

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF012318\
 Data File : BF102344.D
 Acq On : 23 Jan 2018 14:33
 Operator : SJ/JU
 Sample : J1239-17
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-50-02

Quant Time: Jan 24 01:07:20 2018
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF012218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 23 00:43:41 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.72	152	136839	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	569041	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	255479	20.00	ng	0.00
63) Phenanthrene-d10	11.23	188	457511	20.00	ng	0.00
75) Chrysene-d12	13.87	240	389467	20.00	ng	0.00
86) Perylene-d12	15.30	264	290279	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.34	112	1025262	113.61	ng	0.02
7) Phenol-d6	6.36	99	1162519	106.85	ng	0.00
23) Nitrobenzene-d5	7.29	82	858445	86.69	ng	0.00
41) 2,4,6-Tribromophenol	10.55	330	316106	124.87	ng	0.00
44) 2-Fluorobiphenyl	9.07	172	1538532	88.55	ng	0.00
78) Terphenyl-d14	12.83	244	1504983	87.11	ng	0.00
Target Compounds						
10) Phenol	6.37	94	94234	7.933	ng	99
17) 2-Methylphenol	6.97	107	30784	3.747	ng	95
20) 3+4-Methylphenols	7.13	107	88655	9.175	ng	86
27) 2,4-Dimethylphenol	7.66	122	28167	3.637	ng	96
31) Naphthalene	8.03	128	2444842	86.052	ng	99
37) 2-Methylnaphthalene	8.71	142	101274	5.352	ng	98
48) Acenaphthylene	9.61	152	74943	2.709	ng	100
54) Dibenzofuran	9.95	168	51680	2.364	ng	100
57) Fluorene	10.30	166	55479	3.202	ng	98
70) Phenanthrene	11.26	178	93980	3.525	ng	98
72) Carbazole	11.47	167	204782	7.714	ng	97

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

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