

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX091222\
 Data File : VX031289.D
 Acq On : 12 Sep 2022 21:22
 Operator : JC/MD
 Sample : N4298-12 20X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 RE132D7-20220822

Quant Time: Sep 13 04:02:17 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	66824	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.769	114	106576	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.061	117	95196	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	42400	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	80279	50.742	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	101.480%
35) Dibromofluoromethane	5.398	113	53792	46.317	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	92.640%
50) Toluene-d8	8.653	98	172826	40.958	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	81.920%#
62) 4-Bromofluorobenzene	11.086	95	65358	42.266	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	84.540%
Target Compounds						Qvalue
9) 1,1,2-Trichlorotrifluo...	2.319	101	430	0.358	ug/l	96
44) Trichloroethene	7.135	130	37295	28.081	ug/l	97

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX091222\
Data File : VX031289.D
Acq On : 12 Sep 2022 21:22
Operator : JC/MD
Sample : N4298-12 20X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
RE132D7-20220822

Quant Time: Sep 13 04:02:17 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

