

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX011022\
 Data File : VX026407.D
 Acq On : 10 Jan 2022 17:03
 Operator : JC/MD
 Sample : N1046-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 002-35TH-AVE(JAN)

Quant Time: Jan 11 04:56:43 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\624X011022W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
 QLast Update : Mon Jan 10 12:50:55 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Bromochloromethane	4.904	128	20145	30.000	ug/l	0.00
28) 1,4-Difluorobenzene	6.763	114	113287	30.000	ug/l	0.00
57) Chlorobenzene-d5	10.055	117	97396	30.000	ug/l	0.00
System Monitoring Compounds						
27) 1,2-Dichloroethane-d4	5.958	65	49884	30.795	ug/l	0.00
Spiked Amount	30.000	Range	91 - 110	Recovery	=	102.633%
60) 4-Bromofluorobenzene	11.079	95	46861	28.830	ug/l	0.00
Spiked Amount	30.000	Range	63 - 112	Recovery	=	96.100%
63) Toluene-d8	8.653	98	133049	29.736	ug/l	0.00
Spiked Amount	30.000	Range	91 - 112	Recovery	=	99.133%
Target Compounds						
					Qvalue	
15) Acetone	2.386	58	98589	495.998	ug/l	99
16) Carbon Disulfide	2.520	76	3326	1.088	ug/l #	93
30) 2-Butanone	4.568	43	8531	8.396	ug/l #	87
58) 4-Methyl-2-Pentanone	8.580	43	7009	3.858	ug/l	98
62) Toluene	8.720	91	23593	4.537	ug/l	96

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

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