

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF012318\
 Data File : BF102340.D
 Acq On : 23 Jan 2018 12:44
 Operator : SJ/JU
 Sample : J1239-12
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-114767779-01

Manual Integrations
 APPROVED

Sohil
 1/24/2018 2:46:39 PM

Quant Time: Jan 24 11:17:38 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	159807	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	480743	20.00	ng	0.02
38) Acenaphthene-d10	9.75	164	300467	20.00	ng	0.00
63) Phenanthrene-d10	11.24	188	497584	20.00	ng	0.00
75) Chrysene-d12	13.87	240	411868	20.00	ng	0.00
86) Perylene-d12	15.30	264	311856	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.34	112	1071182	101.63	ng	0.02
7) Phenol-d6	6.39	99	1049331	82.58	ng	0.03
23) Nitrobenzene-d5	7.30	82	923736	110.42	ng	0.01
41) 2,4,6-Tribromophenol	10.55	330	343003	115.21	ng	0.00
44) 2-Fluorobiphenyl	9.08	172	1591307	77.87	ng	0.00
78) Terphenyl-d14	12.83	244	1546764	84.66	ng	0.01

Target Compounds

						Qvalue
3) Pyridine	3.22	79	87401m	6.679	ng	
10) Phenol	6.42	94	3516231	253.479	ng	96
17) 2-Methylphenol	6.99	107	914734	95.332	ng	98
20) 3+4-Methylphenols	7.17	107	2829169	250.701	ng	85
27) 2,4-Dimethylphenol	7.68	122	673174	102.890	ng	97
31) Naphthalene	8.06	128	10629725	442.854	ng	97
37) 2-Methylnaphthalene	8.72	142	669631	41.891	ng	99
45) 1,1'-Biphenyl	9.17	154	137143	5.105	ng	97
48) Acenaphthylene	9.61	152	428288	13.164	ng	98
51) Acenaphthene	9.78	154	68731	3.617	ng	96
54) Dibenzofuran	9.96	168	335341	13.040	ng	98
57) Fluorene	10.30	166	305118	14.974	ng	97
70) Phenanthrene	11.26	178	423593	14.609	ng	99
71) Anthracene	11.31	178	103317m	3.491	ng	
72) Carbazole	11.49	167	1076371	37.280	ng	99

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

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