

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP101219\
 Data File : BP000664.D
 Acq On : 12 Oct 2019 13:55
 Operator : JU
 Sample : K5199-11
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BEP78

Quant Time: Oct 12 21:46:57 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	161797	20.00	ng/ul	0.00
18) Naphthalene-d8	8.03	136	632762	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.80	164	357346	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.29	188	656420	20.00	ng/ul	0.00
78) Chrysene-d12	13.97	240	556420	20.00	ng/ul	0.00
86) Perylene-d12	15.45	264	611761	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.36	96	24543	6.29	ng/uL	-0.01
5) Phenol-d5	6.38	99	413002	33.09	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.43	67	224554	35.99	ng/ul	0.00
9) 2-Chlorophenol-d4	6.52	132	357187	33.97	ng/ul	0.00
13) 4-Methylphenol-d8	7.13	113	292779	30.71	ng/ul	0.00
19) Nitrobenzene-d5	7.30	128	169644	35.27	ng/ul	0.00
22) 2-Nitrophenol-d4	7.63	143	183339	35.79	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	7.88	165	322461	32.89	ng/ul	0.00
29) 4-Chloroaniline-d4	8.09	131	396866	36.40	ng/ul	0.00
44) Dimethylphthalate-d6	9.49	166	920007	36.78	ng/ul	0.00
47) Acenaphthylene-d8	9.64	160	1103343	35.86	ng/ul	0.00
52) 4-Nitrophenol-d4	9.91	143	100935	26.20	ng/ul	0.00
58) Fluorene-d10	10.32	176	780461	36.46	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.39	200	81354	26.03	ng/ul	0.00
71) Anthracene-d10	11.35	188	1118272	36.78	ng/ul	0.00
79) Pyrene-d10	12.73	212	1177149	37.41	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.35	264	1119475	37.07	ng/ul	0.00

Target Compounds

					Ovalue
6) Phenol	6.40	94	28889	2.296	ng/ul 98
45) Dimethylphthalate	9.51	163	295440	11.853	ng/ul 99
77) Fluoranthene	12.52	202	40575	1.107	ng/ul 98

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

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