

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM120517\
 Data File : BM013206.D
 Acq On : 06 Dec 2017 01:30
 Operator : SJ/JU
 Sample : I6689-13
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 JHPQ1

Quant Time: Dec 06 03:23:59 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM113017.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 06 01:31:21 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	270948	20.00	ng/ul	0.00
18) Naphthalene-d8	10.48	136	1206967	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.34	164	778807	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.09	188	1874944	20.00	ng/ul	0.00
75) Chrysene-d12	21.27	240	1911082	20.00	ng/ul	0.00
83) Perylene-d12	23.50	264	1504187	20.00	ng/ul	-0.01

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.25	96	22140	4.05	ng/uL	0.00
5) Phenol-d5	6.88	99	706	0.03	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.05	67	4261	0.32	ng/ul	0.00
9) 2-Chlorophenol-d4	7.23	132	3938	0.21	ng/ul	-0.01
13) 4-Methylphenol-d8	8.41	113	1673	0.08	ng/ul	0.00
19) Nitrobenzene-d5	8.86	128	2637	0.28	ng/ul	0.00
22) 2-Nitrophenol-d4	9.57	143	2637	0.26	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.11	165	3319	0.16	ng/ul	0.00
29) 4-Chloroaniline-d4	10.49	131	1544	0.08	ng/ul	-0.14
43) Dimethylphthalate-d6	13.76	166	21447	0.31	ng/ul	0.00
46) Acenaphthylene-d8	14.03	160	24056	0.28	ng/ul	0.00
51) 4-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
57) Fluorene-d10	15.33	176	20006	0.34	ng/ul	-0.01
62) 4,6-Dinitro-2-methylphenol	15.47	200	745	0.06	ng/ul	0.00
70) Anthracene-d10	17.19	188	36299	0.39	ng/ul	0.00
76) Pyrene-d10	19.49	212	31887	0.33	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.36	264	26464	0.35	ng/ul	-0.02

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

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