

Our Ref: EXR/168561 (Ver. 1)
Your Ref: LNO 11202/25
January 14, 2014



ESG 
Environmental **Scientifics** Group
Environmental Chemistry
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Mr A Palliser
ESG Stockport
Unit 5
Crown Industrial Estate
Kenwood Rd
Reddish
Stockport
SK5 6BH

UPM, D.S. screen 2013

For the attention of Mr A Palliser

Dear Mr Palliser

Water Sample Analysis - LNO 11202/26

Samples from the above site have been analysed in accordance with the schedule supplied.
The sample details and the results of analyses for these samples are given in the appended report.

An invoice for this work will follow under a separate cover.

Please be aware that our policy for the retention of paper based laboratory records and analysis reports is 6 years.

The work was carried out in accordance with Environmental Scientifics Group Ltd (Laboratory and Analytical) Standard Terms and Conditions of Contract.

If I can be of any further assistance please do not hesitate to contact me.

Yours sincerely

for ESG

L Bosworth

L Bosworth
Project Co-ordinator
01283 554362

MONITORING RETURNS N. WALTON

	INITIALS	DATE
CHECKED Vs AUTHORISATION	JS	26/2/14
TRACKING	JB	26.2.14
OK FOR PUBLIC REGISTER	JS	26/2/14
FILED TO PPA'S RECORDS	JB	EDPM

TEST REPORT

WATER SAMPLE ANALYSIS



Report No. EXR/168561 (Ver. 1)

ESG Stockport
Unit 5
Crown Industrial Estate
Kenwood Rd
Reddish
Stockport
SK5 6BH

Site: LNO 11202/26

The 3 samples described in this report were registered for analysis by ESG on 11-Dec-2013. This report supersedes any versions previously issued by the laboratory.

The analysis was completed by: 14-Jan-2014

Tests where the accreditation is set to N or No, and any individual data items marked with a * are not UKAS accredited
Any opinions or interpretations expressed herein are outside the scope of any UKAS accreditation held by ESG.

The following tables are contained in this report:

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On behalf of
ESG :
Declan Burns

Operations Director
Laboratory and Analytical Business

Date of Issue: 14-Jan-2014

Tests marked '*' have been subcontracted to another laboratory.

ESG accepts no responsibility for any sampling not carried out by our personnel.

Polycyclic Aromatic Hydrocarbons GC/MS (SIM)

Customer and Site Details:	ESG Stockport: LNO 11202/26		
Sample Details:	HSM 4885 D W	Job Number:	W16_8561
LIMS ID Number:	EX1448492	Date Booked in:	11-Dec-13
QC Batch Number:	130893	Date Extracted:	17-Dec-13
Quantitation File:	Initial Calibration	Date Analysed:	18-Dec-13
Directory:	1713PAHMS14\	Matrix:	Water
Dilution:	1.0	Ext Method:	Bottle

UKAS accredited?: Yes

Target Compounds	CAS #	R.T. (min)	Concentration ug/l	% Fit
Naphthalene	91-20-3	-	< 0.020	-
Acenaphthylene	208-96-8	-	< 0.010	-
Acenaphthene	83-32-9	-	< 0.010	-
Fluorene	86-73-7	-	< 0.010	-
Phenanthrene	85-01-8	-	< 0.010	-
Anthracene	120-12-7	-	< 0.010	-
Fluoranthene	206-44-0	-	< 0.010	-
Pyrene	129-00-0	-	< 0.010	-
Benzo[a]anthracene	56-55-3	-	< 0.010	-
Chrysene	218-01-9	-	< 0.010	-
Benzo[b]fluoranthene	205-99-2	-	< 0.010	-
Benzo[k]fluoranthene	207-08-9	-	< 0.010	-
Benzo[a]pyrene	50-32-8	-	< 0.010	-
Indeno[1,2,3-cd]pyrene	193-39-5	-	< 0.010	-
Dibenzo[a,h]anthracene	53-70-3	-	< 0.010	-
Benzo[g,h,i]perylene	191-24-2	-	< 0.010	-
Total (USEPA16) PAHs	-	-	< 0.170	-

"M" denotes that % fit has been manually interpreted

Internal Standards	% Area
1,4-Dichlorobenzene-d4	NA
Naphthalene-d8	98
Acenaphthene-d10	97
Phenanthrene-d10	98
Chrysene-d12	96
Perylene-d12	92

Surrogates	% Rec
Nitrobenzene-d5	NA
2-Fluorobiphenyl	99
Terphenyl-d14	76

The Total PAH result is the sum of non-rounded individual PAH results and therefore may differ to the sum of the rounded individual PAH results printed above. By convention, where any one or more result is a "less than", the total is expressed as a "less than" and includes the "less than" concentration within the total.

Semi-Volatile Organic Compounds

Customer and Site Details: ESG Stockport: LNO 11202/26
Sample Details: HSM 4885 D W
LIMS ID Number: EX1448492
Job Number: W16_8561

UKAS accredited?: No

Date Booked in: 11-Dec-13
Date Extracted: 18-Dec-13
Date Analysed: 19-Dec-13

Matrix: Water
Ext Method: Sep. Funnel
Operator: JO
Directory/Quant File: 1219_GCC1.DGPC (Y/N)

QC Batch Number: 285

Multiplier: 0.005

Dilution Factor: 1

N

Target Compounds	CAS #	R.T. (min)	Concentration mg/l	% Fit
Phenol	108-95-2	-	< 0.020	-
bis(2-Chloroethyl)ether	111-44-4	-	< 0.005	-
2-Chlorophenol	95-57-8	-	< 0.020	-
1,3-Dichlorobenzene	541-73-1	-	< 0.005	-
1,4-Dichlorobenzene	106-46-7	-	< 0.005	-
Benzyl alcohol	100-51-6	-	< 0.005	-
1,2-Dichlorobenzene	95-50-1	-	< 0.005	-
2-Methylphenol	95-48-7	-	< 0.005	-
bis(2-Chloroisopropyl)ether	108-60-1	-	< 0.005	-
Hexachloroethane	67-72-1	-	< 0.005	-
N-Nitroso-d1-n-propylamine	621-64-7	-	< 0.005	-
3- & 4-Methylphenol	108-39-4/106-44-5	-	< 0.020	-
Nitrobenzene	98-95-3	-	< 0.005	-
Isophorone	78-59-1	-	< 0.005	-
2-Nitrophenol	88-75-5	-	< 0.020	-
2,4-Dimethylphenol	105-67-9	-	< 0.020	-
Benzoic Acid	65-85-0	-	< 0.100	-
bis(2-Chloroethoxy)methane	111-91-1	-	< 0.005	-
2,4-Dichlorophenol	120-83-2	-	< 0.020	-
1,2,4-Trichlorobenzene	120-82-1	-	< 0.005	-
Naphthalene	91-20-3	-	< 0.002	-
4-Chlorophenol	106-48-9	-	< 0.020	-
4-Chloroaniline	106-47-8	-	< 0.005	-
Hexachlorobutadiene	87-68-3	-	< 0.005	-
4-Chloro-3-methylphenol	59-50-7	-	< 0.005	-
2-Methylnaphthalene	91-57-6	-	< 0.002	-
1-Methylnaphthalene	90-12-0	-	< 0.002	-
Hexachlorocyclopentadiene	77-47-4	-	< 0.005	-
2,4,6-Trichlorophenol	88-06-2	-	< 0.020	-
2,4,5-Trichlorophenol	95-95-4	-	< 0.020	-
2-Chloronaphthalene	91-58-7	-	< 0.002	-
Biphenyl	92-52-4	-	< 0.002	-
Diphenyl ether	101-84-8	-	< 0.002	-
2-Nitroaniline	88-74-4	-	< 0.005	-
Acenaphthylene	208-96-8	-	< 0.002	-
Dimethylphthalate	131-11-3	-	< 0.005	-
2,6-Dinitrotoluene	606-20-2	-	< 0.005	-
Acenaphthene	83-32-9	-	< 0.002	-
3-Nitroaniline	99-09-2	-	< 0.005	-

Concentrations are reported on a wet weight basis.

"M" denotes that % fit has been manually interpreted

Target Compounds	CAS #	R.T.	Concentration mg/l	% Fit
2,4-Dinitrophenol	51-28-5	-	< 0.010	-
Dibenzofuran	132-64-9	-	< 0.005	-
4-Nitrophenol	100-02-7	-	< 0.050	-
2,4-Dinitrotoluene	121-14-2	-	< 0.005	-
Fluorene	86-73-7	-	< 0.002	-
Diethylphthalate	84-66-2	-	< 0.005	-
4-Chlorophenyl-phenylether	7005-72-3	-	< 0.005	-
4,6-Dinitro-2-methylphenol	534-52-1	-	< 0.050	-
4-Nitroaniline	100-01-6	-	< 0.005	-
N-Nitrosodiphenylamine	86-30-6	-	< 0.005	-
4-Bromophenyl-phenylether	101-55-3	-	< 0.005	-
Hexachlorobenzene	118-74-1	-	< 0.005	-
Pentachlorophenol	87-86-5	-	< 0.050	-
Phenanthrene	85-01-8	-	< 0.002	-
Anthracene	120-12-7	-	< 0.002	-
Di-n-butylphthalate	84-74-2	-	< 0.005	-
Fluoranthene	206-44-0	-	< 0.002	-
Pyrene	129-00-0	-	< 0.002	-
Butylbenzylphthalate	85-68-7	-	< 0.005	-
Benzo[a]anthracene	56-55-3	-	< 0.002	-
Chrysene	218-01-9	-	< 0.002	-
3,3'-Dichlorobenzidine	91-94-1	-	< 0.020	-
bis(2-Ethylhexyl)phthalate	117-81-7	-	< 0.005	-
Di-n-octylphthalate	117-84-0	-	< 0.002	-
Benzo[b]fluoranthene	205-99-2	-	< 0.002	-
Benzo[k]fluoranthene	207-08-9	-	< 0.002	-
Benzo[a]pyrene	50-32-8	-	< 0.002	-
Indenol 1,2,3-cd]pyrene	193-39-5	-	< 0.002	-
Dibenzol[a,h]anthracene	53-70-3	-	< 0.002	-
Benzo[g,h,i]perylene	191-24-2	-	< 0.002	-

"M" denotes that % fit has been manually interpreted

Internal Standards	% Area
1,4-Dichlorobenzene-d4	48
Naphthalene-d8	45
Acenaphthene-d10	44
Phenanthrene-d10	44
Chrysene-d12	40
Perylene-d12	41

Surrogates	% Rec
2-Fluorophenol	57
Phenol-d5	36
Nitrobenzene-d5	104
2-Fluorobiphenyl	99
2,4,6-Tribromophenol	106
Terphenyl-d14	111

Volatile Organic Compounds by HSA-GCMS

UKAS accredited?: Yes

Customer and Site Details: ESG Stockport: LNO 11202/26

Sample Details: HSM 4885 D W

LIMS ID Number: EX1448492

Job Number: W16_8561

Directory/Quant file:

Date Booked in: 217VOC.MS11\

Date Analysed: 11-Dec-13

Operator: 18-Dec-13

Initial Calibration

Matrix:

Method:

Multiplier:

Position:

Water

Headspace

1

26

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 *	-	< 5	-
Chloroethane	75-00-3	-	< 5	-
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 *	-	< 5	-
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	-	< 5	-
1,3-Dichloropropane	142-28-9	-	< 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	-	< 5	-
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	-

Compounds marked * are not UKAS accredited
"M" denotes that % fit has been manually interpreted

Internal standards	R.T.	Area %	Surrogates	% Rec
Pentafluorobenzene	3.66	108	Dibromofluoromethane	112
1,4-Difluorobenzene	4.01	109	Toluene-d8	98
Chlorobenzene-d5	5.14	111	Bromofluorobenzene	83
1,4-Dichlorobenzene-d4	5.94	92		

Note: Volatile compounds degrade with time, and this may affect the integrity of the data depending on the timescale between sampling and analysis. It is recommended that analysis takes place within 14 days of sampling.

Where individual results are flagged see report notes for status.

Organochlorine Pesticides by GCMS (SIM)

Customer and Site Details:	ESG Stockport: LNO 11202/26		
Sample Details:	HSM 4885 D W	Job Number:	W16_8561
LIMS ID Number:	EX1448492	Date Booked in:	11-Dec-13
QC Batch Number:	130210	Date Extracted:	16-Dec-13
Quantitation File:	12160CPCCC2.D	Date Analysed:	17-Dec-13
Directory:	613PEST.MS9\	Matrix:	Water
Dilution:	1	Ext Method:	Sep. Funnel

UKAS accredited?: No

Target Compounds	CAS #	R.T. (min)	Concentration ug/l	% Fit
1,3,5-Trichlorobenzene	108-70-3	-	< 0.02	-
1,2,3-Trichlorobenzene	87-61-6	-	< 0.01	-
2,6-Dichlorobenzonitrile	1194-65-6	-	< 0.02	-
1,2,3,4-Tetrachlorobenzene	634-66-2	-	< 0.02	-
Pentachlorobenzene	608-93-5	-	< 0.02	-
Tecnazene	117-18-0	-	< 0.01	-
Trifluralin	1582-09-8	-	< 0.02	-
Alpha-HCH	319-84-6	-	< 0.10	-
Hexachlorobenzene	118-74-1	-	< 0.02	-
Beta-HCH	319-85-7	-	< 0.02	-
Gamma-HCH	58-89-9	-	< 0.02	-
Propyzamide	23950-58-5	-	< 0.02	-
Chlorthalonil	1897-45-6	-	< 0.01	-
Triallate	2303-17-5	-	< 0.02	-
Lindane	319-86-8	-	< 0.01	-
Heptachlor	76-44-8	-	< 0.01	-
Aldrin	309-00-2	-	< 0.02	-
Triadimefon	43121-43-3	-	< 0.02	-
Pendimethalin	40487-42-1	-	< 0.02	-
Heptachlorepoxyde	1024-57-3	-	< 0.02	-
Trans-Chlordane	5103-74-2	-	< 0.01	-
Isodrin	465-73-6	-	< 0.02	-
O,P'-DDE	3424-82-6	-	< 0.02	-
Cis-Chlordane	5103-71-9	-	< 0.02	-
Endosulfan I	959-98-8	-	< 0.01	-
P,P'-DDE	72-55-9	-	< 0.01	-
Dieldrin	60-57-1D	-	< 0.01	-
O,P'-DDD	53-19-0	-	< 0.01	-
Endrin	72-20-8	-	< 0.01	-
Endosulfan II	33213-65-9	-	< 0.01	-
P,P'-DDD	72-54-8	-	< 0.01	-
O,P'-DDT	789-02-6	-	< 0.01	-
Endosulfan Sulfate	1031-07-8	-	< 0.01	-
P,P'-DDT	50-29-3	-	< 0.02	-
Endrin Ketone	53494-70-5	-	< 0.02	-
Methoxychlor	72-43-5	-	< 0.01	-
Cis-Permethrin	52645-53-1	-	< 0.01	-
Trans-Permethrin	51877-74-8	-	< 0.01	-

Internal Standards	% Area
Naphthalene-d8	85
Phenanthrene-d10	86
Perylene-d12	81

Surrogates	% Rec
Gamma-HCH-d6	D
P,P'-DDT-d8	78

Where individual results are flagged see report notes for status.

Analytical Report

ESG Environmental Chemistry
P O Box 100
Bretby Business Park
Burton-on-Trent
Staffordshire, DE15 0XD

Report No: 13-34952/1
Date Received: 13/12/2013
Date Tested: 03/01/2014 to 14/01/2014
Date Issued: 14/01/2014
Page: 1 of 1
By email

For the attention of: Louise Smith

1 water sample received from ESG Environmental Chemistry (O/N: 42062 BEC; Project: w168561) in a 1 litre amber glass bottle was analysed as shown below. Analytical methods employed are available on request.

Laboratory reference			237778 W1448492 HSM 4885
chloroalkanes (C10-C13)	[85535-84-8]	ug/l	< 10.0
BDE-3	[101-55-3]	ug/l	< 0.10
4-n-nonylphenol	[104-40-5]	ug/l	< 0.02
nonylphenols	[25154-52-3]	ug/l	< 0.02


Marco Lattughi
Senior Operations Manager

Customer	ESG Stockport
Site	LNO 11202/26
Report No	W168561

Consignment No W63418
Date Logged 11-Dec-2013

Report Due 24-Dec-2013

	VOC HSA-GCMS	>						
SVOCSW	SVOC							
	^Chloroalkanes							
	^ polybrominated diphenyl ethers (PBDEs)							
Sub005	^ phenols (alkylphenols test 20)							
PAHMSW	PAH GC-MS (16)	>						
	Triphenyl Tin							
	Tributyl Tin							
OGSN	Dibutyl Tin							
OCP	Organochlorine Pesticides							
ICPWATVAR	Phosphorus as P (Total) VAR							
CUSTSERV	Report B							
MethodID	Sampled							
ID Number	Description							
EX/1448492	HSM 4885							
EX/1448493	HSM 4886							
EX/1448494	HSM 4887							

Note: For analysis where the scheduled turnaround is greater than the holding time we will do our utmost to prioritise these samples. However, it is possible that samples could become deviant whilst being processed in the laboratory.

In this instance please contact the laboratory immediately should you wish to discuss how you would like us to proceed. If you do not respond within 24 hours, we will proceed as originally requested.

Deviating Sample Key

- A** The sample was received in an inappropriate container for this analysis
B The sample was received without the correct preservation for this analysis
C Headspace present in the sample container
D The sampling date was not supplied so holding time may be compromised - applicable to all analysis
E Sample processing did not commence within the appropriate holding time

Requested Analysis Key

- | | |
|---|---|
| | Analysis Required |
| | Analysis dependant upon trigger result - Note: due date may be affected if triggered |
| | No analysis scheduled |
| A | Analysis Subcontracted - Note: due date may vary |

Where individual results are flagged see report notes for status.

Method Descriptions

Matrix	MethodID	Analysis Basis	Method Description
Water	ICPWATVAR	As Received	Direct determination of Metals and Sulphate in water samples using ICPOES
Water	OCP	As Received	Determination of Organo Chlorine Pesticides by dichloromethane/acetone extraction and GCMS detection
Water	OGSN	As Received	Hexane extraction, sodium tetraethylborate derivitisation and GC-MS quantitation
Water	PAHMSW	As Received	Determination of PolyAromatic Hydrocarbons in water by pentane extraction GCMS quantitation
Water	SubCon*	*	Contact Laboratory for details of the methodology used by the sub-contractor.
Water	SVOCSW	As Received	Determination of Semi Volatile Organic Compounds (SVOC) by DCM extraction followed by GCMS detection
Water	VOCHSAW	As Received	Determination of Volatile Organics Compounds by Headspace GCMS

Report Notes

Generic Notes

Soil/Solid Analysis

Unless stated otherwise,

- Results expressed as mg/kg have been calculated on the basis indicated in the Method Description table.
- All results on MCERTS reports are reported on a 105°C dry weight basis with the exception of pH and conductivity.
- Sulphate analysis not conducted in accordance with BS1377
- Water Soluble Sulphate is on a 2:1 water:soil extract

Waters Analysis

Unless stated otherwise results are expressed as mg/l

Nil: Where "Nil" has been entered against Total Alkalinity or Total Acidity this indicates that a measurement was not required due to the inherent pH of the sample.

Oil analysis specific

Unless stated otherwise,

- Results are expressed as mg/kg
- SG is expressed as g/cm³ @ 15°C

Gas (Tedlar bag) Analysis

Unless stated otherwise, results are expressed as ug/l

Asbestos Analysis

CH Denotes Chrysotile

CR Denotes Crocidolite

AM Denotes Amosite

NAIS No Asbestos Identified in Sample

NADIS No Asbestos Detected In Sample

Symbol Reference

^ Sub-contracted analysis.

\$\$ Unable to analyse due to the nature of the sample

¶ Samples submitted for this analyte were not preserved on site in accordance with laboratory protocols.

This may have resulted in deterioration of the sample(s) during transit to the laboratory.

Consequently the reported data may not represent the concentration of the target analyte present in the sample at the time of sampling

¥ Results for guidance only due to possible interference

& Blank corrected result

I.S Insufficient sample to complete requested analysis

I.S(g) Insufficient sample to re-analyse, results for guidance only

Intf Unable to analyse due to interferences

N.D Not determined

N.Det Not detected

NS Information Not Supplied

Req Analysis requested, see attached sheets for results

P Raised detection limit due to nature of the sample

***** All accreditation has been removed by the laboratory for this result

‡ MCERTS accreditation has been removed for this result

Note: The Laboratory may only claim that data is accredited when all of the requirements of our Quality System have been met. Where these requirements have not been met the laboratory may elect to include the data in its final report and remove the accreditation from individual data items if it believes that the validity of the data has not been affected. If further details are required of the circumstances which have led to the removal of accreditation then please do not hesitate to contact the laboratory.

END OF REPORT

Where individual results are flagged see report notes for status.

