

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081623\
 Data File : VN078909.D
 Acq On : 16 Aug 2023 17:42
 Operator : JC\MD
 Sample : 04022-08
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-24

Quant Time: Aug 17 01:19:28 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	8.236	168	144274	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	289569	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	277102	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	82721	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	124546	55.476	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 110.960%	
35) Dibromofluoromethane	8.171	113	99543	51.068	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 102.140%	
50) Toluene-d8	10.571	98	374297	51.715	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 103.440%	
62) 4-Bromofluorobenzene	12.853	95	108058	45.677	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 91.360%	
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	26619	28.736	ug/l	94
20) Methylene Chloride	5.277	84	15302	7.751	ug/l #	86
25) 2-Butanone	7.494	43	7789	5.245	ug/l #	74
43) Isopropyl Acetate	8.694	43	246515	42.513	ug/l #	68
51) 4-Methyl-2-Pentanone	10.453	43	37123	10.794	ug/l #	74

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

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