

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN080321\
 Data File : VN067990.D
 Acq On : 3 Aug 2021 18:38
 Operator : JC/MD
 Sample : M3242-01
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 257392

Quant Time: Aug 04 03:41:42 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N072921W.M
 Quant Title : SW846 8260
 QLast Update : Thu Jul 29 15:02:35 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.086	168	295784	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	561988	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.746	117	542278	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.677	152	190119	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	230822	55.124	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	110.240%
35) Dibromofluoromethane	8.024	113	174701	53.341	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	106.680%
50) Toluene-d8	10.443	98	716906	47.628	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	95.260%
62) 4-Bromofluorobenzene	12.733	95	231672	42.683	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	85.360%
Target Compounds						
					Qvalue	
16) Acetone	4.291	43	277364	181.539	ug/l	96
25) 2-Butanone	7.348	43	54619	24.641	ug/l	98
29) Tetrahydrofuran	7.689	42	159837	112.608	ug/l	93

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN080321\
Data File : VN067990.D
Acq On : 3 Aug 2021 18:38
Operator : JC/MD
Sample : M3242-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
257392

Quant Time: Aug 04 03:41:42 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N072921W.M
Quant Title : SW846 8260
QLast Update : Thu Jul 29 15:02:35 2021
Response via : Initial Calibration

