

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN011019\
 Data File : VN053336.D
 Acq On : 10 Jan 2019 12:47
 Operator : MD\SY
 Sample : K1095-01
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 9 Sample Multiplier: 28

Instrument :
 MSVOA_N
 ClientSampleId :
 2-AVE-29

Manual Integrations
 APPROVED

MMDadoda
 1/11/2019 10:11:40 AM

Quant Time: Jan 11 06:12:42 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA N\METHODS\624N122818W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
 QLast Update : Fri Dec 28 10:33:50 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Bromochloromethane	7.20	128	148395	30.00	ug/l	0.00
28) 1,4-Difluorobenzene	8.59	114	853359	30.00	ug/l	0.00
57) Chlorobenzene-d5	11.41	117	748513	30.00	ug/l	0.00
System Monitoring Compounds						
27) 1,2-Dichloroethane-d4	8.03	65	286729	30.42	ug/l	0.00
Spiked Amount 30.000	Range 50 - 169		Recovery =	101.40%		
60) 4-Bromofluorobenzene	12.40	95	321366	28.06	ug/l	0.00
Spiked Amount 30.000	Range 56 - 143		Recovery =	93.53%		
63) Toluene-d8	10.09	98	1000432	29.67	ug/l	0.00
Spiked Amount 30.000	Range 66 - 137		Recovery =	98.90%		
Target Compounds						
15) Acetone	3.82	58	5273m	8.229	ug/l	
22) cis-1,2-Dichloroethene	6.83	96	163323	15.867	ug/l	88
25) Chloroform	7.37	83	46838	2.873	ug/l	98
37) Trichloroethene	8.83	130	47805	4.543	ug/l	95
61) Tetrachloroethene	10.63	164	90243	8.158	ug/l	96

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

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