

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101723\
 Data File : BM042320.D
 Acq On : 17 Oct 2023 13:45
 Operator : MA/JU
 Sample : PB156335BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK335

Quant Time: Oct 18 01:57:13 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	156454	20.000	ng/u1	0.00
20) Naphthalene-d8	10.833	136	709905	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.651	164	437816	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.404	188	954280	20.000	ng/u1	0.00
79) Chrysene-d12	21.586	240	862208	20.000	ng/u1	-0.01
88) Perylene-d12	24.062	264	908264	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	31689	6.329	ng/uL	0.00
4) Pyridine-d5	3.804	84	477618	32.207	ng/u1	0.00
7) Phenol-d5	7.157	99	551979	31.835	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.351	67	382924	32.684	ng/u1	0.00
11) 2-Chlorophenol-d4	7.534	132	393132	32.959	ng/u1	0.00
15) 4-Methylphenol-d8	8.716	113	426908	31.089	ng/u1	0.00
21) Nitrobenzene-d5	9.210	128	198648	32.948	ng/u1	0.00
24) 2-Nitrophenol-d4	9.922	143	217170	33.696	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.445	165	368224	31.307	ng/u1	0.00
31) 4-Chloroaniline-d4	10.992	131	578130	31.219	ng/u1	0.00
46) Dimethylphthalate-d6	14.068	166	1279250	33.707	ng/u1	0.00
49) Acenaphthylene-d8	14.357	160	1398493	33.668	ng/u1	0.00
54) 4-Nitrophenol-d4	14.863	143	219434	28.746	ng/u1	-0.01
60) Fluorene-d10	15.645	176	1085401	34.790	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.768	200	160609	24.026	ng/u1	0.00
73) Anthracene-d10	17.504	188	1699842	34.773	ng/u1	0.00
81) Pyrene-d10	19.797	212	1988151	37.834	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.897	264	1795631	36.018	ng/u1	-0.02

Target Compounds Qvalue

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

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