

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX013122\
 Data File : VX026694.D
 Acq On : 31 Jan 2022 20:30
 Operator : JC/MD
 Sample : N1335-11
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MH-4D

Quant Time: Feb 01 06:32:08 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X011122W.M
 Quant Title : SW846 8260
 QLast Update : Tue Jan 11 14:45:13 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	123497	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	209909	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	194412	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	86883	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	80522	46.675	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	93.360%
35) Dibromofluoromethane	5.391	113	65530	48.343	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	96.680%
50) Toluene-d8	8.653	98	250427	51.267	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	102.540%
62) 4-Bromofluorobenzene	11.079	95	93153	53.683	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	107.360%
Target Compounds						
					Qvalue	
16) Acetone	2.386	43	12616	14.340	ug/l	95
20) Methylene Chloride	2.794	84	1562	1.075	ug/l	90
43) Isopropyl Acetate	6.342	43	11416	3.265	ug/l	98

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX013122\
Data File : VX026694.D
Acq On : 31 Jan 2022 20:30
Operator : JC/MD
Sample : N1335-11
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MH-4D

Quant Time: Feb 01 06:32:08 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X011122W.M
Quant Title : SW846 8260
QLast Update : Tue Jan 11 14:45:13 2022
Response via : Initial Calibration

