

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101723\
 Data File : BM042328.D
 Acq On : 17 Oct 2023 19:37
 Operator : MA/JU
 Sample : PB156382TB 10X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLEB382

Quant Time: Oct 18 02:14:27 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM101223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 18 02:04:27 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.010	152	134306	20.000	ng/u1	0.00
20) Naphthalene-d8	10.833	136	608332	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.657	164	370673	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.404	188	799204	20.000	ng/u1	0.00
79) Chrysene-d12	21.586	240	686334	20.000	ng/u1	0.00
88) Perylene-d12	24.056	264	727854	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	31947	7.432	ng/uL	0.00
4) Pyridine-d5	3.804	84	447938	35.187	ng/u1	0.00
7) Phenol-d5	7.157	99	532722	35.791	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.351	67	371442	36.933	ng/u1	0.00
11) 2-Chlorophenol-d4	7.534	132	380806	37.191	ng/u1	0.00
15) 4-Methylphenol-d8	8.716	113	409042	34.701	ng/u1	0.00
21) Nitrobenzene-d5	9.210	128	193070	37.370	ng/u1	0.00
24) 2-Nitrophenol-d4	9.922	143	208017	37.665	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.445	165	350647	34.791	ng/u1	0.00
31) 4-Chloroaniline-d4	10.992	131	570112	35.926	ng/u1	0.00
46) Dimethylphthalate-d6	14.069	166	1230650	38.300	ng/u1	0.00
49) Acenaphthylene-d8	14.357	160	1330590	37.835	ng/u1	0.00
54) 4-Nitrophenol-d4	14.863	143	208085	32.197	ng/u1	0.00
60) Fluorene-d10	15.645	176	1028458	38.936	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.774	200	163844	29.266	ng/u1	0.00
73) Anthracene-d10	17.504	188	1630679	39.831	ng/u1	0.00
81) Pyrene-d10	19.798	212	1896464	45.337	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.897	264	1664625	41.666	ng/u1	0.00

Target Compounds Qvalue

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

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