

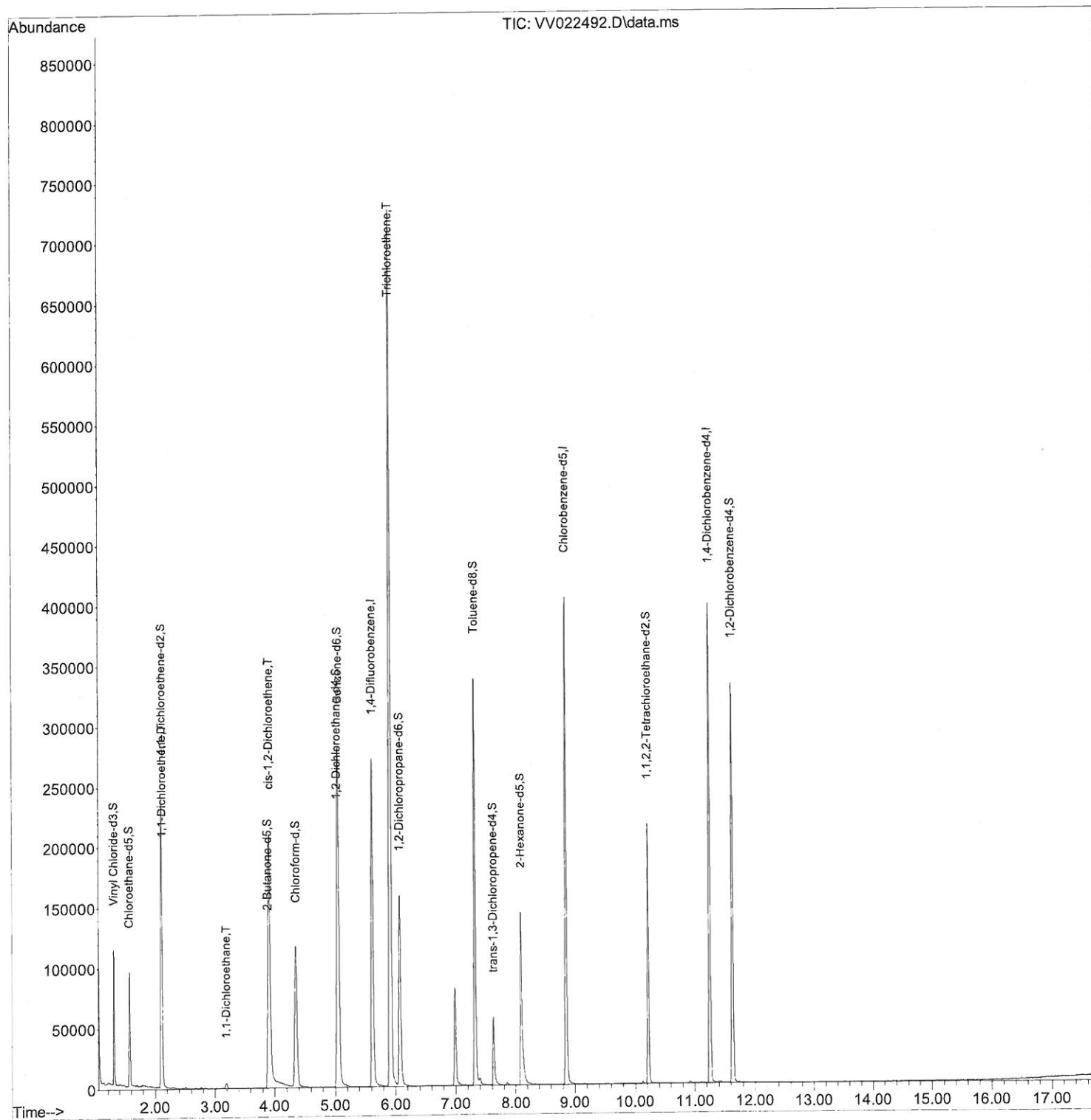
Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV100121\
Data File : VV022492.D
Acq On : 02 Oct 2021 03:15
Operator : SY/MD
Sample : M4024-12
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
COAM0

Manual Integrations
APPROVED

MMDadoda
10/6/2021 2:24:37 PM

Quant Time: Oct 02 03:34:46 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM100121WMA.M
Quant Title : VOC Analysis
QLast Update : Sat Oct 02 01:27:14 2021
Response via : Initial Calibration



Quantitation Report (Qedit)

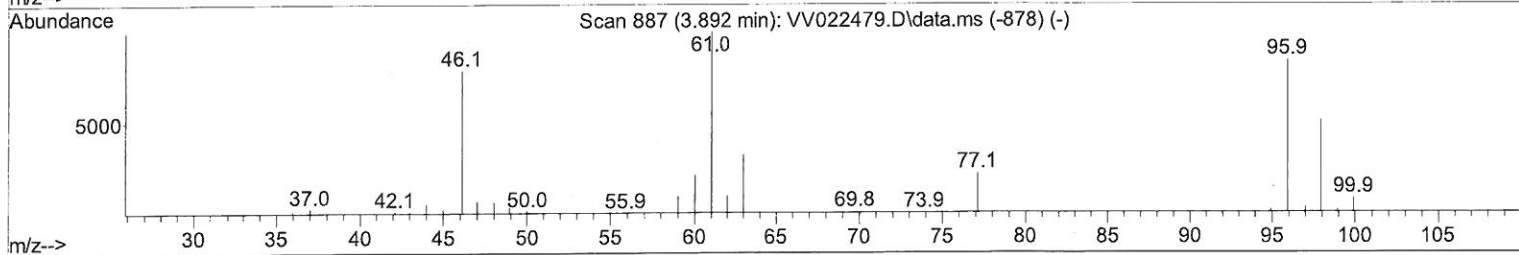
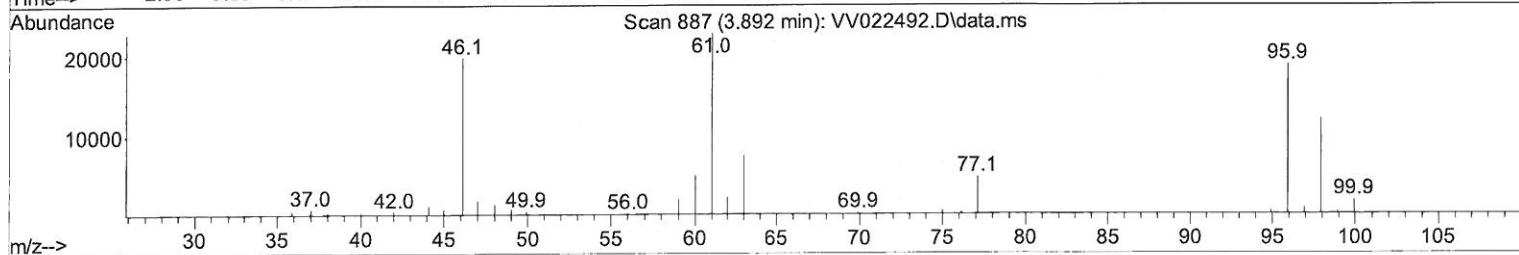
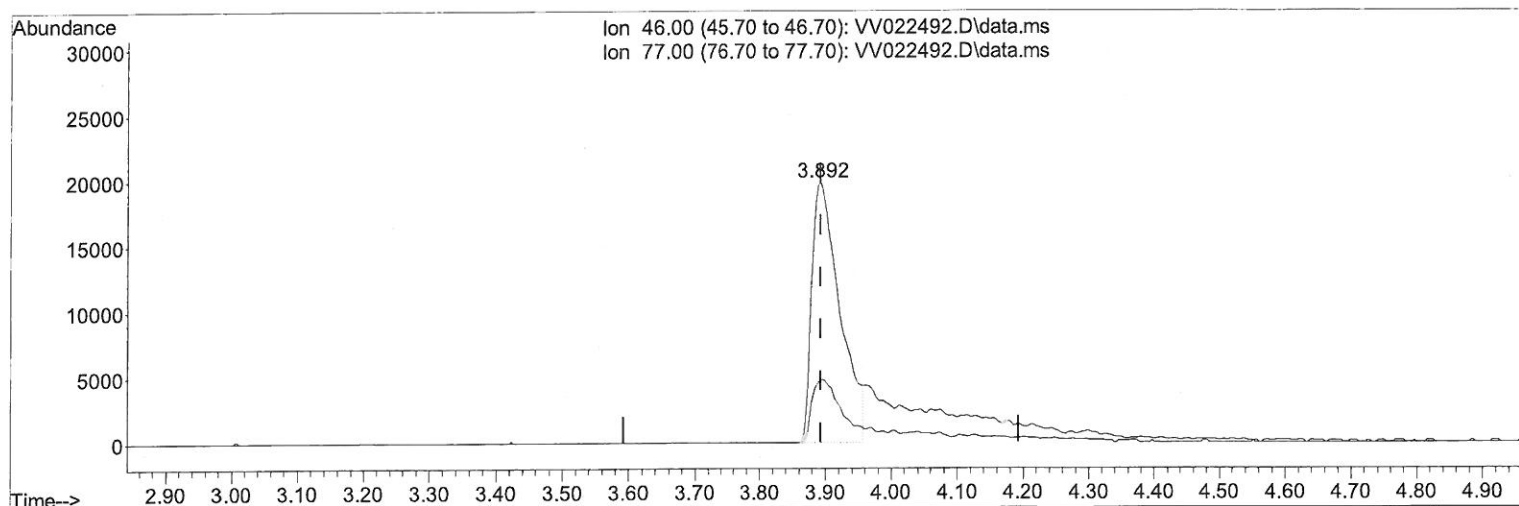
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TIC: VV022492.D\data.ms

(21) 2-Butanone-d5 (S)

3.892min (-0.000) 70.82 ug/L

response 58425

Ion	Exp%	Act%
46.00	100.00	100.00
77.00	26.40	23.96
0.00	0.00	0.00
0.00	0.00	0.00

Quantitation Report (Qedit)

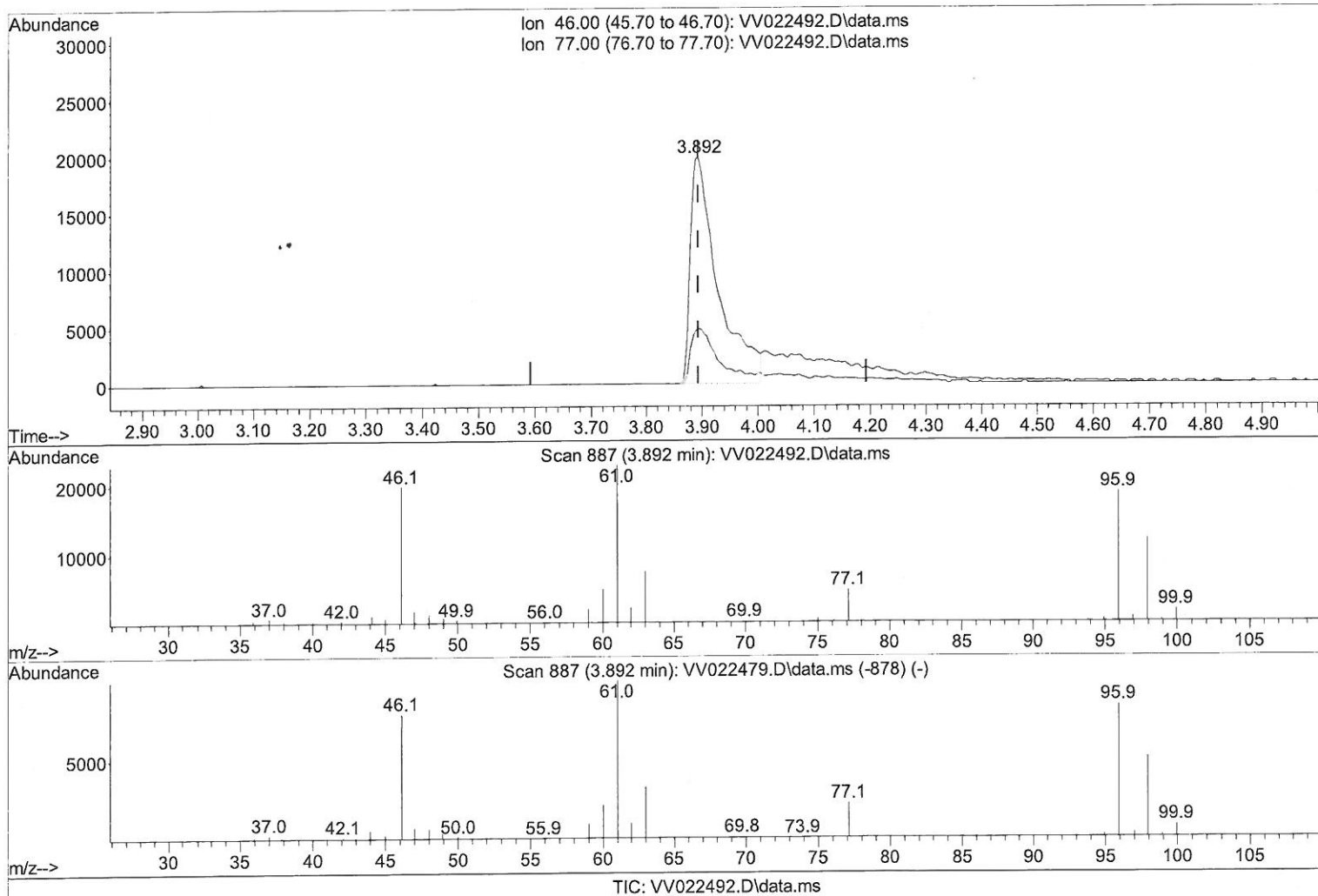
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 Data File : VV022492.D
 Acq On : 02 Oct 2021 03:15
 Operator : SY/MD
 Sample : M4024-12
 Misc : 5.0mL/MSVOA_V/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 COAMO

Manual Integrations
 APPROVED

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 QLast Update : Sat Oct 02 01:27:14 2021
 Response via : Initial Calibration



(21) 2-Butanone-d5 (S)

3.892min (-0.000) 82.85 ug/L m

response 68352

Ion	Exp%	Act%
46.00	100.00	100.00
77.00	26.40	20.48
0.00	0.00	0.00
0.00	0.00	0.00

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 Acq On : 02 Oct 2021 03:15
 Operator : SY/MD
 Sample : M4024-12
 Misc : 5.0mL/MSVOA_V/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampled :
 COAMO

Manual Integrations
 APPROVED

MMDadoda
 10/6/2021 2:24:37 PM

Quant Time: Oct 02 03:34:46 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM100121WMA.M
 Quant Title : VOC Analysis
 QLast Update : Sat Oct 02 01:27:14 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Difluorobenzene	5.619	114	242337	50.000	ug/L	0.00
28) Chlorobenzene-d5	8.853	117	229617	50.000	ug/L	0.00
58) 1,4-Dichlorobenzene-d4	11.252	152	107941	50.000	ug/L	0.00
System Monitoring Compounds						
4) Vinyl Chloride-d3	1.304	65	69790	46.565	ug/L	0.00
Spiked Amount	50.000	Range	60 - 135	Recovery	=	93.140%
7) Chloroethane-d5	1.568	69	57269	48.611	ug/L	0.00
Spiked Amount	50.000	Range	70 - 130	Recovery	=	97.220%
11) 1,1-Dichloroethene-d2	2.108	63	112134	37.945	ug/L	0.00
Spiked Amount	50.000	Range	60 - 125	Recovery	=	75.880%
21) 2-Butanone-d5	3.892	46	68352m	82.852	ug/L	0.00
Spiked Amount	100.000	Range	40 - 130	Recovery	=	82.850%
24) Chloroform-d	4.352	84	123907	45.250	ug/L	0.00
Spiked Amount	50.000	Range	70 - 125	Recovery	=	90.500%
26) 1,2-Dichloroethane-d4	5.034	65	76286	48.025	ug/L	0.00
Spiked Amount	50.000	Range	70 - 125	Recovery	=	96.060%
32) Benzene-d6	5.050	84	253957	47.741	ug/L	0.00
Spiked Amount	50.000	Range	70 - 125	Recovery	=	95.480%
36) 1,2-Dichloropropane-d6	6.072	67	77363	47.242	ug/L	0.00
Spiked Amount	50.000	Range	70 - 120	Recovery	=	94.480%
41) Toluene-d8	7.317	98	225517	46.218	ug/L	0.00
Spiked Amount	50.000	Range	80 - 120	Recovery	=	92.440%
43) trans-1,3-Dichloroprop...	7.625	79	31969	41.261	ug/L	0.00
Spiked Amount	50.000	Range	60 - 125	Recovery	=	82.520%
47) 2-Hexanone-d5	8.091	63	59522	87.741	ug/L	0.00
Spiked Amount	100.000	Range	45 - 130	Recovery	=	87.740%
56) 1,1,2,2-Tetrachloroeth...	10.217	84	100469	44.121	ug/L	0.00
Spiked Amount	50.000	Range	65 - 120	Recovery	=	88.240%
66) 1,2-Dichlorobenzene-d4	11.625	152	86247	48.997	ug/L	0.00
Spiked Amount	50.000	Range	80 - 120	Recovery	=	98.000%
Target Compounds						
12) 1,1-Dichloroethene	2.117	96	22631	13.369	ug/L #	41
19) 1,1-Dichloroethane	3.191	63	4561	1.387	ug/L	95
20) cis-1,2-Dichloroethene	3.912	96	111346	56.244	ug/L	100
34) Trichloroethene	5.915	95	248497	134.239	ug/L	98

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