

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050818\
 Data File : BF105162.D
 Acq On : 8 May 2018 23:25
 Operator : JU/SJ
 Sample : J2133-07
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LOD-MDL-WATER-01-QT2-2018

Quant Time: May 09 09:22:22 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 08 16:33:54 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	201888	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	812306	20.00	ng	0.00
38) Acenaphthene-d10	9.83	164	400138	20.00	ng	0.00
63) Phenanthrene-d10	11.31	188	767686	20.00	ng	0.00
75) Chrysene-d12	14.01	240	668922	20.00	ng	0.00
86) Perylene-d12	15.56	264	595678	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.40	112	1569446	128.07	ng	0.00
7) Phenol-d6	6.42	99	1871420	128.25	ng	0.00
23) Nitrobenzene-d5	7.36	82	1090185	90.96	ng	0.00
41) 2,4,6-Tribromophenol	10.62	330	601881	136.34	ng	0.00
44) 2-Fluorobiphenyl	9.15	172	1980096	76.81	ng	0.00
78) Terphenyl-d14	12.91	244	2219490	77.07	ng	0.00

Target Compounds				Qvalue		
17) 2-Methylphenol	7.03	107	24777	2.207	ng	98
19) n-Nitroso-di-n-propylamine	7.20	70	20623	2.135	ng	91
20) 3+4-Methylphenols	7.19	107	32131	2.436	ng	92
22) Acetophenone	7.19	105	46198	2.440	ng	# 91
25) Isophorone	7.60	82	51935	1.938	ng	97
26) 2-Nitrophenol	7.69	139	9142	11.260	ng	99
27) 2,4-Dimethylphenol	7.72	122	21102	1.772	ng	96
28) bis(2-Chloroethoxy)methane	7.82	93	34106	2.062	ng	95
29) 2,4-Dichlorophenol	7.92	162	21453	1.925	ng	88
30) 1,2,4-Trichlorobenzene	8.01	180	30006	2.335	ng	98
36) 4-Chloro-3-methylphenol	8.61	107	21545	1.857	ng	90
42) 2,4,6-Trichlorophenol	9.06	196	14360	1.848	ng	96
43) 2,4,5-Trichlorophenol	9.09	196	16263	2.003	ng	93
56) 2,4-Dinitrotoluene	10.00	165	12327	5.067	ng	# 96
60) 4-Chlorophenyl-phenylether	10.37	204	31908	2.634	ng	88
61) 4-Nitroaniline	10.37	138	10339	1.678	ng	100
76) Benzidine	12.65	184	11754	0.471	ng	# 90
81) 3,3'-Dichlorobenzidine	13.96	252	19132	1.254	ng	# 90

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

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