

Data Path : U:\HPCHEM1\BNA F\DATA\BF011718\
 Data File : BF102177.D
 Acq On : 18 Jan 2018 4:36
 Operator : SJ/JU
 Sample : J1144-04
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 IGW-4

Quant Time: Jan 18 05:04:56 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	160586	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	617738	20.00	ng	0.00
38) Acenaphthene-d10	9.76	164	256598	20.00	ng	0.00
63) Phenanthrene-d10	11.24	188	370781	20.00	ng	0.00
75) Chrysene-d12	13.86	240	271791	20.00	ng	-0.01
86) Perylene-d12	15.28	264	274707	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	1086359	95.53	ng	0.01
7) Phenol-d6	6.36	99	1286377	94.81	ng	0.00
23) Nitrobenzene-d5	7.29	82	787103	74.88	ng	0.00
41) 2,4,6-Tribromophenol	10.55	330	230785	103.06	ng	0.00
44) 2-Fluorobiphenyl	9.08	172	1255824	76.46	ng	0.00
78) Terphenyl-d14	12.82	244	795522	60.77	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.37	94	56879	3.709	ng	96
31) Naphthalene	8.03	128	671205	21.745	ng	99
37) 2-Methylnaphthalene	8.72	142	660423	32.500	ng	99
49) Dimethylphthalate	9.47	163	115796	5.625	ng	# 98
57) Fluorene	10.30	166	41643	2.591	ng	# 86
70) Phenanthrene	11.26	178	88521	4.087	ng	98

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

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