

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF012318\
 Data File : BF102381.D
 Acq On : 24 Jan 2018 10:11
 Operator : SJ/JU
 Sample : J1239-08DL 10X
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-43-02DL

Quant Time: Jan 24 11:01:40 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	117241	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	491890	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	213671	20.00	ng	0.00
63) Phenanthrene-d10	11.23	188	330271	20.00	ng	0.00
75) Chrysene-d12	13.87	240	241091	20.00	ng	0.00
86) Perylene-d12	15.29	264	215826	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	110925	14.35	ng	0.00
7) Phenol-d6	6.35	99	126333	13.55	ng	0.00
23) Nitrobenzene-d5	7.27	82	92119	10.76	ng	-0.01
41) 2,4,6-Tribromophenol	10.54	330	27206	12.85	ng	0.00
44) 2-Fluorobiphenyl	9.07	172	198345	13.65	ng	-0.01
78) Terphenyl-d14	12.82	244	145462	13.60	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.36	94	360154	35.389	ng	99
17) 2-Methylphenol	6.97	107	94212	13.383	ng	98
20) 3+4-Methylphenols	7.13	107	268128	32.386	ng	83
27) 2,4-Dimethylphenol	7.66	122	47012	7.023	ng	98
31) Naphthalene	8.02	128	1292969	52.647	ng	99
37) 2-Methylnaphthalene	8.71	142	51838	3.169	ng	97
48) Acenaphthylene	9.61	152	48704	2.105	ng	96
72) Carbazole	11.47	167	46956	2.450	ng	96

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

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