

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM013025\
 Data File : BM049526.D
 Acq On : 30 Jan 2025 14:23
 Operator : RC/JU
 Sample : PB166284BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK284

Quant Time: Jan 30 15:09:46 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM012825.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jan 28 14:45:18 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.504	152	271168	20.000	ng/u1	0.00
20) Naphthalene-d8	10.269	136	966233	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.145	164	606502	20.000	ng/u1	0.00
64) Phenanthrene-d10	16.904	188	1177211	20.000	ng/u1	0.00
79) Chrysene-d12	21.133	240	1080408	20.000	ng/u1	0.00
88) Perylene-d12	23.909	264	1140229	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.116	96	42936	5.867	ng/uL	0.00
4) Pyridine-d5	3.504	84	645136	30.002	ng/u1	0.00
7) Phenol-d5	6.728	99	697802	31.712	ng/u1	0.02
9) Bis-(2-Chloroethyl)eth...	6.857	67	449714	29.888	ng/u1	0.00
11) 2-Chlorophenol-d4	7.057	132	559551	32.861	ng/u1	0.00
15) 4-Methylphenol-d8	8.239	113	508583	31.546	ng/u1	0.01
21) Nitrobenzene-d5	8.651	128	239522	32.403	ng/u1	0.00
24) 2-Nitrophenol-d4	9.363	143	273137	33.876	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	9.916	165	514263	31.255	ng/u1	0.01
31) 4-Chloroaniline-d4	10.416	131	606087	28.717	ng/u1	0.00
46) Dimethylphthalate-d6	13.574	166	1365515	30.707	ng/u1	0.00
49) Acenaphthylene-d8	13.839	160	1636895	31.329	ng/u1	0.00
54) 4-Nitrophenol-d4	14.421	143	182471	25.014	ng/u1	0.04
60) Fluorene-d10	15.151	176	1223073	31.275	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.280	200	181359	24.448	ng/u1	0.00
73) Anthracene-d10	16.998	188	1878216	31.950	ng/u1	0.00
81) Pyrene-d10	19.309	212	2122850	33.611	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.721	264	1983523	32.616	ng/u1	0.00

Target Compounds	Qvalue

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

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