

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM121324\
 Data File : BM048979.D
 Acq On : 13 Dec 2024 13:08
 Operator : RC/JU
 Sample : PB165149BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK149

Quant Time: Dec 14 00:29:06 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM121124.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 12 00:53:37 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.569	152	181320	20.000	ng/u1	0.00
20) Naphthalene-d8	10.334	136	690048	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.204	164	371264	20.000	ng/u1	0.00
64) Phenanthrene-d10	16.951	188	688908	20.000	ng/u1	0.00
79) Chrysene-d12	21.174	240	553181	20.000	ng/u1	0.00
88) Perylene-d12	23.980	264	607122	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.152	96	35672	6.924	ng/uL	0.00
4) Pyridine-d5	3.546	84	505080	32.963	ng/u1	0.00
7) Phenol-d5	6.757	99	534936	33.844	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.916	67	377629	35.179	ng/u1	0.00
11) 2-Chlorophenol-d4	7.110	132	414516	36.142	ng/u1	0.00
15) 4-Methylphenol-d8	8.275	113	377583	32.002	ng/u1	0.00
21) Nitrobenzene-d5	8.716	128	184352	37.252	ng/u1	0.00
24) 2-Nitrophenol-d4	9.428	143	161482	35.297	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	9.963	165	328634	33.752	ng/u1	0.00
31) 4-Chloroaniline-d4	10.475	131	487356	32.104	ng/u1	0.00
46) Dimethylphthalate-d6	13.627	166	968025	37.868	ng/u1	0.00
49) Acenaphthylene-d8	13.892	160	1158037	37.111	ng/u1	0.00
54) 4-Nitrophenol-d4	14.410	143	124777	25.475	ng/u1	0.00
60) Fluorene-d10	15.204	176	826424	37.301	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.321	200	76735	21.351	ng/u1	0.00
73) Anthracene-d10	17.051	188	1231993	36.489	ng/u1	0.00
81) Pyrene-d10	19.351	212	1385267	44.604	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.797	264	1186551	38.144	ng/u1	0.00

Target Compounds	Qvalue

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

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