

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092121\
 Data File : VX024384.D
 Acq On : 21 Sep 2021 14:27
 Operator : JC/MD
 Sample : M3782-05
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 DUP02-20210914

Quant Time: Sep 21 15:41:01 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091421W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 21 10:52:48 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	155163	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	272997	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	255453	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	90696	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	111453	50.372	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	= 100.740%		
35) Dibromofluoromethane	5.391	113	89740	49.016	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 98.040%		
50) Toluene-d8	8.653	98	338111	50.694	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	= 101.380%		
62) 4-Bromofluorobenzene	11.085	95	115924	44.662	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 89.320%		
Target Compounds						Qvalue
12) 1,1-Dichloroethene	2.325	96	1086	0.759	ug/l	93
16) Acetone	2.386	43	1701	1.803	ug/l	92
27) cis-1,2-Dichloroethene	4.501	96	1925	1.064	ug/l	98
30) Chloroform	5.099	83	1346	0.422	ug/l	97
44) Trichloroethene	7.135	130	189055	93.469	ug/l	98

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

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