

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN073123\
 Data File : VN078705.D
 Acq On : 31 Jul 2023 15:11
 Operator : JC\MD
 Sample : 03797-05
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 HANOVER-ST-2

Quant Time: Aug 01 01:18:26 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N071023W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jul 10 16:13:53 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.235	168	734854	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.112	114	1331588	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	1228539	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	400256	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	486757	49.409	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	98.820%
35) Dibromofluoromethane	8.177	113	448244	49.589	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	99.180%
50) Toluene-d8	10.570	98	1691163	49.525	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	99.060%
62) 4-Bromofluorobenzene	12.853	95	550047	52.295	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	104.600%
Target Compounds						
					Qvalue	
16) Acetone	4.447	43	180503	58.820	ug/l	99
20) Methylene Chloride	5.294	84	49487	4.510	ug/l	90
25) 2-Butanone	7.494	43	101336	20.978	ug/l #	82
37) Ethyl Acetate	7.577	43	33423	3.072	ug/l	94
43) Isopropyl Acetate	8.694	43	553698	30.910	ug/l #	83
51) 4-Methyl-2-Pentanone	10.453	43	431383	40.487	ug/l	91

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

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