

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV091819\

Quantitation Report (QT Reviewed)

Data File : VV012895.D

Acq On : 18 Sep 2019 14:42

Operator : SY/MD

Sample : K4833-06

Misc : 25.0mL/MSVOA_V/WATER

ALS Vial : 10 Sample Multiplier: 1

Quant Time: Sep 19 03:49:17 2019

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR091719WMA.M

Quant Title : TRACE VOA SOM01.0

QLast Update : Thu Sep 19 01:01:31 2019

Response via : Initial Calibration

Instrument :

MSVOA_V

ClientSampled :

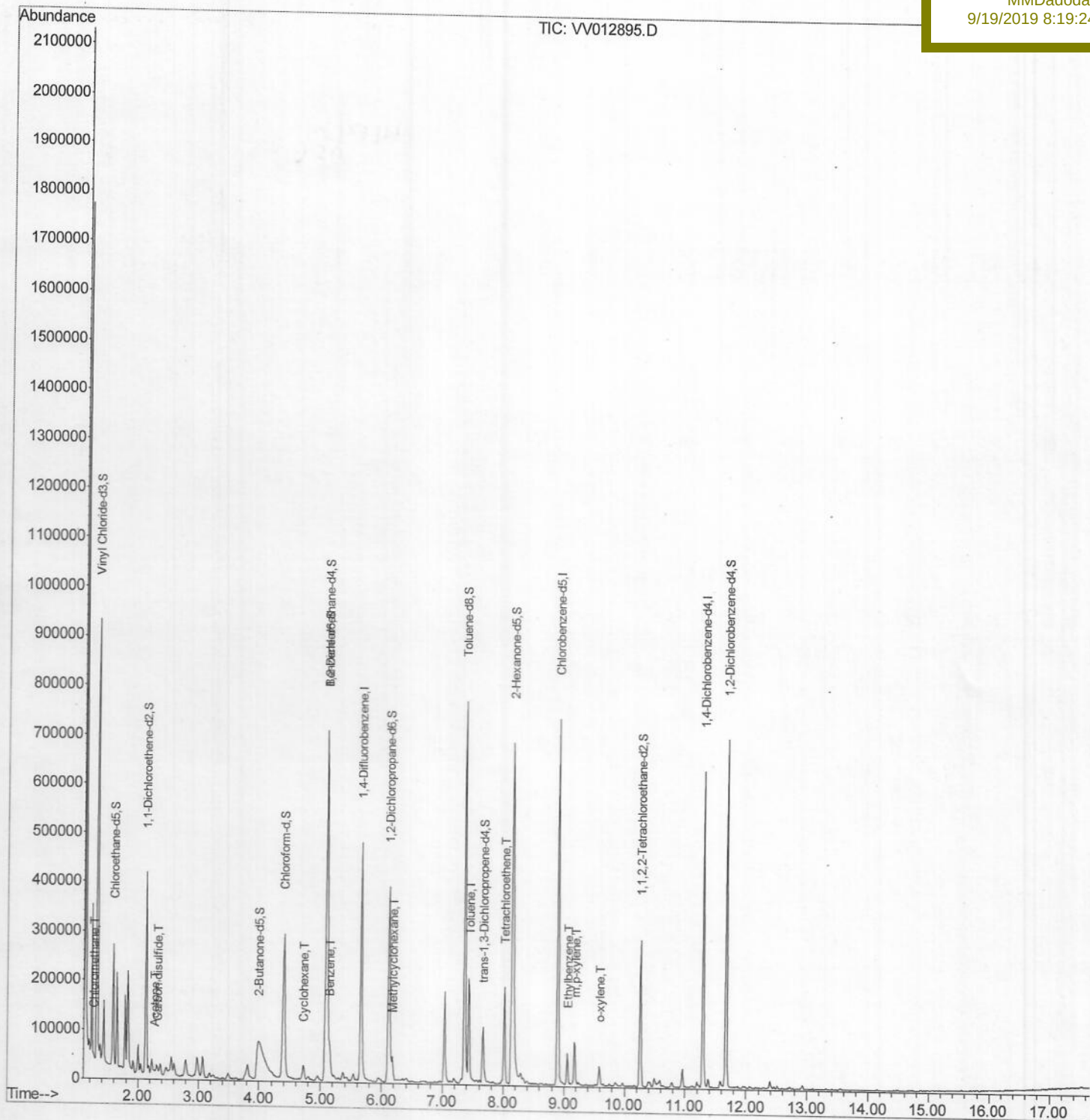
ETOH3

Manual Integrations

APPROVED

MMDadoda

9/19/2019 8:19:24 AM



SOMVTR091719WMA.M Thu Sep 19 03:47:15 2019

Page: 2

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV091819\

Quantitation Report (Qedit)

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Response via : Initial Calibration

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MSVOA_V

ClientSampleId :

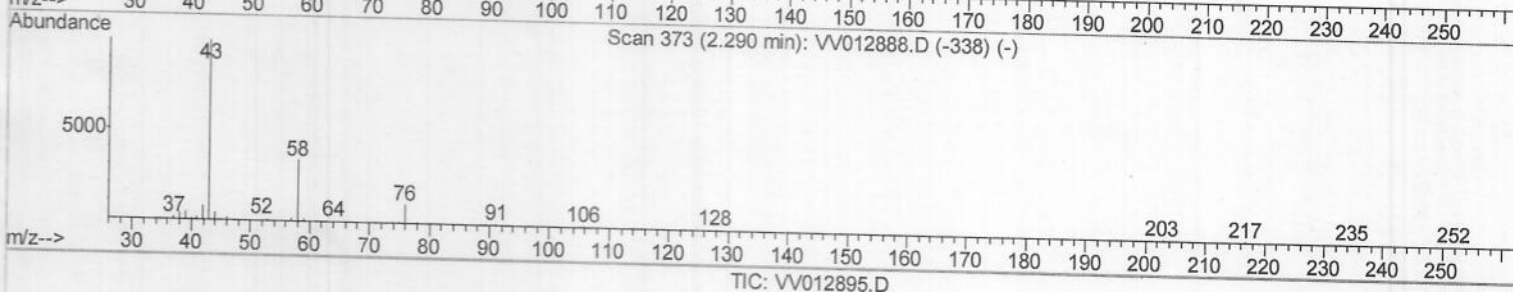
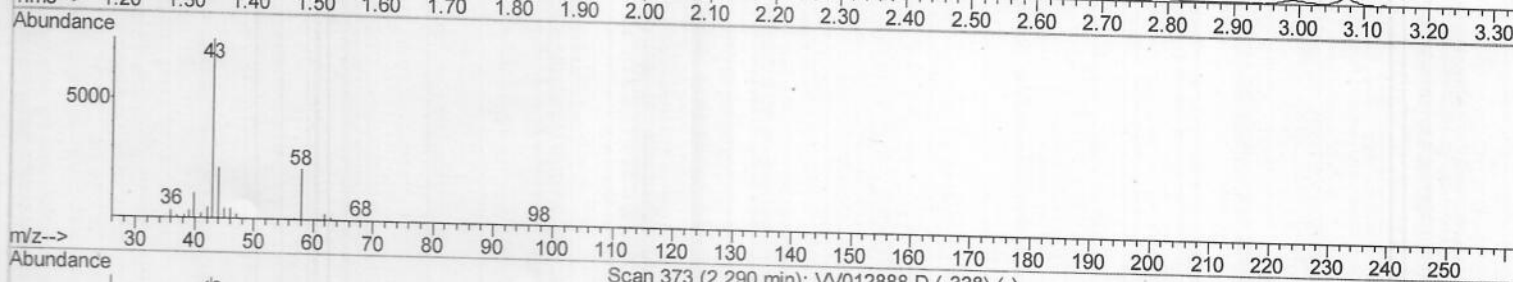
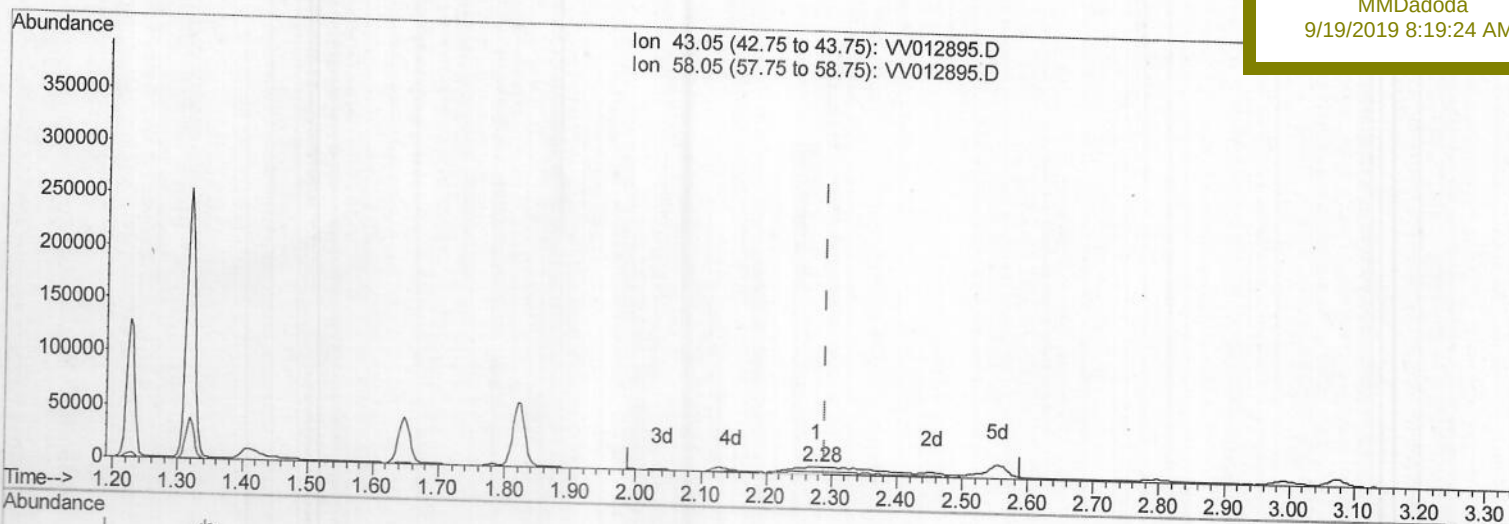
ETOH3

Manual Integrations

APPROVED

MMDadoda

9/19/2019 8:19:24 AM



(13) Acetone (T)

2.277min (-0.013) 4.99ug/L

response 22763

Ion	Exp%	Act%
43.05	100	100
58.05	12.20	2.38
0.00	0.00	0.00
0.00	0.00	0.00

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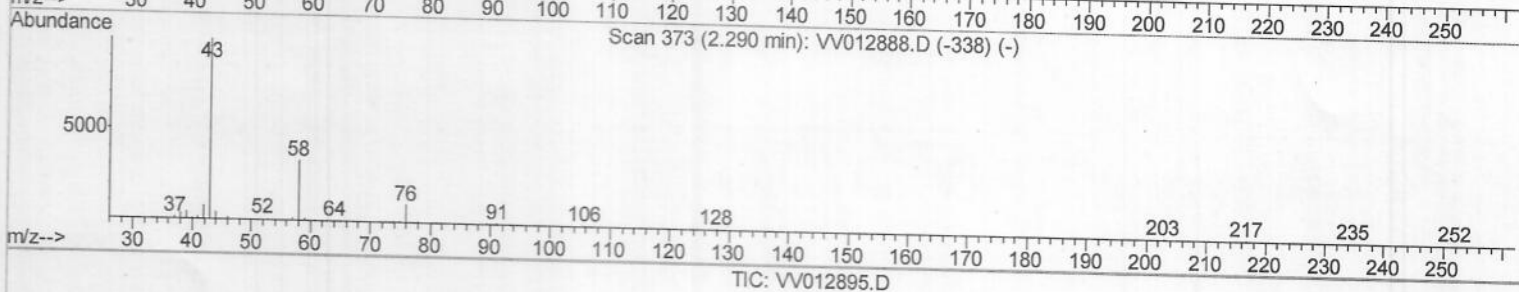
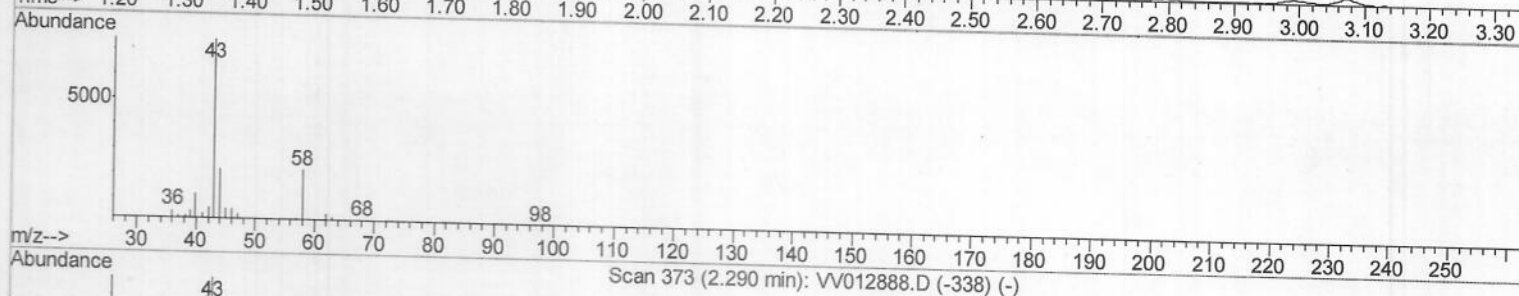
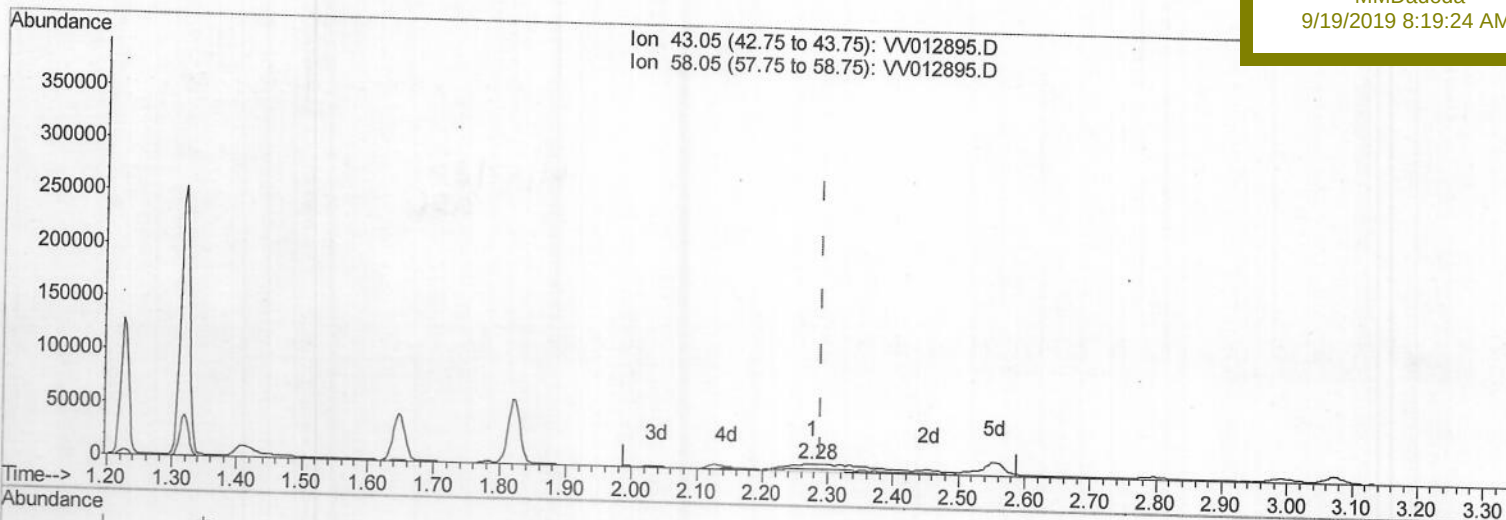
ETOH3

Manual Integrations

APPROVED

MMDadoda

9/19/2019 8:19:24 AM



TIC: VV012895.D

(13) Acetone (T)

2.277min (-0.013) 8.69ug/L m

response 39631

Ion	Exp%	Act%
43.05	100	100
58.05	12.20	1.37
0.00	0.00	0.00
0.00	0.00	0.00

> M.D 09/24/19

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Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Difluorobenzene	5.66	114	416171	5.00	ug/L	0.00
28) Chlorobenzene-d5	8.89	117	402222	5.00	ug/L	0.00
60) 1,4-Dichlorobenzene-d4	11.29	152	166443	5.00	ug/L	0.00

System Monitoring Compounds

4) Vinyl Chloride-d3	1.32	65	167513	4.07	ug/L	0.00
Spiked Amount	5.000	Range	40 - 130	Recovery	=	81.40%
7) Chloroethane-d5	1.58	69	151227	4.66	ug/L	0.00
Spiked Amount	5.000	Range	65 - 130	Recovery	=	93.20%
11) 1,1-Dichloroethene-d2	2.13	63	210122	3.46	ug/L	0.00
Spiked Amount	5.000	Range	60 - 125	Recovery	=	69.20%
20) 2-Butanone-d5	3.99	46	398826	46.24	ug/L	0.00
Spiked Amount	50.000	Range	40 - 130	Recovery	=	92.48%
24) Chloroform-d	4.40	84	304163	4.52	ug/L	0.00
Spiked Amount	5.000	Range	70 - 125	Recovery	=	90.40%
26) 1,2-Dichloroethane-d4	5.08	65	158967	4.84	ug/L	0.00
Spiked Amount	5.000	Range	70 - 130	Recovery	=	96.80%
32) Benzene-d6	5.10	84	636938	4.78	ug/L	0.00
Spiked Amount	5.000	Range	70 - 125	Recovery	=	95.60%
36) 1,2-Dichloropropane-d6	6.12	67	200043	4.93	ug/L	0.00
Spiked Amount	5.000	Range	60 - 140	Recovery	=	98.60%
41) Toluene-d8	7.36	98	516169	4.36	ug/L	0.00
Spiked Amount	5.000	Range	70 - 130	Recovery	=	87.20%
43) trans-1,3-Dichloropropene-	7.66	79	65829	4.51	ug/L	0.00
Spiked Amount	5.000	Range	55 - 130	Recovery	=	90.20%
46) 2-Hexanone-d5	8.14	63	279924	45.64	ug/L	0.00
Spiked Amount	50.000	Range	45 - 130	Recovery	=	91.28%
57) 1,1,2,2-Tetrachloroethane-	10.26	84	143021	4.55	ug/L	0.00
Spiked Amount	5.000	Range	65 - 120	Recovery	=	91.00%
64) 1,2-Dichlorobenzene-d4	11.67	152	184361	5.13	ug/L	0.00
Spiked Amount	5.000	Range	80 - 120	Recovery	=	102.60%

Target Compounds

	R.T.	QIon	Response	Conc	Units	Qvalue
3) Chloromethane	1.26	50	5655	0.195	ug/L	98
13) Acetone	2.28	43	39631m	8.689	ug/L	96
14) Carbon disulfide	2.32	76	10294	0.225	ug/L	96
30) Cyclohexane	4.72	56	13473	0.387	ug/L	59
33) Benzene	5.15	78	55914	0.536	ug/L	100
35) Methylcyclohexane	6.18	83	15160	0.420	ug/L	91
42) Toluene	7.43	91	152893	1.407	ug/L	100
47) Tetrachloroethene	8.02	164	45115	2.088	ug/L	98
52) Ethylbenzene	9.05	91	44239	0.380	ug/L	99
53) m,p-xylene	9.18	106	24272	0.565	ug/L	96
54) o-xylene	9.59	106	10269	0.243	ug/L	98

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