

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042022\
 Data File : VN072080.D
 Acq On : 20 Apr 2022 15:52
 Operator : JC\MD
 Sample : N2480-06
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WC-16-1A-2-GR

Quant Time: Apr 21 03:30:47 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
 Quant Title : SW846 8260
 QLast Update : Thu Mar 31 01:57:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.080	168	702753	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.963	114	1187525	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	1308489	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.668	152	541068	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.433	65	378101	47.129	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	94.260%
35) Dibromofluoromethane	8.016	113	333196	47.682	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	95.360%
50) Toluene-d8	10.439	98	1530293	53.533	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	107.060%
62) 4-Bromofluorobenzene	12.727	95	561344	60.718	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	121.440%
Target Compounds						
					Qvalue	
16) Acetone	4.287	43	51159	14.921	ug/l	95
20) Methylene Chloride	5.092	84	37966	5.254	ug/l	91
39) Methylcyclohexane	9.457	83	33083	2.848	ug/l	99
43) Isopropyl Acetate	8.557	43	68617	3.862	ug/l #	83
80) 1,3,5-Trimethylbenzene	13.051	105	50079	1.386	ug/l	97
84) 1,2,4-Trimethylbenzene	13.362	105	48657	1.380	ug/l	96

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

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