

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX030322\
 Data File : VX027286.D
 Acq On : 03 Mar 2022 22:10
 Operator : JC/MD
 Sample : N1728-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 17 Sample Multiplier: 1

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

Quant Time: Mar 04 02:42:35 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\624X030322W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
 QLast Update : Fri Mar 04 01:00:22 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Bromochloromethane	4.910	128	12545	30.000	ug/l	0.00
28) 1,4-Difluorobenzene	6.769	114	66792	30.000	ug/l	0.00
57) Chlorobenzene-d5	10.055	117	60700	30.000	ug/l	0.00
System Monitoring Compounds						
27) 1,2-Dichloroethane-d4	5.958	65	25484	29.343	ug/l	0.00
Spiked Amount	30.000	Range	91 - 110	Recovery	=	97.800%
60) 4-Bromofluorobenzene	11.085	95	28044	28.519	ug/l	0.00
Spiked Amount	30.000	Range	63 - 112	Recovery	=	95.067%
63) Toluene-d8	8.653	98	78087	29.427	ug/l	0.00
Spiked Amount	30.000	Range	91 - 112	Recovery	=	98.100%
Target Compounds						
					Qvalue	
15) Acetone	2.379	58	7767	64.186	ug/l	98
16) Carbon Disulfide	2.526	76	5887	3.086	ug/l	99
30) 2-Butanone	4.568	43	4241	7.358	ug/l #	77
58) 4-Methyl-2-Pentanone	8.580	43	4594	4.184	ug/l	99
62) Toluene	8.726	91	2943	0.890	ug/l	86
67) m/p-Xylenes	10.305	106	428	0.293	ug/l	94

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX030322\
Data File : VX027286.D
Acq On : 03 Mar 2022 22:10
Operator : JC/MD
Sample : N1728-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 17 Sample Multiplier: 1

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

Quant Time: Mar 04 02:42:35 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\624X030322W.M
Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
QLast Update : Fri Mar 04 01:00:22 2022
Response via : Initial Calibration

