

Data Path : U:\HPCHEM1\BNA_F\DATA\BF011818\
 Data File : BF102209.D
 Acq On : 19 Jan 2018 00:22
 Operator : SJ/JU
 Sample : J1187-05
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :

Quant Time: Jan 19 02:46:33 2018
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF010918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 17 15:28:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.72	152	150957	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	582897	20.00	ng	0.00
38) Acenaphthene-d10	9.76	164	223457	20.00	ng	0.00
63) Phenanthrene-d10	11.24	188	327596	20.00	ng	0.00
75) Chrysene-d12	13.87	240	286792	20.00	ng	0.00
86) Perylene-d12	15.29	264	252762	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.34	112	1175618	109.97	ng	0.02
7) Phenol-d6	6.36	99	1348126	105.70	ng	0.00
23) Nitrobenzene-d5	7.29	82	838605	84.54	ng	0.00
41) 2,4,6-Tribromophenol	10.55	330	222512	114.10	ng	0.00
44) 2-Fluorobiphenyl	9.08	172	1264021	88.37	ng	0.00
78) Terphenyl-d14	12.82	244	849686	61.51	ng	0.00

Target Compounds

						Qvalue
9) Benzaldehyde	6.49	77	25027	2.826	ng	81
10) Phenol	6.37	94	61447	4.263	ng	97
19) n-Nitroso-di-n-propylamine	7.29	70	111633	13.355	ng	# 80
25) Isophorone	7.29	82	831470	42.423	ng	# 64
49) Dimethylphthalate	9.47	163	209487	11.686	ng	99
56) 2,4-Dinitrotoluene	9.76	165	28012	6.242	ng	# 23
64) 4,6-Dinitro-2-methylphenol	10.55	198	564	5.942	ng	# 1
66) 4-Bromophenyl-phenylether	10.55	248	14570	4.212	ng	# 1

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

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