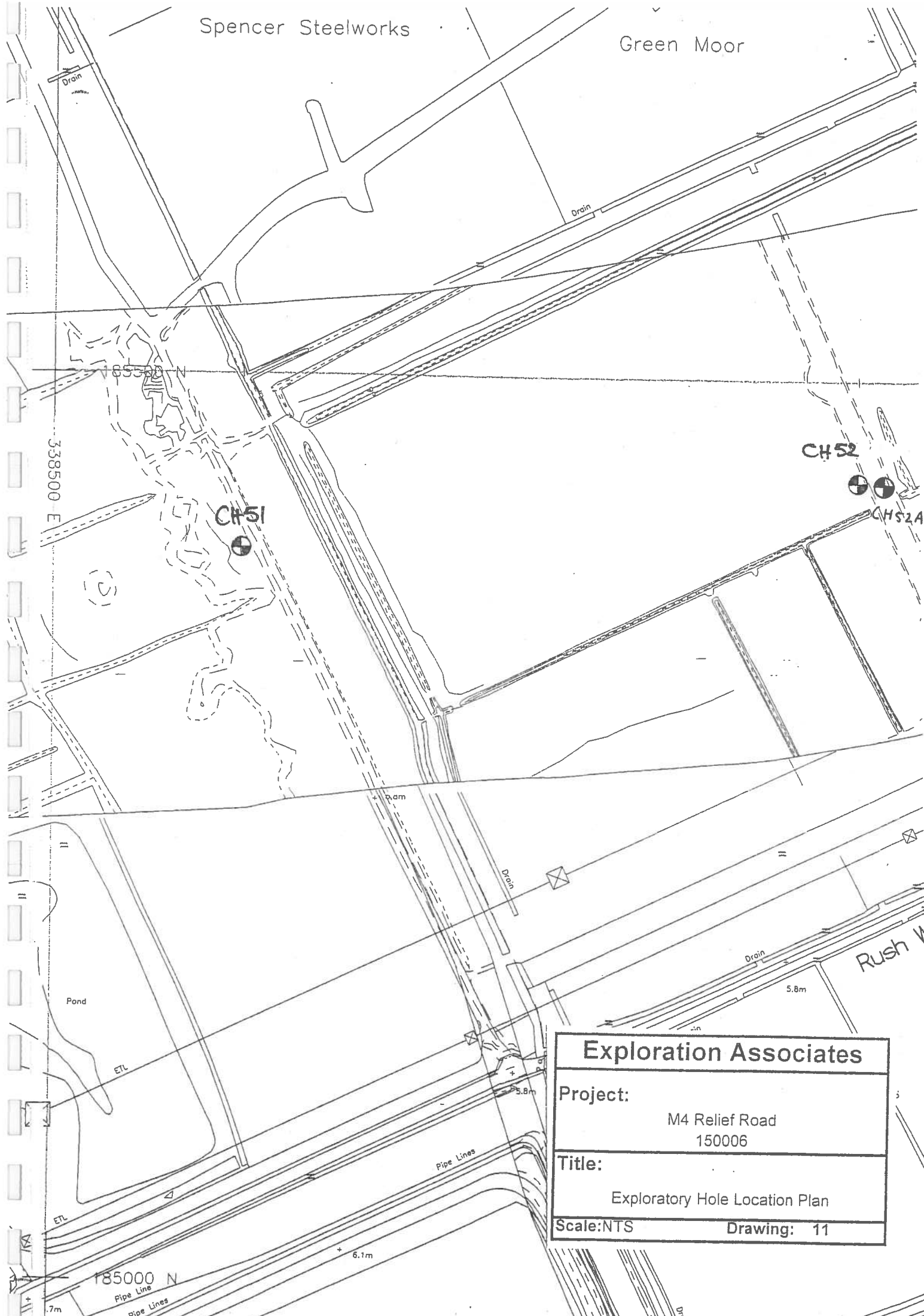
Gludge Bed

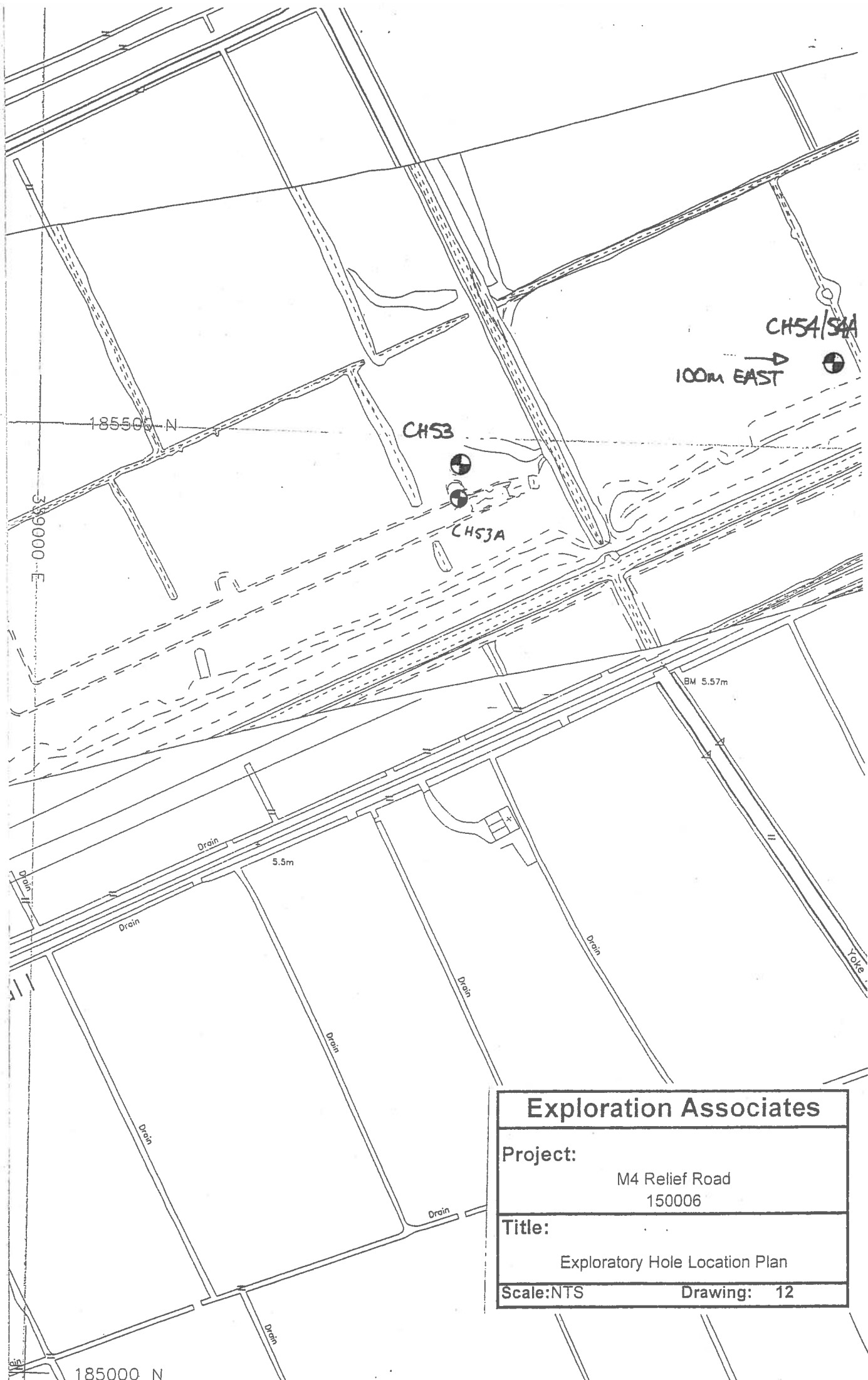


Exploration Associates	
Project:	M4 Relief Road 150006
Title:	Exploratory Hole Location Plan
Scale: NTS	Drawing: 10

Spencer Steelworks

Green Moor





Exploration Associates

Project:

M4 Relief Road
150006

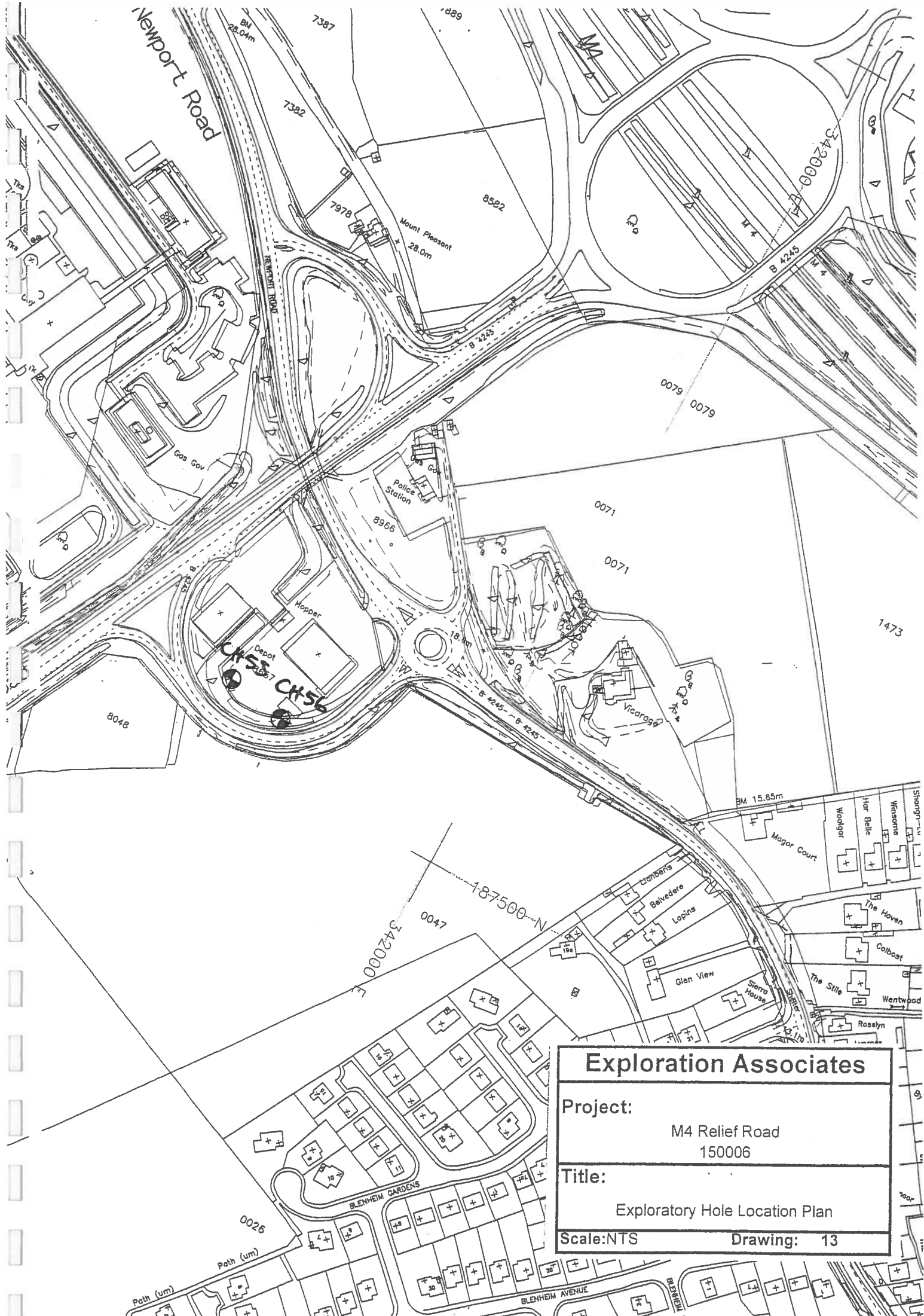
Title:

Exploratory Hole Location Plan

Scale: NTS

Drawing: 12

185000 N



Exploration Associates

Project:

M4 Relief Road
150006

Title:

Exploratory Hole Location Plan

Scale: NTS

Drawing: 13

General Notes

These notes, which accompany the ground investigation report, are intended to assist the user of the information contained in the report. They point out some inevitable shortcomings of any ground investigation and do not constitute a disclaimer of responsibility for the results obtained by Exploration Associates Limited.

1. The information in this report is based on the ground conditions encountered during the ground investigation work and the results of any field and laboratory testing. The exploratory records describe the ground conditions at their specific locations and should not be regarded as representative of the ground as a whole.
2. Ground investigations are performed by the company in general accordance with the recommendations in BS 5930 (1981) "Code of Practice for Site Investigations". The testing of soils, rocks and aggregates generally follow the recommendations of BS 1377 (1990) "Methods of test for soils for Civil Engineering Purposes", the International Society of Rock Mechanics (Brown, 1981) "Rock characterisation, testing and monitoring, ISRM suggested methods", and BS 812 (1975) "Methods of sampling and testing of mineral aggregates, sands and filters", respectively.
3. The primary purpose of ground investigation boreholes and trial pits is to probe the stratified sequences of soil and/or rock. From the results of these probings no conclusions should be drawn concerning the presence of size, lithological nature and numbers per unit volume of ground of cobbles and boulders in soil types such as glacial till (boulder clay).
4. When cable percussion boring techniques are used in superficial and drift deposits some mixing of thin-layered soils inevitably occurs. If strong randomly-occurring pieces of rock are encountered in soil material then the rock may be either pushed aside or penetrated and broken up in which case the arisings that are recovered may not be indicative of the nature of the material in situ.
5. Rotary drilling techniques may sometimes be used for drilling through superficial deposits and rocks in order to provide a very general indication of the nature of the ground. Where open-hole methods have been used for the ground investigation the description of the ground is based on the cuttings recovered from the flushing medium and the rate of progress in advancing the hole. Descriptions of strata and the depths of changes in strata may not be accurate under these conditions.
6. Groundwater conditions noted during boring may be subject to change through seasonal and/or other effects such as, for example, boring and constructional excavation. When a groundwater inflow is encountered during boring, work on the hole is suspended, typically for 20 minutes, and any change in level is recorded. The groundwater level recorded on resumption of boring may not be the natural pre-boring standing water level. When piezometers are installed in boreholes the reported groundwater levels may also be subject to variation due to seasonal and/or other effects.
7. The factual information contained within the ground investigation report should not be used for any purpose other than for the development project for which it was prepared unless a check has been carried out on its applicability. Where the ground investigation report contains an interpretation of the factual information that interpretation must be considered in the context of the stated development proposals and should not be used in any other context.
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