

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP101619\
 Data File : BP000782.D
 Acq On : 15 Oct 2019 21:48
 Operator : JU
 Sample : K5178-07
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BEPFO

Quant Time: Oct 16 05:51:17 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP100919MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 16 05:11:27 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	219839	20.00	ng/ul	0.00
18) Naphthalene-d8	8.03	136	857240	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.80	164	483542	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.29	188	916897	20.00	ng/ul	0.00
78) Chrysene-d12	13.97	240	726666	20.00	ng/ul	0.00
86) Perylene-d12	15.45	264	701492	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.36	96	20755	3.92	ng/uL	0.00
5) Phenol-d5	6.39	99	363812	21.45	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.43	67	191712	22.61	ng/ul	0.00
9) 2-Chlorophenol-d4	6.52	132	317406	22.22	ng/ul	0.00
13) 4-Methylphenol-d8	7.13	113	271560	20.96	ng/ul	0.00
19) Nitrobenzene-d5	7.30	128	151365	23.23	ng/ul	0.00
22) 2-Nitrophenol-d4	7.63	143	163303	23.53	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	7.88	165	292628	22.03	ng/ul	0.00
29) 4-Chloroaniline-d4	8.09	131	306069	20.72	ng/ul	0.00
44) Dimethylphthalate-d6	9.49	166	836239	24.70	ng/ul	0.00
47) Acenaphthylene-d8	9.64	160	993931	23.87	ng/ul	0.00
52) 4-Nitrophenol-d4	9.91	143	93457	17.93	ng/ul	0.00
58) Fluorene-d10	10.32	176	703532	24.29	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.39	200	67466	15.45	ng/ul	0.00
71) Anthracene-d10	11.35	188	1032264	24.31	ng/ul	0.00
79) Pyrene-d10	12.73	212	1075726	26.18	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.36	264	847930	24.48	ng/ul	0.00

Target Compounds

				Ovalue	
6) Phenol	6.40	94	46494	2.720	ng/ul 99
45) Dimethylphthalate	9.51	163	450597	13.360	ng/ul 99
84) Bis(2-ethylhexyl)phthalate	13.96	149	85736	3.061	ng/ul# 98

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

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