

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM111124\
 Data File : BM048579.D
 Acq On : 11 Nov 2024 10:52
 Operator : RC/JU
 Sample : PB164746BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK746

Quant Time: Nov 11 23:14:55 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM110724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 07 22:50:20 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.663	152	136111	20.000	ng/u1	0.00
20) Naphthalene-d8	10.445	136	553138	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.310	164	337002	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.057	188	677038	20.000	ng/u1	0.00
79) Chrysene-d12	21.280	240	652442	20.000	ng/u1	0.00
88) Perylene-d12	24.191	264	773102	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.122	96	19617	5.746	ng/uL	0.00
4) Pyridine-d5	3.540	84	262079	27.853	ng/u1	0.00
7) Phenol-d5	6.840	99	291704	28.153	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.998	67	176297	28.328	ng/u1	0.00
11) 2-Chlorophenol-d4	7.198	132	251923	29.352	ng/u1	0.00
15) 4-Methylphenol-d8	8.375	113	227477	27.921	ng/u1	0.00
21) Nitrobenzene-d5	8.816	128	116056	28.948	ng/u1	0.00
24) 2-Nitrophenol-d4	9.539	143	124640	29.229	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.081	165	205362	25.924	ng/u1	0.00
31) 4-Chloroaniline-d4	10.592	131	313356	27.849	ng/u1	0.00
46) Dimethylphthalate-d6	13.727	166	632208	29.388	ng/u1	0.00
49) Acenaphthylene-d8	13.998	160	742944	29.292	ng/u1	0.00
54) 4-Nitrophenol-d4	14.551	143	74391	18.618	ng/u1	0.02
60) Fluorene-d10	15.304	176	552961	29.447	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.439	200	64490	18.935	ng/u1	0.00
73) Anthracene-d10	17.157	188	867540	30.280	ng/u1	0.00
81) Pyrene-d10	19.451	212	1014779	30.992	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.991	264	1081349	29.936	ng/u1	0.00

Target Compounds	Qvalue

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

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