

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060622\
 Data File : VX029245.D
 Acq On : 06 Jun 2022 17:54
 Operator : JC/MD
 Sample : N3160-06
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-6

Quant Time: Jun 07 06:05:54 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	305878	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.763	114	573429	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	552275	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	265823	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	209481	60.733	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	= 121.460%	
35) Dibromofluoromethane	5.391	113	189526	49.759	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	= 99.520%	
50) Toluene-d8	8.653	98	678499	48.706	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	= 97.420%	
62) 4-Bromofluorobenzene	11.079	95	277523	51.537	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	= 103.080%	
Target Compounds						
					Qvalue	
3) Chloromethane	1.295	50	1035	0.372	ug/l	# 81
16) Acetone	2.380	43	12277	9.039	ug/l	91
20) Methylene Chloride	2.788	84	834	0.263	ug/l	# 86
25) 2-Butanone	4.568	43	4884	2.229	ug/l	# 77
30) Chloroform	5.099	83	122948	21.645	ug/l	98
40) Benzene	6.044	78	4514	0.311	ug/l	91
52) Toluene	8.726	92	3097	0.319	ug/l	96

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

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