

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101424\
 Data File : BM048085.D
 Acq On : 16 Oct 2024 13:15
 Operator : RC/JU
 Sample : PB163982BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK982

Quant Time: Oct 16 23:33:28 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 10 05:10:14 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.775	152	277096	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.569	136	1164513	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.416	164	720113	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.157	188	1401283	20.000	ng/u1	0.00
79) Chrysene-d12	21.386	240	1251430	20.000	ng/u1	0.00
88) Perylene-d12	24.380	264	1389792	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.240	96	30978	3.613	ng/uL	0.00
4) Pyridine-d5	3.652	84	442986	17.644	ng/u1	0.00
7) Phenol-d5	6.951	99	525599	20.059	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.110	67	314937	16.938	ng/u1	0.00
11) 2-Chlorophenol-d4	7.316	132	453824	23.806	ng/u1	0.00
15) 4-Methylphenol-d8	8.487	113	417516	21.432	ng/u1	0.00
21) Nitrobenzene-d5	8.934	128	215049	23.194	ng/u1	0.00
24) 2-Nitrophenol-d4	9.657	143	241139	27.139	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.198	165	409140	23.051	ng/u1	0.00
31) 4-Chloroaniline-d4	10.704	131	587900	21.424	ng/u1	-0.01
46) Dimethylphthalate-d6	13.822	166	1170143	22.413	ng/u1	-0.01
49) Acenaphthylene-d8	14.110	160	1397976	22.584	ng/u1	0.00
54) 4-Nitrophenol-d4	14.633	143	177485	17.075	ng/u1	0.01
60) Fluorene-d10	15.410	176	1025181	22.670	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.533	200	116051	14.152	ng/u1	0.00
73) Anthracene-d10	17.257	188	1559529	22.634	ng/u1	0.00
81) Pyrene-d10	19.545	212	1737124	24.458	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.174	264	1711055	23.729	ng/u1	0.00

Target Compounds Qvalue

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

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