

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070621\
 Data File : VX023152.D
 Acq On : 06 Jul 2021 17:40
 Operator : JC/MD
 Sample : M2923-07
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-12A-070121

Quant Time: Jul 07 02:54:24 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X070221W.M
 Quant Title : SW846 8260
 QLast Update : Fri Jul 02 14:29:42 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	259619	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	428708	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	380029	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	158065	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	182067	52.179	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	104.360%
35) Dibromofluoromethane	5.397	113	137080	49.516	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	99.040%
50) Toluene-d8	8.653	98	512847	49.662	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	99.320%
62) 4-Bromofluorobenzene	11.085	95	184566	46.295	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	92.580%
Target Compounds						
						Qvalue
16) Acetone	2.386	43	4116	2.982	ug/l	# 81
27) cis-1,2-Dichloroethene	4.501	96	10015	3.308	ug/l	96
44) Trichloroethene	7.129	130	59287	18.835	ug/l	99
64) Tetrachloroethene	9.275	164	992	0.327	ug/l	88

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

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