

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM013025\
 Data File : BM049547.D
 Acq On : 31 Jan 2025 04:57
 Operator : RC/JU
 Sample : PB166286BL
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK286

Quant Time: Jan 31 09:05:13 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	7.510	152	254017	20.000	ng/u1	0.00
20) Naphthalene-d8	10.269	136	957517	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.151	164	645324	20.000	ng/u1	0.00
64) Phenanthrene-d10	16.904	188	1400671	20.000	ng/u1	0.00
79) Chrysene-d12	21.139	240	1502058	20.000	ng/u1	0.00
88) Perylene-d12	23.921	264	1610011	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.110	96	36513	5.326	ng/uL	-0.01
4) Pyridine-d5	3.499	84	572655	28.429	ng/u1	-0.01
7) Phenol-d5	6.722	99	659087	31.974	ng/u1	0.01
9) Bis-(2-Chloroethyl)eth...	6.857	67	401136	28.459	ng/u1	0.00
11) 2-Chlorophenol-d4	7.057	132	525536	32.947	ng/u1	0.00
15) 4-Methylphenol-d8	8.234	113	487877	32.305	ng/u1	0.00
21) Nitrobenzene-d5	8.651	128	231269	31.571	ng/u1	0.00
24) 2-Nitrophenol-d4	9.363	143	266512	33.355	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	9.916	165	505371	30.994	ng/u1	0.01
31) 4-Chloroaniline-d4	10.416	131	600591	28.716	ng/u1	0.00
46) Dimethylphthalate-d6	13.574	166	1417816	29.965	ng/u1	0.00
49) Acenaphthylene-d8	13.839	160	1680676	30.231	ng/u1	0.00
54) 4-Nitrophenol-d4	14.421	143	217636	28.039	ng/u1	0.04
60) Fluorene-d10	15.151	176	1282091	30.812	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.286	200	181608	20.576	ng/u1	0.00
73) Anthracene-d10	17.004	188	2111800	30.192	ng/u1	0.00
81) Pyrene-d10	19.309	212	2589553	29.491	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.733	264	2674498	31.146	ng/u1	0.00

Target Compounds Qvalue

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

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