

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN020724\
 Data File : VN080946.D
 Acq On : 07 Feb 2024 19:15
 Operator : JC\MD
 Sample : P1372-01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 CG-MW-63

Quant Time: Feb 08 00:31:05 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N020624W.M
 Quant Title : SW846 8260
 QLast Update : Wed Feb 07 06:22:23 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.229	168	256724	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	573896	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	671116	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	292692	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	186741	57.661	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 115.320%	
35) Dibromofluoromethane	8.171	113	177436	57.618	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 115.240%	
50) Toluene-d8	10.570	98	648171	58.190	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 116.380%#	
62) 4-Bromofluorobenzene	12.853	95	281715	69.619	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 139.240%#	
Target Compounds						
						Qvalue
16) Acetone	4.436	43	6646	6.237	ug/l	99
27) cis-1,2-Dichloroethene	7.494	96	12030	3.351	ug/l	81
30) Chloroform	7.971	83	5174	0.850	ug/l	88
44) Trichloroethene	9.353	130	7974	1.883	ug/l	92
64) Tetrachloroethene	11.106	164	130780	25.817	ug/l	89

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

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