

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101323\
 Data File : BM042286.D
 Acq On : 13 Oct 2023 16:44
 Operator : MA/JU
 Sample : PB156222BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK222

Quant Time: Oct 14 02:05:33 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM101223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 12 15:38:21 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.010	152	130887	20.000	ng/ul	0.00
20) Naphthalene-d8	10.833	136	586740	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.657	164	368253	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.409	188	797472	20.000	ng/ul	0.00
79) Chrysene-d12	21.592	240	720908	20.000	ng/ul	0.00
88) Perylene-d12	24.074	264	810177	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.363	96	27682	6.608	ng/uL	0.00
4) Pyridine-d5	3.799	84	418193	33.708	ng/ul	0.00
7) Phenol-d5	7.151	99	475979	32.814	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.345	67	337893	34.475	ng/ul	0.00
11) 2-Chlorophenol-d4	7.528	132	343137	34.387	ng/ul	0.00
15) 4-Methylphenol-d8	8.710	113	371750	32.361	ng/ul	0.00
21) Nitrobenzene-d5	9.204	128	177428	35.606	ng/ul	0.00
24) 2-Nitrophenol-d4	9.922	143	191629	35.974	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.439	165	323298	33.258	ng/ul	0.00
31) 4-Chloroaniline-d4	10.992	131	526170	34.377	ng/ul	0.00
46) Dimethylphthalate-d6	14.069	166	1169158	36.625	ng/ul	0.00
49) Acenaphthylene-d8	14.357	160	1245904	35.660	ng/ul	0.00
54) 4-Nitrophenol-d4	14.868	143	198632	30.936	ng/ul	0.00
60) Fluorene-d10	15.645	176	969318	36.938	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.774	200	154208	27.604	ng/ul	0.00
73) Anthracene-d10	17.509	188	1517246	37.140	ng/ul	0.00
81) Pyrene-d10	19.803	212	1738582	39.569	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.909	264	1682465	37.834	ng/ul	0.00

Target Compounds Qvalue

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

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