

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF012318\
 Data File : BF102338.D
 Acq On : 23 Jan 2018 11:50
 Operator : SJ/JU
 Sample : J1239-10
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-096613-02

Manual Integrations
 APPROVED

Sohil
 1/24/2018 2:46:33 PM

Quant Time: Jan 24 00:32:29 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	131429	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	539245	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	244679	20.00	ng	0.00
63) Phenanthrene-d10	11.24	188	433672	20.00	ng	0.00
75) Chrysene-d12	13.87	240	371724	20.00	ng	0.00
86) Perylene-d12	15.30	264	280284	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.34	112	983828	113.50	ng	0.02
7) Phenol-d6	6.37	99	1036742	99.21	ng	0.02
23) Nitrobenzene-d5	7.29	82	808631	86.18	ng	0.00
41) 2,4,6-Tribromophenol	10.55	330	307208	126.71	ng	0.00
44) 2-Fluorobiphenyl	9.08	172	1420239	85.35	ng	0.00
78) Terphenyl-d14	12.83	244	1415257	85.83	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.39	94	1546941	135.595	ng	98
17) 2-Methylphenol	6.99	107	629858	79.816	ng	99
20) 3+4-Methylphenols	7.15	107	1653003	178.105	ng	87
27) 2,4-Dimethylphenol	7.66	122	427822	58.296	ng	97
31) Naphthalene	8.03	128	4023968	149.458	ng	98
37) 2-Methylnaphthalene	8.71	142	435135	24.268	ng	98
45) 1,1'-Biphenyl	9.17	154	65380	2.988	ng	95
48) Acenaphthylene	9.61	152	160973	6.076	ng	98
54) Dibenzofuran	9.96	168	76774	3.666	ng	99
57) Fluorene	10.30	166	129652	7.813	ng	99
70) Phenanthrene	11.26	178	273363	10.817	ng	98
71) Anthracene	11.31	178	58454m	2.266	ng	
72) Carbazole	11.47	167	520276	20.675	ng	99

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

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