

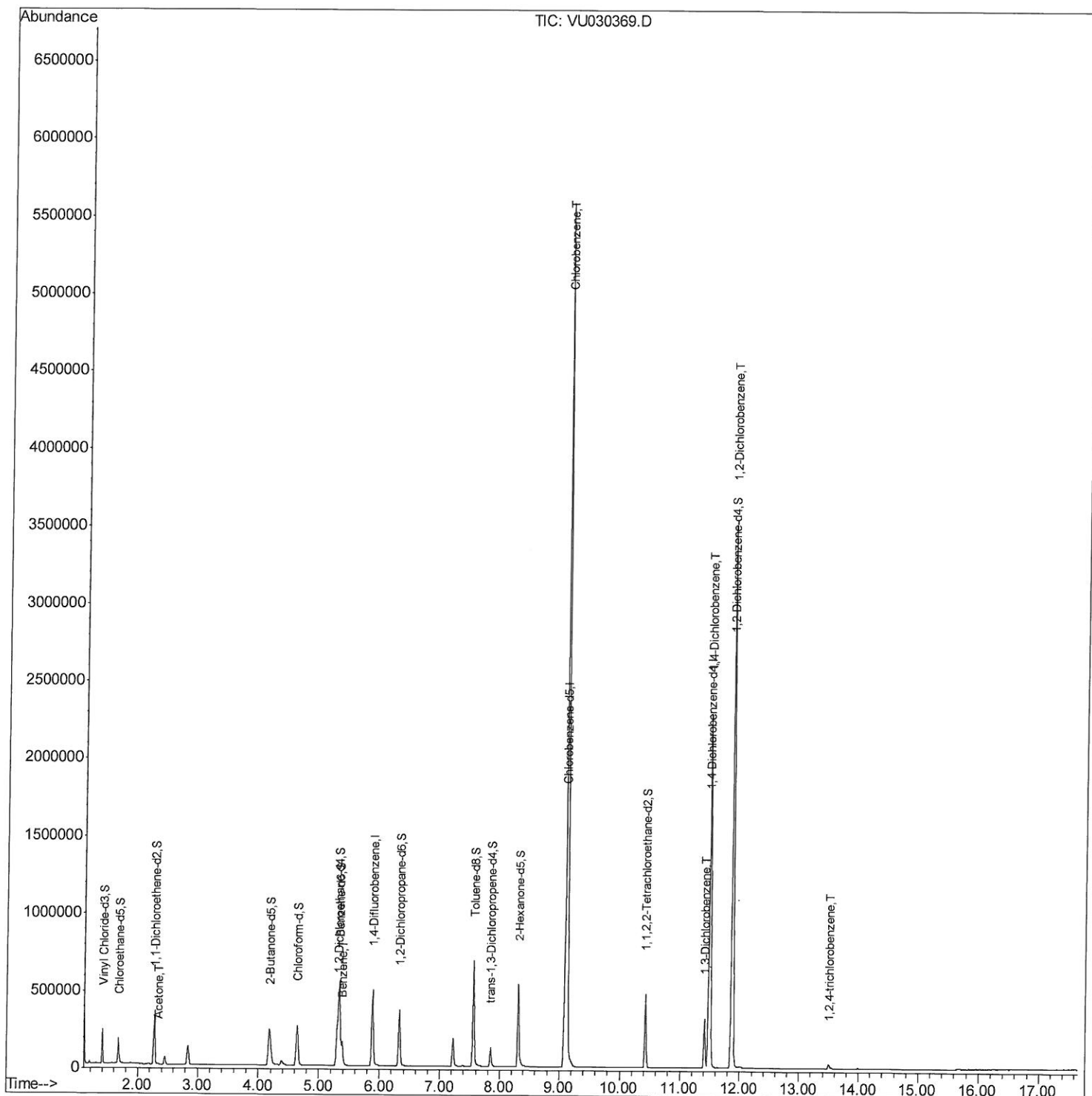
Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU032419\
Quantitation Report (QT Reviewed)
Data File : VU030369.D
Acq On : 24 Mar 2019 14:39
Operator : JC/SP
Sample : K2016-03
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
C0980

Manual Integrations
APPROVED

MMDadoda
3/25/2019 12:35:52 PM

Quant Time: Mar 25 04:19:08 2019
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM032319WMA.M
Quant Title : VOC Analysis
QLast Update : Mon Mar 25 03:20:51 2019
Response via : Initial Calibration



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU032419\
Quantitation Report (Qedit)

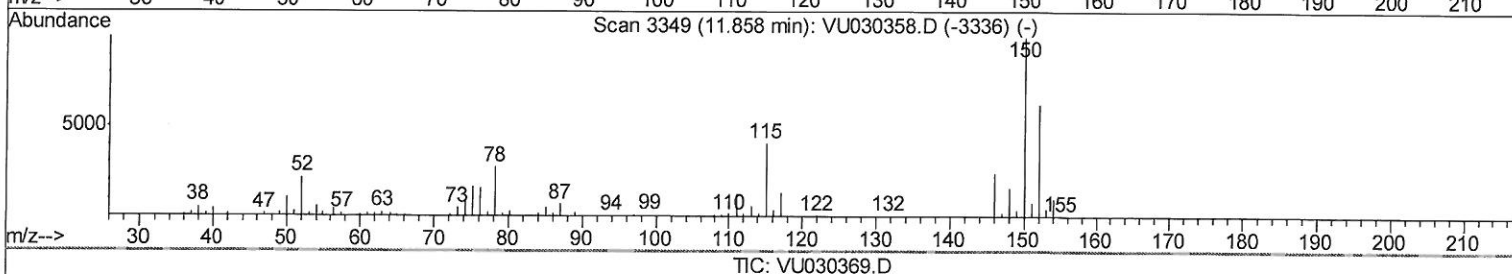
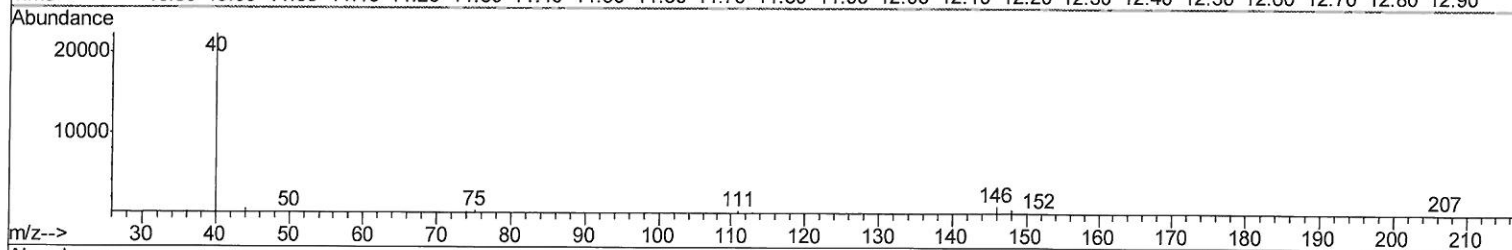
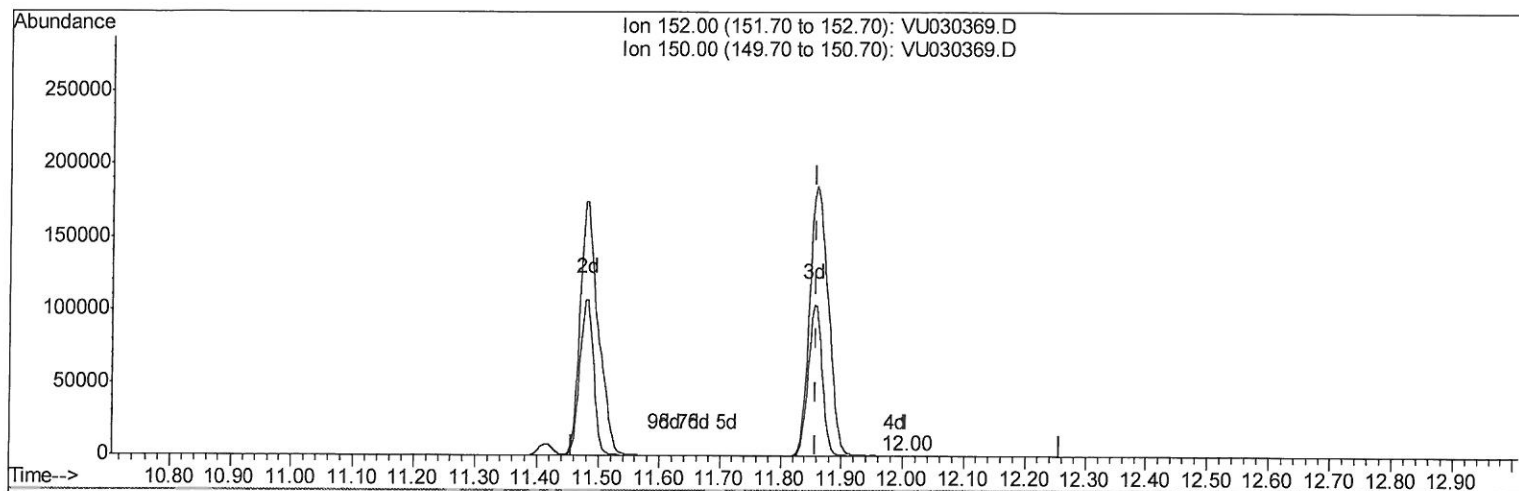
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(64) 1,2-Dichlorobenzene-d4 (S)

12.003min (+0.145) 0.01ug/L

response 44

Ion	Exp%	Act%
152.00	100	100
150.00	173.80	127.27
0.00	0.00	0.00
0.00	0.00	0.00

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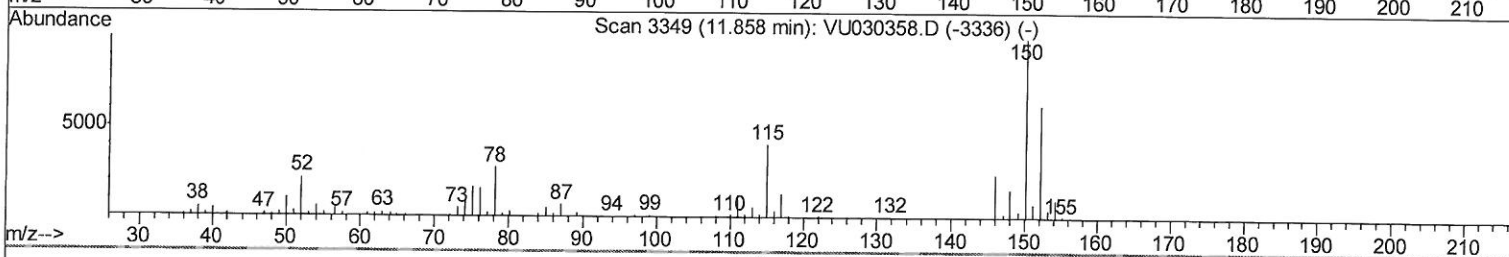
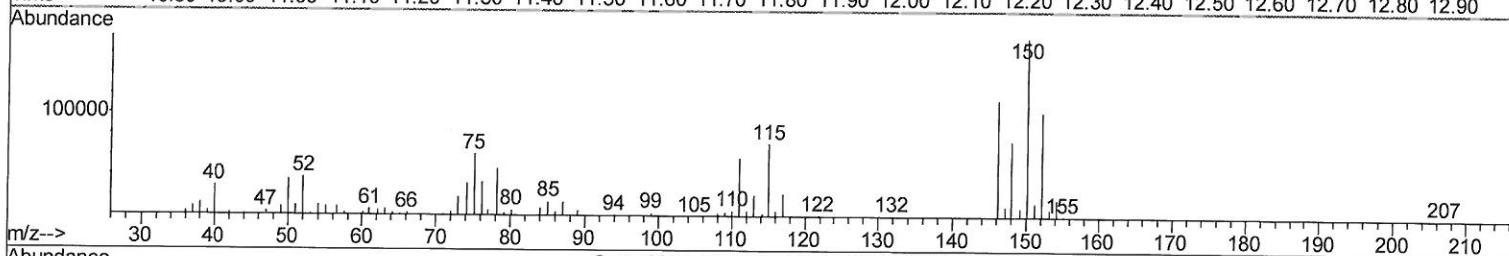
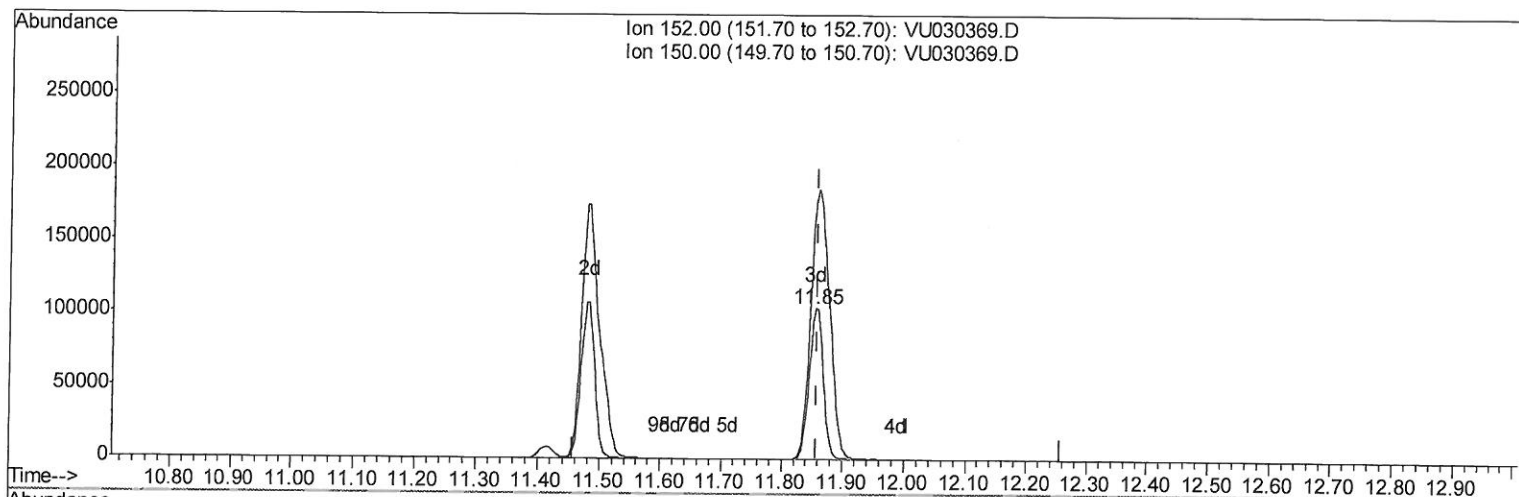
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Manual Integrations
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Quant Title : VOC Analysis
QLast Update : Mon Mar 25 03:20:51 2019
Response via : Initial Calibration



TIC: VU030369.D

(64) 1,2-Dichlorobenzene-d4 (S)

11.855min (-0.003) 49.41ug/L m

response 167935

Ion	Exp%	Act%
152.00	100	100
150.00	173.80	0.03#
0.00	0.00	0.00
0.00	0.00	0.00

MMD
3/27/19

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU032419\

Quantitation Report (QT Reviewed)

Data File : VU030369.D

Acq On : 24 Mar 2019 14:39

Operator : JC/SP

Sample : K2016-03

Misc : 5.0mL/MSVOA_U/WATER

ALS Vial : 13 Sample Multiplier: 1

Instrument :

MSVOA_U

ClientSampled :

C0980

Manual Integrations

APPROVED

MMDadoda

3/25/2019 12:35:52 PM

Quant Time: Mar 25 04:19:08 2019

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMULM032319WMA.M

Quant Title : VOC Analysis

QLast Update : Mon Mar 25 03:20:51 2019

Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Difluorobenzene	5.88	114	408735	50.00	ug/L	0.00
28) Chlorobenzene-d5	9.09	117	390294	50.00	ug/L	0.00
60) 1,4-Dichlorobenzene-d4	11.48	152	169645	50.00	ug/L	0.00

System Monitoring Compounds

4) Vinyl Chloride-d3	1.40	65	139961	42.55	ug/L	0.00
Spiked Amount	50.000	Range	60 - 135	Recovery	=	85.10%
7) Chloroethane-d5	1.68	69	117020	44.88	ug/L	0.00
Spiked Amount	50.000	Range	70 - 130	Recovery	=	89.76%
11) 1,1-Dichloroethene-d2	2.27	63	203771	32.66	ug/L	0.00
Spiked Amount	50.000	Range	60 - 125	Recovery	=	65.32%
21) 2-Butanone-d5	4.17	46	279165	90.28	ug/L	0.00
Spiked Amount	100.000	Range	40 - 130	Recovery	=	90.28%
24) Chloroform-d	4.64	84	262765	46.36	ug/L	0.00
Spiked Amount	50.000	Range	70 - 125	Recovery	=	92.72%
26) 1,2-Dichloroethane-d4	5.31	65	173951	46.90	ug/L	0.00
Spiked Amount	50.000	Range	70 - 125	Recovery	=	93.80%
32) Benzene-d6	5.34	84	536319	47.92	ug/L	0.00
Spiked Amount	50.000	Range	70 - 125	Recovery	=	95.84%
36) 1,2-Dichloropropane-d6	6.33	67	192526	50.28	ug/L	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	100.56%
41) Toluene-d8	7.56	98	470762	46.83	ug/L	0.00
Spiked Amount	50.000	Range	80 - 120	Recovery	=	93.66%
43) trans-1,3-Dichloropropene-	7.85	79	74198	41.21	ug/L	0.00
Spiked Amount	50.000	Range	60 - 125	Recovery	=	82.42%
47) 2-Hexanone-d5	8.31	63	189015	91.08	ug/L	0.00
Spiked Amount	100.000	Range	45 - 130	Recovery	=	91.08%
57) 1,1,2,2-Tetrachloroethane-	10.43	84	249189	45.47	ug/L	0.00
Spiked Amount	50.000	Range	65 - 120	Recovery	=	90.94%
64) 1,2-Dichlorobenzene-d4	11.85	152	167935m	49.41	ug/L	0.00
Spiked Amount	50.000	Range	80 - 120	Recovery	=	98.82%

7 MLO
3/27/19

Target Compounds

						Qvalue
13) Acetone	2.32	43	4332	1.385	ug/L	98
33) Benzene	5.39	78	142293	10.710	ug/L	100
51) Chlorobenzene	9.12	112	2854011	341.521	ug/L	98
62) 1,3-Dichlorobenzene	11.41	146	123741	22.304	ug/L	99
63) 1,4-Dichlorobenzene	11.50	146	852165	149.205	ug/L	99
65) 1,2-Dichlorobenzene	11.87	146	1305744	232.763	ug/L	99
68) 1,2,4-trichlorobenzene	13.51	180	12022	3.331	ug/L	93

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