

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM101424\
 Data File : BM048078.D
 Acq On : 16 Oct 2024 07:37
 Operator : RC/JU
 Sample : PB164081BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK081

Quant Time: Oct 16 08:12:23 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	7.781	152	524659	20.000	ng/u1	0.00
20) Naphthalene-d8	10.569	136	2210358	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.416	164	1387813	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.162	188	2736026	20.000	ng/u1	0.00
79) Chrysene-d12	21.392	240	2356142	20.000	ng/u1	0.00
88) Perylene-d12	24.385	264	2579191	20.000	ng/u1	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.240	96	57951	3.569	ng/uL	0.00
4) Pyridine-d5	3.652	84	821430	17.280	ng/u1	0.00
7) Phenol-d5	6.951	99	985898	19.872	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.110	67	595566	16.917	ng/u1	0.00
11) 2-Chlorophenol-d4	7.316	132	845423	23.422	ng/u1	0.00
15) 4-Methylphenol-d8	8.492	113	783595	21.244	ng/u1	0.00
21) Nitrobenzene-d5	8.934	128	417976	23.750	ng/u1	0.00
24) 2-Nitrophenol-d4	9.657	143	471647	27.966	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.198	165	768962	22.824	ng/u1	0.00
31) 4-Chloroaniline-d4	10.710	131	1112215	21.354	ng/u1	0.00
46) Dimethylphthalate-d6	13.827	166	2261289	22.474	ng/u1	0.00
49) Acenaphthylene-d8	14.110	160	2645930	22.179	ng/u1	0.00
54) 4-Nitrophenol-d4	14.627	143	382314	19.084	ng/u1	0.00
60) Fluorene-d10	15.410	176	1918946	22.018	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.533	200	313925	19.607	ng/u1	0.00
73) Anthracene-d10	17.256	188	2888402	21.470	ng/u1	0.00
81) Pyrene-d10	19.550	212	3258318	24.366	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.185	264	3084117	23.047	ng/u1	0.02

Target Compounds Qvalue

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

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