

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_F\DATA\BF012318\
 Data File : BF102339.D
 Acq On : 23 Jan 2018 12:17
 Operator : SJ/JU
 Sample : J1239-11
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-50-01

Manual Integrations
 APPROVED

Sohil
 1/24/2018 2:46:36 PM

Quant Time: Jan 24 11:14:10 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	157266	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	531768	20.00	ng	0.02
38) Acenaphthene-d10	9.75	164	303408	20.00	ng	0.00
63) Phenanthrene-d10	11.24	188	520418	20.00	ng	0.00
75) Chrysene-d12	13.87	240	419441	20.00	ng	0.00
86) Perylene-d12	15.30	264	308275	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	1264598	121.92	ng	0.02
7) Phenol-d6	6.40	99	1171932	93.72	ng	0.04
23) Nitrobenzene-d5	7.30	82	1079588	116.67	ng	0.02
41) 2,4,6-Tribromophenol	10.55	330	391821	130.33	ng	0.00
44) 2-Fluorobiphenyl	9.08	172	1841839	89.26	ng	0.00
78) Terphenyl-d14	12.83	244	1706205	91.70	ng	0.01
Target Compounds						
3) Pyridine	3.25	79	136481m	10.598	ng	Qvalue
10) Phenol	6.43	94	5943388	435.371	ng	96
17) 2-Methylphenol	7.00	107	1686308	178.584	ng	98
20) 3+4-Methylphenols	7.19	107	4586988	413.034	ng	86
27) 2,4-Dimethylphenol	7.69	122	1299934	179.621	ng	93
31) Naphthalene	8.05	128	8470937	319.052	ng	97
37) 2-Methylnaphthalene	8.72	142	563244	31.854	ng	100
45) 1,1'-Biphenyl	9.17	154	121798	4.489	ng	98
48) Acenaphthylene	9.62	152	946154	28.800	ng	99
51) Acenaphthene	9.79	154	51038	2.660	ng	99
54) Dibenzofuran	9.96	168	302678	11.656	ng	100
57) Fluorene	10.30	166	288749	14.033	ng	96
70) Phenanthrene	11.26	178	359621	11.859	ng	99
71) Anthracene	11.31	178	89497m	2.891	ng	
72) Carbazole	11.48	167	1036493	34.324	ng	100

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

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