

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN083122\
 Data File : VN074218.D
 Acq On : 31 Aug 2022 16:48
 Operator : JC\MD
 Sample : N4455-01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 KTS2-WC-L-MH-02

Quant Time: Sep 01 00:27:17 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	184623	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	337186	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	324317	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.615	152	151819	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.375	65	154228	53.851	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 107.700%		
35) Dibromofluoromethane	7.951	113	126717	57.753	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 115.500%		
50) Toluene-d8	10.380	98	412674	47.759	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 95.520%		
62) 4-Bromofluorobenzene	12.674	95	157988	60.124	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 120.240%		
Target Compounds						
						Qvalue
3) Chloromethane	2.269	50	20748	7.711	ug/l	90
16) Acetone	4.228	43	45754	23.908	ug/l	98
18) Methyl Acetate	4.792	43	11420	0.343	ug/l #	79
25) 2-Butanone	7.257	43	11289	3.688	ug/l #	70
40) Benzene	8.386	78	6813	0.609	ug/l	100
52) Toluene	10.439	92	1574	0.239	ug/l #	71

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

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