

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP100919\
 Data File : BP000579.D
 Acq On : 09 Oct 2019 19:29
 Operator : JU
 Sample : K5201-05
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AB9

Quant Time: Oct 10 01:21:47 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	234970	20.00	ng/ul	0.00
18) Naphthalene-d8	8.03	136	907408	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.80	164	531474	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.29	188	981077	20.00	ng/ul	0.00
78) Chrysene-d12	13.97	240	863121	20.00	ng/ul	0.00
86) Perylene-d12	15.44	264	946185	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	2.36	96	28392	5.01	ng/uL	0.00
5) Phenol-d5	6.39	99	111079	6.13	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.43	67	241712	26.67	ng/ul	0.00
9) 2-Chlorophenol-d4	6.52	132	326153	21.36	ng/ul	0.00
13) 4-Methylphenol-d8	7.13	113	190482	13.76	ng/ul	0.00
19) Nitrobenzene-d5	7.31	128	191588	27.78	ng/ul	0.00
22) 2-Nitrophenol-d4	7.63	143	207416	28.24	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	7.89	165	354266	25.20	ng/ul	0.00
29) 4-Chloroaniline-d4	8.10	131	21931	1.40	ng/ul	0.00
44) Dimethylphthalate-d6	9.49	166	1143603	30.74	ng/ul	0.00
47) Acenaphthylene-d8	9.65	160	1306442	28.55	ng/ul	0.00
52) 4-Nitrophenol-d4	9.91	143	34819	6.08	ng/ul	0.00
58) Fluorene-d10	10.32	176	952821	29.93	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.39	200	116792	25.00	ng/ul	0.00
71) Anthracene-d10	11.35	188	1434119	31.56	ng/ul	0.00
79) Pyrene-d10	12.73	212	1584072	32.46	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.35	264	1563799	33.48	ng/ul	0.00
Target Compounds						
6) Phenol	6.40	94	27479	1.504	ng/ul	96
45) Dimethylphthalate	9.51	163	134464	3.627	ng/ul	100

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

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