

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX111122\
 Data File : VX032775.D
 Acq On : 11 Nov 2022 15:24
 Operator : JC/MD
 Sample : N5559-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 001-WILLETS-PT-BLVD(NOV)

Quant Time: Nov 14 05:13:17 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\624X111022W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
 QLast Update : Fri Nov 11 05:24:06 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Bromochloromethane	4.903	128	10230	30.000	ug/l	0.00
28) 1,4-Difluorobenzene	6.763	114	72695	30.000	ug/l	0.00
57) Chlorobenzene-d5	10.055	117	78546	30.000	ug/l	0.00
System Monitoring Compounds						
27) 1,2-Dichloroethane-d4	5.958	65	31131	31.544	ug/l	0.00
Spiked Amount	30.000	Range	91 - 110	Recovery	=	105.133%
60) 4-Bromofluorobenzene	11.079	95	41330	27.508	ug/l	0.00
Spiked Amount	30.000	Range	63 - 112	Recovery	=	91.700%
63) Toluene-d8	8.653	98	78883	29.759	ug/l	0.00
Spiked Amount	30.000	Range	91 - 112	Recovery	=	99.200%
Target Compounds						
					Qvalue	
15) Acetone	2.386	58	28620	208.266	ug/l	81
16) Carbon Disulfide	2.508	76	954	0.385	ug/l #	91
25) Chloroform	5.092	83	965	0.435	ug/l #	61
30) 2-Butanone	4.568	43	12202	16.231	ug/l #	85
58) 4-Methyl-2-Pentanone	8.580	43	2589	1.684	ug/l #	87
62) Toluene	8.720	91	14606	2.859	ug/l	95
82) p-Isopropyltoluene	12.012	119	4236	0.870	ug/l	94

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

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