

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP100919\
 Data File : BP000603.D
 Acq On : 10 Oct 2019 05:30
 Operator : JU
 Sample : K5097-10
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 JLME6

Quant Time: Oct 10 05:58:17 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP100919MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 09 17:20:45 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	240217	20.00	ng/ul	0.00
18) Naphthalene-d8	8.03	136	923829	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.80	164	532582	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.29	188	978481	20.00	ng/ul	0.00
78) Chrysene-d12	13.96	240	843983	20.00	ng/ul	0.00
86) Perylene-d12	15.44	264	924775	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.36	96	32244	5.57	ng/uL	0.00
5) Phenol-d5	6.39	99	544545	29.39	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.43	67	282230	30.47	ng/ul	0.00
9) 2-Chlorophenol-d4	6.52	132	466192	29.86	ng/ul	0.00
13) 4-Methylphenol-d8	7.13	113	435731	30.78	ng/ul	0.00
19) Nitrobenzene-d5	7.31	128	220251	31.36	ng/ul	0.00
22) 2-Nitrophenol-d4	7.63	143	240042	32.10	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	7.88	165	426846	29.82	ng/ul	0.00
29) 4-Chloroaniline-d4	8.10	131	585538	36.79	ng/ul	0.00
44) Dimethylphthalate-d6	9.49	166	1167068	31.30	ng/ul	0.00
47) Acenaphthylene-d8	9.64	160	1404095	30.62	ng/ul	0.00
52) 4-Nitrophenol-d4	9.90	143	146049	25.43	ng/ul	0.00
58) Fluorene-d10	10.32	176	1002166	31.42	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.39	200	101341	21.75	ng/ul	0.00
71) Anthracene-d10	11.35	188	1448079	31.95	ng/ul	0.00
79) Pyrene-d10	12.73	212	1483701	31.09	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.35	264	1431562	31.36	ng/ul	0.00

Target Compounds

				Ovalue	
6) Phenol	6.40	94	82570	4.420	ng/ul 99
45) Dimethylphthalate	9.51	163	336900	9.069	ng/ul 100

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

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