

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110824\
 Data File : BM048570.D
 Acq On : 09 Nov 2024 17:52
 Operator : RC/JU
 Sample : PB164726BL
 Misc :
 ALS Vial : 52 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK726

Quant Time: Nov 09 22:42:31 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	226811	20.000	ng/u1	0.00
20) Naphthalene-d8	10.445	136	932810	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.310	164	575562	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.057	188	1151800	20.000	ng/u1	0.00
79) Chrysene-d12	21.280	240	1065207	20.000	ng/u1	0.00
88) Perylene-d12	24.186	264	1245926	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.128	96	29788	5.236	ng/uL	0.00
4) Pyridine-d5	3.546	84	417132	26.604	ng/u1	0.00
7) Phenol-d5	6.846	99	500373	28.981	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.998	67	296885	28.628	ng/u1	0.00
11) 2-Chlorophenol-d4	7.198	132	428980	29.994	ng/u1	0.00
15) 4-Methylphenol-d8	8.381	113	396854	29.232	ng/u1	0.00
21) Nitrobenzene-d5	8.816	128	201238	29.765	ng/u1	0.00
24) 2-Nitrophenol-d4	9.539	143	224735	31.251	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.081	165	385990	28.893	ng/u1	0.00
31) 4-Chloroaniline-d4	10.592	131	515020	27.141	ng/u1	0.00
46) Dimethylphthalate-d6	13.722	166	1097841	29.881	ng/u1	0.00
49) Acenaphthylene-d8	13.998	160	1294273	29.879	ng/u1	0.00
54) 4-Nitrophenol-d4	14.533	143	172897	25.336	ng/u1	0.00
60) Fluorene-d10	15.304	176	956632	29.828	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.439	200	138821	23.958	ng/u1	0.00
73) Anthracene-d10	17.157	188	1491813	30.607	ng/u1	0.00
81) Pyrene-d10	19.451	212	1739596	32.541	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.991	264	1822630	31.309	ng/u1	0.00

Target Compounds	Qvalue

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

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