

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP100919\
 Data File : BP000599.D
 Acq On : 10 Oct 2019 03:30
 Operator : JU
 Sample : PB123576BL
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK76

Quant Time: Oct 10 03:53:05 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	223108	20.00	ng/ul	0.00
18) Naphthalene-d8	8.03	136	873062	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.80	164	502351	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.29	188	937629	20.00	ng/ul	0.00
78) Chrysene-d12	13.96	240	802860	20.00	ng/ul	0.00
86) Perylene-d12	15.44	264	893773	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.36	96	37620	6.99	ng/uL	-0.01
5) Phenol-d5	6.39	99	614429	35.70	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.43	67	308309	35.83	ng/ul	0.00
9) 2-Chlorophenol-d4	6.52	132	512980	35.38	ng/ul	0.00
13) 4-Methylphenol-d8	7.13	113	470279	35.77	ng/ul	0.00
19) Nitrobenzene-d5	7.31	128	242662	36.56	ng/ul	0.00
22) 2-Nitrophenol-d4	7.64	143	259781	36.76	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	7.88	165	462691	34.21	ng/ul	0.00
29) 4-Chloroaniline-d4	8.10	131	653890	43.47	ng/ul	0.00
44) Dimethylphthalate-d6	9.49	166	1284218	36.52	ng/ul	0.00
47) Acenaphthylene-d8	9.65	160	1529856	35.36	ng/ul	0.00
52) 4-Nitrophenol-d4	9.90	143	191361	35.33	ng/ul	0.00
58) Fluorene-d10	10.32	176	1082999	35.99	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.39	200	122169	27.36	ng/ul	0.00
71) Anthracene-d10	11.35	188	1575359	36.27	ng/ul	0.00
79) Pyrene-d10	12.73	212	1624648	35.79	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.35	264	1592636	36.09	ng/ul	0.00

Target Compounds Ovalue

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

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