

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN100518\
 Data File : VN051688.D
 Acq On : 5 Oct 2018 15:41
 Operator : MD\SY
 Sample : J5239-02
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 17 Sample Multiplier: 28

Instrument :
 MSVOA_N
 ClientSampleId :
 002M-35TH-AVE(AUG)

Quant Time: Oct 05 16:39:09 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	82519	30.00	ug/l	0.00
28) 1,4-Difluorobenzene	8.59	114	460915	30.00	ug/l	0.00
57) Chlorobenzene-d5	11.41	117	375059	30.00	ug/l	0.00

System Monitoring Compounds

27) 1,2-Dichloroethane-d4	8.03	65	169255	29.55	ug/l	0.00
Spiked Amount	30.000	Range	50 - 169	Recovery	=	98.50%
60) 4-Bromofluorobenzene	12.40	95	134622	23.52	ug/l	0.00
Spiked Amount	30.000	Range	56 - 143	Recovery	=	78.40%
63) Toluene-d8	10.09	98	535017	32.13	ug/l	0.00
Spiked Amount	30.000	Range	66 - 137	Recovery	=	107.10%

Target Compounds

					Ovalue
15) Acetone	3.82	58	194677	656.356	ug/l 82
25) Chloroform	7.37	83	150502	16.902	ug/l 98
41) Bromodichloromethane	9.41	83	14785	2.110	ug/l # 91
62) Toluene	10.16	91	1219939	64.211	ug/l 100
82) p-Isopropyltoluene	13.29	119	49205	2.415	ug/l 98

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

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