

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM122022\
 Data File : BM038163.D
 Acq On : 21 Dec 2022 09:44
 Operator : CG/JU
 Sample : PB149779BL
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Dec 21 23:30:15 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 21 00:11:56 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.039	152	157396	20.000	ng/ul	0.00
20) Naphthalene-d8	10.857	136	693369	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.668	164	442180	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.409	188	959876	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.574	240	751519	20.000	ng/ul	0.00
88) Perylene-d12	24.027	264	722771	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.399	96	19909	5.743	ng/uL	0.00
4) Pyridine-d5	3.828	84	282808	27.502	ng/ul	0.00
7) Phenol-d5	7.198	99	369600	28.897	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.369	67	210148	26.983	ng/ul	0.00
11) 2-Chlorophenol-d4	7.569	132	330394	31.490	ng/ul	0.00
15) 4-Methylphenol-d8	8.751	113	318505	31.902	ng/ul	0.00
21) Nitrobenzene-d5	9.216	128	164135	29.898	ng/ul	0.00
24) 2-Nitrophenol-d4	9.939	143	192894	33.529	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.475	165	348469	31.889	ng/ul	0.00
31) 4-Chloroaniline-d4	10.998	131	476443	31.586	ng/ul	0.00
46) Dimethylphthalate-d6	14.080	166	1033697	32.790	ng/ul	0.00
49) Acenaphthylene-d8	14.363	160	1210044	31.406	ng/ul	0.00
54) 4-Nitrophenol-d4	14.868	143	173574	32.307	ng/ul	0.00
60) Fluorene-d10	15.657	176	915036	32.580	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.774	200	121131	23.604	ng/ul	0.00
73) Anthracene-d10	17.504	188	1475800	32.943	ng/ul	0.00
81) Pyrene-d10	19.792	212	1639641	36.320	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.868	264	1216100	33.686	ng/ul	0.00

Target Compounds Qvalue

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

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