

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP101219\
 Data File : BP000670.D
 Acq On : 12 Oct 2019 16:19
 Operator : JU
 Sample : PB123733BL
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK33

Quant Time: Oct 12 22:53:56 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	185371	20.00	ng/ul	0.00
18) Naphthalene-d8	8.03	136	718172	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.80	164	406453	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.29	188	761857	20.00	ng/ul	0.00
78) Chrysene-d12	13.97	240	662378	20.00	ng/ul	0.00
86) Perylene-d12	15.45	264	748993	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	2.35	96	31354	7.02	ng/uL	-0.02
5) Phenol-d5	6.38	99	508611	35.57	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.43	67	258364	36.14	ng/ul	0.00
9) 2-Chlorophenol-d4	6.52	132	422547	35.07	ng/ul	0.00
13) 4-Methylphenol-d8	7.13	113	396848	36.33	ng/ul	0.00
19) Nitrobenzene-d5	7.30	128	196393	35.97	ng/ul	0.00
22) 2-Nitrophenol-d4	7.63	143	211746	36.42	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	7.88	165	371994	33.43	ng/ul	0.00
29) 4-Chloroaniline-d4	8.09	131	525150	42.44	ng/ul	0.00
44) Dimethylphthalate-d6	9.49	166	1040480	36.57	ng/ul	0.00
47) Acenaphthylene-d8	9.64	160	1234055	35.26	ng/ul	0.00
52) 4-Nitrophenol-d4	9.90	143	152483	34.79	ng/ul	0.00
58) Fluorene-d10	10.32	176	867076	35.62	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.39	200	92889	25.60	ng/ul	0.00
71) Anthracene-d10	11.35	188	1298837	36.81	ng/ul	0.00
79) Pyrene-d10	12.73	212	1363382	36.40	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.36	264	1370189	37.06	ng/ul	0.00

Target Compounds Ovalue

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

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