

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012725\
 Data File : BM049452.D
 Acq On : 27 Jan 2025 17:23
 Operator : RC/JU
 Sample : PB166251BL
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK251

Quant Time: Jan 27 23:39:27 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	164384	20.000	ng/u1	0.00
20) Naphthalene-d8	10.275	136	576712	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.157	164	370092	20.000	ng/u1	0.00
64) Phenanthrene-d10	16.904	188	774804	20.000	ng/u1	-0.01
79) Chrysene-d12	21.139	240	745214	20.000	ng/u1	0.00
88) Perylene-d12	23.915	264	817078	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.116	96	26735	6.729	ng/uL	0.00
4) Pyridine-d5	3.504	84	394064	31.156	ng/u1	0.00
7) Phenol-d5	6.710	99	407239	29.512	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	6.863	67	275006	29.930	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.057	132	320315	30.344	ng/u1	0.00
15) 4-Methylphenol-d8	8.228	113	295777	27.641	ng/u1	-0.01
21) Nitrobenzene-d5	8.657	128	132915	30.223	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.375	143	140164	28.510	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	9.910	165	288426	28.750	ng/u1	0.00
31) 4-Chloroaniline-d4	10.422	131	349621	26.485	ng/u1	0.00
46) Dimethylphthalate-d6	13.574	166	843059	30.606	ng/u1	-0.01
49) Acenaphthylene-d8	13.845	160	968702	30.806	ng/u1	0.00
54) 4-Nitrophenol-d4	14.386	143	85162	18.670	ng/u1	0.00
60) Fluorene-d10	15.157	176	755174	31.735	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.286	200	85140	17.242	ng/u1	0.00
73) Anthracene-d10	17.004	188	1200529	31.344	ng/u1	-0.01
81) Pyrene-d10	19.309	212	1421783	31.349	ng/u1	-0.01
92) Benzo(a)pyrene-d12	23.727	264	1331898	30.366	ng/u1	-0.02

Target Compounds Qvalue

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

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