

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102021\
 Data File : VN069066.D
 Acq On : 20 Oct 2021 17:01
 Operator : JC/MD
 Sample : M4249-11
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP2-(D)

Quant Time: Oct 20 23:15:14 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N101221W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 13 05:44:40 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.091	168	354806	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	620870	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.749	117	604331	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.678	152	218861	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	231399	49.465	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	98.940%
35) Dibromofluoromethane	8.024	113	183832	50.891	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	101.780%
50) Toluene-d8	10.443	98	784243	54.260	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	108.520%
62) 4-Bromofluorobenzene	12.734	95	274045	50.255	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	100.520%
Target Compounds						
					Qvalue	
16) Acetone	4.293	43	11907	6.821	ug/l	98
18) Methyl Acetate	4.873	43	8788	1.327	ug/l	100
20) Methylene Chloride	5.101	84	15449	3.845	ug/l	92
43) Isopropyl Acetate	8.574	43	62601	6.478	ug/l #	78

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

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