

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062422\
 Data File : VX029782.D
 Acq On : 24 Jun 2022 17:53
 Operator : JC/MD
 Sample : N3473-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BP-MW17411-GW-305-307

Quant Time: Jun 25 02:05:31 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jun 20 01:31:01 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	242081	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	390218	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	347183	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	150211	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	133091	41.533	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	83.060%
35) Dibromofluoromethane	5.391	113	117059	46.032	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	92.060%
50) Toluene-d8	8.653	98	441819	47.558	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	95.120%
62) 4-Bromofluorobenzene	11.085	95	159034	45.394	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	90.780%
Target Compounds						
					Qvalue	
12) 1,1-Dichloroethene	2.319	96	2933	1.196	ug/l	95
16) Acetone	2.386	43	3309	2.591	ug/l	95
19) Methyl tert-butyl Ether	3.123	73	3538	0.416	ug/l	98
24) 1,1-Dichloroethane	3.611	63	32127	6.820	ug/l	98
27) cis-1,2-Dichloroethene	4.501	96	1966	0.628	ug/l	84
30) Chloroform	5.099	83	1621	0.325	ug/l #	77

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

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