

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM112219\
 Data File : BM023881.D
 Acq On : 23 Nov 2019 05:44
 Operator : JU
 Sample : K5854-02
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AB7

Quant Time: Nov 25 01:32:24 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111119MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Nov 25 01:09:42 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.50	152	379185	20.00	ng/ul	0.00
18) Naphthalene-d8	10.28	136	1568489	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.17	164	967019	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.92	188	2046933	20.00	ng/ul	0.00
78) Chrysene-d12	21.13	240	1925160	20.00	ng/ul	0.00
86) Perylene-d12	23.28	264	2418435	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.05	96	24360	2.65	ng/uL	0.00
5) Phenol-d5	6.70	99	511677	14.80	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.85	67	286666	14.43	ng/ul	0.00
9) 2-Chlorophenol-d4	7.05	132	413216	16.50	ng/ul	0.00
13) 4-Methylphenol-d8	8.23	113	363325	13.18	ng/ul	0.00
19) Nitrobenzene-d5	8.66	128	197898	17.63	ng/ul	0.00
22) 2-Nitrophenol-d4	9.37	143	225372	19.10	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.92	165	409496	15.22	ng/ul	0.00
29) 4-Chloroaniline-d4	10.43	131	436513	14.77	ng/ul	0.00
44) Dimethylphthalate-d6	13.59	166	1264971	16.77	ng/ul	0.00
47) Acenaphthylene-d8	13.86	160	1529008	17.77	ng/ul	0.00
52) 4-Nitrophenol-d4	14.41	143	138667	10.77	ng/ul	0.01
58) Fluorene-d10	15.17	176	1117481	16.78	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.31	200	65898	5.57	ng/ul	0.00
71) Anthracene-d10	17.02	188	1705170	17.56	ng/ul	0.00
79) Pyrene-d10	19.33	212	1950140	19.43	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.14	264	2314452	18.19	ng/ul	0.00

Target Compounds						Qvalue
6) Phenol	6.73	94	129604	3.640	ng/ul	98
45) Dimethylphthalate	13.63	163	278682	3.771	ng/ul	99

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

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