

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030723\
 Data File : BM038891.D
 Acq On : 07 Mar 2023 11:42
 Operator : CG/JU
 Sample : PB151199BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK199

Quant Time: Mar 07 12:21:04 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM022823.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Mar 06 23:57:30 2023
 Response via : Initial Calibration

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

Internal Standards					
1) 1,4-Dichlorobenzene-d4	8.004	152	249028	20.000 ng/u1	0.00
20) Naphthalene-d8	10.816	136	1114819	20.000 ng/u1	0.00
38) Acenaphthene-d10	14.639	164	696553	20.000 ng/u1	0.00
64) Phenanthrene-d10	17.380	188	1484176	20.000 ng/u1	0.00
79) Chrysene-d12	21.562	240	1422316	20.000 ng/u1	0.00
88) Perylene-d12	24.015	264	1736989	20.000 ng/u1	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.375	96	40603	6.088 ng/uL	0.00
4) Pyridine-d5	3.799	84	571578	31.172 ng/u1	0.00
7) Phenol-d5	7.151	99	675098	28.916 ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.328	67	440589	29.535 ng/u1	0.00
11) 2-Chlorophenol-d4	7.528	132	533707	30.678 ng/u1	0.00
15) 4-Methylphenol-d8	8.704	113	521611	28.359 ng/u1	0.00
21) Nitrobenzene-d5	9.169	128	259031	33.097 ng/u1	0.00
24) 2-Nitrophenol-d4	9.892	143	258767	35.286 ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.434	165	485957	28.517 ng/u1	0.00
31) 4-Chloroaniline-d4	10.951	131	786101	27.675 ng/u1	0.00
46) Dimethylphthalate-d6	14.051	166	1605234	29.754 ng/u1	0.00
49) Acenaphthylene-d8	14.333	160	1826979	28.974 ng/u1	0.00
54) 4-Nitrophenol-d4	14.822	143	263842	24.430 ng/u1	0.00
60) Fluorene-d10	15.627	176	1410405	29.642 ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.739	200	174246	22.583 ng/u1	0.00
73) Anthracene-d10	17.480	188	2235755	30.550 ng/u1	0.00
81) Pyrene-d10	19.768	212	2595130	33.174 ng/u1	0.00
92) Benzo(a)pyrene-d12	23.856	264	2780644	31.766 ng/u1	0.00

Target Compounds Qvalue

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

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