

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080924\
 Data File : BM047087.D
 Acq On : 08 Aug 2024 15:25
 Operator : RC/JU
 Sample : PB162538BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK538

Quant Time: Aug 08 23:18:52 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Aug 07 04:48:25 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.387	152	230343	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.145	136	887976	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.039	164	541951	20.000	ng/u1	-0.01
64) Phenanthrene-d10	16.798	188	1078386	20.000	ng/u1	#-0.02
79) Chrysene-d12	21.033	240	901192	20.000	ng/u1	#-0.02
88) Perylene-d12	23.738	264	1024483	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.016	96	34908	6.004	ng/uL	0.00
4) Pyridine-d5	3.405	84	479248	30.219	ng/u1	0.00
7) Phenol-d5	6.593	99	561174	32.071	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	6.745	67	369743	32.114	ng/u1	-0.01
11) 2-Chlorophenol-d4	6.928	132	446843	30.147	ng/u1	-0.01
15) 4-Methylphenol-d8	8.104	113	428987	30.948	ng/u1	-0.02
21) Nitrobenzene-d5	8.534	128	219532	30.809	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.245	143	237774	30.557	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	9.781	165	429953	28.779	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.292	131	596373	30.373	ng/u1	-0.01
46) Dimethylphthalate-d6	13.469	166	1208947	29.175	ng/u1	-0.01
49) Acenaphthylene-d8	13.727	160	1410222	30.488	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.280	143	208889	26.072	ng/u1	-0.02
60) Fluorene-d10	15.045	176	1031391	29.096	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.186	200	162070	23.788	ng/u1	-0.01
73) Anthracene-d10	16.898	188	1585376	31.011	ng/u1	-0.01
81) Pyrene-d10	19.209	212	1694430	32.845	ng/u1	-0.02
92) Benzo(a)pyrene-d12	23.556	264	1665956	30.882	ng/u1	-0.02

Target Compounds Qvalue

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

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