

Data Path : Z:\HPCHEM1\BNA E\DATA\BE051518\
 Data File : BE096581.D
 Acq On : 15 May 2018 16:58
 Operator : SJ/JU
 Sample : J2604-13DL 100X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 PT-BNA-SoilDL

Quant Time: May 16 05:59:38 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE051518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 15 13:02:03 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.34	152	653	0.40	ng	0.00
7) Naphthalene-d8	11.17	136	2764	0.40	ng	-0.01
13) Acenaphthene-d10	14.92	164	1693	0.40	ng	0.00
19) Phenanthrene-d10	17.64	188	3854	0.40	ng	0.00
27) Chrysene-d12	21.73	240	3706	0.40	ng	0.00
34) Perylene-d12	24.33	264	3542	0.40	ng	0.00

System Monitoring Compounds						
4) 2-Fluorophenol	5.84	112	2525	1.40	ng	-0.01
5) Phenol-d6	7.49	99	3733	1.46	ng	-0.01
8) Nitrobenzene-d5	9.52	82	2673	0.91	ng	-0.01
11) 2-Methylnaphthalene-d10	12.71	152	5	0.00	ng	-0.01
14) 2,4,6-Tribromophenol	16.40	330	1061	1.59	ng	0.00
15) 2-Fluorobiphenyl	13.56	172	6424	0.88	ng	0.00
25) Fluoranthene-d10	19.62	212	32	0.00	ng	0.00
29) Terphenyl-d14	20.18	244	8300	0.99	ng	0.00

Target Compounds				Qvalue		
10) Hexachlorobutadiene	11.48	225	2863	1.761	ng	94
12) 2-Methylnaphthalene	12.79	142	3605	0.723	ng	93
16) Acenaphthylene	14.64	152	14100	1.596	ng	99
17) Acenaphthene	14.99	154	6529	1.212	ng	93
18) Fluorene	15.96	166	11538	1.722	ng	# 77
20) 4-Bromophenyl-phenylether	16.82	248	4784	2.046	ng	# 83
21) Hexachlorobenzene	16.95	284	4058	1.665	ng	# 81
22) Pentachlorophenol	17.30	266	2010	2.744	ng	# 74
23) Phenanthrene	17.67	178	16593	1.533	ng	99
26) Fluoranthene	19.65	202	15608	1.206	ng	# 97
30) Benzo(a)anthracene	21.72	228	11618	1.020	ng	97
32) Bis(2-ethylhexyl)phthalate	21.58	149	51826	2.125	ng	100
33) Indeno(1,2,3-cd)pyrene	27.10	276	11470	1.009	ng	# 93
36) Benzo(k)fluoranthene	23.58	252	11147	0.964	ng	# 94
37) Benzo(a)pyrene	24.21	252	13845	1.326	ng	# 92
38) Dibenzo(a,h)anthracene	27.11	278	10526	1.102	ng	# 95
39) Benzo(a,h,i)perylene	27.96	276	11528	1.153	ng	# 89

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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