

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101122\
 Data File : VX032069.D
 Acq On : 12 Oct 2022 08:11
 Operator : JC/MD
 Sample : N5066-09
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 46 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 B-4(11-12)

Quant Time: Oct 12 13:27:07 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100622W.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 06 14:28:46 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	77934	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	119121	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	130071	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	75386	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	85053	52.844	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	105.680%
35) Dibromofluoromethane	5.385	113	65246	48.388	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	96.780%
50) Toluene-d8	8.653	98	250445	51.617	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	103.240%
62) 4-Bromofluorobenzene	11.079	95	84639	49.344	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	98.680%
Target Compounds						
					Qvalue	
16) Acetone	2.386	43	22227	40.634	ug/l	97
20) Methylene Chloride	2.788	84	7305	5.082	ug/l	94
37) Ethyl Acetate	4.715	43	13826	8.213	ug/l #	87
43) Isopropyl Acetate	6.342	43	57152	22.377	ug/l	99

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101122\
Data File : VX032069.D
Acq On : 12 Oct 2022 08:11
Operator : JC/MD
Sample : N5066-09
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B-4(11-12)

Quant Time: Oct 12 13:27:07 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100622W.M
Quant Title : SW846 8260
QLast Update : Thu Oct 06 14:28:46 2022
Response via : Initial Calibration

