

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM112119\
 Data File : BM023857.D
 Acq On : 22 Nov 2019 08:32
 Operator : JU
 Sample : PB124901BL
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK01

Quant Time: Nov 22 12:07:50 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111119MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 21 16:09:32 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.51	152	453460	20.00	ng/ul	-0.01
18) Naphthalene-d8	10.29	136	1757705	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.17	164	1018664	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.93	188	1923709	20.00	ng/ul	0.00
78) Chrysene-d12	21.13	240	1660181	20.00	ng/ul	0.00
86) Perylene-d12	23.29	264	2212322	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.08	96	103	0.01	ng/uL	0.02
5) Phenol-d5	0.00	99	0	0.00	ng/ul	
7) Bis-(2-Chloroethyl)ether-d	6.86	67	72	0.00	ng/ul	-0.01
9) 2-Chlorophenol-d4	0.00	132	0	0.00	ng/ul	
13) 4-Methylphenol-d8	0.00	113	0	0.00	ng/ul	
19) Nitrobenzene-d5	0.00	128	0	0.00	ng/ul	
22) 2-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
26) 2,4-Dichlorophenol-d3	0.00	165	0	0.00	ng/ul	
29) 4-Chloroaniline-d4	10.29	131	2855	0.09	ng/ul	-0.15
44) Dimethylphthalate-d6	0.00	166	0	0.00	ng/ul	
47) Acenaphthylene-d8	0.00	160	0	0.00	ng/ul	
52) 4-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
58) Fluorene-d10	0.00	176	0	0.00	ng/ul	
63) 4,6-Dinitro-2-methylphenol	0.00	200	0	0.00	ng/ul	
71) Anthracene-d10	16.93	188	1923709	21.08	ng/ul	-0.10
79) Pyrene-d10	19.33	212	1232	0.01	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.29	264	2188077	18.80	ng/ul	0.14

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

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