

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062921\
 Data File : VX023009.D
 Acq On : 29 Jun 2021 17:29
 Operator : JC/MD
 Sample : M2875-18
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-9

Quant Time: Jun 30 01:45:11 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062321W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 23 19:12:26 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	55664	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.769	114	116395	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.061	117	100428	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	38747	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	52928	52.394	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 104.780%	
35) Dibromofluoromethane	5.391	113	35306	47.767	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 95.540%	
50) Toluene-d8	8.653	98	140952	48.658	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 97.320%	
62) 4-Bromofluorobenzene	11.085	95	48902	45.233	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 90.460%	
Target Compounds						
					Qvalue	
16) Acetone	2.386	43	3071	6.857	ug/l	90
30) Chloroform	5.098	83	2720	1.774	ug/l	98
64) Tetrachloroethene	9.281	164	2434	3.449	ug/l	# 81
67) Ethyl Benzene	10.195	91	5908	1.517	ug/l	95
68) m/p-Xylenes	10.305	106	4106	2.839	ug/l	90
69) o-Xylene	10.646	106	854	0.607	ug/l	86
73) Isopropylbenzene	10.969	105	1002	0.306	ug/l	80
78) n-propylbenzene	11.311	91	1295	0.326	ug/l	95
80) 1,3,5-Trimethylbenzene	11.457	105	1105	0.410	ug/l	84
84) 1,2,4-Trimethylbenzene	11.756	105	8461	3.123	ug/l	93
95) Naphthalene	13.780	128	13365	5.991	ug/l	99

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

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