

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100724\
 Data File : BM047913.D
 Acq On : 08 Oct 2024 14:29
 Operator : RC/JU
 Sample : PB163712BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK712

Quant Time: Oct 08 23:38:59 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 04 08:32:31 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.793	152	295965	20.000	ng/u1	0.00
20) Naphthalene-d8	10.586	136	1162907	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.433	164	720105	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.168	188	1467576	20.000	ng/u1	0.00
79) Chrysene-d12	21.398	240	1350928	20.000	ng/u1	0.00
88) Perylene-d12	24.397	264	1502640	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.252	96	33462	3.654	ng/uL	0.00
4) Pyridine-d5	3.663	84	476653	17.775	ng/u1	0.00
7) Phenol-d5	6.963	99	528864	18.897	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.122	67	340339	17.138	ng/u1	0.00
11) 2-Chlorophenol-d4	7.328	132	442742	21.744	ng/u1	0.00
15) 4-Methylphenol-d8	8.498	113	409953	19.702	ng/u1	0.00
21) Nitrobenzene-d5	8.945	128	206631	22.316	ng/u1	0.00
24) 2-Nitrophenol-d4	9.669	143	218358	24.609	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.210	165	391952	22.113	ng/u1	0.00
31) 4-Chloroaniline-d4	10.722	131	570270	20.810	ng/u1	0.00
46) Dimethylphthalate-d6	13.839	166	1166290	22.339	ng/u1	0.00
49) Acenaphthylene-d8	14.122	160	1342576	21.689	ng/u1	0.00
54) 4-Nitrophenol-d4	14.627	143	175732	16.906	ng/u1	0.00
60) Fluorene-d10	15.421	176	999941	22.112	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.539	200	136416	15.884	ng/u1	0.00
73) Anthracene-d10	17.268	188	1590243	22.037	ng/u1	0.00
81) Pyrene-d10	19.556	212	1812593	23.641	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.191	264	1809442	23.209	ng/u1	0.00

Target Compounds Qvalue

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

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