

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090922\
 Data File : VX031260.D
 Acq On : 09 Sep 2022 16:31
 Operator : JC/MD
 Sample : N4495-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 ETGI-345

Quant Time: Sep 12 00:58:51 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X090722W.M
 Quant Title : SW846 8260
 QLast Update : Wed Sep 07 23:22:41 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	52130	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.769	114	93324	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	83396	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	34999	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	76038	61.609	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	123.220%
35) Dibromofluoromethane	5.391	113	48232	47.427	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	94.860%
50) Toluene-d8	8.653	98	159550	43.181	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	86.360%
62) 4-Bromofluorobenzene	11.079	95	64156	47.380	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	94.760%
Target Compounds						
						Qvalue
16) Acetone	2.380	43	28643	67.026	ug/l	# 87
20) Methylene Chloride	2.788	84	2417	2.406	ug/l	# 86
43) Isopropyl Acetate	6.342	43	5186	2.434	ug/l	# 90

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090922\
Data File : VX031260.D
Acq On : 09 Sep 2022 16:31
Operator : JC/MD
Sample : N4495-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
ETGI-345

Quant Time: Sep 12 00:58:51 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X090722W.M
Quant Title : SW846 8260
QLast Update : Wed Sep 07 23:22:41 2022
Response via : Initial Calibration

