

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM101718\
 Data File : BM017184.D
 Acq On : 18 Oct 2018 09:46
 Operator : MJ/SJ
 Sample : J5164-01
 Misc : LOD 2 PPM
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 LOD-MDL-SOIL-01-QT4-2018

Quant Time: Oct 18 14:40:29 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM101718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 17 14:23:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	163612	20.00	ng	0.00
21) Naphthalene-d8	10.45	136	686365	20.00	ng	0.00
39) Acenaphthene-d10	14.31	164	391138	20.00	ng	0.00
64) Phenanthrene-d10	17.06	188	915113	20.00	ng	-0.01
76) Chrysene-d12	21.25	240	1062635	20.00	ng	-0.01
87) Perylene-d12	23.46	264	1082718	20.00	ng	-0.02

System Monitoring Compounds						
5) 2-Fluorophenol	5.28	112	1540033	159.20	ng	0.00
7) Phenol-d6	6.85	99	2010217	152.58	ng	0.00
23) Nitrobenzene-d5	8.82	82	1195733	101.89	ng	-0.01
42) 2,4,6-Tribromophenol	15.80	330	884038	167.13	ng	-0.01
45) 2-Fluorobiphenyl	12.93	172	2863033	97.39	ng	0.00
79) Terphenyl-d14	19.70	244	4559035	96.70	ng	0.00

Target Compounds				Qvalue		
17) 2-Methylphenol	8.12	107	18289	2.010	ng	92
19) n-Nitroso-di-n-propylamine	8.47	70	16844	1.942	ng	# 96
20) 3+4-Methylphenols	8.44	107	24876	1.998	ng	95
22) Acetophenone	8.49	105	36074	2.073	ng	# 96
25) Isophorone	9.39	82	43899	2.265	ng	97
26) 2-Nitrophenol	9.57	139	9119	7.086	ng	95
27) 2,4-Dimethylphenol	9.63	122	20207	2.358	ng	97
28) bis(2-Chloroethoxy)methane	9.87	93	29119	2.135	ng	98
29) 2,4-Dichlorophenol	10.10	162	18001	1.809	ng	99
30) 1,2,4-Trichlorobenzene	10.31	180	25948	2.174	ng	96
36) 4-Chloro-3-methylphenol	11.75	107	19865	1.771	ng	99
43) 2,4,6-Trichlorophenol	12.73	196	14289	1.787	ng	94
44) 2,4,5-Trichlorophenol	12.80	196	15042	1.754	ng	95
57) 2,4-Dinitrotoluene	14.69	165	12820	1.436	ng	# 91
61) 4-Chlorophenyl-phenylether	15.36	204	35529	2.148	ng	92
62) 4-Nitroaniline	15.40	138	9215	5.789	ng	# 84
77) Benzidine	19.32	184	23837	0.885	ng	95
82) 3,3'-Dichlorobenzidine	21.18	252	28940	1.299	ng	# 99

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

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