

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF012318\
 Data File : BF102384.D
 Acq On : 24 Jan 2018 11:32
 Operator : SJ/JU
 Sample : J1239-15DL 10X
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-42-01DL

Quant Time: Jan 24 12:30:56 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	121746	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	501393	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	210124	20.00	ng	0.00
63) Phenanthrene-d10	11.23	188	306321	20.00	ng	0.00
75) Chrysene-d12	13.87	240	243244	20.00	ng	0.00
86) Perylene-d12	15.29	264	217385	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	106287	13.24	ng	0.00
7) Phenol-d6	6.35	99	120210	12.42	ng	0.00
23) Nitrobenzene-d5	7.27	82	86322	9.89	ng	-0.01
41) 2,4,6-Tribromophenol	10.54	330	23464	11.27	ng	0.00
44) 2-Fluorobiphenyl	9.07	172	181033	12.67	ng	-0.01
78) Terphenyl-d14	12.82	244	129681	12.02	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.36	94	485379	45.929	ng	99
17) 2-Methylphenol	6.97	107	117057	16.013	ng	99
20) 3+4-Methylphenols	7.13	107	354397	41.222	ng	87
27) 2,4-Dimethylphenol	7.66	122	69789	10.227	ng	98
31) Naphthalene	8.02	128	1274747	50.921	ng	100
37) 2-Methylnaphthalene	8.71	142	37203	2.231	ng	97
48) Acenaphthylene	9.61	152	58963	2.592	ng	97
72) Carbazole	11.47	167	51475	2.896	ng	99

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

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