

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP101919\
 Data File : BP000919.D
 Acq On : 19 Oct 2019 17:03
 Operator : JU
 Sample : K5354-15
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BFBB8

Quant Time: Oct 19 22:48:24 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP100919MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Oct 19 22:14:11 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.74	152	117788	20.00	ng/ul	0.00
18) Naphthalene-d8	8.02	136	449147	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.79	164	257728	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.28	188	500648	20.00	ng/ul	0.00
78) Chrysene-d12	13.96	240	470672	20.00	ng/ul	0.00
86) Perylene-d12	15.43	264	536167	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.35	96	14030	4.94	ng/uL	0.00
5) Phenol-d5	6.39	99	55579	6.12	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.42	67	134387	29.58	ng/ul	0.00
9) 2-Chlorophenol-d4	6.51	132	195776	25.58	ng/ul	0.00
13) 4-Methylphenol-d8	7.13	113	113493	16.35	ng/ul	0.00
19) Nitrobenzene-d5	7.30	128	115183	33.74	ng/ul	0.00
22) 2-Nitrophenol-d4	7.62	143	125340	34.47	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	7.87	165	210661	30.28	ng/ul	0.00
29) 4-Chloroaniline-d4	8.09	131	18111	2.34	ng/ul	0.00
44) Dimethylphthalate-d6	9.48	166	618574	34.29	ng/ul	0.00
47) Acenaphthylene-d8	9.63	160	738584	33.28	ng/ul	0.00
52) 4-Nitrophenol-d4	9.92	143	15047	5.41	ng/ul	0.00
58) Fluorene-d10	10.31	176	529026	34.27	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.39	200	59687	25.03	ng/ul	0.00
71) Anthracene-d10	11.34	188	842228	36.32	ng/ul	0.00
79) Pyrene-d10	12.72	212	921652	34.63	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.34	264	976931	36.91	ng/ul	0.00

Target Compounds				Ovalue
45) Dimethylphthalate	9.50	163	42107	2.342 ng/ul 99

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

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