

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121624\
 Data File : VN085211.D
 Acq On : 16 Dec 2024 11:57
 Operator : JC\MD
 Sample : P5276-03
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 001-WILLETS-PT-BLVD(DEC)

Quant Time: Dec 16 12:22:32 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\624N120624W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
 QLast Update : Fri Dec 06 22:50:42 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Bromochloromethane	7.812	128	27333	30.000	ug/l	0.00
28) 1,4-Difluorobenzene	9.100	114	145796	30.000	ug/l	0.00
57) Chlorobenzene-d5	11.865	117	134362	30.000	ug/l	0.00
System Monitoring Compounds						
27) 1,2-Dichloroethane-d4	8.576	65	65807	30.343	ug/l	0.00
Spiked Amount	30.000	Range	91 - 110	Recovery	=	101.133%
60) 4-Bromofluorobenzene	12.847	95	59091	26.580	ug/l	0.00
Spiked Amount	30.000	Range	63 - 112	Recovery	=	88.600%
63) Toluene-d8	10.565	98	178120	28.261	ug/l	0.00
Spiked Amount	30.000	Range	91 - 112	Recovery	=	94.200%
Target Compounds						
					Qvalue	
15) Acetone	4.424	58	89517	405.114	ug/l	100
30) 2-Butanone	7.488	43	18689	17.311	ug/l #	98
62) Toluene	10.629	91	591725	75.514	ug/l	99
82) p-Isopropyltoluene	13.723	119	3174	0.494	ug/l	96

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

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