

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101724\
 Data File : BM048113.D
 Acq On : 17 Oct 2024 19:44
 Operator : RC/JU
 Sample : PB164043BL
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK043

Quant Time: Oct 17 23:25:16 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM101724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 17 15:44:59 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.769	152	225925	20.000	ng/u1	0.00
20) Naphthalene-d8	10.563	136	880004	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.404	164	526779	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.145	188	1023645	20.000	ng/u1	0.00
79) Chrysene-d12	21.374	240	925949	20.000	ng/u1	0.00
88) Perylene-d12	24.362	264	1068396	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	38423	6.458	ng/uL	0.00
4) Pyridine-d5	3.651	84	517554	32.329	ng/u1	0.00
7) Phenol-d5	6.945	99	589698	32.223	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.104	67	358033	32.936	ng/u1	0.00
11) 2-Chlorophenol-d4	7.304	132	515180	32.547	ng/u1	0.00
15) 4-Methylphenol-d8	8.481	113	461785	32.421	ng/u1	0.00
21) Nitrobenzene-d5	8.928	128	235986	31.778	ng/u1	0.00
24) 2-Nitrophenol-d4	9.645	143	264312	32.165	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.186	165	426022	29.436	ng/u1	0.00
31) 4-Chloroaniline-d4	10.698	131	628628	33.186	ng/u1	0.00
46) Dimethylphthalate-d6	13.816	166	1239422	32.395	ng/u1	0.00
49) Acenaphthylene-d8	14.098	160	1452751	31.934	ng/u1	0.00
54) 4-Nitrophenol-d4	14.621	143	174774	24.714	ng/u1	0.00
60) Fluorene-d10	15.398	176	1061814	32.507	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.527	200	173833	26.255	ng/u1	0.00
73) Anthracene-d10	17.245	188	1618200	33.332	ng/u1	0.00
81) Pyrene-d10	19.533	212	1799769	31.809	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.156	264	1868274	32.740	ng/u1	0.00

Target Compounds Qvalue

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

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