

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN082622\
 Data File : VN074140.D
 Acq On : 26 Aug 2022 16:58
 Operator : JC\MD
 Sample : N4297-01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW202D-20220819

Quant Time: Aug 27 00:21:33 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N081522W.M
 Quant Title : SW846 8260
 QLast Update : Fri Aug 26 06:44:47 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.016	168	574680	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	871203	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	736914	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.616	152	317570	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.375	65	298743	53.372	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 106.740%	
35) Dibromofluoromethane	7.951	113	262463	50.942	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 101.880%	
50) Toluene-d8	10.381	98	973344	48.779	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 97.560%	
62) 4-Bromofluorobenzene	12.674	95	299640	45.247	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 90.500%	
Target Compounds						
					Qvalue	
7) Trichlorofluoromethane	3.322	101	11619	1.385	ug/l	90
12) 1,1-Dichloroethene	4.116	96	3391	0.748	ug/l #	70
16) Acetone	4.240	43	6513	2.466	ug/l #	86
19) Methyl tert-butyl Ether	5.534	73	4625	0.283	ug/l	95
24) 1,1-Dichloroethane	6.310	63	18673	2.106	ug/l #	90
27) cis-1,2-Dichloroethene	7.251	96	23904	3.877	ug/l	86
30) Chloroform	7.745	83	3451	0.348	ug/l	99
42) 1,2-Dichloroethane	8.469	62	5544	0.714	ug/l	98
44) Trichloroethene	9.151	130	138421	21.219	ug/l	98
64) Tetrachloroethene	10.916	164	37214	6.398	ug/l	98

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

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