

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM101519\
 Data File : BM023160.D
 Acq On : 15 Oct 2019 11:24
 Operator : JU
 Sample : K5177-09
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BEP98

Quant Time: Oct 15 16:48:17 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM101419MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 15 02:28:10 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	262295	20.00	ng/ul	0.00
18) Naphthalene-d8	10.47	136	1125380	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.33	164	731409	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.07	188	1601347	20.00	ng/ul	0.00
78) Chrysene-d12	21.26	240	1840867	20.00	ng/ul	0.00
86) Perylene-d12	23.47	264	2237855	20.00	ng/ul	-0.02

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.20	96	21204	3.56	ng/uL	0.00
5) Phenol-d5	6.86	99	342282	14.98	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.03	67	237529	18.38	ng/ul	0.00
9) 2-Chlorophenol-d4	7.22	132	286892	16.41	ng/ul	0.00
13) 4-Methylphenol-d8	8.39	113	175406	9.56	ng/ul	0.00
19) Nitrobenzene-d5	8.83	128	132332	17.38	ng/ul	0.00
22) 2-Nitrophenol-d4	9.55	143	120176	16.58	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.09	165	296466	16.08	ng/ul	0.00
29) 4-Chloroaniline-d4	10.60	131	185366	8.26	ng/ul	0.00
44) Dimethylphthalate-d6	13.74	166	1119898	19.98	ng/ul	0.00
47) Acenaphthylene-d8	14.02	160	1292298	19.66	ng/ul	0.00
52) 4-Nitrophenol-d4	14.51	143	96063	10.59	ng/ul	-0.01
58) Fluorene-d10	15.33	176	962409	20.15	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.43	200	60820	9.58	ng/ul	-0.01
71) Anthracene-d10	17.17	188	1514849	19.62	ng/ul	-0.01
79) Pyrene-d10	19.46	212	1923048	21.57	ng/ul	-0.01
90) Benzo(a)pyrene-d12	23.34	264	2369747	20.89	ng/ul	-0.01

Target Compounds

					Ovalue
6) Phenol	6.88	94	56355	2.392	ng/ul 96
45) Dimethylphthalate	13.79	163	395261	7.045	ng/ul 99

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

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