

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF012318\
 Data File : BF102342.D
 Acq On : 23 Jan 2018 13:39
 Operator : SJ/JU
 Sample : J1239-15
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-42-01

Manual Integrations
 APPROVED

Sohil
 1/24/2018 2:46:45 PM

Quant Time: Jan 24 11:34:14 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	163563	20.00	ng	0.00
21) Naphthalene-d8	8.01	136	571763	20.00	ng	0.01
38) Acenaphthene-d10	9.75	164	309266	20.00	ng	0.00
63) Phenanthrene-d10	11.23	188	519604	20.00	ng	0.00
75) Chrysene-d12	13.87	240	402238	20.00	ng	0.00
86) Perylene-d12	15.30	264	301346	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	1233981	114.39	ng	0.02
7) Phenol-d6	6.38	99	1195795	91.95	ng	0.02
23) Nitrobenzene-d5	7.30	82	1020115	102.53	ng	0.01
41) 2,4,6-Tribromophenol	10.55	330	361312	117.91	ng	0.00
44) 2-Fluorobiphenyl	9.08	172	1766015	83.96	ng	0.00
78) Terphenyl-d14	12.83	244	1597755	89.55	ng	0.00
Target Compounds						
3) Pyridine	3.24	79	99294m	7.413	ng	Qvalue
10) Phenol	6.42	94	4397202	309.708	ng	96
17) 2-Methylphenol	6.99	107	1036109	105.502	ng	98
20) 3+4-Methylphenols	7.17	107	3238030	280.342	ng	87
27) 2,4-Dimethylphenol	7.67	122	854729	109.843	ng	96
31) Naphthalene	8.05	128	7875740	275.884	ng	98
37) 2-Methylnaphthalene	8.72	142	477988	25.142	ng	97
45) 1,1'-Biphenyl	9.17	154	95378	3.449	ng	96
48) Acenaphthylene	9.62	152	785003	23.442	ng	98
51) Acenaphthene	9.78	154	65759	3.362	ng	96
54) Dibenzofuran	9.96	168	249915	9.442	ng	98
57) Fluorene	10.30	166	227110	10.828	ng	96
70) Phenanthrene	11.26	178	253698	8.379	ng	99
72) Carbazole	11.47	167	803463	26.649	ng	99

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

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