

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX032322\
 Data File : VX027673.D
 Acq On : 23 Mar 2022 17:24
 Operator : JC/MD
 Sample : N1905-16
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 WC-14A-1A-1

Quant Time: Mar 23 18:38:30 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X031722W.M
 Quant Title : SW846 8260
 QLast Update : Thu Mar 17 15:02:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	344857	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	557530	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	541118	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	230727	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	201926	47.531	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	95.060%
35) Dibromofluoromethane	5.391	113	182974	51.010	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	102.020%
50) Toluene-d8	8.653	98	670911	49.427	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	98.860%
62) 4-Bromofluorobenzene	11.085	95	233004	46.702	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	93.400%
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	45964	26.544	ug/l	99
20) Methylene Chloride	2.794	84	11500	3.089	ug/l	99
25) 2-Butanone	4.562	43	13740	5.274	ug/l	90
43) Isopropyl Acetate	6.348	43	25077	3.190	ug/l	97

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX032322\
Data File : VX027673.D
Acq On : 23 Mar 2022 17:24
Operator : JC/MD
Sample : N1905-16
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
WC-14A-1A-1

Quant Time: Mar 23 18:38:30 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X031722W.M
Quant Title : SW846 8260
QLast Update : Thu Mar 17 15:02:59 2022
Response via : Initial Calibration

