

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM122022\
 Data File : BM038182.D
 Acq On : 21 Dec 2022 21:52
 Operator : CG/JU
 Sample : N6032-10
 Misc :
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Dec 21 23:34:17 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	8.040	152	160278	20.000	ng/ul	0.00
20) Naphthalene-d8	10.857	136	678338	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.669	164	408506	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.404	188	856755	20.000	ng/ul	0.00
79) Chrysene-d12	21.574	240	630890	20.000	ng/ul	0.00
88) Perylene-d12	24.021	264	629728	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.358	96	740113	209.665	ng/uL	-0.04
4) Pyridine-d5	3.840	84	62618	5.980	ng/ul	0.00
7) Phenol-d5	7.199	99	68479	5.258	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.363	67	197425	24.894	ng/ul	0.00
11) 2-Chlorophenol-d4	7.569	132	153718	14.387	ng/ul	0.00
15) 4-Methylphenol-d8	8.751	113	107971	10.620	ng/ul	0.00
21) Nitrobenzene-d5	9.210	128	158770	29.562	ng/ul	0.00
24) 2-Nitrophenol-d4	9.934	143	113234	20.118	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.475	165	188807	17.661	ng/ul	0.00
31) 4-Chloroaniline-d4	10.992	131	299542	20.298	ng/ul	-0.01
46) Dimethylphthalate-d6	14.075	166	890219	30.567	ng/ul	0.00
49) Acenaphthylene-d8	14.363	160	1122944	31.548	ng/ul	0.00
54) 4-Nitrophenol-d4	14.869	143	24534	4.943	ng/ul	0.00
60) Fluorene-d10	15.657	176	889963	34.299	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.774	200	52488	11.459	ng/ul	0.00
73) Anthracene-d10	17.504	188	1500018	37.514	ng/ul	0.00
81) Pyrene-d10	19.792	212	1551604	40.942	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.862	264	1120649	35.629	ng/ul	0.00

Target Compounds Qvalue

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

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