

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN080823\
 Data File : VN078798.D
 Acq On : 08 Aug 2023 13:35
 Operator : JC\MD
 Sample : 03913-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 72-11933

Quant Time: Aug 09 02:48:25 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N080723W.M
 Quant Title : SW846 8260
 QLast Update : Tue Aug 08 05:03:52 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	189653	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.112	114	361170	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	322349	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	88021	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	146560	51.548	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	103.100%
35) Dibromofluoromethane	8.177	113	121067	49.637	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	99.280%
50) Toluene-d8	10.571	98	454644	50.653	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	101.300%
62) 4-Bromofluorobenzene	12.853	95	119142	44.416	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	88.840%
Target Compounds						
						Qvalue
16) Acetone	4.453	43	42399	38.758	ug/l #	86
20) Methylene Chloride	5.283	84	17632	6.508	ug/l #	85
25) 2-Butanone	7.506	43	13918	7.513	ug/l #	77
43) Isopropyl Acetate	8.694	43	80638	11.888	ug/l #	77
51) 4-Methyl-2-Pentanone	10.453	43	49433	11.898	ug/l #	85

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

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