

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM013123\
 Data File : BM038639.D
 Acq On : 01 Feb 2023 03:46
 Operator : CG/JU
 Sample : PB150491BL
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK491

Quant Time: Feb 01 05:37:57 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM013023.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jan 31 01:02:54 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.951	152	44104	20.000	ng/ul	0.00
20) Naphthalene-d8	10.757	136	127438	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.586	164	102507	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.327	188	236768	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.503	240	206081	20.000	ng/ul	# 0.00
88) Perylene-d12	23.909	264	293728	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.334	96	5812	4.699	ng/uL	0.00
4) Pyridine-d5	3.763	84	73281	23.659	ng/ul	0.00
7) Phenol-d5	7.116	99	76578	26.012	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	46265	26.924	ng/ul	0.00
11) 2-Chlorophenol-d4	7.481	132	68057	27.876	ng/ul	0.00
15) 4-Methylphenol-d8	8.669	113	61622	26.234	ng/ul	0.00
21) Nitrobenzene-d5	9.128	128	30257	31.322	ng/ul	0.00
24) 2-Nitrophenol-d4	9.845	143	43338	33.972	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.381	165	94403	33.437	ng/ul	0.00
31) 4-Chloroaniline-d4	10.910	131	75370	27.470	ng/ul	0.00
46) Dimethylphthalate-d6	13.998	166	266816	31.510	ng/ul	0.00
49) Acenaphthylene-d8	14.280	160	281252	31.212	ng/ul	0.00
54) 4-Nitrophenol-d4	14.798	143	23147	19.760	ng/ul	0.00
60) Fluorene-d10	15.574	176	242312	31.999	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.704	200	52042	25.678	ng/ul	0.00
73) Anthracene-d10	17.427	188	375502	31.507	ng/ul	0.00
81) Pyrene-d10	19.715	212	407185	38.346	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.750	264	495043	32.835	ng/ul	-0.01

Target Compounds Qvalue

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

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