

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX091721\
 Data File : VX024332.D
 Acq On : 18 Sep 2021 01:23
 Operator : JC/MD
 Sample : M3782-09
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW181S-20210914

Quant Time: Sep 20 02:43:56 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091421W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 14 12:18:04 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	135431	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.775	114	208945	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.061	117	188423	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	79014	50.000	ug/l	0.00

System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	96846	50.148	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 100.300%	
35) Dibromofluoromethane	5.397	113	80674	57.572	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 115.140%	
50) Toluene-d8	8.653	98	264510	51.816	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 103.640%	
62) 4-Bromofluorobenzene	11.085	95	86957	43.772	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 87.540%	

Target Compounds			Qvalue			
3) Chloromethane	1.300	50	538	0.396	ug/l #	86
12) 1,1-Dichloroethene	2.325	96	863	0.691	ug/l	84
16) Acetone	2.386	43	1963	2.384	ug/l #	70
19) Methyl tert-butyl Ether	3.129	73	1404	0.306	ug/l	92
24) 1,1-Dichloroethane	3.617	63	2304	0.881	ug/l #	92
27) cis-1,2-Dichloroethene	4.495	96	2406	1.524	ug/l	85
30) Chloroform	5.098	83	2155	0.775	ug/l	97
32) 1,1,1-Trichloroethane	5.397	97	1292	0.519	ug/l #	44
44) Trichloroethene	7.129	130	2963	1.914	ug/l	58
64) Tetrachloroethene	9.281	164	2605	1.785	ug/l	91

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

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