

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012225\
 Data File : BM049410.D
 Acq On : 23 Jan 2025 09:02
 Operator : RC/JU
 Sample : PB166176BL
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK176

Quant Time: Jan 23 22:08:06 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	7.516	152	341919	20.000	ng/ul	0.00
20) Naphthalene-d8	10.280	136	1250411	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.157	164	765209	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.909	188	1460113	20.000	ng/ul	0.00
79) Chrysene-d12	21.139	240	1266724	20.000	ng/ul	0.00
88) Perylene-d12	23.915	264	1314737	20.000	ng/ul	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.122	96	52519	6.355	ng/uL	0.00
4) Pyridine-d5	3.510	84	765871	29.112	ng/ul	0.00
7) Phenol-d5	6.710	99	858148	29.898	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.869	67	569285	29.787	ng/ul	0.00
11) 2-Chlorophenol-d4	7.057	132	691413	31.490	ng/ul	0.00
15) 4-Methylphenol-d8	8.228	113	636643	28.604	ng/ul	-0.01
21) Nitrobenzene-d5	8.663	128	320042	33.564	ng/ul	0.00
24) 2-Nitrophenol-d4	9.375	143	343479	32.223	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.910	165	649931	29.880	ng/ul	0.00
31) 4-Chloroaniline-d4	10.422	131	791170	27.642	ng/ul	0.00
46) Dimethylphthalate-d6	13.580	166	1815248	31.873	ng/ul	0.00
49) Acenaphthylene-d8	13.845	160	2164086	33.285	ng/ul	0.00
54) 4-Nitrophenol-d4	14.386	143	205248	21.763	ng/ul	0.00
60) Fluorene-d10	15.157	176	1590606	32.329	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.286	200	137815	14.810	ng/ul	0.00
73) Anthracene-d10	17.009	188	2426056	33.612	ng/ul	0.00
81) Pyrene-d10	19.315	212	2670632	34.642	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.727	264	2398243	33.981	ng/ul	-0.02

Target Compounds Qvalue

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

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