

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM122022\
 Data File : BM038167.D
 Acq On : 21 Dec 2022 12:10
 Operator : CG/JU
 Sample : N6014-13
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :

Quant Time: Dec 21 23:30:43 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	202408	20.000	ng/ul	0.00
20) Naphthalene-d8	10.857	136	858688	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.669	164	515739	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.410	188	1067040	20.000	ng/ul	0.00
79) Chrysene-d12	21.574	240	775186	20.000	ng/ul	0.00
88) Perylene-d12	24.027	264	758581	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	16628	3.730	ng/uL	0.00
4) Pyridine-d5	3.834	84	206147	15.589	ng/ul	0.00
7) Phenol-d5	7.198	99	300979	18.299	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.369	67	175237	17.497	ng/ul	0.00
11) 2-Chlorophenol-d4	7.569	132	269032	19.939	ng/ul	0.00
15) 4-Methylphenol-d8	8.751	113	244999	19.082	ng/ul	0.00
21) Nitrobenzene-d5	9.216	128	132923	19.551	ng/ul	0.00
24) 2-Nitrophenol-d4	9.939	143	151739	21.297	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.475	165	279307	20.639	ng/ul	0.00
31) 4-Chloroaniline-d4	10.998	131	336345	18.005	ng/ul	0.00
46) Dimethylphthalate-d6	14.080	166	798069	21.705	ng/ul	0.00
49) Acenaphthylene-d8	14.363	160	923693	20.555	ng/ul	0.00
54) 4-Nitrophenol-d4	14.869	143	112947	18.024	ng/ul	0.00
60) Fluorene-d10	15.657	176	697910	21.305	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.774	200	62650	10.982	ng/ul	0.00
73) Anthracene-d10	17.504	188	1044482	20.974	ng/ul	0.00
81) Pyrene-d10	19.792	212	1091414	23.438	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.868	264	795781	21.003	ng/ul	0.00

Target Compounds Qvalue

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

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