

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM031624\
 Data File : BM044641.D
 Acq On : 15 Mar 2024 20:05
 Operator : MA/JU
 Sample : P1759-08
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 E0738

Quant Time: Mar 15 21:44:22 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030524.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Mar 15 11:12:49 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.698	152	181563	20.000	ng/ul	0.00
20) Naphthalene-d8	10.475	136	750607	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.339	164	449509	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.098	188	918277	20.000	ng/ul	0.00
79) Chrysene-d12	21.292	240	837139	20.000	ng/ul	0.00
88) Perylene-d12	23.539	264	976110	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.257	96	14016	2.618	ng/uL	0.00
4) Pyridine-d5	3.657	84	208812	13.253	ng/ul	0.00
7) Phenol-d5	6.869	99	278750	14.933	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.045	67	189592	16.443	ng/ul	0.00
11) 2-Chlorophenol-d4	7.228	132	213386	15.556	ng/ul	0.00
15) 4-Methylphenol-d8	8.398	113	220885	15.067	ng/ul	0.00
21) Nitrobenzene-d5	8.857	128	110551	17.491	ng/ul	0.00
24) 2-Nitrophenol-d4	9.569	143	117050	17.178	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.098	165	206111	17.774	ng/ul	0.00
31) 4-Chloroaniline-d4	10.628	131	318904	16.558	ng/ul	0.00
46) Dimethylphthalate-d6	13.763	166	716571	20.322	ng/ul	0.00
49) Acenaphthylene-d8	14.033	160	774451	19.204	ng/ul	0.00
54) 4-Nitrophenol-d4	14.557	143	110119	15.059	ng/ul	0.00
60) Fluorene-d10	15.339	176	613087	20.690	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.468	200	87769	14.724	ng/ul	0.00
73) Anthracene-d10	17.198	188	937396	21.356	ng/ul	0.00
81) Pyrene-d10	19.498	212	1099842	21.421	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.397	264	1103271	21.504	ng/ul	0.00

Target Compounds Qvalue

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

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