

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN101821\
 Data File : VN068988.D
 Acq On : 18 Oct 2021 18:55
 Operator : JC/MD
 Sample : M4243-01
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 T3-1-TOP-GRAB

Quant Time: Oct 19 06:39:25 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	8.088	168	326172	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	567378	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.747	117	537432	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.678	152	204343	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	214115	49.789	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	99.580%
35) Dibromofluoromethane	8.024	113	169392	51.314	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	102.620%
50) Toluene-d8	10.443	98	709989	53.754	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	107.500%
62) 4-Bromofluorobenzene	12.734	95	252189	50.607	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	101.220%
Target Compounds						
					Qvalue	
16) Acetone	4.293	43	23604	14.708	ug/l	100
20) Methylene Chloride	5.101	84	13066	3.537	ug/l	98
43) Isopropyl Acetate	8.566	43	44968	5.092	ug/l #	83

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN101821\
Data File : VN068988.D
Acq On : 18 Oct 2021 18:55
Operator : JC/MD
Sample : M4243-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
T3-1-TOP-GRAB

Quant Time: Oct 19 06:39:25 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

