

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101724\
 Data File : BM048109.D
 Acq On : 17 Oct 2024 16:27
 Operator : RC/JU
 Sample : PB164148BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK148

Quant Time: Oct 17 23:40:24 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	7.775	152	222943	20.000	ng/u1	0.00
20) Naphthalene-d8	10.563	136	851497	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.410	164	506346	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.145	188	988052	20.000	ng/u1	0.00
79) Chrysene-d12	21.374	240	908589	20.000	ng/u1	0.00
88) Perylene-d12	24.356	264	1025923	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	38307	6.524	ng/uL	0.00
4) Pyridine-d5	3.651	84	489725	31.000	ng/u1	0.00
7) Phenol-d5	6.945	99	558972	30.953	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.104	67	340238	31.718	ng/u1	0.00
11) 2-Chlorophenol-d4	7.304	132	488179	31.254	ng/u1	0.00
15) 4-Methylphenol-d8	8.481	113	439764	31.288	ng/u1	0.00
21) Nitrobenzene-d5	8.928	128	229311	31.913	ng/u1	0.00
24) 2-Nitrophenol-d4	9.645	143	250892	31.554	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.192	165	406192	29.006	ng/u1	0.00
31) 4-Chloroaniline-d4	10.698	131	594417	32.430	ng/u1	0.00
46) Dimethylphthalate-d6	13.816	166	1169270	31.794	ng/u1	0.00
49) Acenaphthylene-d8	14.098	160	1380837	31.578	ng/u1	0.00
54) 4-Nitrophenol-d4	14.621	143	162940	23.971	ng/u1	0.00
60) Fluorene-d10	15.398	176	1002555	31.931	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.527	200	166058	25.984	ng/u1	0.00
73) Anthracene-d10	17.245	188	1538227	32.826	ng/u1	0.00
81) Pyrene-d10	19.539	212	1718830	30.959	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.156	264	1783308	32.544	ng/u1	0.00

Target Compounds Qvalue

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

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