

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX042624\
 Data File : VX041264.D
 Acq On : 26 Apr 2024 17:13
 Operator : JC/MD
 Sample : P2271-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 RT-3860

Quant Time: Apr 27 02:21:11 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X042224W.M
 Quant Title : SW846 8260
 QLast Update : Tue Apr 23 05:30:28 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	135421	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.757	114	197284	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	189758	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	93660	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	74056	57.448	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	=	114.900%	
35) Dibromofluoromethane	5.385	113	62861	52.382	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	104.760%	
50) Toluene-d8	8.647	98	229012	56.562	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	=	113.120%#	
62) 4-Bromofluorobenzene	11.079	95	84114	56.285	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	=	112.580%	
Target Compounds						
					Qvalue	
16) Acetone	2.386	43	35047	74.690	ug/l	93
18) Methyl Acetate	2.715	43	2794	2.013	ug/l	94
20) Methylene Chloride	2.788	84	4065	3.547	ug/l	95
43) Isopropyl Acetate	6.342	43	49664	18.301	ug/l	98

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX042624\
Data File : VX041264.D
Acq On : 26 Apr 2024 17:13
Operator : JC/MD
Sample : P2271-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
RT-3860

Quant Time: Apr 27 02:21:11 2024
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X042224W.M
Quant Title : SW846 8260
QLast Update : Tue Apr 23 05:30:28 2024
Response via : Initial Calibration

