

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX042221\
 Data File : VX021823.D
 Acq On : 22 Apr 2021 17:42
 Operator : JC/MD
 Sample : M2138-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 152140-MW-5D

Quant Time: Apr 23 07:44:53 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X042021W.M
 Quant Title : SW846 8260
 QLast Update : Tue Apr 20 13:46:22 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.629	168	158753	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.830	114	287128	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.092	117	305238	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.061	152	164598	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.031	65	90958	50.51	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	101.02%
35) Dibromofluoromethane	5.464	113	93259	50.79	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	101.58%
50) Toluene-d8	8.695	98	352534	49.93	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	99.86%
62) 4-Bromofluorobenzene	11.116	95	153447	58.20	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	116.40%
Target Compounds						
					Qvalue	
16) Acetone	2.422	43	2722	4.40	ug/l	97
27) cis-1,2-Dichloroethene	4.562	96	3732	1.95	ug/l	84
44) Trichloroethene	7.190	130	5539	2.51	ug/l	99
64) Tetrachloroethene	9.317	164	215274	90.20	ug/l	96
90) Hexachloroethane	12.579	117	1610	0.83	ug/l	99

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

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