

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN101518\
 Data File : VN051867.D
 Acq On : 15 Oct 2018 12:02
 Operator : MD\SY
 Sample : J5512-02
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 7 Sample Multiplier: 28

Instrument :
 MSVOA_N
 ClientSampleId :
 002M-35TH-AVE-SEP

Quant Time: Oct 16 01:16:09 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\624N101218W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
 QLast Update : Fri Oct 12 10:23:44 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	7.20	128	116505	30.00	ug/l	0.00
28) 1,4-Difluorobenzene	8.59	114	649953	30.00	ug/l	0.00
57) Chlorobenzene-d5	11.41	117	544880	30.00	ug/l	0.00

System Monitoring Compounds

27) 1,2-Dichloroethane-d4	8.03	65	216295	30.51	ug/l	0.00
Spiked Amount	30.000	Range	50 - 169	Recovery	=	101.70%
60) 4-Bromofluorobenzene	12.40	95	199122	25.34	ug/l	0.00
Spiked Amount	30.000	Range	56 - 143	Recovery	=	84.47%
63) Toluene-d8	10.09	98	737924	31.19	ug/l	0.00
Spiked Amount	30.000	Range	66 - 137	Recovery	=	103.97%

Target Compounds

						Qvalue
15) Acetone	3.82	58	366197	995.08	ug/l	96
25) Chloroform	7.37	83	237228	20.18	ug/l	94
30) 2-Butanone	6.85	43	11042	6.31	ug/l #	86
41) Bromodichloromethane	9.40	83	15678	1.75	ug/l	95
58) 4-Methyl-2-Pentanone	9.99	43	20473	5.61	ug/l #	88
62) Toluene	10.16	91	5713674	211.08	ug/l	99
82) p-Isopropyltoluene	13.29	119	150727	5.31	ug/l	97

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

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