

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF012318\
 Data File : BF102387.D
 Acq On : 24 Jan 2018 12:54
 Operator : SJ/JU
 Sample : J1239-05
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-14R30-01

Manual Integrations
 APPROVED

Sohil
 1/24/2018 2:47:48 PM

Quant Time: Jan 24 13:36:14 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	133325	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	542409	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	233756	20.00	ng	0.00
63) Phenanthrene-d10	11.23	188	346125	20.00	ng	0.00
75) Chrysene-d12	13.87	240	253001	20.00	ng	0.00
86) Perylene-d12	15.29	264	234666	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	1012436	115.14	ng	0.01
7) Phenol-d6	6.36	99	1060714	100.06	ng	0.00
23) Nitrobenzene-d5	7.28	82	799253	84.68	ng	0.00
41) 2,4,6-Tribromophenol	10.54	330	246875	106.59	ng	0.00
44) 2-Fluorobiphenyl	9.07	172	1397709	87.92	ng	0.00
78) Terphenyl-d14	12.82	244	1010024	90.00	ng	0.00

Target Compounds				Qvalue		
10) Phenol	6.37	94	36570	3.160	ng	90
17) 2-Methylphenol	6.97	107	25821	3.226	ng	# 92
20) 3+4-Methylphenols	7.13	107	70274	7.464	ng	87
27) 2,4-Dimethylphenol	7.67	122	20631	2.795	ng	97
31) Naphthalene	8.03	128	3050143	112.628	ng	99
37) 2-Methylnaphthalene	8.71	142	286060	15.861	ng	99
45) 1,1'-Biphenyl	9.17	154	56519	2.704	ng	100
48) Acenaphthylene	9.61	152	185428	7.326	ng	99
51) Acenaphthene	9.78	154	89274	6.039	ng	98
54) Dibenzofuran	9.95	168	42879	2.143	ng	99
57) Fluorene	10.29	166	89936	5.673	ng	99
70) Phenanthrene	11.26	178	152193	7.546	ng	98
72) Carbazole	11.47	167	149680	7.453	ng	98

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

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