

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101723\
 Data File : BM042321.D
 Acq On : 17 Oct 2023 15:11
 Operator : MA/JU
 Sample : PB156336BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK336

Quant Time: Oct 18 01:58:43 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	136108	20.000	ng/u1	0.00
20) Naphthalene-d8	10.834	136	606767	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.657	164	368264	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.404	188	790565	20.000	ng/u1	0.00
79) Chrysene-d12	21.586	240	691533	20.000	ng/u1	-0.01
88) Perylene-d12	24.056	264	766563	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	28304	6.498	ng/uL	0.00
4) Pyridine-d5	3.805	84	419757	32.536	ng/u1	0.00
7) Phenol-d5	7.157	99	474288	31.443	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.351	67	336406	33.006	ng/u1	0.00
11) 2-Chlorophenol-d4	7.534	132	334211	32.208	ng/u1	0.00
15) 4-Methylphenol-d8	8.716	113	364392	30.504	ng/u1	0.00
21) Nitrobenzene-d5	9.210	128	167758	32.554	ng/u1	0.00
24) 2-Nitrophenol-d4	9.922	143	180991	32.856	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.445	165	304607	30.301	ng/u1	0.00
31) 4-Chloroaniline-d4	10.992	131	490533	30.991	ng/u1	0.00
46) Dimethylphthalate-d6	14.069	166	1073743	33.635	ng/u1	0.00
49) Acenaphthylene-d8	14.357	160	1170635	33.505	ng/u1	0.00
54) 4-Nitrophenol-d4	14.863	143	181318	28.239	ng/u1	-0.01
60) Fluorene-d10	15.645	176	904177	34.455	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.768	200	132866	23.992	ng/u1	0.00
73) Anthracene-d10	17.504	188	1407484	34.755	ng/u1	0.00
81) Pyrene-d10	19.798	212	1612574	38.260	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.897	264	1520226	36.131	ng/u1	-0.02

Target Compounds	Qvalue

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

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