

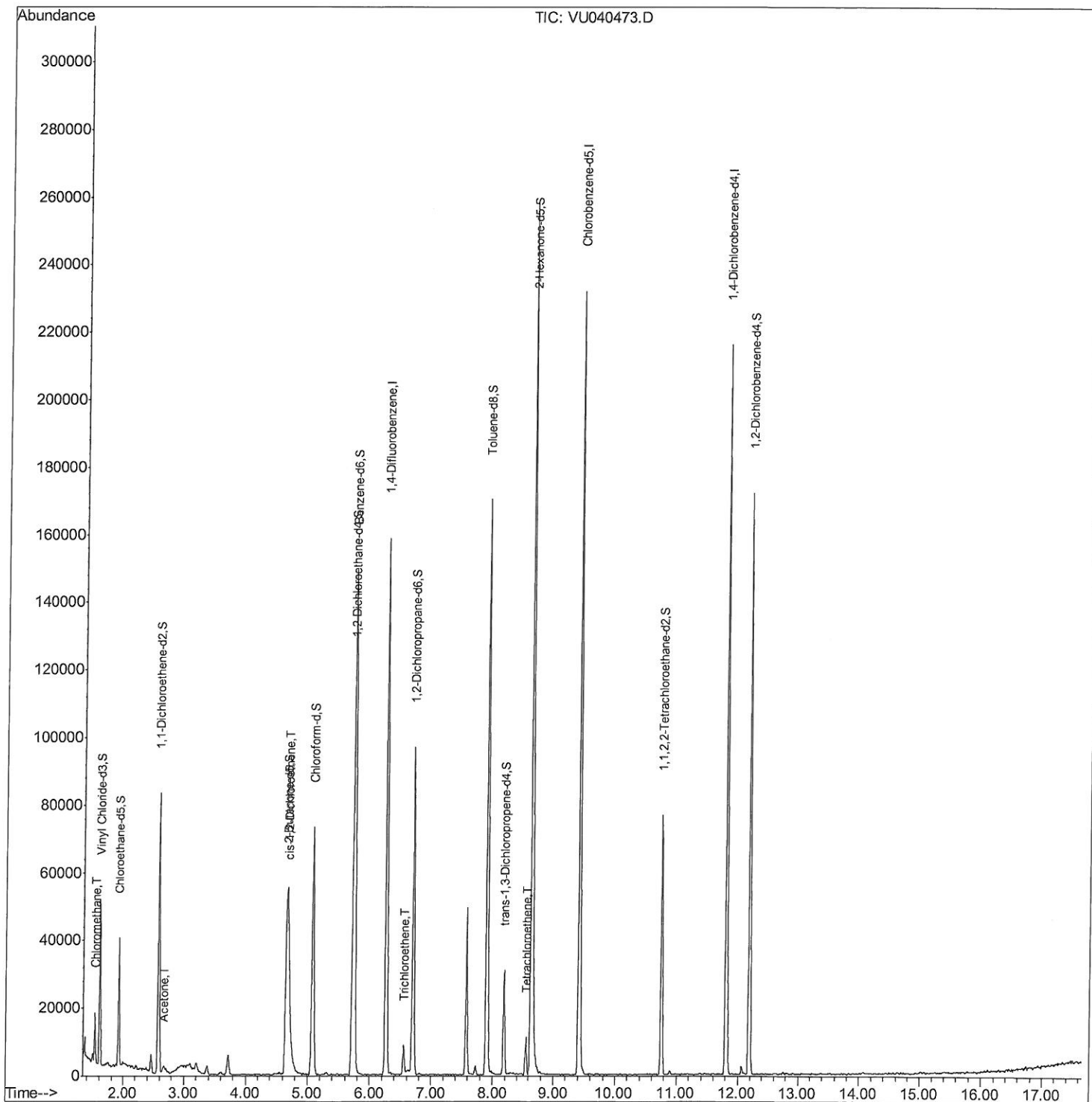
Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU100220\
Data File : VU040473.D
Acq On : 02 Oct 2020 20:35
Operator : SY/MD
Sample : L4185-14
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
C0AR6

Manual Integrations
APPROVED

MMDadoda
10/6/2020 9:04:19 AM

Quant Time: Oct 03 03:24:15 2020
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR092920WMA.M
Quant Title : TRACE VOA SOM01.0
QLast Update : Sat Oct 03 00:09:31 2020
Response via : Initial Calibration



Quantitation Report (Qedit)

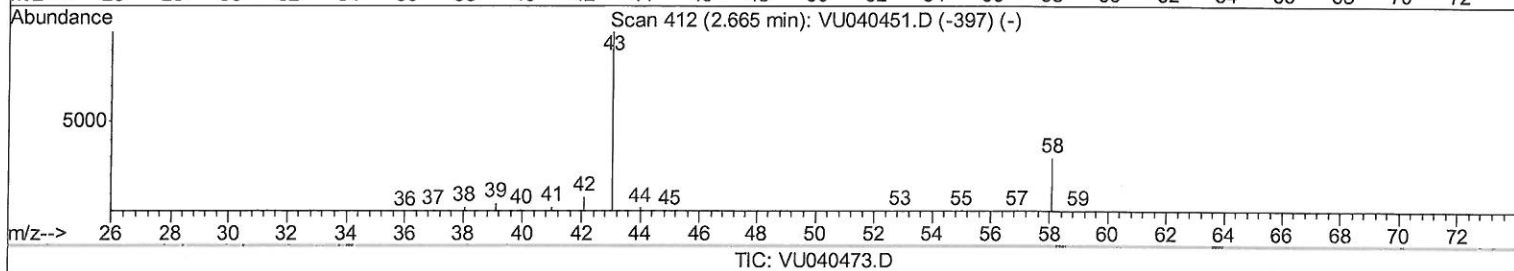
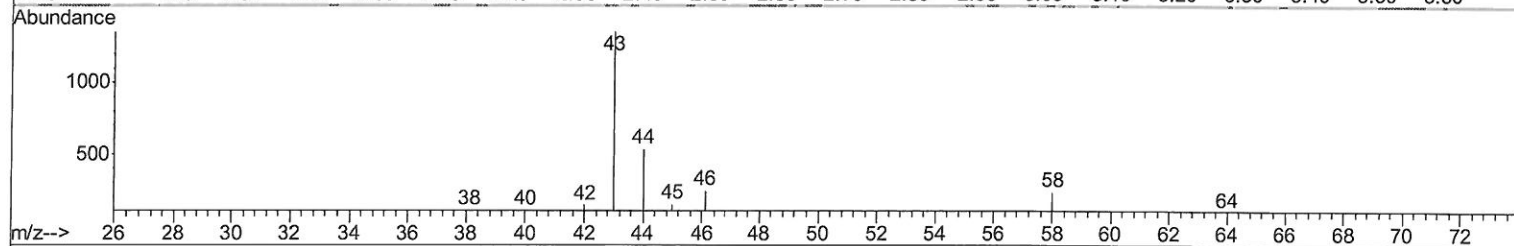
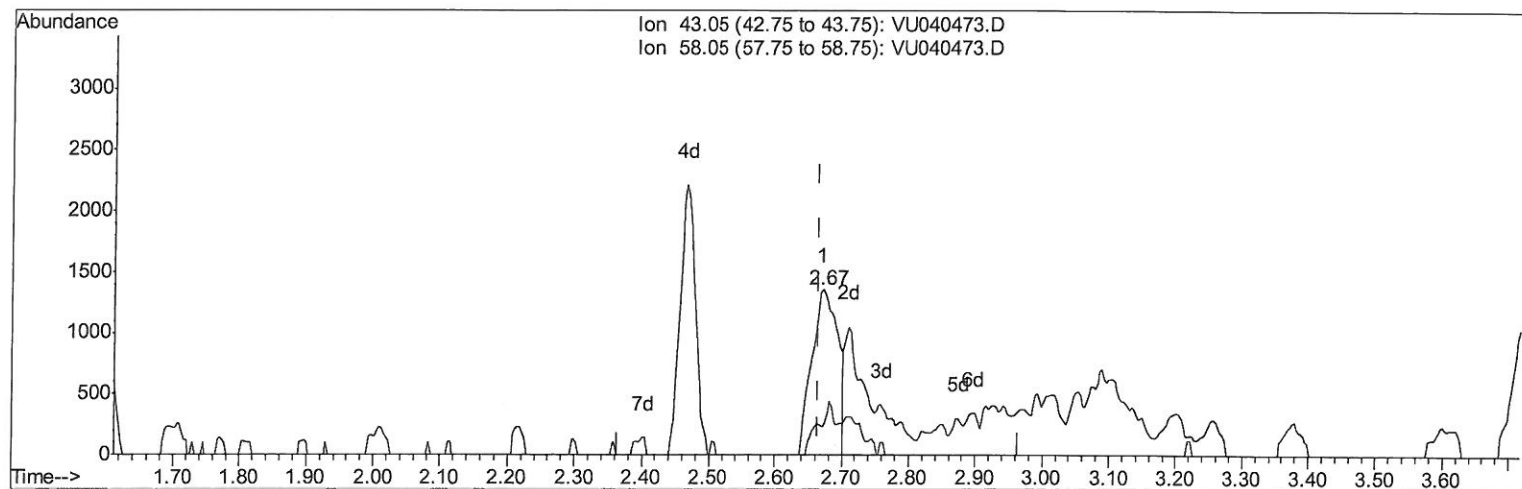
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(13) Acetone (T)

2.671min (+0.007) 1.47ug/L

response 3614

Ion	Exp%	Act%
43.05	100	100
58.05	30.90	8.74
0.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

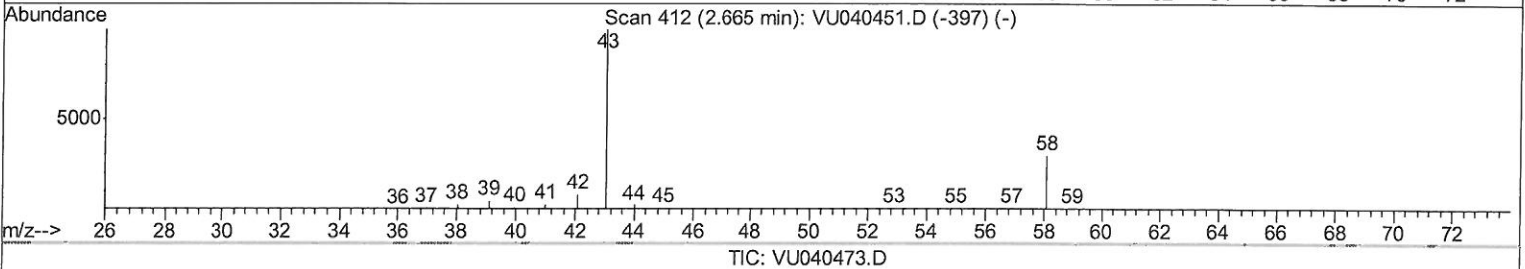
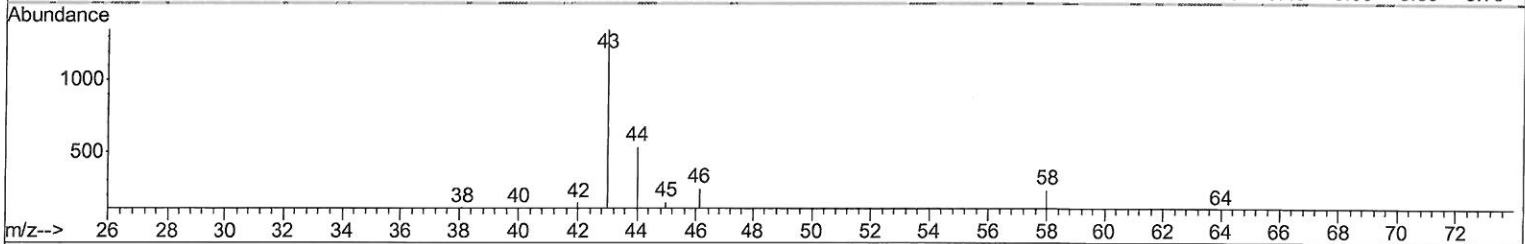
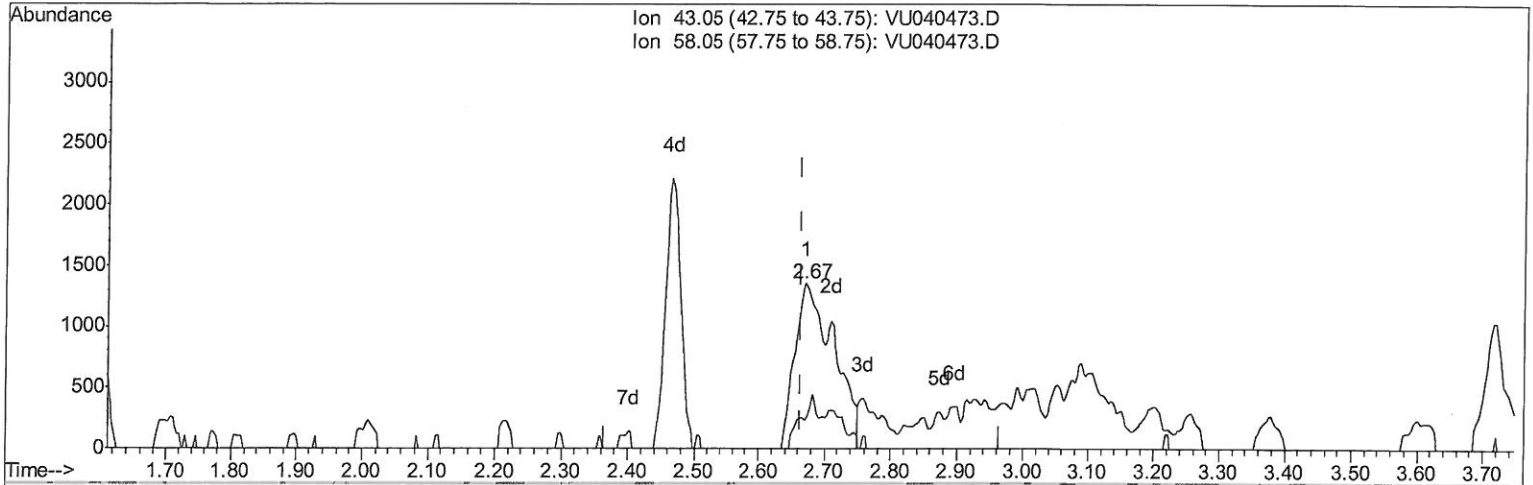
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Quant Time: Oct 03 00:20:35 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR092920WMA.M
 Quant Title : TRACE VOA SOM01.0
 QLast Update : Sat Oct 03 00:09:31 2020
 Response via : Initial Calibration



(13) Acetone (T)

2.671min (+0.007) 2.26ug/L m

response 5547

Ion	Exp%	Act%
43.05	100	100
58.05	30.90	5.70
0.00	0.00	0.00
0.00	0.00	0.00

MD
10/08/20

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU100220\
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 Operator : SY/MD
 Sample : L4185-14
 Misc : 25.0mL/MSVOA U/WATER
 ALS Vial : 26 Sample Multiplier: 1

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Manual Integrations
 APPROVED

MMDadoda
 10/6/2020 9:04:19 AM

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 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR092920WMA.M
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Difluorobenzene	6.26	114	119722	5.00	ug/L	0.00
28) Chlorobenzene-d5	9.42	117	120394	5.00	ug/L	0.00
60) 1,4-Dichlorobenzene-d4	11.81	152	52398	5.00	ug/L	0.00

System Monitoring Compounds

4) Vinyl Chloride-d3	1.60	65	31176	5.26	ug/L	0.00
Spiked Amount 5.000	Range 40 - 130		Recovery =	105.20%		
7) Chloroethane-d5	1.92	69	26189	5.05	ug/L	0.00
Spiked Amount 5.000	Range 65 - 130		Recovery =	101.00%		
11) 1,1-Dichloroethene-d2	2.58	63	50702	3.56	ug/L	0.00
Spiked Amount 5.000	Range 60 - 125		Recovery =	71.20%		
20) 2-Butanone-d5	4.66	46	134702	45.75	ug/L	0.00
Spiked Amount 50.000	Range 40 - 130		Recovery =	91.50%		
24) Chloroform-d	5.08	84	64269	4.56	ug/L	0.00
Spiked Amount 5.000	Range 70 - 125		Recovery =	91.20%		
26) 1,2-Dichloroethane-d4	5.72	65	41838	4.87	ug/L	0.00
Spiked Amount 5.000	Range 70 - 130		Recovery =	97.40%		
32) Benzene-d6	5.75	84	127399	5.26	ug/L	0.00
Spiked Amount 5.000	Range 70 - 125		Recovery =	105.20%		
36) 1,2-Dichloropropane-d6	6.70	67	43088	4.87	ug/L	0.00
Spiked Amount 5.000	Range 60 - 140		Recovery =	97.40%		
41) Toluene-d8	7.91	98	106532	5.07	ug/L	0.00
Spiked Amount 5.000	Range 70 - 130		Recovery =	101.40%		
43) trans-1,3-Dichloropropene-	8.19	79	17076	4.73	ug/L	0.00
Spiked Amount 5.000	Range 55 - 130		Recovery =	94.60%		
46) 2-Hexanone-d5	8.64	63	87556	41.20	ug/L	0.00
Spiked Amount 50.000	Range 45 - 130		Recovery =	82.40%		
57) 1,1,2,2-Tetrachloroethane-	10.75	84	36429	4.23	ug/L	0.00
Spiked Amount 5.000	Range 65 - 120		Recovery =	84.60%		
64) 1,2-Dichlorobenzene-d4	12.19	152	41673	5.25	ug/L	0.00
Spiked Amount 5.000	Range 80 - 120		Recovery =	105.00%		

Target Compounds

					Ovalue
3) Chloromethane	1.53	50	9638	0.819 ug/L	95
13) Acetone	2.67	43	5547m	2.261 ug/L	93
22) cis-1,2-Dichloroethene	4.69	96	2541	0.244 ug/L	96
34) Trichloroethene	6.56	95	2844	0.265 ug/L	96
47) Tetrachloroethene	8.56	164	2502	0.329 ug/L	87

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