

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU102620\
 Data File : VU040948.D
 Acq On : 26 Oct 2020 16:47
 Operator : SY/MD
 Sample : L4492-01
 Misc : 25.0mL/MSVOA U/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 C0AM9

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.1

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR102620WMA.M

Title : TRACE VOA SOM01.0

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.362	3	7	27	rVB	768089	731981	100.00%	17.965%
2	1.494	41	48	56	rBV	55329	52852	7.22%	1.297%
3	1.594	62	79	92	rBV2	329071	552489	75.48%	13.560%
4	1.665	98	101	110	rVB3	16059	15798	2.16%	0.388%
5	1.739	118	124	138	rVB	58514	75142	10.27%	1.844%
6	1.916	172	179	197	rVB3	47397	75607	10.33%	1.856%
7	2.002	200	206	220	rVB2	35135	50743	6.93%	1.245%
8	2.163	248	256	264	rBV	79550	109358	14.94%	2.684%
9	2.215	265	272	287	rVB3	39132	67604	9.24%	1.659%
10	2.327	301	307	325	rVB3	15448	33685	4.60%	0.827%
11	2.424	327	337	347	rBV4	26811	41015	5.60%	1.007%
12	2.578	373	385	398	rBV	38578	61594	8.41%	1.512%
13	2.700	405	423	433	rBV4	3535	9742	1.33%	0.239%
14	2.986	499	512	524	rBV5	12268	37355	5.10%	0.917%
15	3.198	573	578	594	rVB3	4685	9352	1.28%	0.230%
16	3.607	690	705	718	rBV6	11008	25373	3.47%	0.623%
17	3.874	774	788	804	rBV2	11355	30307	4.14%	0.744%
18	4.214	875	894	904	rBV7	4035	9406	1.29%	0.231%
19	4.668	1018	1035	1088	rBV2	30806	166708	22.77%	4.091%
20	5.079	1149	1163	1182	rVB	40864	92537	12.64%	2.271%
21	5.739	1349	1368	1381	rBV3	75417	202676	27.69%	4.974%
22	6.179	1489	1505	1518	rBV3	5690	14518	1.98%	0.356%
23	6.263	1518	1531	1558	rVB	78072	145700	19.90%	3.576%
24	6.703	1655	1668	1683	rBV	52782	99859	13.64%	2.451%
25	7.574	1927	1939	1953	rBV2	23516	40518	5.54%	0.994%
26	7.906	2029	2042	2054	rBV	78611	133491	18.24%	3.276%
27	7.976	2054	2064	2074	rVB	11915	19263	2.63%	0.473%
28	8.185	2119	2129	2144	rBV	15550	25391	3.47%	0.623%
29	8.642	2256	2271	2308	rBV	176117	343437	46.92%	8.429%
30	9.420	2501	2513	2538	rBV	115357	185252	25.31%	4.547%
31	9.693	2589	2598	2606	rBV3	6658	10309	1.41%	0.253%
32	10.237	2757	2767	2778	rBV	9168	14467	1.98%	0.355%
33	10.754	2917	2928	2942	rVB2	46944	73977	10.11%	1.816%
34	11.465	3140	3149	3160	rBV3	8821	13217	1.81%	0.324%

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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.1 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR102620WMA.M
Title : TRACE VOA SOM01.0

35	11.568	3169	3181	3199	rVB2	15098	23946	3.27%	0.588%
36	11.806	3244	3255	3269	rVB	110696	175593	23.99%	4.310%
37	12.057	3323	3333	3347	rBV5	6382	12728	1.74%	0.312%
38	12.188	3363	3374	3385	rVB	91155	145461	19.87%	3.570%
39	15.211	4303	4314	4327	rBV	50981	81342	11.11%	1.996%
40	15.397	4361	4372	4384	rBV	41622	64744	8.85%	1.589%

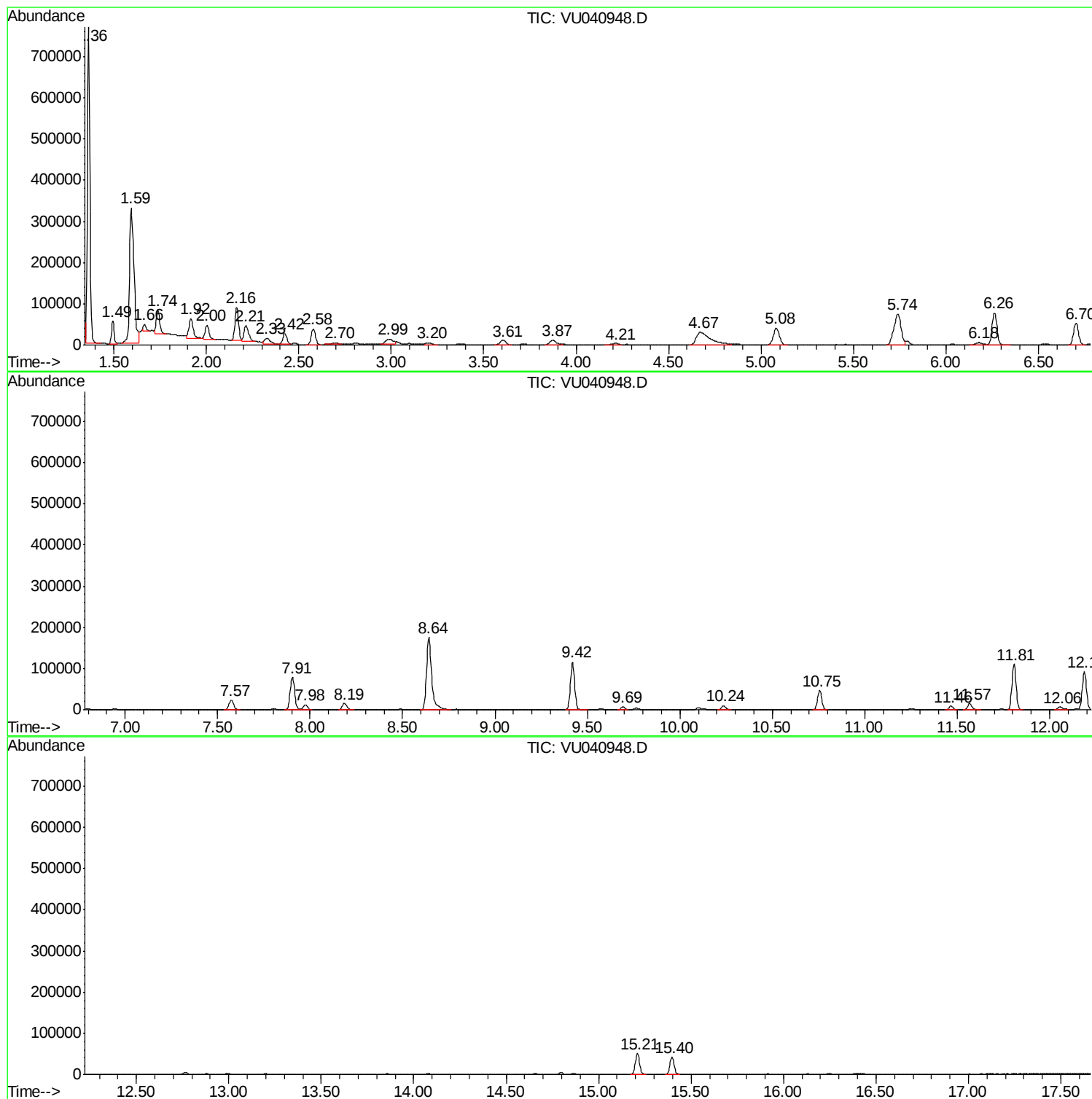
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR102620WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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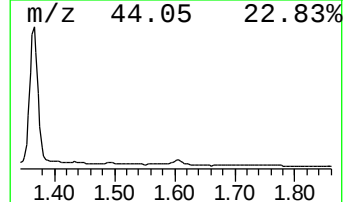
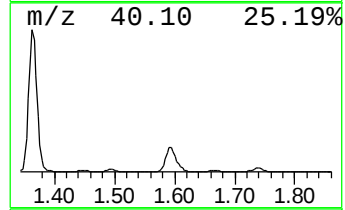
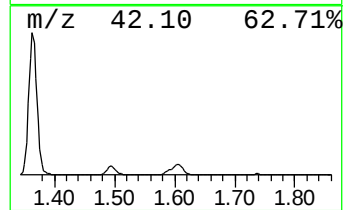
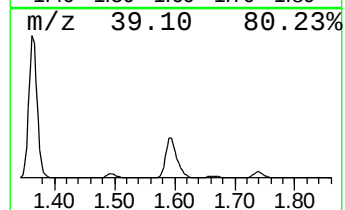
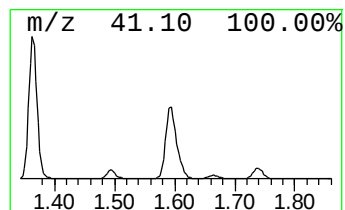
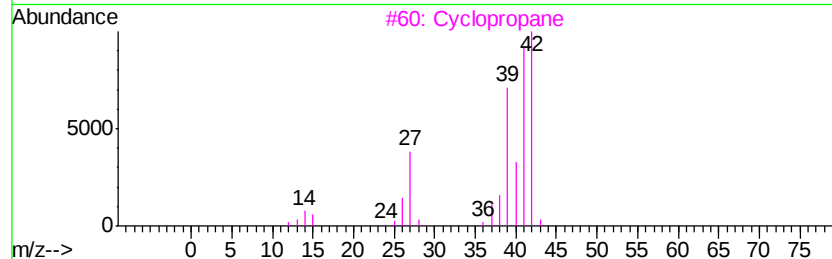
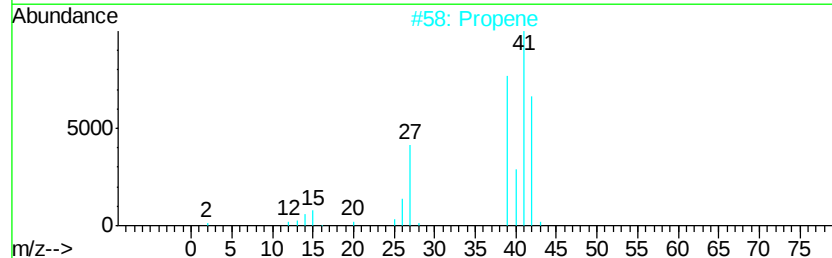
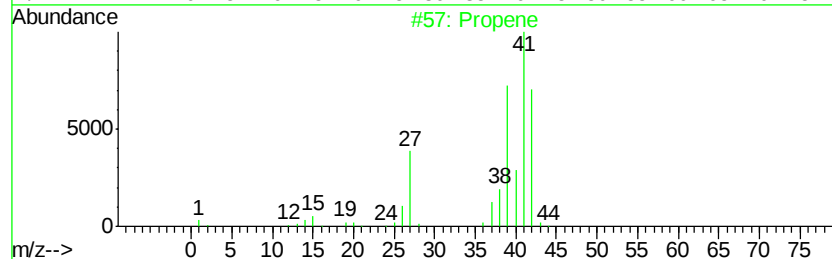
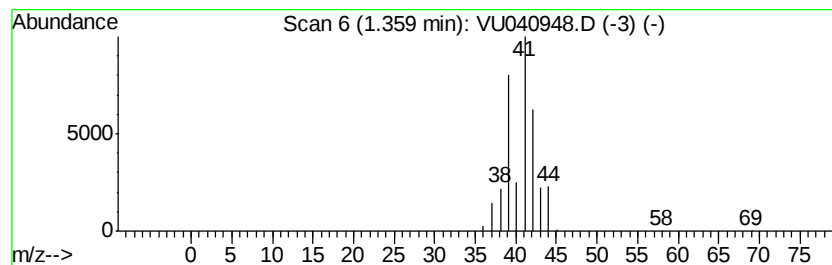
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.36	25.12 ug/L	731981	1,4-Difluorobenzene	6.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propene	42	C3H6	000115-07-1	90
2		Propene	42	C3H6	000115-07-1	42
3		Cyclopropane	42	C3H6	000075-19-4	9
4		Cyclopropene	40	C3H4	002781-85-3	9
5		Cyclopropane	42	C3H6	000075-19-4	7



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Instrument :
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ClientSampleId :
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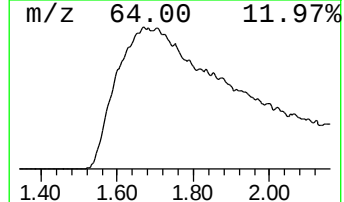
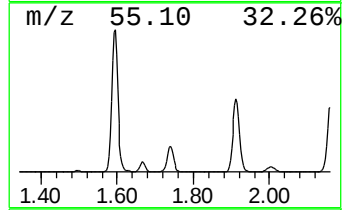
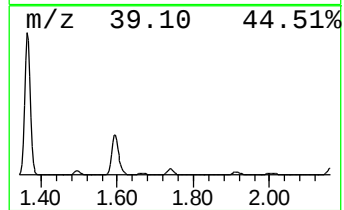
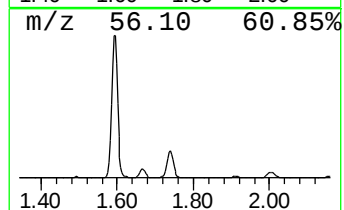
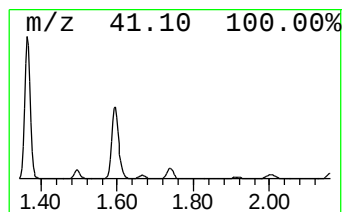
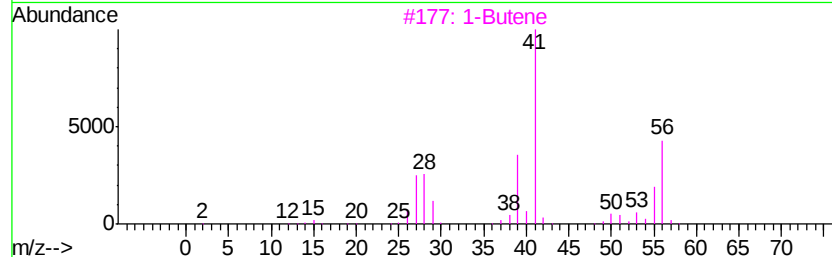
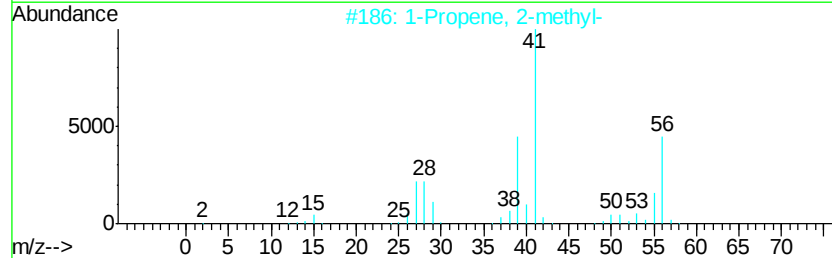
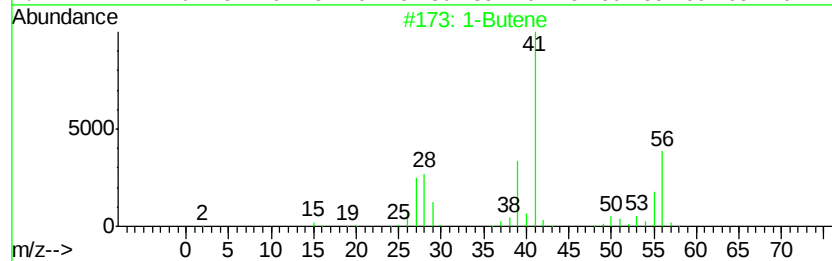
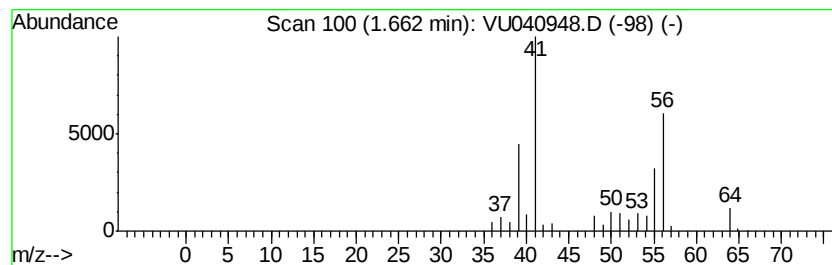
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR102620WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.66	0.54 ug/L	15798	1,4-Difluorobenzene	6.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Butene	56	C4H8	000106-98-9	64
2		1-Propene, 2-methyl-	56	C4H8	000115-11-7	59
3		1-Butene	56	C4H8	000106-98-9	59
4		1-Propene, 2-methyl-	56	C4H8	000115-11-7	59
5		2-Butene, (Z)-	56	C4H8	000590-18-1	58



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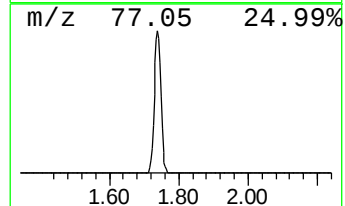
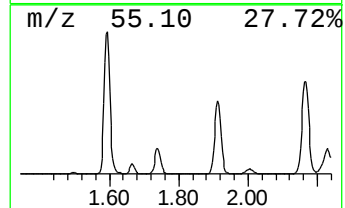
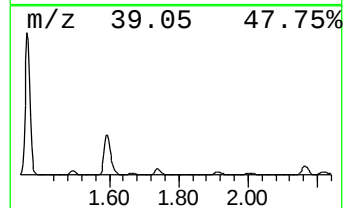
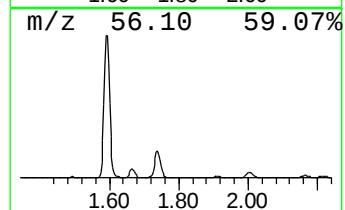
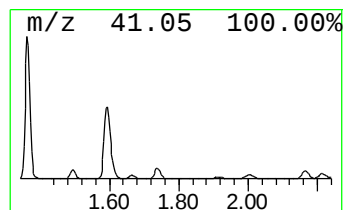
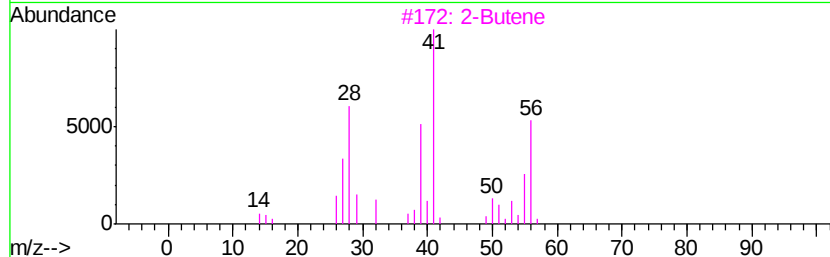
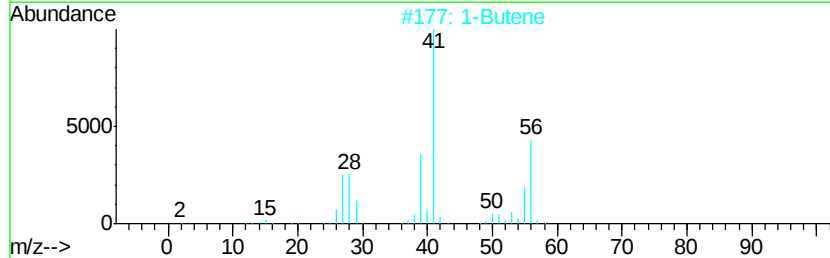
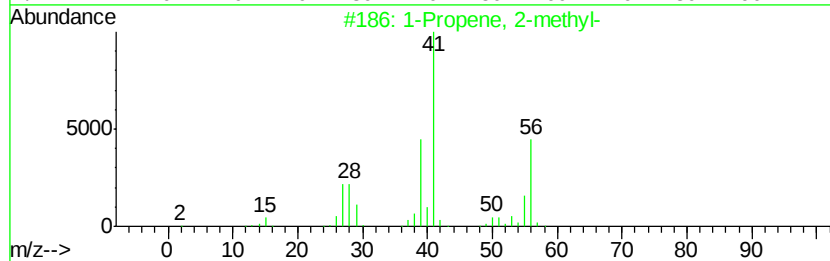
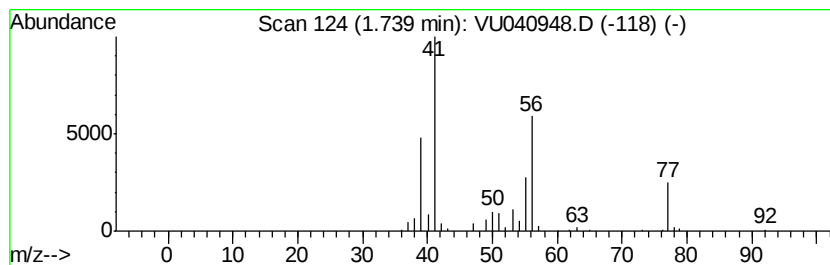
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.74	2.58 ug/L	75142	1,4-Difluorobenzene	6.26

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Propene, 2-methyl-	56	C4H8	000115-11-7	62
2		1-Butene	56	C4H8	000106-98-9	58
3		2-Butene	56	C4H8	000107-01-7	53
4		1-Propene, 2-methyl-	56	C4H8	000115-11-7	52
5		2-Butene, (Z)-	56	C4H8	000590-18-1	52



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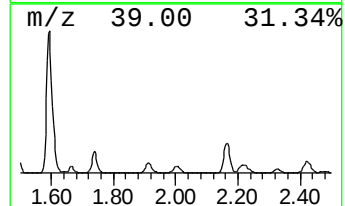
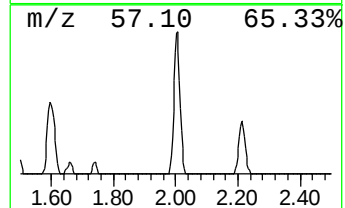
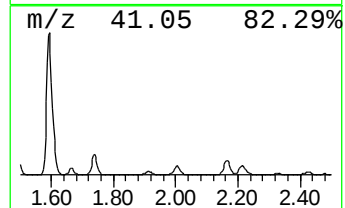
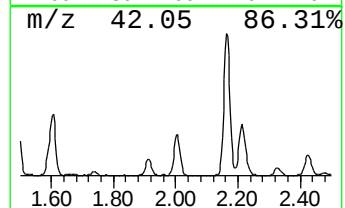
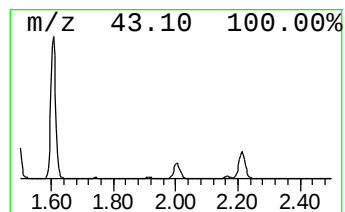
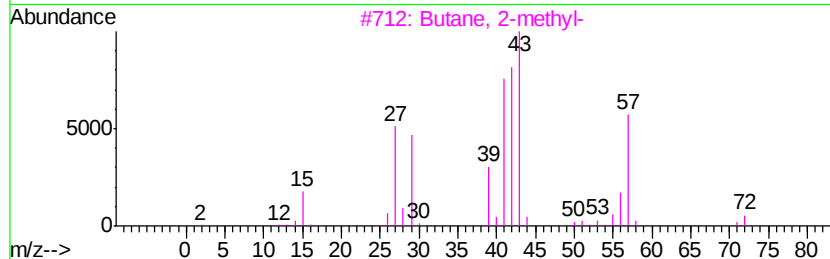
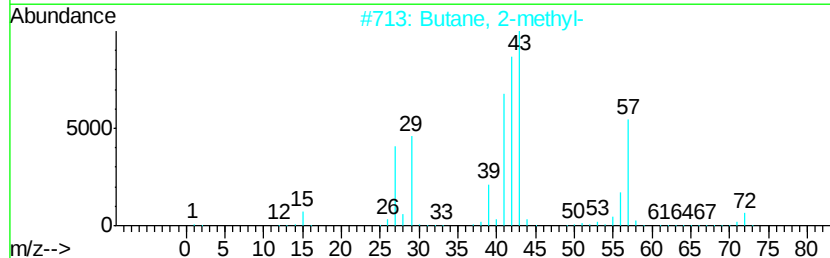
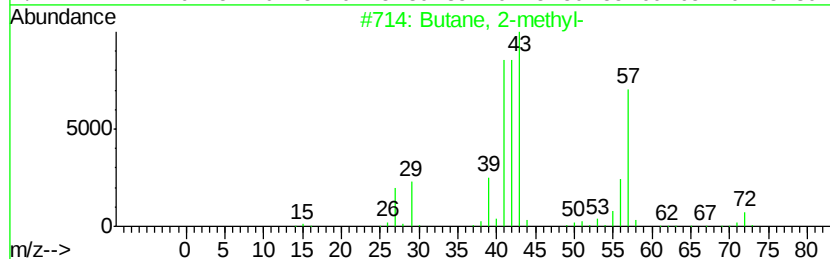
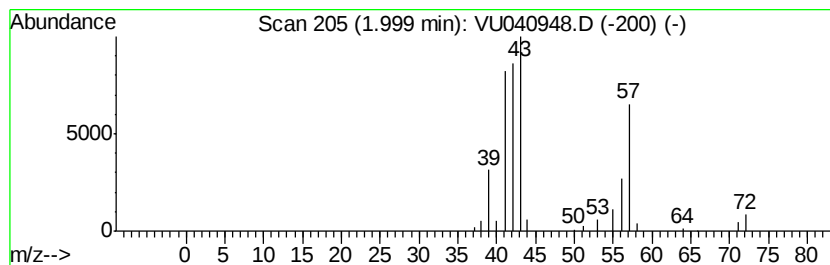
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.00	1.74 ug/L	50743	1,4-Difluorobenzene	6.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methyl-	72	C5H12	000078-78-4	91
2		Butane, 2-methyl-	72	C5H12	000078-78-4	86
3		Butane, 2-methyl-	72	C5H12	000078-78-4	86
4		Pentane	72	C5H12	000109-66-0	53
5		3-Buten-1-ol	72	C4H8O	000627-27-0	47



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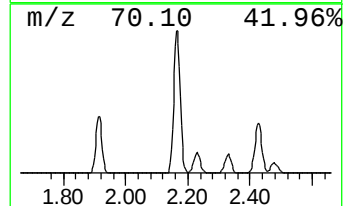
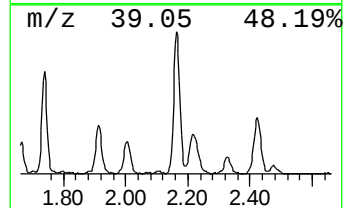
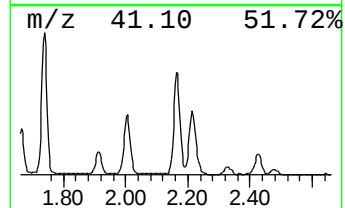
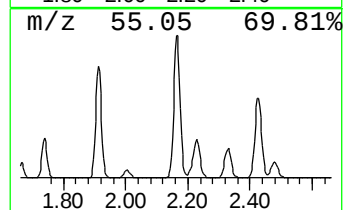
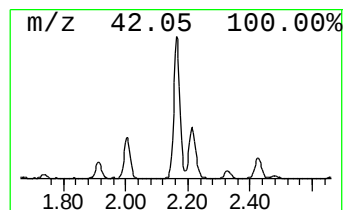
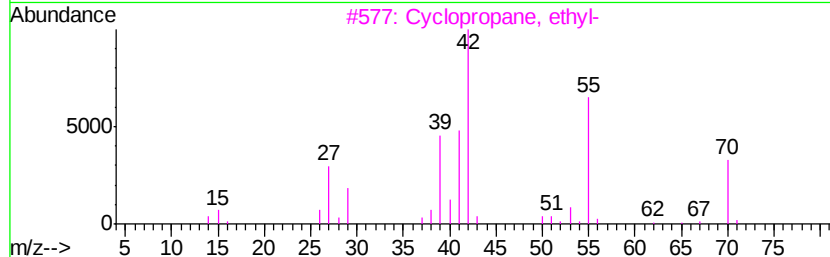
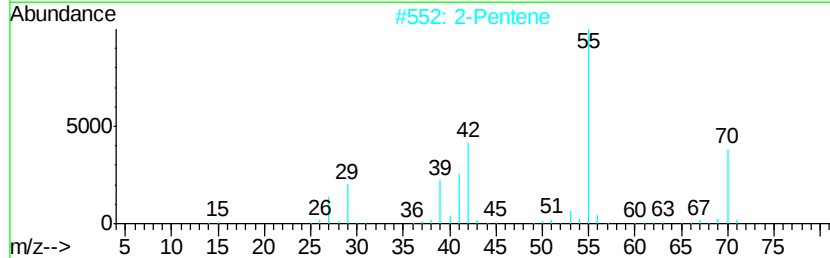
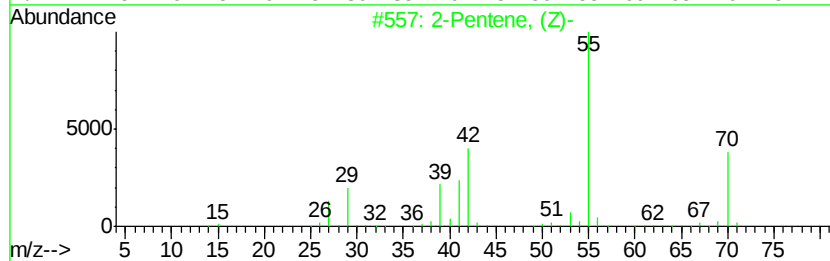
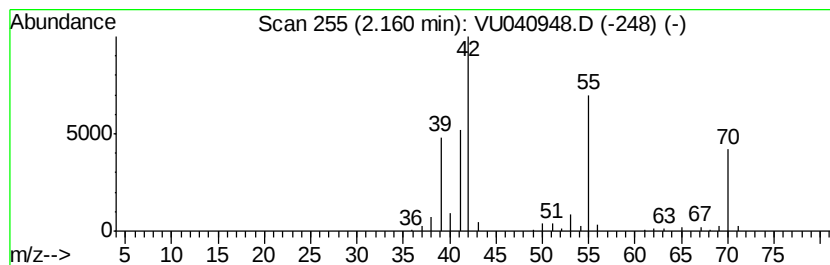
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR102620WMA.M
Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.16	3.75 ug/L	109358	1,4-Difluorobenzene	6.26

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentene, (Z)-		70	C5H10	000627-20-3	87
2	2-Pentene		70	C5H10	000109-68-2	87
3	Cyclopropane, ethyl-		70	C5H10	001191-96-4	86
4	Cyclopropane, ethyl-		70	C5H10	001191-96-4	70
5	1-Pentene		70	C5H10	000109-67-1	70



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU102620\
Data File : VU040948.D
Acq On : 26 Oct 2020 16:47
Operator : SY/MD
Sample : L4492-01
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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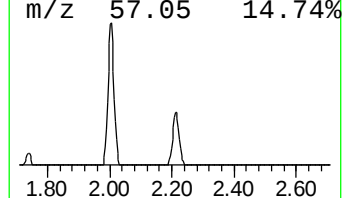
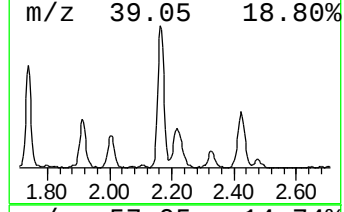
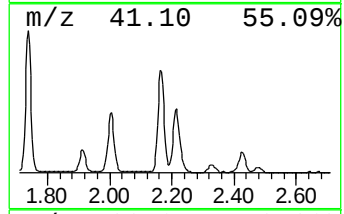
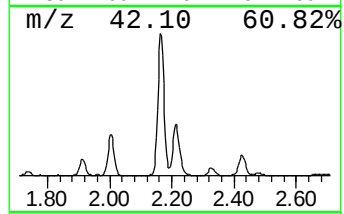
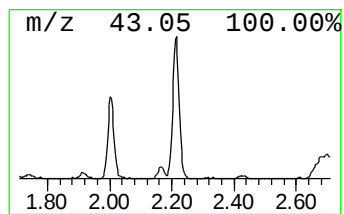
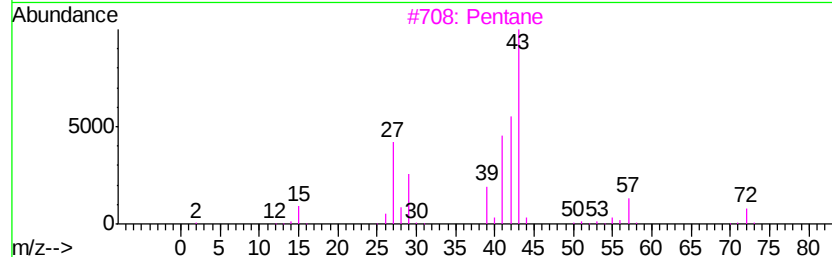
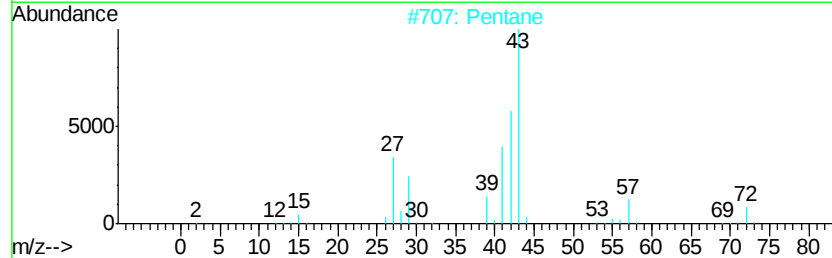
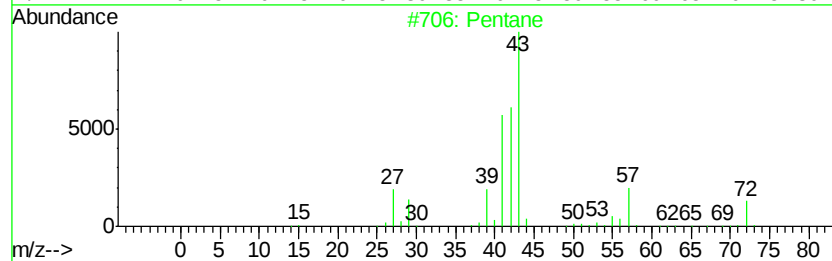
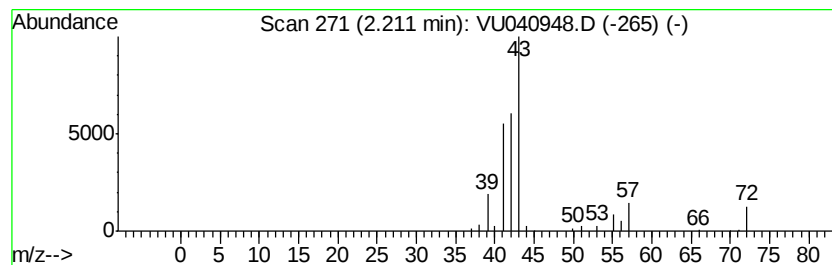
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.21	2.32 ug/L	67604	1,4-Difluorobenzene	6.26

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Pentane	72	C5H12	000109-66-0	86
2		Pentane	72	C5H12	000109-66-0	64
3		Pentane	72	C5H12	000109-66-0	64
4		Pentane	72	C5H12	000109-66-0	45
5		Oxirane, ethyl-	72	C4H8O	000106-88-7	42



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU102620\
Data File : VU040948.D
Acq On : 26 Oct 2020 16:47
Operator : SY/MD
Sample : L4492-01
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampled :
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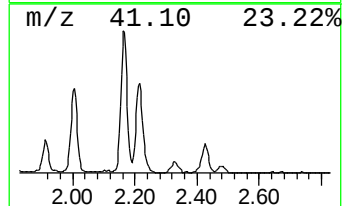
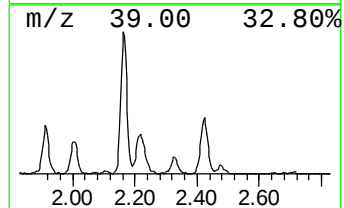
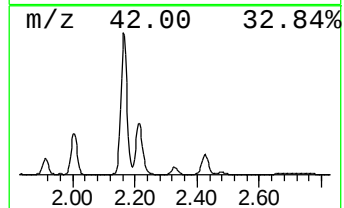
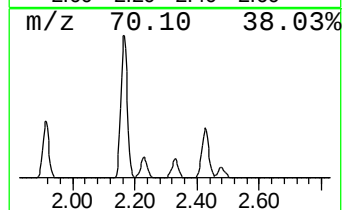
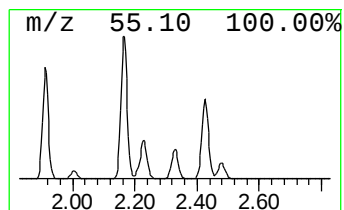
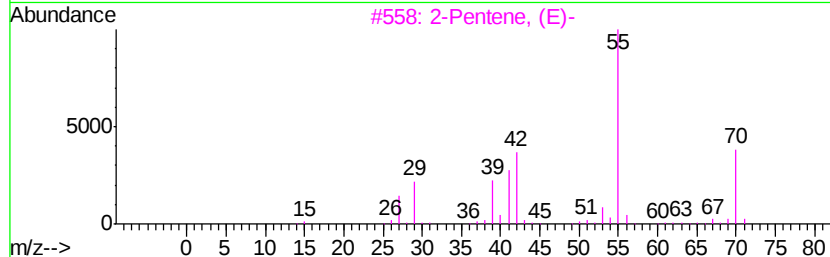
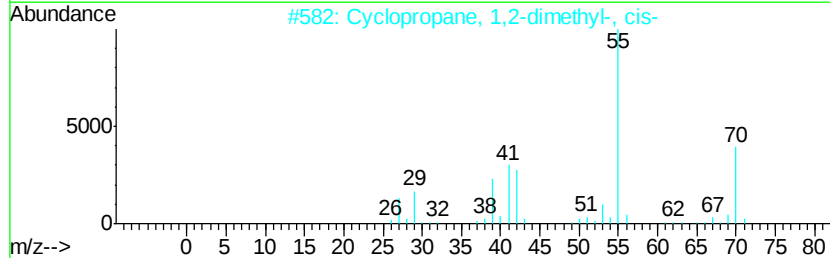
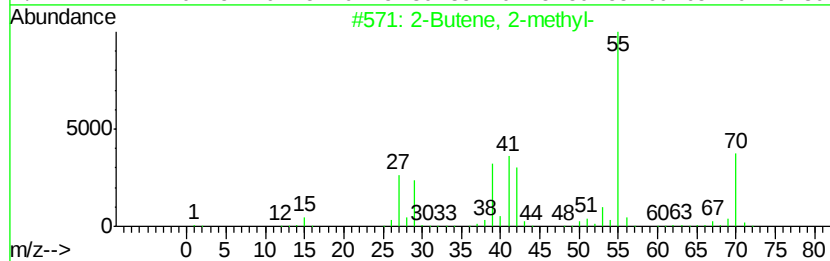
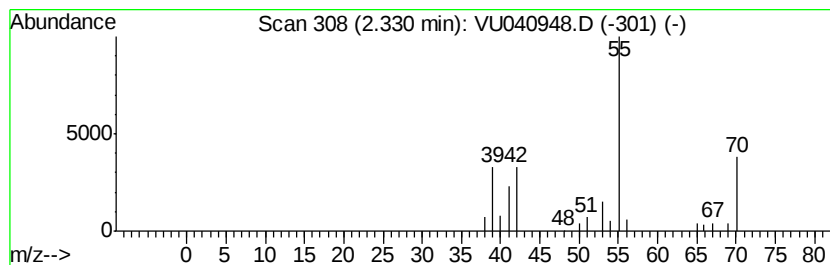
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.33	1.16 ug/L	33685	1,4-Difluorobenzene	6.26

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Butene, 2-methyl-	70	C5H10	000513-35-9	86
2			Cyclopropane, 1,2-dimethyl-, cis-	70	C5H10	000930-18-7	86
3			2-Pentene, (E)-	70	C5H10	000646-04-8	86
4			2-Pentene, (E)-	70	C5H10	000646-04-8	83
5			1-Butene, 3-methyl-	70	C5H10	000563-45-1	80



Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU102620\
Data File : VU040948.D
Acq On : 26 Oct 2020 16:47
Operator : SY/MD
Sample : L4492-01
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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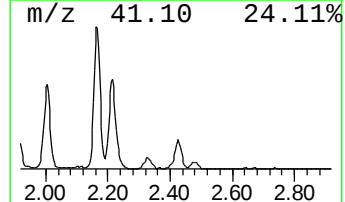
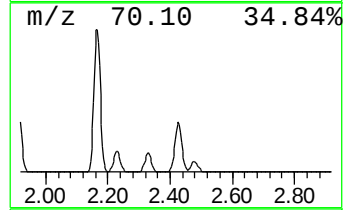
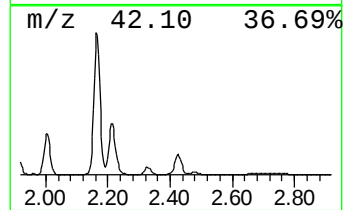
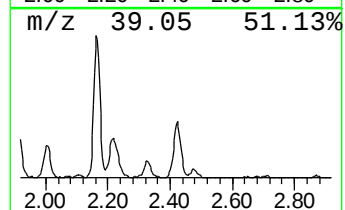
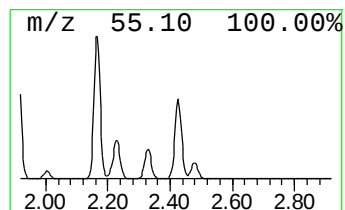
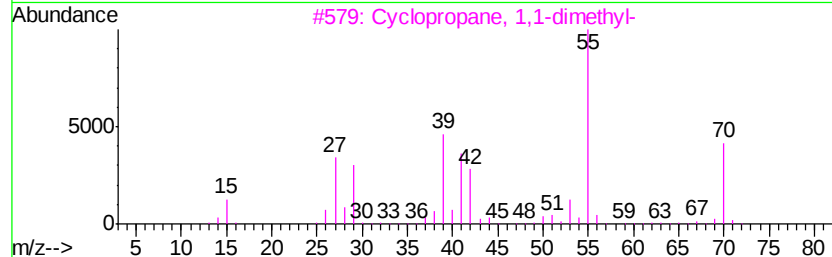
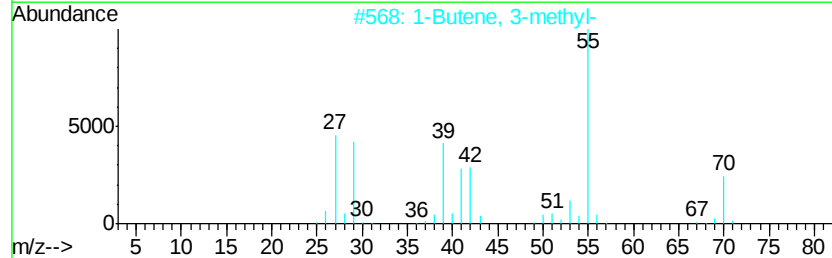
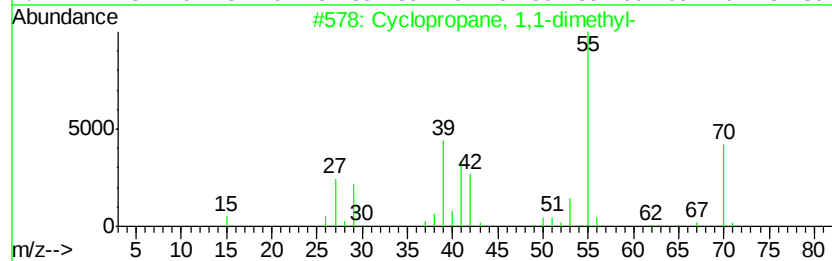
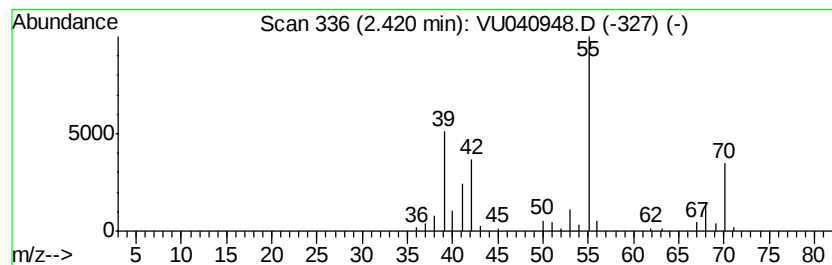
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 (DEL) Alkane: Cyclic2.42 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.42	1.41 ug/L	41015	1,4-Difluorobenzene	6.26

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopropane, 1,1-dimethyl-	70	C5H10	001630-94-0	80
2			1-Butene, 3-methyl-	70	C5H10	000563-45-1	72
3			Cyclopropane, 1,1-dimethyl-	70	C5H10	001630-94-0	72
4			2-Pentene	70	C5H10	000109-68-2	53
5			2-Butene, 2-methyl-	70	C5H10	000513-35-9	53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU102620\
Data File : VU040948.D
Acq On : 26 Oct 2020 16:47
Operator : SY/MD
Sample : L4492-01
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampled :
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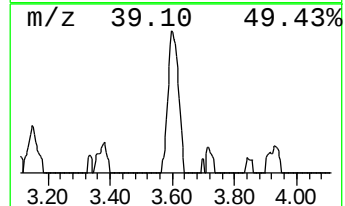
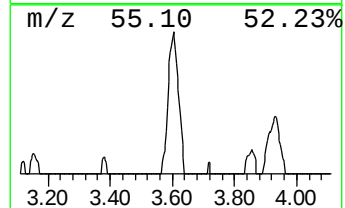
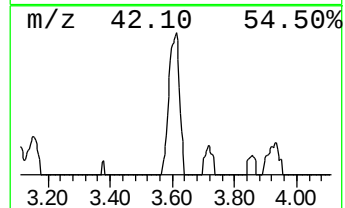
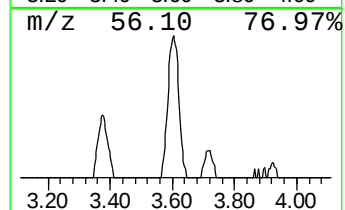
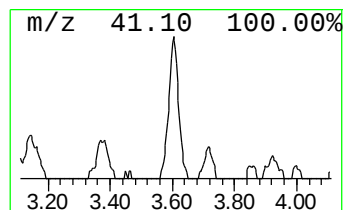
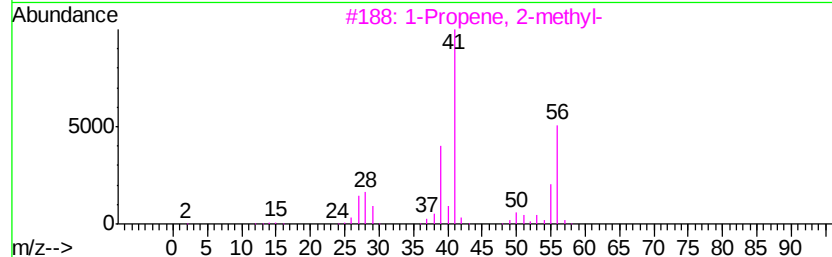
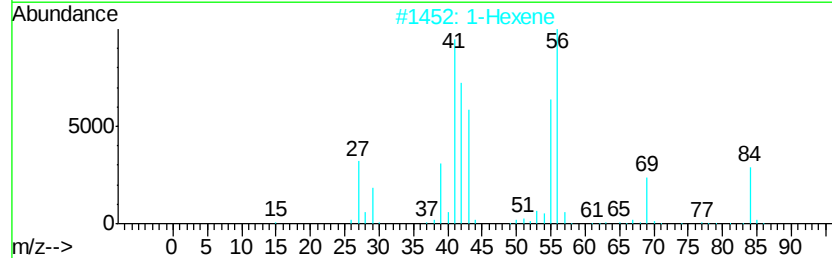
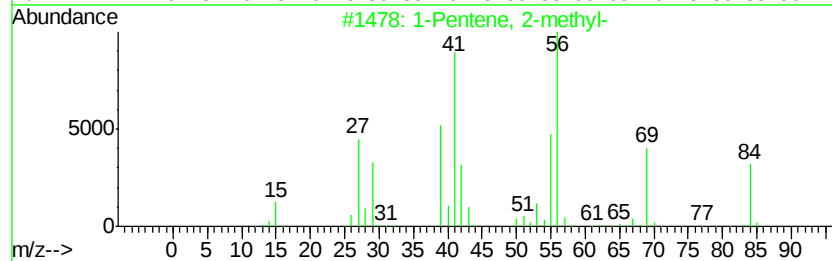
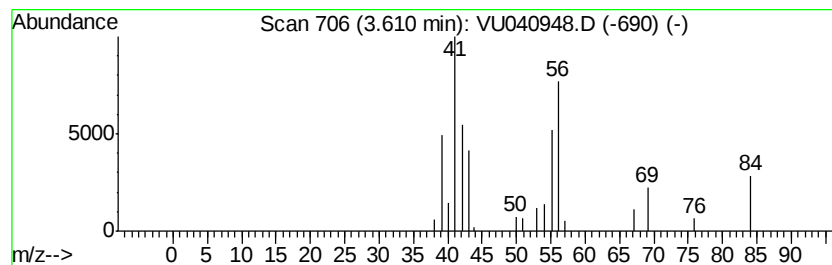
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.61	0.87 ug/L	25373	1,4-Difluorobenzene	6.26

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Pentene, 2-methyl-	84	C6H12	000763-29-1	59
2	1-Hexene	84	C6H12	000592-41-6	53
3	1-Propene, 2-methyl-	56	C4H8	000115-11-7	52
4	3-Hexene, (E)-	84	C6H12	013269-52-8	52
5	3-Hexene, (Z)-	84	C6H12	007642-09-3	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU102620\
Data File : VU040948.D
Acq On : 26 Oct 2020 16:47
Operator : SY/MD
Sample : L4492-01
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampled :
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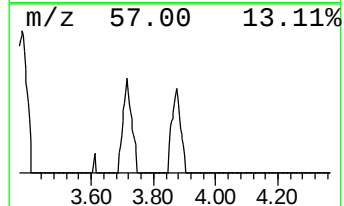
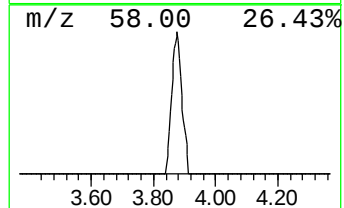
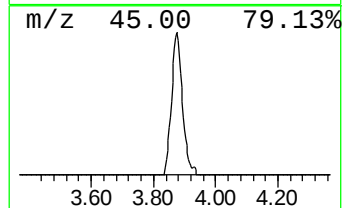
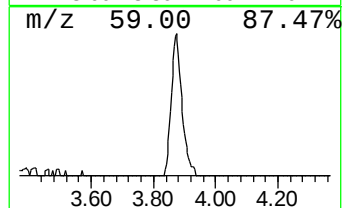
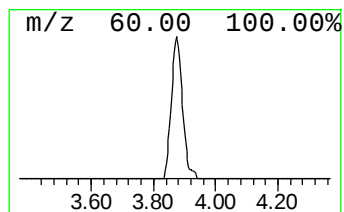
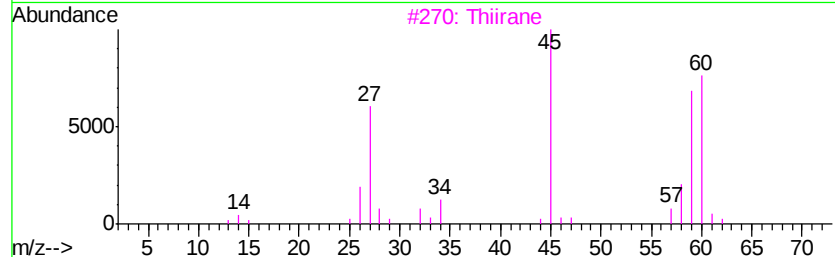
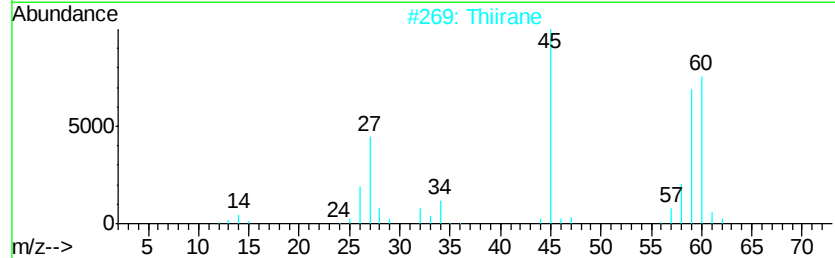
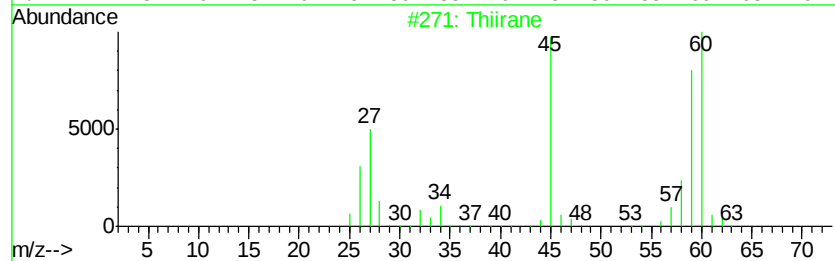
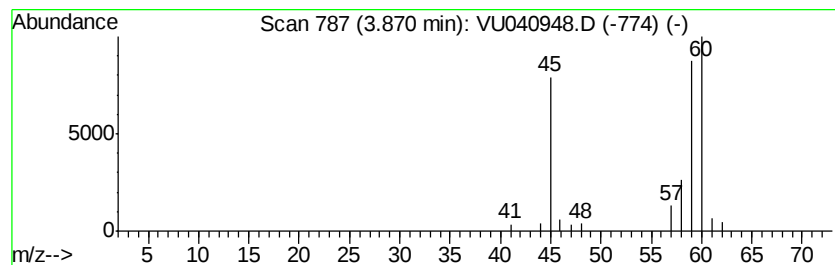
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Thiirane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.87	1.04 ug/L	30307	1,4-Difluorobenzene	6.26

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Thiirane	60	C2H4S	000420-12-2	90
2			Thiirane	60	C2H4S	000420-12-2	72
3			Thiirane	60	C2H4S	000420-12-2	49
4			Silane, dimethyl-	60	C2H8Si	001111-74-6	28
5			Hydrazine, 1,1-dimethyl-	60	C2H8N2	000057-14-7	9



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU102620\
Data File : VU040948.D
Acq On : 26 Oct 2020 16:47
Operator : SY/MD
Sample : L4492-01
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ALS Vial : 17 Sample Multiplier: 1

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ClientSampleId :
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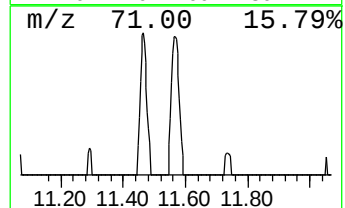
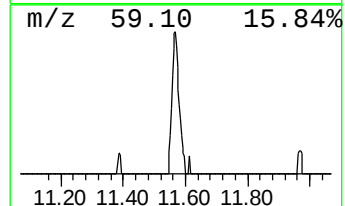
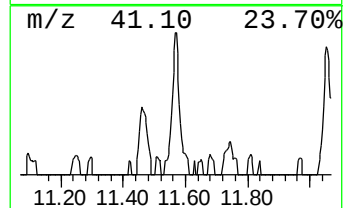
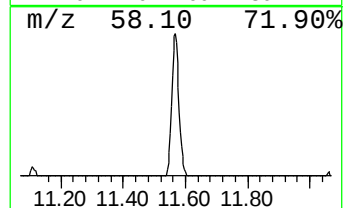
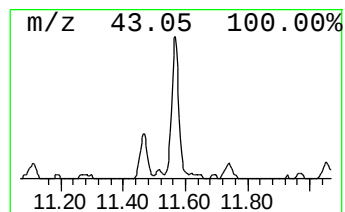
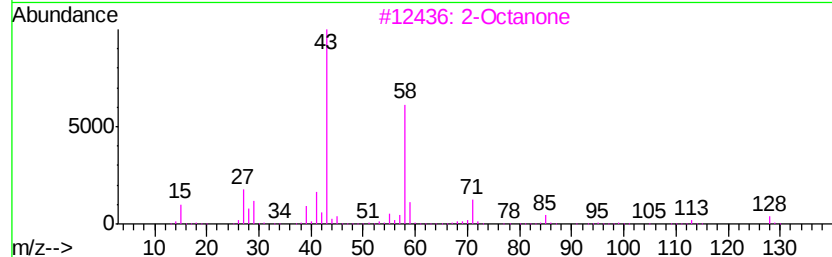
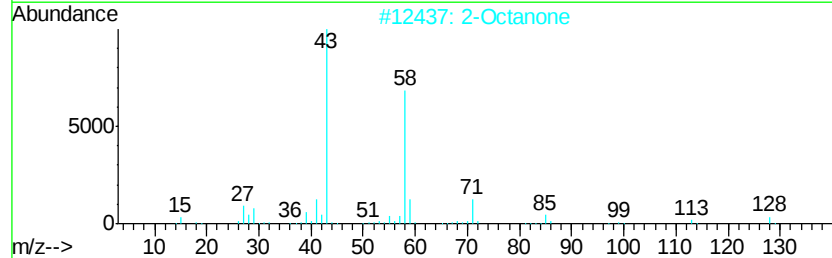
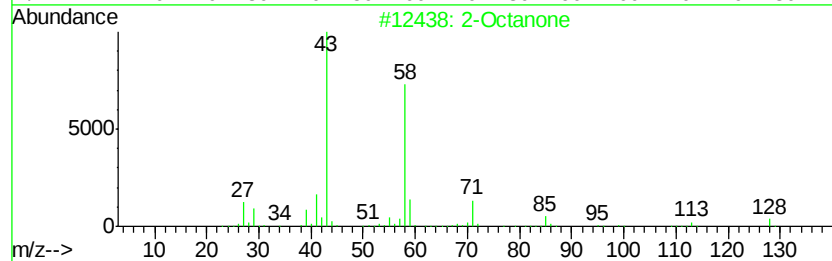
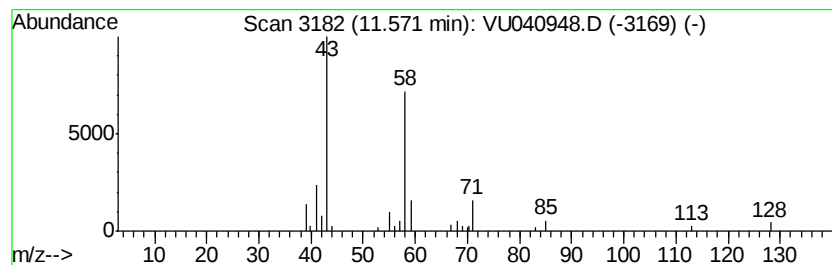
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 2-Octanone Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.57	0.68 ug/L	23946	1,4-Dichlorobenzene-d4	11.81

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Octanone	128	C8H16O	000111-13-7	91
2			2-Octanone	128	C8H16O	000111-13-7	91
3			2-Octanone	128	C8H16O	000111-13-7	86
4			2-Octanone	128	C8H16O	000111-13-7	80
5			2-Heptanone	114	C7H14O	000110-43-0	72



Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU102620\
Data File : VU040948.D
Acq On : 26 Oct 2020 16:47
Operator : SY/MD
Sample : L4492-01
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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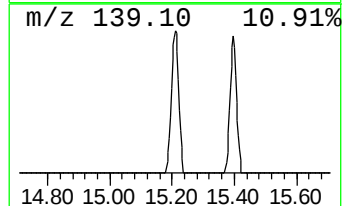
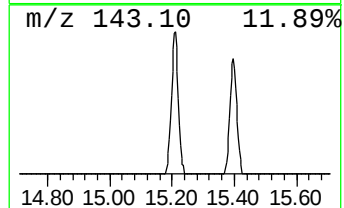
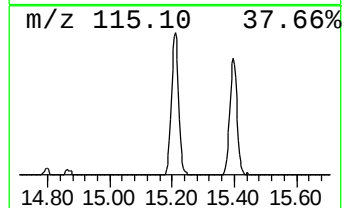
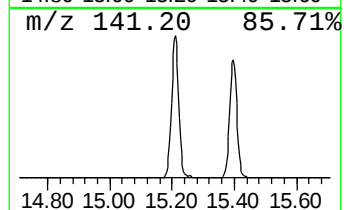
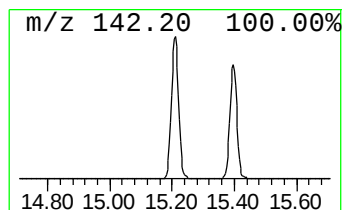
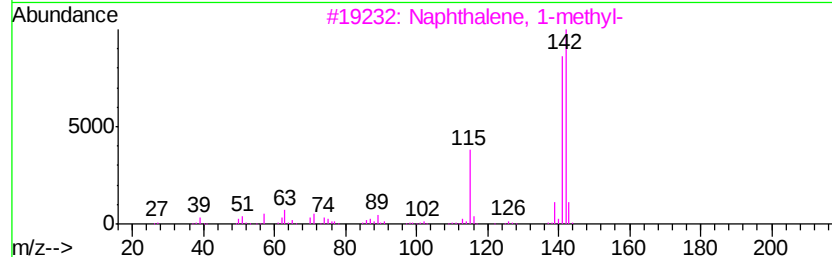
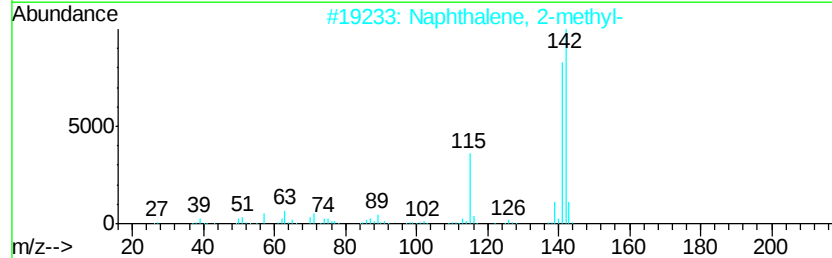
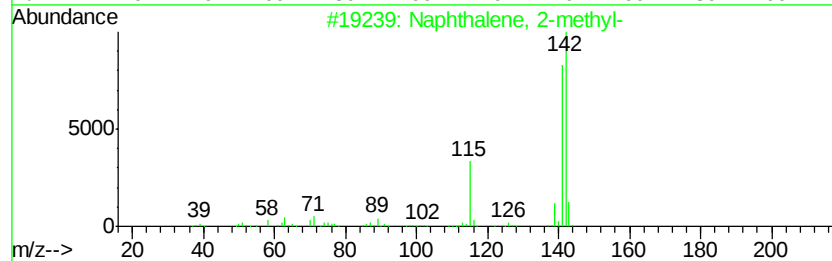
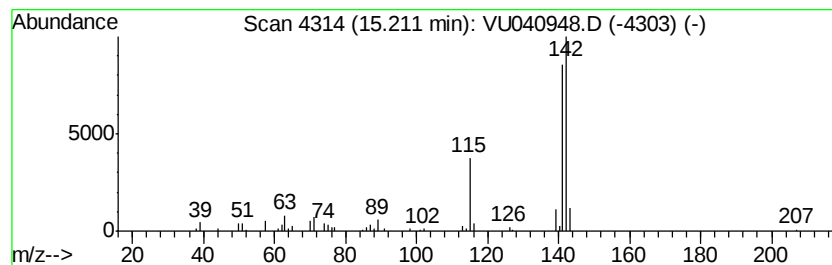
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Naphthalene, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.21	2.32 ug/L	81342	1,4-Dichlorobenzene-d4	11.81

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2		Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
4		1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	94
5		Naphthalene, 1-methyl-	142	C11H10	000090-12-0	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU102620\
Data File : VU040948.D
Acq On : 26 Oct 2020 16:47
Operator : SY/MD
Sample : L4492-01
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
C0AM9

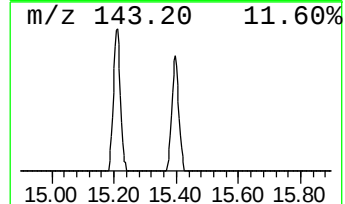
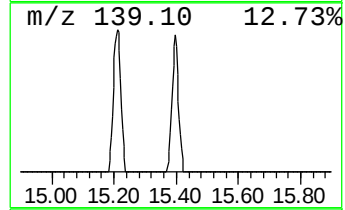
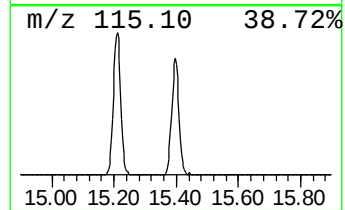
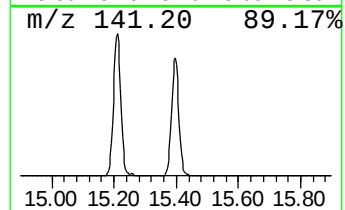
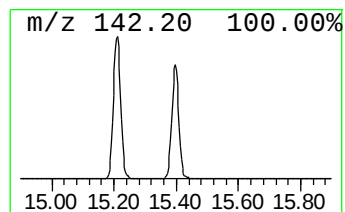
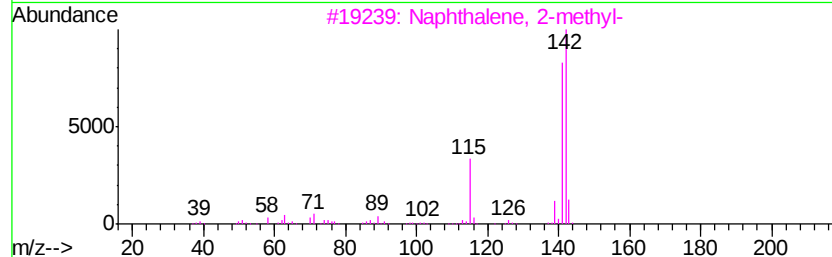
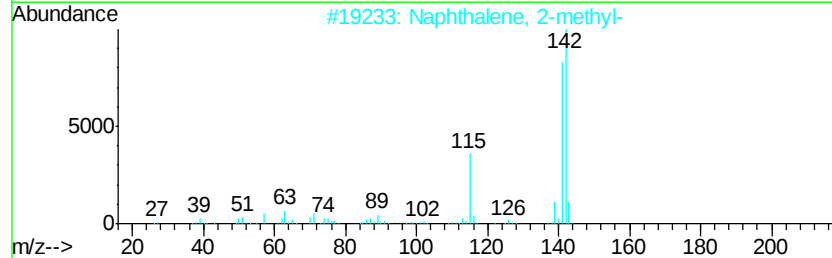
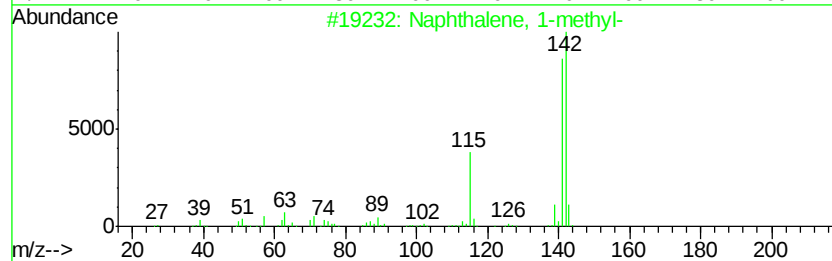
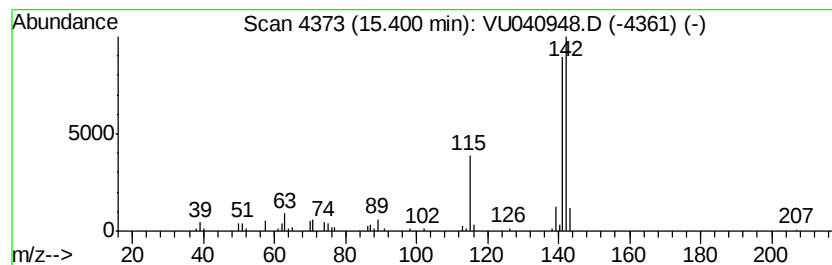
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Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 Naphthalene, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.40	1.84 ug/L	64744	1,4-Dichlorobenzene-d4	11.81

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
4			Benzocycloheptatriene	142	C11H10	000264-09-5	94
5			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	94



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU102620\
 Data File : VU040948.D
 Acq On : 26 Oct 2020 16:47
 Operator : SY/MD
 Sample : L4492-01
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 C0AM9

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR102620WMA.M
 Quant Title : TRACE VOA SOM01.0

TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: Str...	1.36	25.1	ug/L	731981	1	6.26	145700	5.0
(DEL) Alkane: Str...	1.66	0.5	ug/L	15798	1	6.26	145700	5.0
(DEL) Alkane: Str...	1.74	2.6	ug/L	75142	1	6.26	145700	5.0
(DEL) Alkane: Str...	2.00	1.7	ug/L	50743	1	6.26	145700	5.0
(DEL) Alkane: Str...	2.16	3.8	ug/L	109358	1	6.26	145700	5.0
(DEL) Alkane: Str...	2.21	2.3	ug/L	67604	1	6.26	145700	5.0
(DEL) Alkane: Str...	2.33	1.2	ug/L	33685	1	6.26	145700	5.0
(DEL) Alkane: Cyc...	2.42	1.4	ug/L	41015	1	6.26	145700	5.0
(DEL) Alkane: Str...	3.61	0.9	ug/L	25373	1	6.26	145700	5.0
Thiirane	3.87	1.0	ug/L	30307	1	6.26	145700	5.0
2-Octanone	11.57	0.7	ug/L	23946	3	11.81	175593	5.0
Naphthalene, 2-me...	15.21	2.3	ug/L	81342	3	11.81	175593	5.0
Naphthalene, 1-me...	15.40	1.8	ug/L	64744	3	11.81	175593	5.0