

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN100518\
 Data File : VN051687.D
 Acq On : 5 Oct 2018 15:16
 Operator : MD\SY
 Sample : J5239-01
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 16 Sample Multiplier: 28

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

Quant Time: Oct 05 16:38:27 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA N\METHODS\624N100518W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
 QLast Update : Fri Oct 05 11:11:26 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	7.20	128	80810	30.00	ug/l	0.00
28) 1,4-Difluorobenzene	8.59	114	463050	30.00	ug/l	0.00
57) Chlorobenzene-d5	11.41	117	378792	30.00	ug/l	0.00

System Monitoring Compounds

27) 1,2-Dichloroethane-d4	8.03	65	165290	29.47	ug/l	0.00
Spiked Amount	30.000	Range	50 - 169	Recovery	=	98.23%
60) 4-Bromofluorobenzene	12.40	95	135654	23.47	ug/l	0.00
Spiked Amount	30.000	Range	56 - 143	Recovery	=	78.23%
63) Toluene-d8	10.09	98	536188	31.88	ug/l	0.00
Spiked Amount	30.000	Range	66 - 137	Recovery	=	106.27%

Target Compounds

					Ovalue
15) Acetone	3.82	58	272717	938.914	ug/l 79
25) Chloroform	7.38	83	130075	14.917	ug/l 94
41) Bromodichloromethane	9.41	83	12858	1.826	ug/l 89
62) Toluene	10.16	91	1945572	101.396	ug/l 99
82) p-Isopropyltoluene	13.29	119	78501	3.815	ug/l 98

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

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