

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF012318\
 Data File : BF102385.D
 Acq On : 24 Jan 2018 11:59
 Operator : SJ/JU
 Sample : J1239-16DL 20X
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-43-01DL

Quant Time: Jan 24 12:37:27 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	6.71	152	115127	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	463931	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	191744	20.00	ng	0.00
63) Phenanthrene-d10	11.23	188	281915	20.00	ng	0.00
75) Chrysene-d12	13.87	240	210895	20.00	ng	0.00
86) Perylene-d12	15.29	264	204660	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	51784	6.82	ng	0.00
7) Phenol-d6	6.35	99	60460	6.60	ng	0.00
23) Nitrobenzene-d5	7.27	82	43978	5.45	ng	-0.01
41) 2,4,6-Tribromophenol	10.54	330	9974	5.25	ng	0.00
44) 2-Fluorobiphenyl	9.07	172	93235	7.15	ng	-0.01
78) Terphenyl-d14	12.82	244	60136	6.43	ng	0.00
Target Compounds						
10) Phenol	6.36	94	622557	62.296	ng	99
17) 2-Methylphenol	6.97	107	123069	17.804	ng	97
20) 3+4-Methylphenols	7.13	107	385119	47.371	ng	85
27) 2,4-Dimethylphenol	7.66	122	66196	10.484	ng	96
31) Naphthalene	8.02	128	978255	42.233	ng	99
72) Carbazole	11.47	167	38654	2.363	ng	96

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

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