

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM111919\
 Data File : BM023824.D
 Acq On : 20 Nov 2019 08:07
 Operator : JU
 Sample : PB124868BL
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK68

Quant Time: Nov 20 10:24:02 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM11119MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 20 06:55:09 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.52	152	394669	20.00	ng/ul	0.00
18) Naphthalene-d8	10.29	136	1512516	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.17	164	888671	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.93	188	1632402	20.00	ng/ul	0.00
78) Chrysene-d12	21.13	240	1240553	20.00	ng/ul	0.00
86) Perylene-d12	23.29	264	1658362	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.06	96	303	0.03	ng/uL	0.00
5) Phenol-d5	0.00	99	0	0.00	ng/ul	
7) Bis-(2-Chloroethyl)ether-d	6.88	67	145	0.01	ng/ul	0.00
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	10.30	131	900	0.03	ng/ul	-0.14
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	1632402	21.08	ng/ul	-0.10
79) Pyrene-d10	19.34	212	3223	0.05	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.29	264	1642042	18.82	ng/ul	0.13

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

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