

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN100824\
 Data File : VN084348.D
 Acq On : 08 Oct 2024 14:53
 Operator : JC\MD
 Sample : P4334-01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 13 Sample Multiplier: 1

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

Quant Time: Oct 09 02:30:43 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\624N100824W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
 QLast Update : Wed Oct 09 02:21:08 2024
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By :John Carlone 10/09/2024
 Supervised By :Mahesh Dadoda 10/09/2024

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

Internal Standards						
1) Bromochloromethane	7.818	128	31390	30.000	ug/l	0.00
28) 1,4-Difluorobenzene	9.100	114	165437	30.000	ug/l	0.00
57) Chlorobenzene-d5	11.865	117	154322	30.000	ug/l	0.00
System Monitoring Compounds						
27) 1,2-Dichloroethane-d4	8.576	65	79279	31.371	ug/l	0.00
Spiked Amount 30.000	Range 91 - 110		Recovery = 104.567%			
60) 4-Bromofluorobenzene	12.847	95	69842	27.422	ug/l	0.00
Spiked Amount 30.000	Range 63 - 112		Recovery = 91.400%			
63) Toluene-d8	10.565	98	210498	29.444	ug/l	0.00
Spiked Amount 30.000	Range 91 - 112		Recovery = 98.133%			
Target Compounds						
15) Acetone	4.424	58	309723	986.299	ug/l	97
25) Chloroform	7.971	83	5224	1.320	ug/l	98
30) 2-Butanone	7.482	43	36140	24.994	ug/l	99
58) 4-Methyl-2-Pentanone	10.441	43	12613	4.387	ug/l	97
62) Toluene	10.629	91	495692	54.868	ug/l	99
82) p-Isopropyltoluene	13.729	119	10396	1.363	ug/l	98
91) Naphthalene	15.647	128	7120m	1.013	ug/l	

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

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