

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM112219\
 Data File : BM023871.D
 Acq On : 22 Nov 2019 23:46
 Operator : JU
 Sample : PB124904BL
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK04

Quant Time: Nov 23 05:24:49 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111119MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Nov 23 05:21:37 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.50	152	323988	20.00	ng/ul	0.00
18) Naphthalene-d8	10.28	136	1345567	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.17	164	840655	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.92	188	1811607	20.00	ng/ul	0.00
78) Chrysene-d12	21.13	240	1808664	20.00	ng/ul	0.00
86) Perylene-d12	23.27	264	2155955	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	3.05	96	54931	6.99	ng/uL	0.00
5) Phenol-d5	6.70	99	1058920	35.85	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.85	67	595910	35.12	ng/ul	0.00
9) 2-Chlorophenol-d4	7.05	132	843018	39.40	ng/ul	0.00
13) 4-Methylphenol-d8	8.22	113	822546	34.91	ng/ul	0.00
19) Nitrobenzene-d5	8.66	128	403754	41.92	ng/ul	0.00
22) 2-Nitrophenol-d4	9.37	143	464959	45.93	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.92	165	838390	36.32	ng/ul	0.00
29) 4-Chloroaniline-d4	10.43	131	1070796	42.25	ng/ul	0.00
44) Dimethylphthalate-d6	13.59	166	2467555	37.64	ng/ul	0.00
47) Acenaphthylene-d8	13.86	160	3036278	40.58	ng/ul	0.00
52) 4-Nitrophenol-d4	14.40	143	417041	37.27	ng/ul	0.00
58) Fluorene-d10	15.17	176	2183394	37.72	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.31	200	277547	26.49	ng/ul	0.00
71) Anthracene-d10	17.02	188	3488142	40.59	ng/ul	0.00
79) Pyrene-d10	19.33	212	3826291	40.57	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.14	264	4522170	39.86	ng/ul	0.00

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

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