

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN083122\
 Data File : VN074217.D
 Acq On : 31 Aug 2022 16:23
 Operator : JC\MD
 Sample : N4456-01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 KTS2-GW-MH-02-5-20

Quant Time: Sep 01 00:26:54 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N083022W.M
 Quant Title : SW846 8260
 QLast Update : Wed Aug 31 07:02:40 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.016	168	190353	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	353408	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	344339	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.615	152	155490	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.369	65	157587	53.368	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 106.740%		
35) Dibromofluoromethane	7.951	113	129737	56.415	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 112.820%		
50) Toluene-d8	10.380	98	420138	46.391	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 92.780%		
62) 4-Bromofluorobenzene	12.668	95	163242	59.272	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 118.540%		
Target Compounds						
						Qvalue
3) Chloromethane	2.269	50	21315	7.683	ug/l #	84
16) Acetone	4.228	43	37967	19.242	ug/l	96
17) Carbon Disulfide	4.463	76	3532	0.556	ug/l #	90
25) 2-Butanone	7.257	43	10639	3.371	ug/l	96
40) Benzene	8.386	78	6092	0.520	ug/l #	90
52) Toluene	10.445	92	1791	0.259	ug/l #	24

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

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