

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP100919\
 Data File : BP000585.D
 Acq On : 09 Oct 2019 21:54
 Operator : JU
 Sample : K5100-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 ETJY1

Quant Time: Oct 10 01:22:40 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	243065	20.00	ng/ul	0.00
18) Naphthalene-d8	8.03	136	949918	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.80	164	554808	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.29	188	1030282	20.00	ng/ul	0.00
78) Chrysene-d12	13.96	240	906551	20.00	ng/ul	0.00
86) Perylene-d12	15.44	264	1004591	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.36	96	29744	5.08	ng/uL	0.00
5) Phenol-d5	6.39	99	118123	6.30	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.43	67	251739	26.86	ng/ul	0.00
9) 2-Chlorophenol-d4	6.52	132	341417	21.61	ng/ul	0.00
13) 4-Methylphenol-d8	7.13	113	210392	14.69	ng/ul	0.00
19) Nitrobenzene-d5	7.31	128	199813	27.67	ng/ul	0.00
22) 2-Nitrophenol-d4	7.63	143	215235	27.99	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	7.88	165	364836	24.79	ng/ul	0.00
29) 4-Chloroaniline-d4	8.10	131	35507	2.17	ng/ul	0.00
44) Dimethylphthalate-d6	9.49	166	1138861	29.32	ng/ul	0.00
47) Acenaphthylene-d8	9.65	160	1343457	28.12	ng/ul	0.00
52) 4-Nitrophenol-d4	9.91	143	37557	6.28	ng/ul	0.00
58) Fluorene-d10	10.32	176	976133	29.37	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.39	200	130072	26.51	ng/ul	0.00
71) Anthracene-d10	11.35	188	1544006	32.35	ng/ul	0.00
79) Pyrene-d10	12.73	212	1680368	32.78	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.35	264	1647242	33.21	ng/ul	0.00

Target Compounds					Ovalue	
45) Dimethylphthalate	9.51	163	63478	1.640	ng/ul	99

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

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