

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072721\
 Data File : VX023478.D
 Acq On : 27 Jul 2021 15:37
 Operator : JC/MD
 Sample : M3146-12
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 152140-DDC-3-PS

Quant Time: Jul 28 00:51:46 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	166086	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	284219	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	241507	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	94584	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.970	65	110350	53.263	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	= 106.520%		
35) Dibromofluoromethane	5.397	113	90563	50.962	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 101.920%		
50) Toluene-d8	8.653	98	332085	50.863	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	= 101.720%		
62) 4-Bromofluorobenzene	11.085	95	102519	46.177	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 92.360%		
Target Compounds						
						Qvalue
16) Acetone	2.386	43	3650	3.591	ug/l	92
20) Methylene Chloride	2.800	84	843	0.426	ug/l	89
27) cis-1,2-Dichloroethene	4.495	96	1090	0.541	ug/l	91
30) Chloroform	5.099	83	2897	0.879	ug/l	85
64) Tetrachloroethene	9.275	164	82651	43.001	ug/l	96
90) Hexachloroethane	12.542	117	1771	3.492	ug/l	89

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

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