

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP100919\
 Data File : BP000597.D
 Acq On : 10 Oct 2019 02:42
 Operator : JU
 Sample : PB123660BL
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK60

Quant Time: Oct 10 03:43:20 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	219792	20.00	ng/ul	0.00
18) Naphthalene-d8	8.03	136	851871	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.80	164	492327	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.29	188	929292	20.00	ng/ul	0.00
78) Chrysene-d12	13.96	240	824803	20.00	ng/ul	0.00
86) Perylene-d12	15.44	264	917604	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.37	96	7158	1.35	ng/uL	0.00
5) Phenol-d5	6.39	99	121314	7.16	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.43	67	241748	28.52	ng/ul	0.00
9) 2-Chlorophenol-d4	6.52	132	336359	23.55	ng/ul	0.00
13) 4-Methylphenol-d8	7.13	113	211940	16.36	ng/ul	0.00
19) Nitrobenzene-d5	7.31	128	194191	29.99	ng/ul	0.00
22) 2-Nitrophenol-d4	7.63	143	208257	30.20	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	7.88	165	357747	27.11	ng/ul	0.00
29) 4-Chloroaniline-d4	8.10	131	37139	2.53	ng/ul	0.00
44) Dimethylphthalate-d6	9.49	166	1142074	33.14	ng/ul	0.00
47) Acenaphthylene-d8	9.65	160	1312700	30.96	ng/ul	0.00
52) 4-Nitrophenol-d4	9.90	143	36896	6.95	ng/ul	0.00
58) Fluorene-d10	10.32	176	991617	33.63	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.39	200	126119	28.50	ng/ul	0.00
71) Anthracene-d10	11.35	188	1652789	38.40	ng/ul	0.00
79) Pyrene-d10	12.73	212	1739571	37.30	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.35	264	1742446	38.46	ng/ul	0.00

Target Compounds

					Ovalue
6) Phenol	6.40	94	22199	1.299 ng/ul	97
45) Dimethylphthalate	9.51	163	84936	2.473 ng/ul	99

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

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