

Data Path : Z:\voasrv\HPCHEM1\MSVOA V\Data\VV041319\

Data File : VV010326.D

Acq On : 13 Apr 2019 18:43

Operator : SY/MD

Sample : K2297-04MS

Misc : 25mL/MSVOA V/WATER

ALS Vial : 11 Sample Multiplier: 1

Quant Time: Apr 14 01:49:52 2019

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR041319WMA.M

Quant Title : TRACE VOA SOM01.0

QLast Update : Sat Apr 13 14:26:57 2019

Response via : Initial Calibration

Instrument :

MSVOA_V

ClientSampleId :

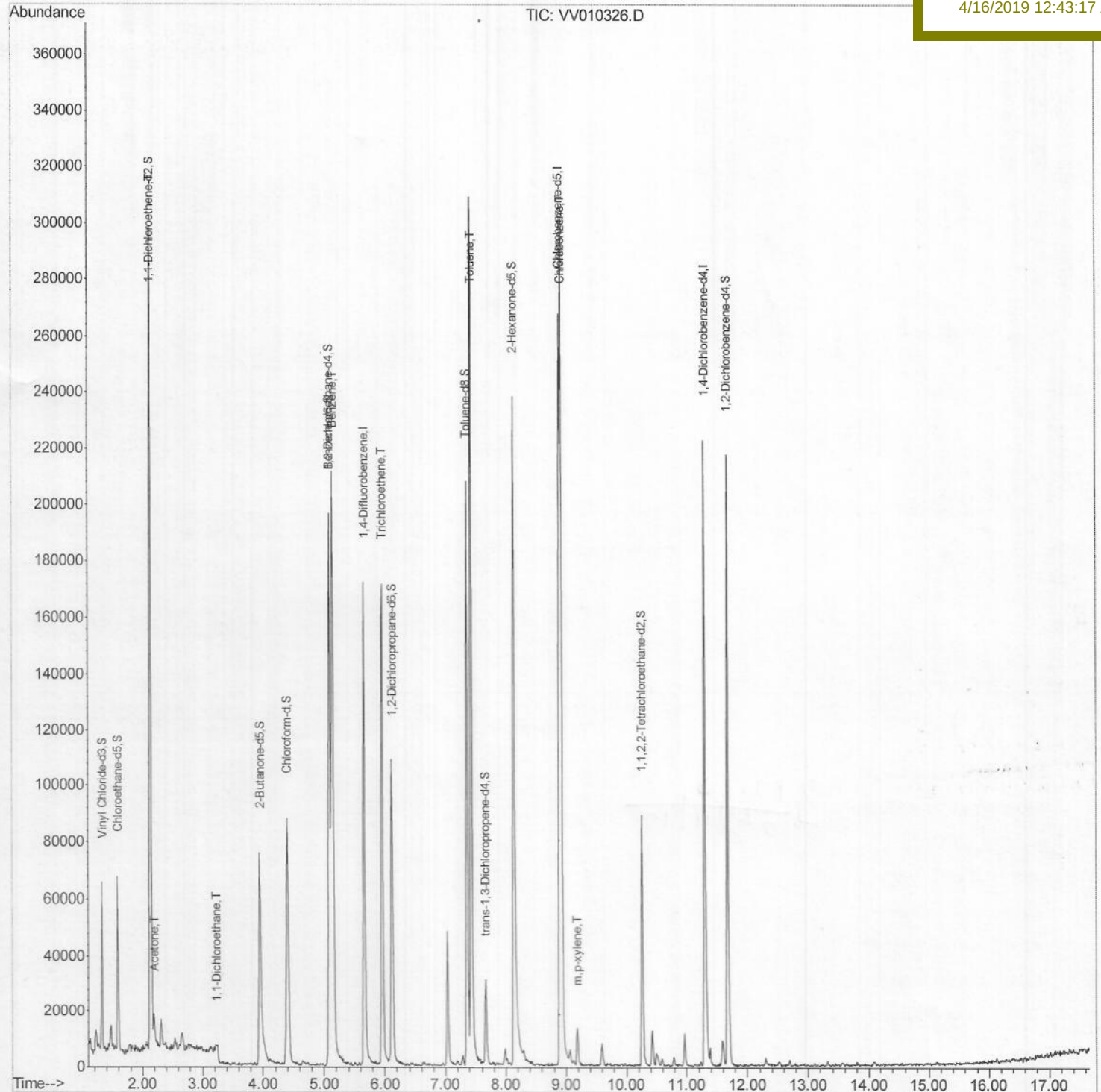
E3YC4MS

Manual Integrations

APPROVED

MMDadoda

4/16/2019 12:43:17 AM



Quantitation Report (Qedit)

Data Path : Z:\voasrv\HPCHEM1\MSVOA V\Data\VV041319\

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Response via : Initial Calibration

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MSVOA_V

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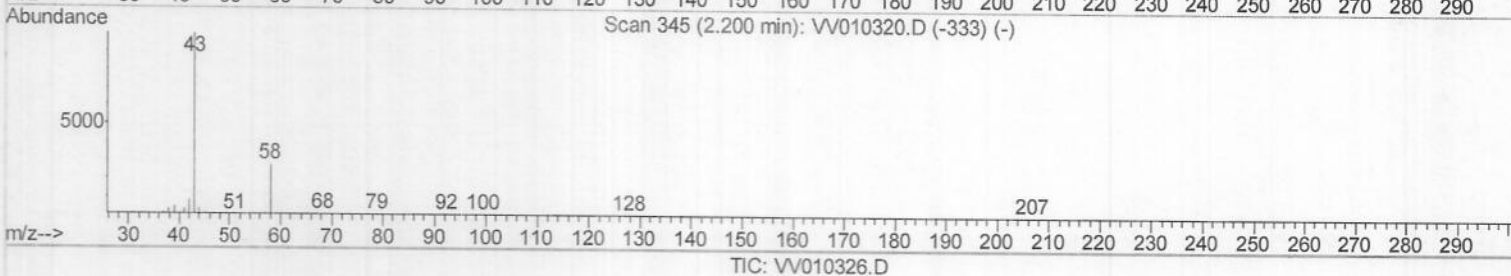
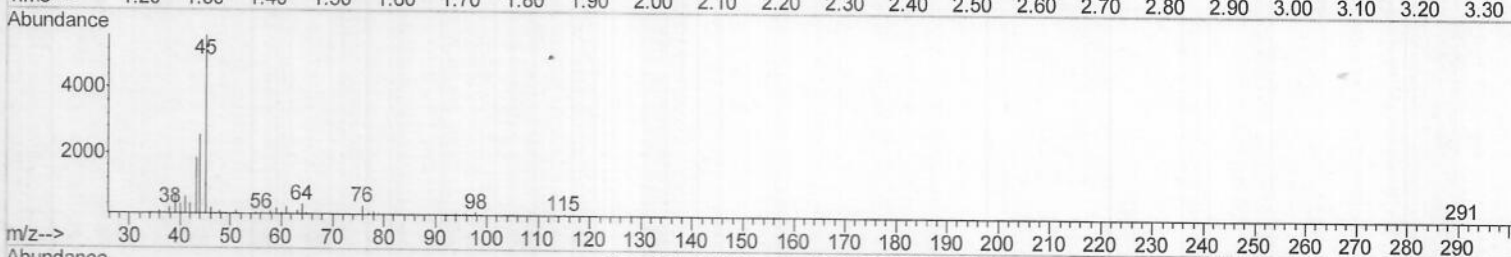
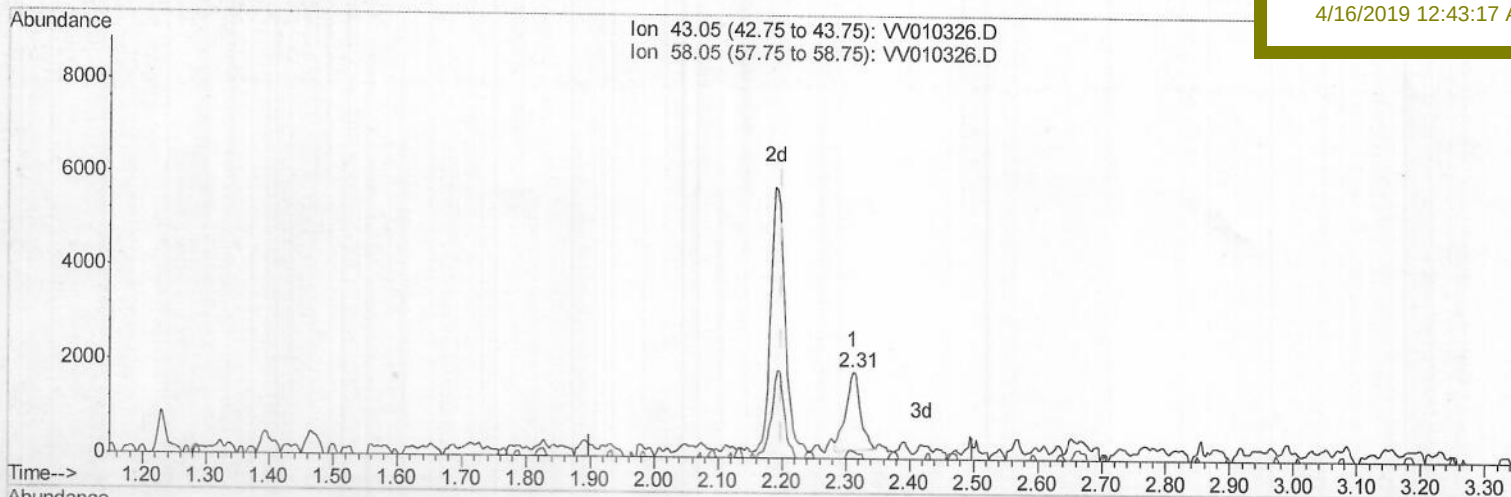
E3YC4MS

Manual Integrations

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4/16/2019 12:43:17 AM



(13) Acetone (T)

2.312min (+0.113) 1.53ug/L

response 2617

Ion	Exp%	Act%
43.05	100	100
58.05	11.50	8.29
0.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

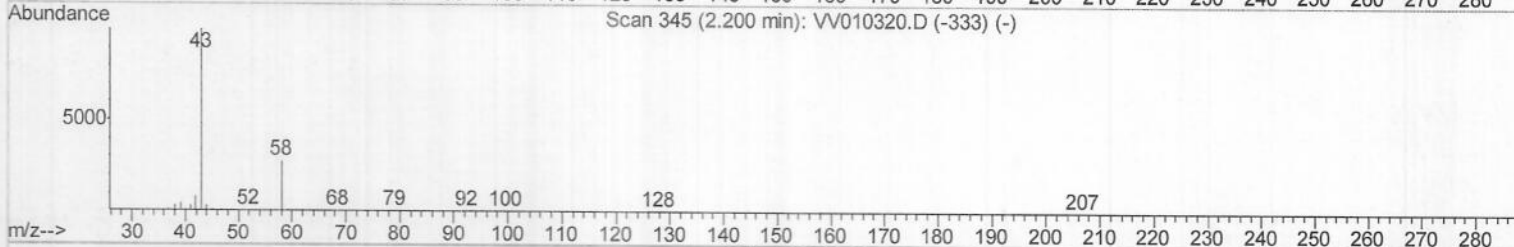
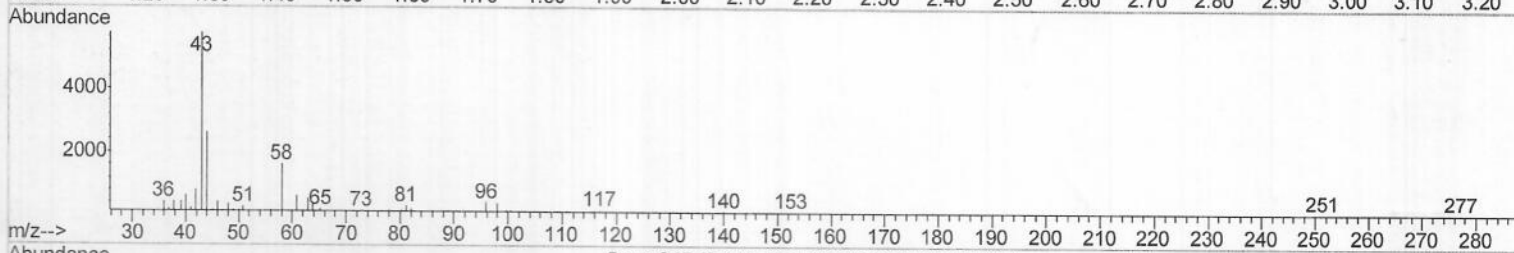
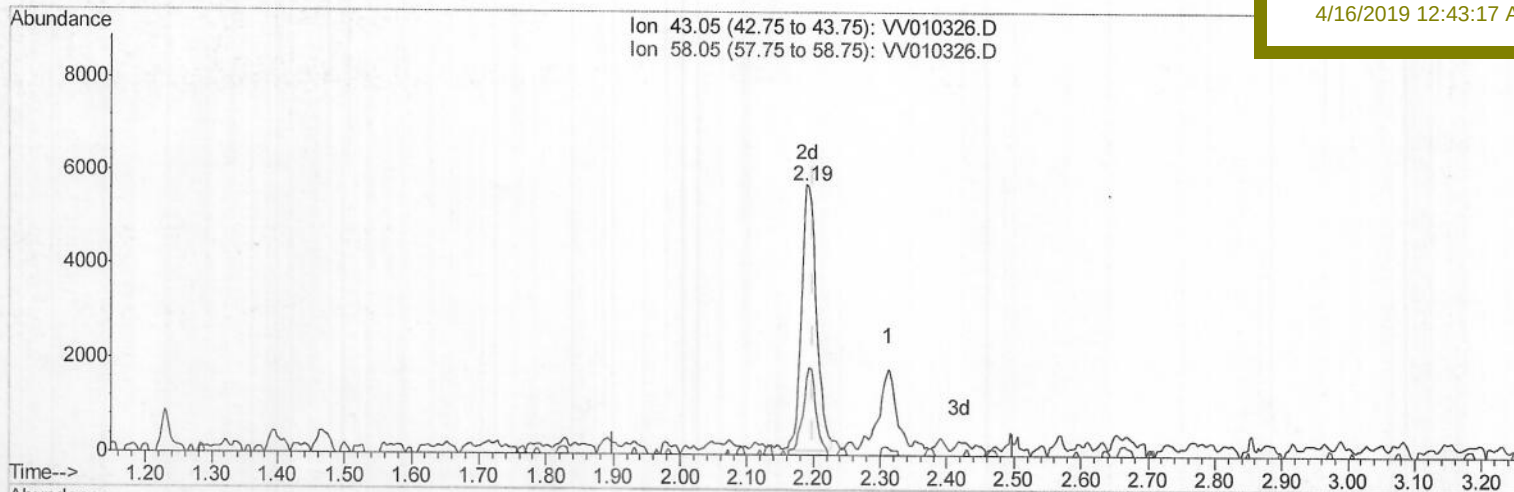
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 Data File : VV010326.D
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 Operator : SY/MD
 Sample : K2297-04MS
 Misc : 25mL/MSVOA V/WATER
 ALS Vial : 11 Sample Multiplier: 1

Quant Time: Apr 14 01:14:44 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR041319WMA.M
 Quant Title : TRACE VOA SOM01.0
 QLast Update : Sat Apr 13 14:26:57 2019
 Response via : Initial Calibration

Instrument :
 MSVOA_V
ClientSampleId :
 E3YC4MS

Manual Integrations
APPROVED

MMDadoda
 4/16/2019 12:43:17 AM



TIC: VV010326.D

(13) Acetone (T)

2.190min (-0.010) 5.10ug/L m

response 8740

Handwritten: } M.D.
 04/16/19

Ion	Exp%	Act%
43.05	100	100
58.05	11.50	2.48
0.00	0.00	0.00
0.00	0.00	0.00

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Acq On : 13 Apr 2019 18:43

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Quant Title : TRACE VOA SOM01.0

QLast Update : Sat Apr 13 14:26:57 2019

Response via : Initial Calibration

Instrument :

MSVOA_V

ClientSampleId :

E3YC4MS

Manual Integrations

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4/16/2019 12:43:17 AM

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev (Min)
1) 1,4-Difluorobenzene	5.67	114	149893	5.00	ug/L	0.00
28) Chlorobenzene-d5	8.90	117	146576	5.00	ug/L	0.00
60) 1,4-Dichlorobenzene-d4	11.30	152	67817	5.00	ug/L	0.00

System Monitoring Compounds

4) Vinyl Chloride-d3	1.32	65	32538	4.08	ug/L	0.00
Spiked Amount	5.000	Range	40 - 130	Recovery	=	81.60%
7) Chloroethane-d5	1.58	69	36724	4.19	ug/L	0.00
Spiked Amount	5.000	Range	65 - 130	Recovery	=	83.80%
11) 1,1-Dichloroethene-d2	2.14	63	99709	4.48	ug/L	0.00
Spiked Amount	5.000	Range	60 - 125	Recovery	=	89.60%
20) 2-Butanone-d5	3.95	46	105937	52.05	ug/L	0.00
Spiked Amount	50.000	Range	40 - 130	Recovery	=	104.10%
24) Chloroform-d	4.40	84	84729	4.60	ug/L	0.00
Spiked Amount	5.000	Range	70 - 125	Recovery	=	92.00%
26) 1,2-Dichloroethane-d4	5.09	65	44384	4.77	ug/L	0.00
Spiked Amount	5.000	Range	70 - 130	Recovery	=	95.40%
32) Benzene-d6	5.10	84	159482	4.69	ug/L	0.00
Spiked Amount	5.000	Range	70 - 125	Recovery	=	93.80%
36) 1,2-Dichloropropane-d6	6.12	67	49485	4.63	ug/L	0.00
Spiked Amount	5.000	Range	60 - 140	Recovery	=	92.60%
41) Toluene-d8	7.36	98	139516	4.26	ug/L	0.00
Spiked Amount	5.000	Range	70 - 130	Recovery	=	85.20%
43) trans-1,3-Dichloropropene-	7.67	79	16685	4.89	ug/L	0.00
Spiked Amount	5.000	Range	55 - 130	Recovery	=	97.80%
46) 2-Hexanone-d5	8.14	63	90030	58.22	ug/L	0.00
Spiked Amount	50.000	Range	45 - 130	Recovery	=	116.44%
57) 1,1,2,2-Tetrachloroethane-	10.26	84	39769	4.85	ug/L	0.00
Spiked Amount	5.000	Range	65 - 120	Recovery	=	97.00%
64) 1,2-Dichlorobenzene-d4	11.68	152	61379	5.02	ug/L	0.00
Spiked Amount	5.000	Range	80 - 120	Recovery	=	100.40%

Target Compounds

					Ovalue	
12) 1,1-Dichloroethene	2.14	96	60346	4.936	ug/L	86
13) Acetone	2.19	43	8740m	5.102	ug/L	
19) 1,1-Dichloroethane	3.23	63	2611	0.112	ug/L #	83
33) Benzene	5.15	78	206872	5.031	ug/L	100
34) Trichloroethene	5.96	95	57258	4.822	ug/L	94
42) Toluene	7.43	91	226789	4.934	ug/L	96
51) Chlorobenzene	8.93	112	158334	5.104	ug/L	96
53) m,p-xylene	9.19	106	4490	0.231	ug/L	90

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

3 M.D.
04/16/19