

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP102219\
 Data File : BP000944.D
 Acq On : 22 Oct 2019 15:59
 Operator : JU
 Sample : K5345-19
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 ETOL3

Quant Time: Oct 23 05:36:57 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP100919MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 23 05:34:10 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.74	152	130175	20.00	ng/ul	0.00
18) Naphthalene-d8	8.02	136	485508	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.79	164	278055	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.28	188	515786	20.00	ng/ul	0.00
78) Chrysene-d12	13.96	240	454610	20.00	ng/ul	0.00
86) Perylene-d12	15.43	264	498825	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.35	96	10167	3.24	ng/uL	0.00
5) Phenol-d5	6.38	99	181054	18.03	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.42	67	92752	18.48	ng/ul	0.00
9) 2-Chlorophenol-d4	6.51	132	164268	19.42	ng/ul	0.00
13) 4-Methylphenol-d8	7.13	113	153938	20.07	ng/ul	0.00
19) Nitrobenzene-d5	7.30	128	76954	20.85	ng/ul	0.00
22) 2-Nitrophenol-d4	7.62	143	83144	21.16	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	7.88	165	150997	20.08	ng/ul	0.00
29) 4-Chloroaniline-d4	8.09	131	111103	13.28	ng/ul	0.00
44) Dimethylphthalate-d6	9.48	166	447942	23.01	ng/ul	0.00
47) Acenaphthylene-d8	9.63	160	512468	21.40	ng/ul	0.00
52) 4-Nitrophenol-d4	9.91	143	51301	17.11	ng/ul	0.00
58) Fluorene-d10	10.30	176	372312	22.36	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.39	200	37007	15.07	ng/ul	0.00
71) Anthracene-d10	11.34	188	570742	23.89	ng/ul	0.00
79) Pyrene-d10	12.72	212	607413	23.63	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.35	264	609661	24.76	ng/ul	0.00

Target Compounds

				Ovalue	
6) Phenol	6.39	94	16614	1.641	ng/ul# 90
45) Dimethylphthalate	9.50	163	155047	7.994	ng/ul 99

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

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