

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090122\
 Data File : VX031038.D
 Acq On : 01 Sep 2022 22:29
 Operator : JC/MD
 Sample : N4297-22
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 RW9-MW01D2-20220821

Quant Time: Sep 02 02:56:55 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X090122W.M
 Quant Title : SW846 8260
 QLast Update : Fri Sep 02 02:49:21 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	219001	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	357209	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	319699	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	142382	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	157992	56.068	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 112.140%	
35) Dibromofluoromethane	5.391	113	155642	59.706	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 119.420%	
50) Toluene-d8	8.653	98	552552	57.594	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 115.180%#	
62) 4-Bromofluorobenzene	11.085	95	185377	54.707	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 109.420%	
Target Compounds						
						Qvalue
16) Acetone	2.391	43	3574	3.285	ug/l #	78
29) Tetrahydrofuran	5.019	42	13524	12.491	ug/l #	77
44) Trichloroethene	7.135	130	2416	0.728	ug/l	99

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

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