

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091324\
 Data File : BM047511.D
 Acq On : 13 Sep 2024 09:29
 Operator : RC/JU
 Sample : PB163349BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK349

Quant Time: Sep 14 05:21:26 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091024.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Sep 11 00:35:12 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.863	152	403239	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.663	136	1801777	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.492	164	1128216	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.227	188	2188412	20.000	ng/u1	-0.01
79) Chrysene-d12	21.451	240	1495886	20.000	ng/u1	-0.02
88) Perylene-d12	24.486	264	1404087	20.000	ng/u1	-0.04
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	55619	5.563	ng/uL	0.00
4) Pyridine-d5	3.710	84	805182	25.206	ng/u1	0.00
7) Phenol-d5	7.016	99	903354	23.899	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.193	67	605666	25.743	ng/u1	0.00
11) 2-Chlorophenol-d4	7.393	132	730887	25.389	ng/u1	-0.01
15) 4-Methylphenol-d8	8.557	113	728647	24.525	ng/u1	-0.01
21) Nitrobenzene-d5	9.016	128	377400	27.123	ng/u1	0.00
24) 2-Nitrophenol-d4	9.739	143	348162	26.486	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.275	165	680420	25.457	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.787	131	1039441	23.761	ng/u1	0.00
46) Dimethylphthalate-d6	13.898	166	2184713	26.346	ng/u1	-0.01
49) Acenaphthylene-d8	14.186	160	2622208	26.205	ng/u1	0.00
54) 4-Nitrophenol-d4	14.663	143	325741	19.826	ng/u1	-0.01
60) Fluorene-d10	15.480	176	1896675	26.469	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.586	200	184950	15.987	ng/u1	-0.01
73) Anthracene-d10	17.321	188	2908558	27.600	ng/u1	-0.01
81) Pyrene-d10	19.604	212	3075781	35.280	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.280	264	1990131	26.673	ng/u1	-0.03

Target Compounds Qvalue

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

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