

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF012318\
 Data File : BF102343.D
 Acq On : 23 Jan 2018 14:06
 Operator : SJ/JU
 Sample : J1239-16
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-43-01

Manual Integrations
 APPROVED

Sohil
 1/24/2018 2:46:48 PM

Quant Time: Jan 24 01:02:00 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	144560	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	449960	20.00	ng	0.02
38) Acenaphthene-d10	9.75	164	280040	20.00	ng	0.00
63) Phenanthrene-d10	11.24	188	485722	20.00	ng	0.00
75) Chrysene-d12	13.87	240	393062	20.00	ng	0.00
86) Perylene-d12	15.30	264	295133	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	1201037	125.97	ng	0.02
7) Phenol-d6	6.41	99	963878	83.86	ng	0.05
23) Nitrobenzene-d5	7.31	82	994097	126.96	ng	0.02
41) 2,4,6-Tribromophenol	10.55	330	341585	123.10	ng	0.00
44) 2-Fluorobiphenyl	9.08	172	1683871	88.41	ng	0.00
78) Terphenyl-d14	12.83	244	1616999	92.74	ng	0.00
Target Compounds						Qvalue
3) Pyridine	3.20	79	233734	19.744	ng	99
10) Phenol	6.46	94	9066956	722.560	ng	96
17) 2-Methylphenol	7.01	107	1918294	221.008	ng	97
20) 3+4-Methylphenols	7.20	107	5572991	545.925	ng	87
27) 2,4-Dimethylphenol	7.70	122	1434766	234.297	ng	92
31) Naphthalene	8.06	128	9769582	434.864	ng	98
37) 2-Methylnaphthalene	8.72	142	644947	43.107	ng	99
45) 1,1'-Biphenyl	9.17	154	124025	4.953	ng	97
48) Acenaphthylene	9.62	152	932903	30.766	ng	100
54) Dibenzofuran	9.96	168	316912	13.223	ng	98
57) Fluorene	10.30	166	276292	14.548	ng	# 96
70) Phenanthrene	11.26	178	326747	11.544	ng	99
71) Anthracene	11.31	178	83835m	2.902	ng	
72) Carbazole	11.48	167	1127146	39.992	ng	99

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

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