

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072822\
 Data File : VX030286.D
 Acq On : 28 Jul 2022 21:40
 Operator : JC/MD
 Sample : N3695-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 EB-02-07112022

Quant Time: Jul 29 06:01:05 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072522W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jul 25 16:49:17 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	139443	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	245633	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	251154	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	137440	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	146782	52.371	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 104.740%	
35) Dibromofluoromethane	5.391	113	117319	49.963	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 99.920%	
50) Toluene-d8	8.653	98	440379	48.428	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 96.860%	
62) 4-Bromofluorobenzene	11.085	95	158113	46.851	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 93.700%	
Target Compounds						
					Qvalue	
16) Acetone	2.379	43	10579	7.373	ug/l	99
25) 2-Butanone	4.586	43	4352	1.805	ug/l #	73
30) Chloroform	5.105	83	1217	0.272	ug/l	88
52) Toluene	8.726	92	2051	0.316	ug/l	96
64) Tetrachloroethene	9.281	164	25357	9.378	ug/l	96

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

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