

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090622\
 Data File : VX031147.D
 Acq On : 06 Sep 2022 22:40
 Operator : JC/MD
 Sample : N4479-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 SB-8

Quant Time: Sep 07 01:15:55 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X090122W.M
 Quant Title : SW846 8260
 QLast Update : Fri Sep 02 02:49:21 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	213658	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	334013	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	297638	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	133493	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	137758	50.110	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 100.220%		
35) Dibromofluoromethane	5.391	113	124533	51.090	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 102.180%		
50) Toluene-d8	8.653	98	438086	48.834	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 97.660%		
62) 4-Bromofluorobenzene	11.079	95	158462	50.012	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 100.020%		
Target Compounds						
					Qvalue	
16) Acetone	2.386	43	19951	18.799	ug/l	95
20) Methylene Chloride	2.794	84	14525	3.591	ug/l #	81
37) Ethyl Acetate	4.721	43	22325	7.526	ug/l #	89
43) Isopropyl Acetate	6.348	43	100973	21.779	ug/l #	92

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090622\
Data File : VX031147.D
Acq On : 06 Sep 2022 22:40
Operator : JC/MD
Sample : N4479-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SB-8

Quant Time: Sep 07 01:15:55 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X090122W.M
Quant Title : SW846 8260
QLast Update : Fri Sep 02 02:49:21 2022
Response via : Initial Calibration

