

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN030922\
 Data File : VN071208.D
 Acq On : 09 Mar 2022 14:33
 Operator : JC\MD
 Sample : N1830-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 11986

Quant Time: Mar 10 01:12:18 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N030722W.M
 Quant Title : SW846 8260
 QLast Update : Tue Mar 08 06:52:46 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.086	168	836806	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.963	114	1344120	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	1254349	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.674	152	415600	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.433	65	514923	44.898	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	89.800%
35) Dibromofluoromethane	8.022	113	394038	46.457	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	92.920%
50) Toluene-d8	10.439	98	1685383	49.352	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	98.700%
62) 4-Bromofluorobenzene	12.727	95	573499	49.593	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	99.180%
Target Compounds						
					Qvalue	
16) Acetone	4.287	43	67495	12.095	ug/l	97
18) Methyl Acetate	4.857	43	9182	1.077	ug/l #	81
20) Methylene Chloride	5.104	84	29519	3.254	ug/l	91
25) 2-Butanone	7.351	43	11841	1.317	ug/l	99
43) Isopropyl Acetate	8.557	43	75381	2.695	ug/l #	90

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

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