

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM010224\
 Data File : BM043699.D
 Acq On : 02 Jan 2024 18:06
 Operator : MA/JU
 Sample : PB158101BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK101

Quant Time: Jan 02 23:11:32 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121523.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 29 22:31:57 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.957	152	214966	20.000	ng/u1	0.00
20) Naphthalene-d8	10.769	136	869488	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.598	164	467707	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.345	188	973810	20.000	ng/u1	0.00
79) Chrysene-d12	21.521	240	872063	20.000	ng/u1	0.00
88) Perylene-d12	23.938	264	1027053	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	32216	4.871	ng/uL	0.00
4) Pyridine-d5	3.793	84	436214	22.915	ng/u1	0.00
7) Phenol-d5	7.128	99	554951	22.694	ng/u1	0.01
9) Bis-(2-Chloroethyl)eth...	7.298	67	335051	21.634	ng/u1	0.00
11) 2-Chlorophenol-d4	7.487	132	430918	23.079	ng/u1	0.00
15) 4-Methylphenol-d8	8.675	113	411290	21.674	ng/u1	0.00
21) Nitrobenzene-d5	9.134	128	203543	24.196	ng/u1	0.00
24) 2-Nitrophenol-d4	9.851	143	211293	24.309	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.392	165	363749	23.706	ng/u1	0.01
31) 4-Chloroaniline-d4	10.916	131	547980	24.275	ng/u1	0.00
46) Dimethylphthalate-d6	14.016	166	941379	22.359	ng/u1	0.00
49) Acenaphthylene-d8	14.292	160	1210112	24.229	ng/u1	0.00
54) 4-Nitrophenol-d4	14.816	143	199896	25.085	ng/u1	0.02
60) Fluorene-d10	15.586	176	859359	24.683	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.716	200	96979	15.058	ng/u1	0.00
73) Anthracene-d10	17.439	188	1364018	24.976	ng/u1	0.00
81) Pyrene-d10	19.733	212	1505901	21.873	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.780	264	1565621	24.702	ng/u1	0.00

Target Compounds Qvalue

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

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