

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN071823\
 Data File : VN078567.D
 Acq On : 18 Jul 2023 17:40
 Operator : JC\MD
 Sample : 03646-05RE
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 YARD-1RE

Quant Time: Jul 19 03:14:32 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	400526	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.112	114	741823	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	731631	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	259490	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	269620	50.213	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 100.420%	
35) Dibromofluoromethane	8.177	113	269858	53.589	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 107.180%	
50) Toluene-d8	10.576	98	906057	47.628	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 95.260%	
62) 4-Bromofluorobenzene	12.853	95	312120	53.266	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 106.540%	
Target Compounds						
					Qvalue	
16) Acetone	4.447	43	69346	41.460	ug/l	97
20) Methylene Chloride	5.289	84	18505	2.700	ug/l	95
25) 2-Butanone	7.494	43	43546	16.539	ug/l #	76
37) Ethyl Acetate	7.565	43	14053	2.318	ug/l #	89
43) Isopropyl Acetate	8.700	43	352099	35.283	ug/l #	78
51) 4-Methyl-2-Pentanone	10.453	43	143592	24.191	ug/l	90

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

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