

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072221\
 Data File : VX023395.D
 Acq On : 23 Jul 2021 00:29
 Operator : JC/MD
 Sample : M3145-14
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 152140-MW-3D(OFFSITE)

Quant Time: Jul 23 06:11:54 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072221W.M
 Quant Title : SW846 8260
 QLast Update : Fri Jul 23 06:05:38 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	170011	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	282400	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	244026	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	88552	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.971	65	107030	50.468	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 100.940%	
35) Dibromofluoromethane	5.397	113	87591	49.607	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 99.220%	
50) Toluene-d8	8.653	98	336916	51.936	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 103.880%	
62) 4-Bromofluorobenzene	11.085	95	101983	46.232	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 92.460%	
Target Compounds						
					Qvalue	
16) Acetone	2.386	43	2527	2.429	ug/l #	83
27) cis-1,2-Dichloroethene	4.495	96	39993	19.406	ug/l	91
30) Chloroform	5.105	83	1483	0.440	ug/l	99
44) Trichloroethene	7.135	130	10005	4.808	ug/l	90
64) Tetrachloroethene	9.275	164	15621	8.043	ug/l	93

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

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