

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN083022\
 Data File : VN074195.D
 Acq On : 31 Aug 2022 03:50
 Operator : JC\MD
 Sample : N4421-05
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 33 Sample Multiplier: 1

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

Quant Time: Aug 31 07:07:43 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

Manual Integrations
 APPROVED

Reviewed By : John Carlone 08/31/2022
 Supervised By : Semsettin Yesilyurt 08/31/2022

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

Internal Standards						
1) Pentafluorobenzene	8.022	168	168238	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	317228	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	316587	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.615	152	131530	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.369	65	125892	48.238	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	96.480%		
35) Dibromofluoromethane	7.957	113	106779	51.727	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	103.460%		
50) Toluene-d8	10.380	98	371173	45.658	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	91.320%		
62) 4-Bromofluorobenzene	12.674	95	134289	54.320	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	108.640%		
Target Compounds						
					Qvalue	
16) Acetone	4.228	43	46157	26.468	ug/l	96
17) Carbon Disulfide	4.469	76	6707m	1.195	ug/l	
20) Methylene Chloride	5.010	84	18309	5.881	ug/l #	78
37) Ethyl Acetate	7.333	43	43703	7.802	ug/l #	93
43) Isopropyl Acetate	8.492	43	172568	22.917	ug/l #	84
84) 1,2,4-Trimethylbenzene	13.304	105	18145	2.052	ug/l	100

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

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