

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM120417\
 Data File : BM013178.D
 Acq On : 04 Dec 2017 21:57
 Operator : SJ/JU
 Sample : I6646-18
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0244

Quant Time: Dec 05 06:07:39 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM113017.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Dec 05 04:53:13 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	284976	20.00	ng/ul	0.00
18) Naphthalene-d8	10.49	136	1229529	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.34	164	713304	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.09	188	1598305	20.00	ng/ul	0.00
75) Chrysene-d12	21.27	240	1414619	20.00	ng/ul	0.00
83) Perylene-d12	23.50	264	1071099	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.25	96	17128	2.98	ng/uL	0.00
5) Phenol-d5	6.88	99	427898	17.35	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.05	67	273241	19.43	ng/ul	0.00
9) 2-Chlorophenol-d4	7.24	132	371745	19.29	ng/ul	0.00
13) 4-Methylphenol-d8	8.41	113	214208	10.17	ng/ul	0.00
19) Nitrobenzene-d5	8.86	128	199975	20.51	ng/ul	0.00
22) 2-Nitrophenol-d4	9.57	143	207989	19.93	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.11	165	362730	16.87	ng/ul	0.00
29) 4-Chloroaniline-d4	10.62	131	334716	16.94	ng/ul	0.00
43) Dimethylphthalate-d6	13.76	166	1207389	19.02	ng/ul	0.00
46) Acenaphthylene-d8	14.03	160	1525128	19.47	ng/ul	0.00
51) 4-Nitrophenol-d4	14.55	143	168914	15.32	ng/ul	0.00
57) Fluorene-d10	15.34	176	1052397	19.28	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.46	200	133699	13.30	ng/ul	0.00
70) Anthracene-d10	17.19	188	1594581	19.92	ng/ul	0.00
76) Pyrene-d10	19.49	212	1669788	23.69	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.36	264	1096755	20.10	ng/ul	0.00

Target Compounds

					Ovalue	
6) Phenol	6.91	94	44309	1.818	ng/ul#	84
44) Dimethylphthalate	13.80	163	203111	3.346	ng/ul#	98
49) Acenaphthene	14.40	153	67686	1.371	ng/ul	99
58) Fluorene	15.39	166	90068	1.494	ng/ul	89
69) Phenanthrene	17.13	178	618032	6.806	ng/ul	99
74) Fluoranthene	19.14	202	421168	3.785	ng/ul#	93
77) Pyrene	19.51	202	459309	5.344	ng/ul#	90
80) Benzo(a)anthracene	21.26	228	142702	1.590	ng/ul#	91
82) Chrysene	21.31	228	153342	1.842	ng/ul#	95
85) Benzo(b)fluoranthene	22.84	252	140759	1.949	ng/ul#	83
88) Benzo(a)pyrene	23.40	252	88427	1.363	ng/ul#	89

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

Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA M\DATA\BM120417\
Data File : BM013178.D
Acq On : 04 Dec 2017 21:57
Operator : SJ/JU
Sample : I6646-18
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0244

Quant Time: Dec 05 06:07:39 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM113017.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Dec 05 04:53:13 2017
Response via : Initial Calibration

