

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN112524\
 Data File : VN085026.D
 Acq On : 25 Nov 2024 13:14
 Operator : JC\MD
 Sample : P4967-01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 001-WILLETS-PT-BLVD

Manual Integrations
 APPROVED

Reviewed By :John Carlone 11/25/2024
 Supervised By :Mahesh Dadoda 11/26/2024

Quant Time: Nov 25 15:14:10 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\624N103124W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
 QLast Update : Fri Nov 01 02:38:04 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Bromochloromethane	7.806	128	32568	30.000	ug/l	0.00
28) 1,4-Difluorobenzene	9.094	114	164674	30.000	ug/l	0.00
57) Chlorobenzene-d5	11.865	117	145379	30.000	ug/l	0.00
System Monitoring Compounds						
27) 1,2-Dichloroethane-d4	8.571	65	74069	28.289	ug/l	0.00
Spiked Amount 30.000	Range 91 - 110		Recovery =	94.300%		
60) 4-Bromofluorobenzene	12.847	95	64237	25.662	ug/l	0.00
Spiked Amount 30.000	Range 63 - 112		Recovery =	85.533%		
63) Toluene-d8	10.565	98	195474	28.403	ug/l	0.00
Spiked Amount 30.000	Range 91 - 112		Recovery =	94.667%		
Target Compounds						
15) Acetone	4.424	58	2744m	11.034	ug/l	Qvalue
25) Chloroform	7.959	83	4251	1.066	ug/l	87
41) Bromodichloromethane	9.882	83	3850	1.301	ug/l #	68
53) Dibromochloromethane	11.353	129	5331	2.460	ug/l	86
56) Bromoform	12.576	173	3101	2.135	ug/l #	98
62) Toluene	10.624	91	154315	18.637	ug/l	100
82) p-Isopropyltoluene	13.729	119	1861	0.285	ug/l	97

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

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