

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081823\
 Data File : VN078956.D
 Acq On : 18 Aug 2023 20:49
 Operator : JC\MD
 Sample : 04042-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-376

Quant Time: Aug 18 23:49:35 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N081523W.M
 Quant Title : SW846 8260
 QLast Update : Wed Aug 16 02:42:02 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	202889	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	396544	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	384310	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	119305	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	163536	51.799	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	103.600%
35) Dibromofluoromethane	8.171	113	134487	50.382	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	100.760%
50) Toluene-d8	10.571	98	492761	49.716	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	99.440%
62) 4-Bromofluorobenzene	12.853	95	149602	46.178	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	92.360%
Target Compounds						
						Qvalue
16) Acetone	4.436	43	31188	23.941	ug/l #	84
20) Methylene Chloride	5.277	84	8648	3.115	ug/l #	81
25) 2-Butanone	7.500	43	18073	8.653	ug/l #	74
43) Isopropyl Acetate	8.694	43	236529	29.787	ug/l #	70
51) 4-Methyl-2-Pentanone	10.447	43	69181	14.689	ug/l #	79

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

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