

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012225\
 Data File : BM049369.D
 Acq On : 21 Jan 2025 17:44
 Operator : RC/JU
 Sample : PB166156BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK156

Quant Time: Jan 22 00:22:22 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM011325.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Jan 13 16:56:06 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.522	152	168732	20.000	ng/ul	0.00
20) Naphthalene-d8	10.281	136	589973	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.163	164	343327	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.910	188	661412	20.000	ng/ul	0.00
79) Chrysene-d12	21.139	240	580212	20.000	ng/ul	0.00
88) Perylene-d12	23.921	264	633377	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.122	96	28154	6.904	ng/uL	0.00
4) Pyridine-d5	3.510	84	415925	32.037	ng/ul	0.00
7) Phenol-d5	6.716	99	413379	29.185	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.869	67	290796	30.832	ng/ul	0.00
11) 2-Chlorophenol-d4	7.063	132	333632	30.791	ng/ul	0.00
15) 4-Methylphenol-d8	8.234	113	298195	27.149	ng/ul	0.00
21) Nitrobenzene-d5	8.663	128	137910	30.654	ng/ul	0.00
24) 2-Nitrophenol-d4	9.381	143	144372	28.705	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.916	165	283553	27.629	ng/ul	0.00
31) 4-Chloroaniline-d4	10.428	131	344553	25.514	ng/ul	0.00
46) Dimethylphthalate-d6	13.580	166	780610	30.548	ng/ul	0.00
49) Acenaphthylene-d8	13.845	160	947669	32.486	ng/ul	0.00
54) 4-Nitrophenol-d4	14.392	143	62906	14.866	ng/ul	0.00
60) Fluorene-d10	15.163	176	714630	32.373	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.286	200	76545	18.159	ng/ul	0.00
73) Anthracene-d10	17.010	188	1086234	33.222	ng/ul	0.00
81) Pyrene-d10	19.315	212	1226317	34.729	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.733	264	1090012	32.059	ng/ul	-0.01

Target Compounds Qvalue

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

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