

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060622\
 Data File : VX029243.D
 Acq On : 06 Jun 2022 17:07
 Operator : JC/MD
 Sample : N3160-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-4

Quant Time: Jun 07 06:05:28 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X053122W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 01 04:45:42 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	291000	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.763	114	545608	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	538165	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	252250	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	202387	61.676	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	= 123.360%	
35) Dibromofluoromethane	5.385	113	185243	51.115	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	= 102.220%	
50) Toluene-d8	8.653	98	654631	49.389	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	= 98.780%	
62) 4-Bromofluorobenzene	11.085	95	262074	51.150	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	= 102.300%	
Target Compounds						
					Qvalue	
3) Chloromethane	1.288	50	987	0.373	ug/l	# 87
16) Acetone	2.380	43	13944	10.791	ug/l	98
20) Methylene Chloride	2.788	84	819	0.272	ug/l	# 91
25) 2-Butanone	4.586	43	7011	3.364	ug/l	# 66
30) Chloroform	5.099	83	125154	23.160	ug/l	95
40) Benzene	6.050	78	5078	0.368	ug/l	# 87
52) Toluene	8.720	92	3959	0.429	ug/l	96

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

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