

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100724\
 Data File : BM047932.D
 Acq On : 09 Oct 2024 04:17
 Operator : RC/JU
 Sample : PB163712BL
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK712

Quant Time: Oct 09 07:21:27 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	7.792	152	264321	20.000	ng/u1	0.00
20) Naphthalene-d8	10.586	136	1047844	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.427	164	643924	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.168	188	1308123	20.000	ng/u1	0.00
79) Chrysene-d12	21.392	240	1221964	20.000	ng/u1	0.00
88) Perylene-d12	24.385	264	1393242	20.000	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.246	96	31495	3.851	ng/uL	0.00
4) Pyridine-d5	3.663	84	436568	18.229	ng/u1	0.00
7) Phenol-d5	6.963	99	489836	19.598	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.122	67	302612	17.062	ng/u1	0.00
11) 2-Chlorophenol-d4	7.328	132	412567	22.688	ng/u1	0.00
15) 4-Methylphenol-d8	8.498	113	380849	20.494	ng/u1	0.00
21) Nitrobenzene-d5	8.945	128	190127	22.789	ng/u1	0.00
24) 2-Nitrophenol-d4	9.669	143	202036	25.270	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.204	165	368527	23.074	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.716	131	536543	21.730	ng/u1	-0.01
46) Dimethylphthalate-d6	13.833	166	1069609	22.911	ng/u1	-0.01
49) Acenaphthylene-d8	14.121	160	1256767	22.705	ng/u1	0.00
54) 4-Nitrophenol-d4	14.627	143	163181	17.556	ng/u1	0.00
60) Fluorene-d10	15.421	176	939639	23.237	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.533	200	119369	15.594	ng/u1	0.00
73) Anthracene-d10	17.262	188	1484831	23.084	ng/u1	-0.01
81) Pyrene-d10	19.551	212	1693016	24.412	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.180	264	1756401	24.297	ng/u1	-0.01

Target Compounds Qvalue

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

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