

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN120722\
 Data File : VN075621.D
 Acq On : 07 Dec 2022 21:08
 Operator : JC\MD
 Sample : N5897-17
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 2025

Quant Time: Dec 08 00:24:19 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N112122W.M
 Quant Title : SW846 8260
 QLast Update : Wed Nov 23 00:18:03 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	185285	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	321019	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	289255	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	131399	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	55314	55.615	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	111.240%
35) Dibromofluoromethane	8.177	113	69254	54.917	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	109.840%
50) Toluene-d8	10.571	98	148684	52.348	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	104.700%
62) 4-Bromofluorobenzene	12.853	95	108935	51.962	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	103.920%
Target Compounds						
						Qvalue
16) Acetone	4.448	43	26257	28.303	ug/l	98
17) Carbon Disulfide	4.718	76	6036	1.354	ug/l #	74
20) Methylene Chloride	5.283	84	10746	4.550	ug/l #	84
37) Ethyl Acetate	7.571	43	23656	9.553	ug/l	97
43) Isopropyl Acetate	8.694	43	131896	33.184	ug/l #	87

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

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