

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN083022\
 Data File : VN074194.D
 Acq On : 31 Aug 2022 03:25
 Operator : JC\MD
 Sample : N4421-01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WC-PALM-SP-01

Quant Time: Aug 31 07:07:33 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	171370	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	320964	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	330190	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.616	152	146742	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.369	65	132925	50.002	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 100.000%		
35) Dibromofluoromethane	7.951	113	108511	51.955	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 103.900%		
50) Toluene-d8	10.381	98	372338	45.269	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 90.540%		
62) 4-Bromofluorobenzene	12.674	95	137813	55.097	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 110.200%		
Target Compounds						
					Qvalue	
16) Acetone	4.222	43	82579	46.488	ug/l	94
20) Methylene Chloride	5.022	84	17592	5.417	ug/l #	70
25) 2-Butanone	7.269	43	13587	4.781	ug/l #	78
51) 4-Methyl-2-Pentanone	10.263	43	12303	2.225	ug/l #	81

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

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