

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM112219\
 Data File : BM023862.D
 Acq On : 22 Nov 2019 17:09
 Operator : JU
 Sample : PB124750BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK50

Quant Time: Nov 22 17:42:04 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM111119MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 22 15:55:14 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.51	152	420768	20.00	ng/ul	0.00
18) Naphthalene-d8	10.29	136	1688374	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.17	164	1029709	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.93	188	2235609	20.00	ng/ul	0.00
78) Chrysene-d12	21.13	240	2121008	20.00	ng/ul	0.00
86) Perylene-d12	23.28	264	2620596	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	63283	6.20	ng/uL	0.00
5) Phenol-d5	6.70	99	1172426	30.57	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.86	67	661856	30.03	ng/ul	0.00
9) 2-Chlorophenol-d4	7.05	132	947569	34.10	ng/ul	0.00
13) 4-Methylphenol-d8	8.23	113	922874	30.16	ng/ul	0.00
19) Nitrobenzene-d5	8.66	128	452904	37.48	ng/ul	0.00
22) 2-Nitrophenol-d4	9.38	143	528057	41.57	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.92	165	927516	32.02	ng/ul	0.00
29) 4-Chloroaniline-d4	10.43	131	1213302	38.15	ng/ul	0.00
44) Dimethylphthalate-d6	13.59	166	2779778	34.62	ng/ul	0.00
47) Acenaphthylene-d8	13.86	160	3318420	36.21	ng/ul	0.00
52) 4-Nitrophenol-d4	14.40	143	468632	34.19	ng/ul	0.00
58) Fluorene-d10	15.17	176	2413251	34.03	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.31	200	330533	25.56	ng/ul	0.00
71) Anthracene-d10	17.02	188	3795596	35.79	ng/ul	0.00
79) Pyrene-d10	19.33	212	4184038	37.83	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.15	264	4967942	36.03	ng/ul	0.00

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

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