

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101722\
 Data File : VX032179.D
 Acq On : 18 Oct 2022 01:11
 Operator : JC/MD
 Sample : N5159-12
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 SB-43

Quant Time: Oct 18 04:34:36 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X101722W.M
 Quant Title : SW846 8260
 QLast Update : Tue Oct 18 04:18:35 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	123232	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	176548	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	202901	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	91483	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	84031	42.814	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	85.620%
35) Dibromofluoromethane	5.385	113	71721	59.538	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	119.080%
50) Toluene-d8	8.647	98	270353	63.985	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	127.980%#
62) 4-Bromofluorobenzene	11.079	95	100643	62.561	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	125.120%#
Target Compounds						
					Qvalue	
16) Acetone	2.386	43	17930	37.994	ug/l	95
20) Methylene Chloride	2.794	84	7206	4.966	ug/l	89
37) Ethyl Acetate	4.715	43	19782	11.311	ug/l	99
43) Isopropyl Acetate	6.342	43	69449	26.351	ug/l	98

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101722\
Data File : VX032179.D
Acq On : 18 Oct 2022 01:11
Operator : JC/MD
Sample : N5159-12
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 34 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
SB-43

Quant Time: Oct 18 04:34:36 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X101722W.M
Quant Title : SW846 8260
QLast Update : Tue Oct 18 04:18:35 2022
Response via : Initial Calibration

