LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV052020\
Data File : VV016136.D
Acq On : 20 May 2020 17:02
Operator : SY/MD
Sample : L2725-07
Misc : 25.0mL/MSVOA V/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : F9B17

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_V\METHOD\SOMVTR050720WMA.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	1.316	63	70	78	rVB	54520	58433	2.38%	0.434%
2	1.579	145	152	167	rVB	64826	79701	3.24%	0.592%
3	1.984	269	278	286	rBV2	18491	25482	1.04%	0.189%
4	2.042	286	296	311	rVB	151263	215817	8.77%	1.603%
5	2.126	311	322	340	rVB	105059	164178	6.67%	1.219%
6	2.788	512	528	556	rVB	1174664	2460317	100.00%	18.273%
7	3.299	673	687	703	rBV3	23612	52893	2.15%	0.393%
8	3.396	703	717	737	rVB3	14767	35863	1.46%	0.266%
9	3.608	769	783	798	rBV2	33609	79544	3.23%	0.591%
10	3.878	848	867	873	rBV2	18551	42397	1.72%	0.315%
11	3.930	873	883	925	rVB	62852	214257	8.71%	1.591%
12	4.380	1004	1023	1041	rBV	80386	218604	8.89%	1.624%
13	4.489	1043	1057	1082	rVB	106880	265048	10.77%	1.969%
14	5.077	1223	1240	1264	rBV2	194894	492530	20.02%	3.658%
15	5.258	1273	1296	1345	rBV2	461010	2210194	89.83%	16.415%
16	5.643	1404	1416	1436	rBV	196097	455491	18.51%	3.383%
17	6.097	1539	1557	1573	rBV	119213	245135	9.96%	1.821%
18	6.827	1765	1784	1794	rBV2	44710	91874	3.73%	0.682%
19	6.881	1794	1801	1815	rVB	16031	28835	1.17%	0.214%
20	7.010	1832	1841	1854	rVV2	54111	99979	4.06%	0.743%
21	7.100	1855	1869	1882	rVV	310147	680401	27.66%	5.053%
22	7.174	1882	1892	1930	rVV	328332	792060	32.19%	5.883%
23	7.338	1931	1943	1961	rVB	243862	423043	17.19%	3.142%
24	7.524	1985	2001	2022	rBV	632025	1375367	55.90%	10.215%
25	7.647	2031	2039	2054	rVB	33513	60212	2.45%	0.447%
26	8.113	2174	2184	2199	rBV	300447	566168	23.01%	4.205%
27	8.341	2234	2255	2277	rVB7	32849	104411	4.24%	0.775%
28	8.833	2390	2408	2412	rBV5	37390	80409	3.27%	0.597%
29	8.875	2412	2421	2431	rVV	301102	486367	19.77%	3.612%
30	8.929	2431	2438	2445	rVV	45332	74137	3.01%	0.551%
31	8.978	2445	2453	2479	rVB	135961	257147	10.45%	1.910%
32	9.225	2519	2530	2547	rVB2	23000	40708	1.65%	0.302%
33	10.238	2835	2845	2857	rVB	87511	139772	5.68%	1.038%
34	11.273	3154	3167	3192	rBV	297385	469576	19.09%	3.488%

LSC Area Percent Report

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Misc : 25.0mL/MSVOA V/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : F9B17

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_V\METHOD\SOMVTR050720WMA.M
Title : TRACE VOA SOM01.0

35	11.646	3271	3283	3299	rBV	239150	378089	15.37%	2.808%

Sum of corrected areas: 13464439

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV052020\
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Acq On : 20 May 2020 17:02
Operator : SY/MD
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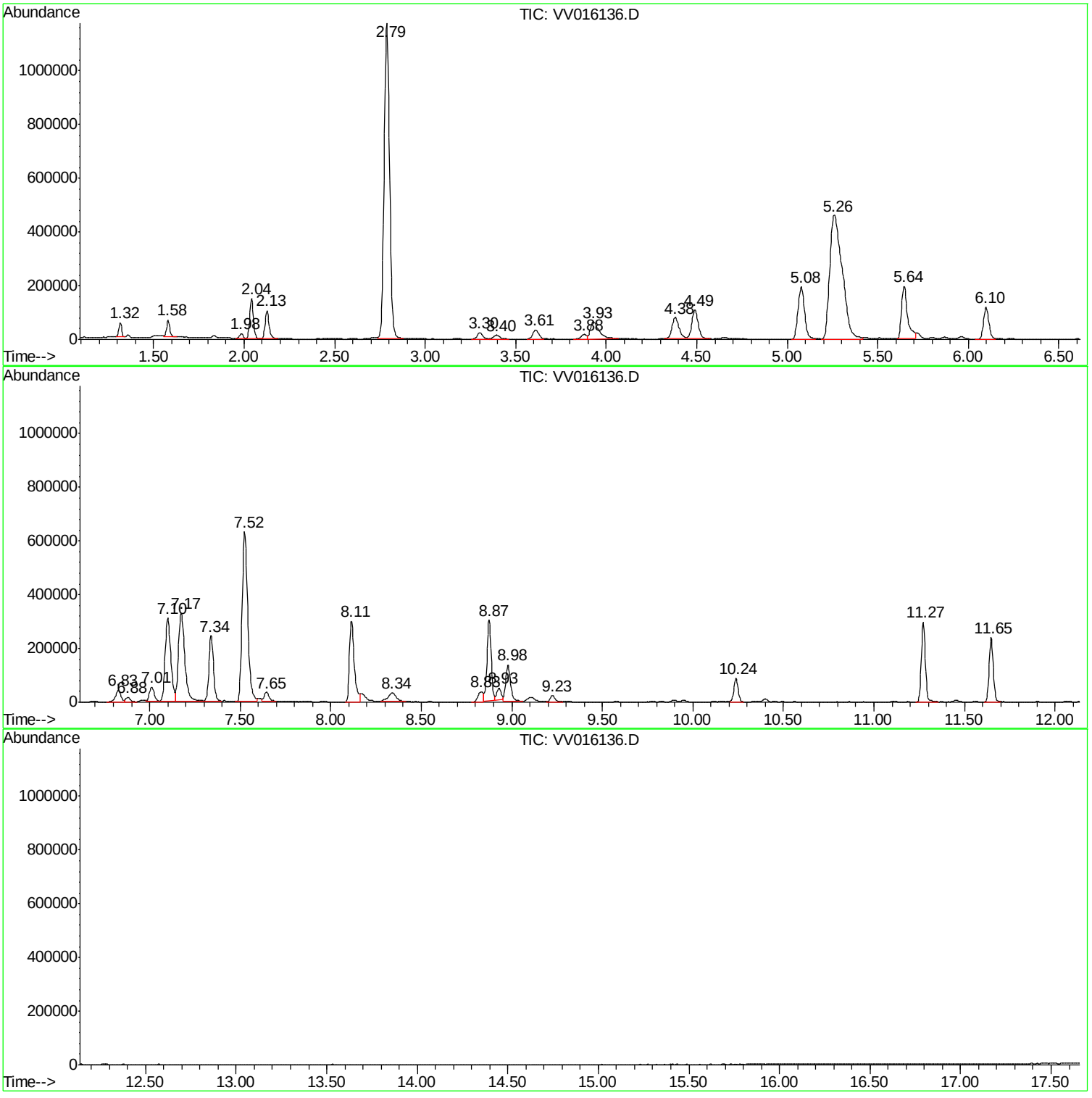
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR050720WMA.M
Quant Title : TRACE VOA SOM01.0

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



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Data File : VV016136.D
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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR050720WMA.M
Quant Title : TRACE VOA SOM01.0

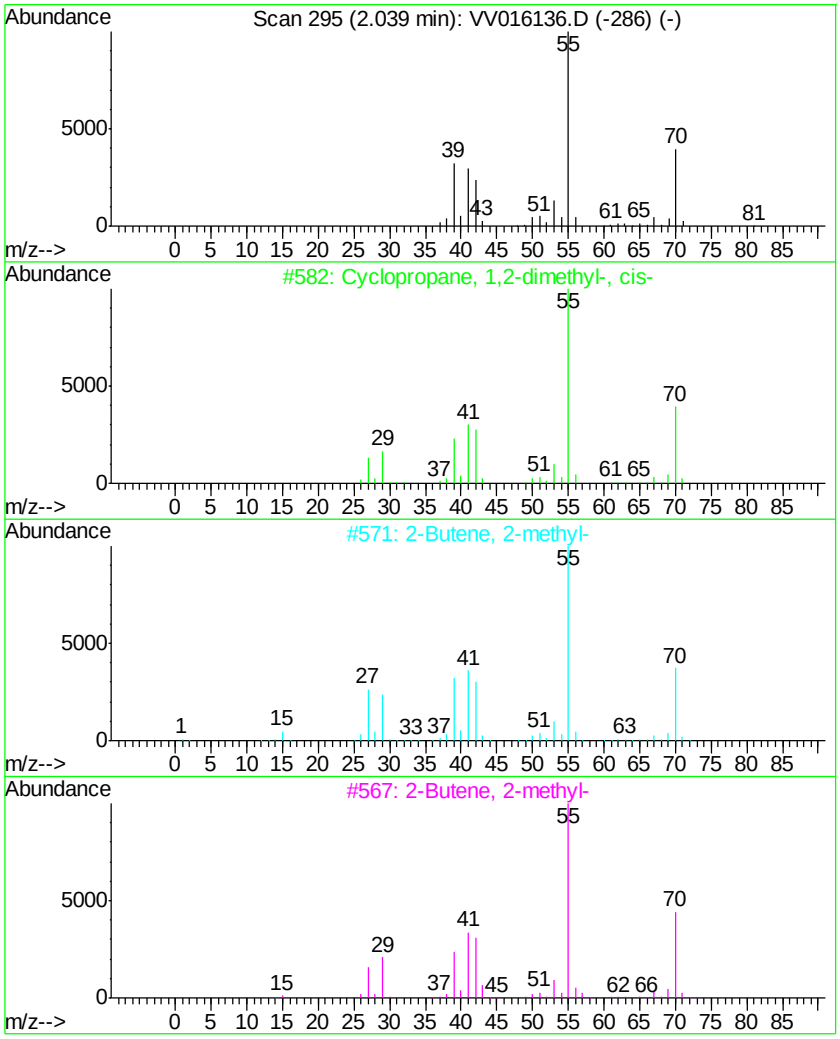
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

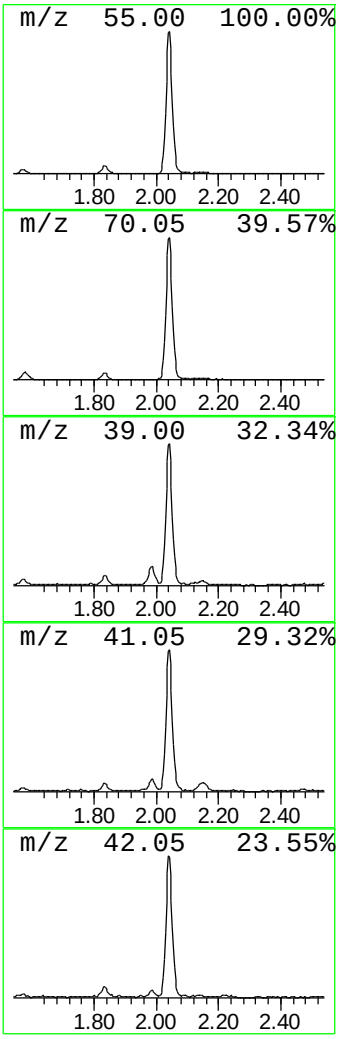
****

Peak Number 1 (DEL) Alkane: Cyclic2.04 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.04	2.37 ug/L	215817	1,4-Difluorobenzene	5.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopropane, 1,2-dimethyl-, cis-	70	C5H10	000930-18-7	91
2	2-Butene, 2-methyl-	70	C5H10	000513-35-9	91
3	2-Butene, 2-methyl-	70	C5H10	000513-35-9	90
4	1-Butene, 3-methyl-	70	C5H10	000563-45-1	87
5	2-Pentene, (E)-	70	C5H10	000646-04-8	87





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV052020\
Data File : VV016136.D
Acq On : 20 May 2020 17:02
Operator : SY/MD
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Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : F9B17

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR050720WMA.M
Quant Title : TRACE VOA SOM01.0

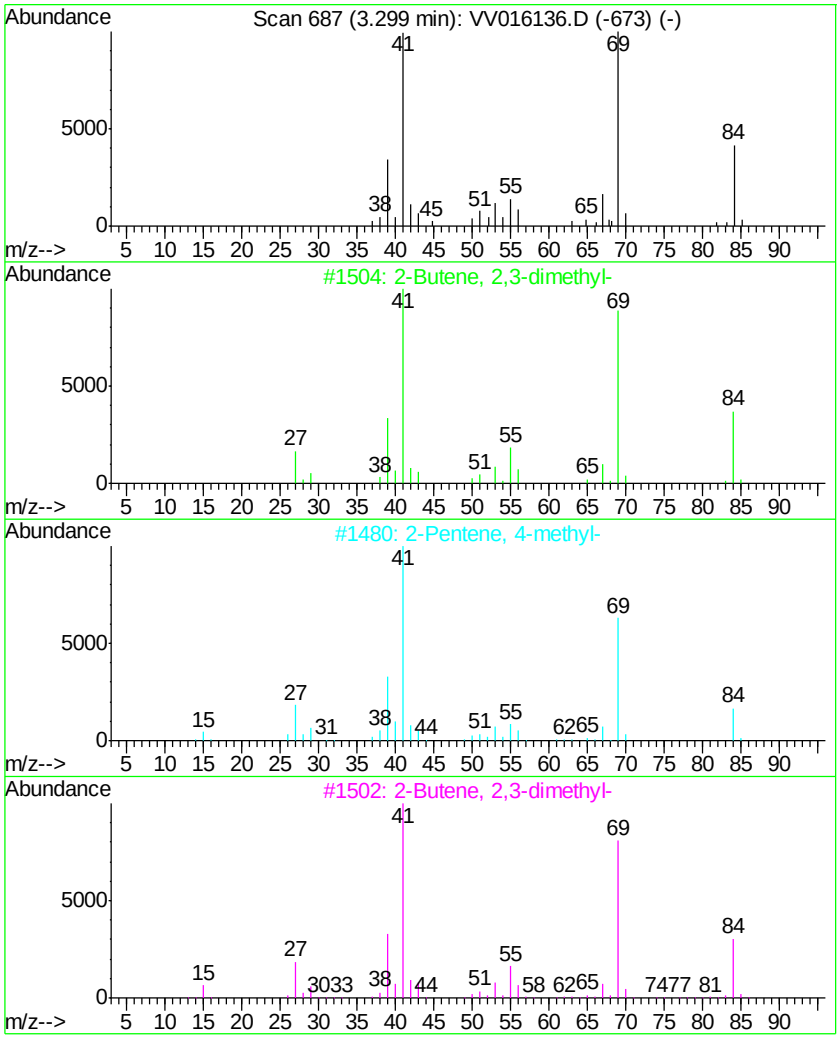
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

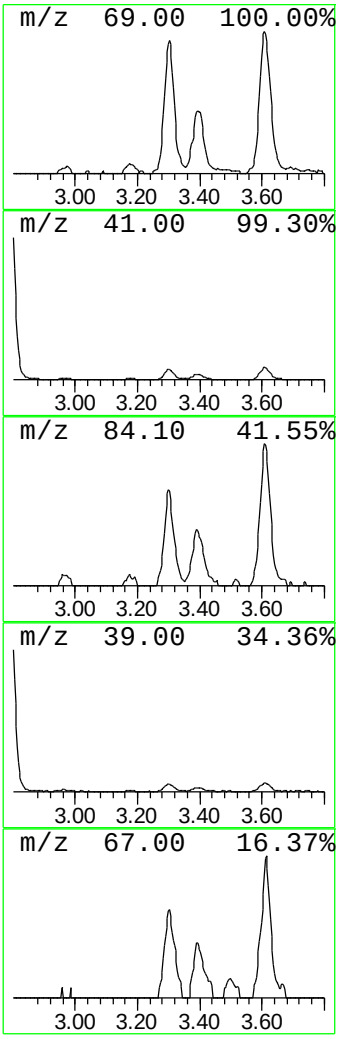
****

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.30	0.58 ug/L	52893	1,4-Difluorobenzene	5.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butene, 2,3-dimethyl-	84	C6H12	000563-79-1	91
2	2-Pentene, 4-methyl-	84	C6H12	004461-48-7	91
3	2-Butene, 2,3-dimethyl-	84	C6H12	000563-79-1	91
4	2-Pentene, 4-methyl-, (Z)-	84	C6H12	000691-38-3	91
5	2-Pentene, 4-methyl-, (Z)-	84	C6H12	000691-38-3	91





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV052020\
Data File : VV016136.D
Acq On : 20 May 2020 17:02
Operator : SY/MD
Sample : L2725-07
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : F9B17

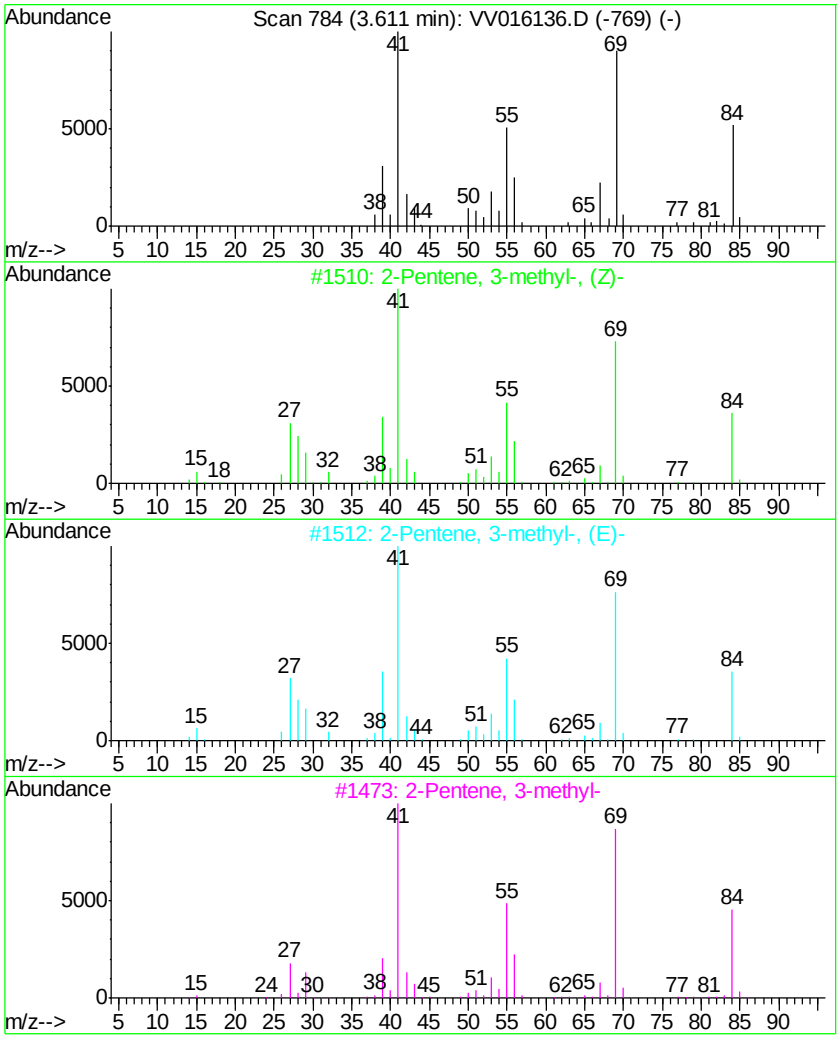
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR050720WMA.M
Quant Title : TRACE VOA SOM01.0

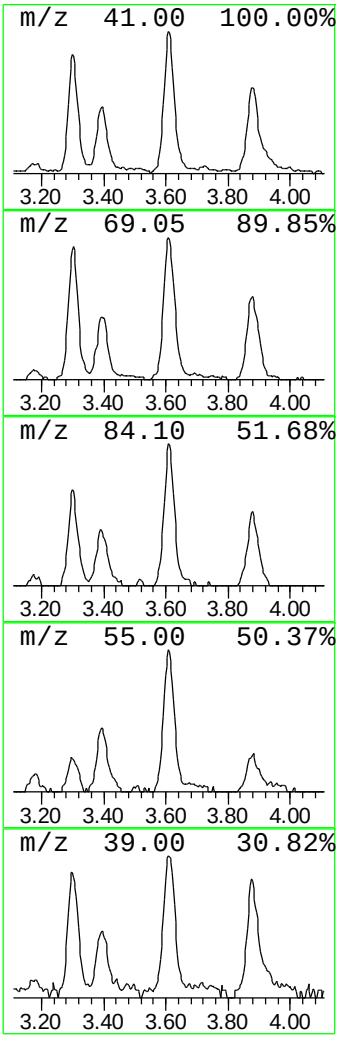
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.61	0.87 ug/L	79544	1,4-Difluorobenzene	5.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentene, 3-methyl-, (Z)-	84	C6H12	000922-62-3	94
2	2-Pentene, 3-methyl-, (E)-	84	C6H12	000616-12-6	94
3	2-Pentene, 3-methyl-	84	C6H12	000922-61-2	90
4	2-Pentene, 3-methyl-, (Z)-	84	C6H12	000922-62-3	90
5	2-Pentene, 3-methyl-, (E)-	84	C6H12	000616-12-6	90





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV052020\
Data File : VV016136.D
Acq On : 20 May 2020 17:02
Operator : SY/MD
Sample : L2725-07
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : F9B17

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR050720WMA.M
Quant Title : TRACE VOA SOM01.0

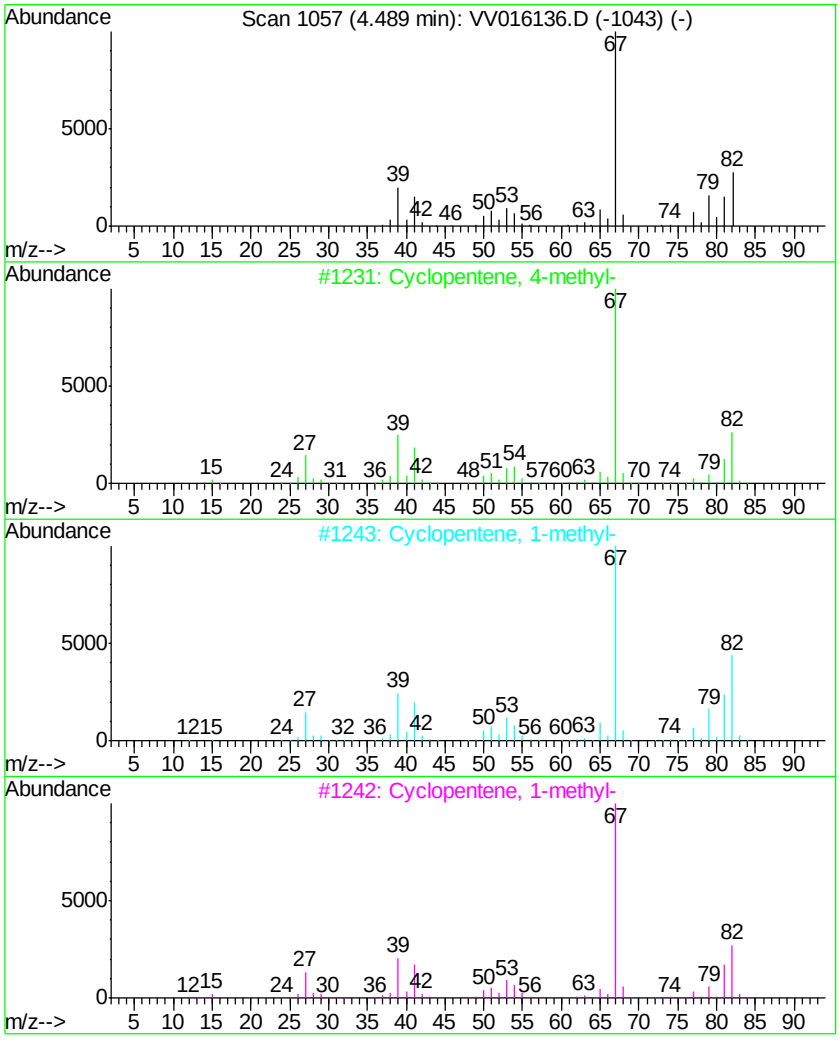
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

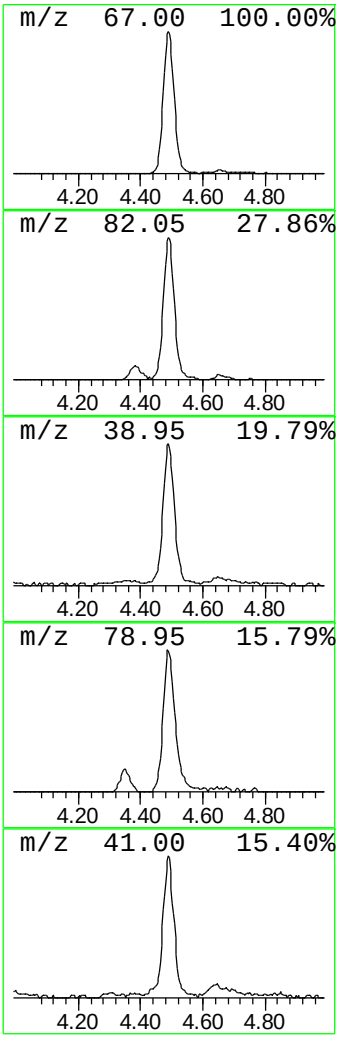
****

Peak Number 4 Cyclopentene, 4-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.49	2.91 ug/L	265048	1,4-Difluorobenzene	5.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentene, 4-methyl-	82	C6H10	001759-81-5	87
2	Cyclopentene, 1-methyl-	82	C6H10	000693-89-0	87
3	Cyclopentene, 1-methyl-	82	C6H10	000693-89-0	87
4	Cyclopentene, 3-methyl-	82	C6H10	001120-62-3	86
5	Cyclopentene, 1-methyl-	82	C6H10	000693-89-0	83





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV052020\
Data File : VV016136.D
Acq On : 20 May 2020 17:02
Operator : SY/MD
Sample : L2725-07
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : F9B17

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR050720WMA.M
Quant Title : TRACE VOA SOM01.0

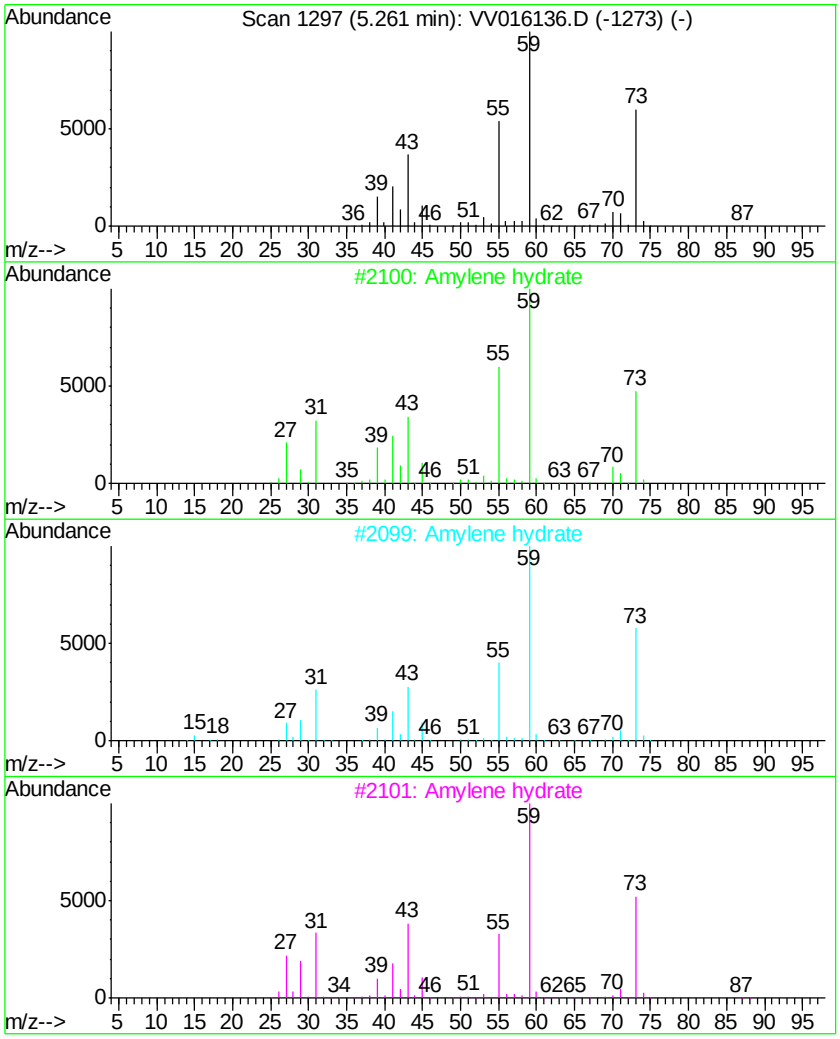
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

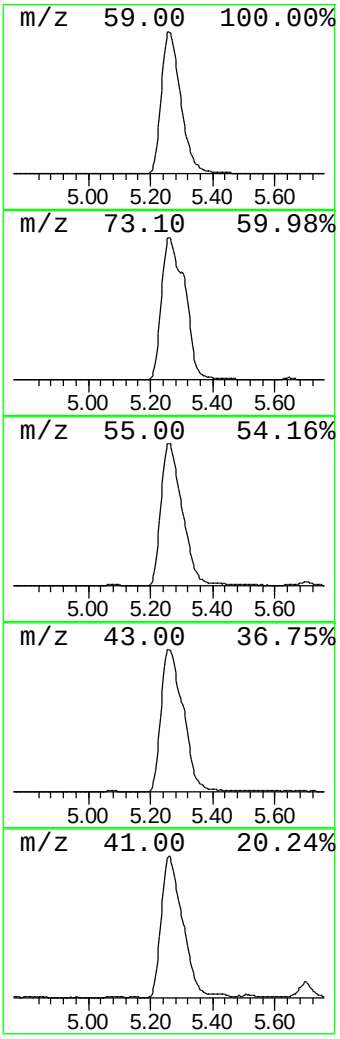
****

Peak Number 5 Amylene hydrate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.26	24.26 ug/L	2210190	1,4-Difluorobenzene	5.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Amylene hydrate	88	C5H12O	000075-85-4	83
2	Amylene hydrate	88	C5H12O	000075-85-4	78
3	Amylene hydrate	88	C5H12O	000075-85-4	72
4	Propanoic acid, 2-hydroxy-2-methyl-	104	C4H8O3	000594-61-6	53
5	3-Pentanol, 2-methyl-	102	C6H14O	000565-67-3	39





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ALS Vial : 15 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : F9B17

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR050720WMA.M
Quant Title : TRACE VOA SOM01.0

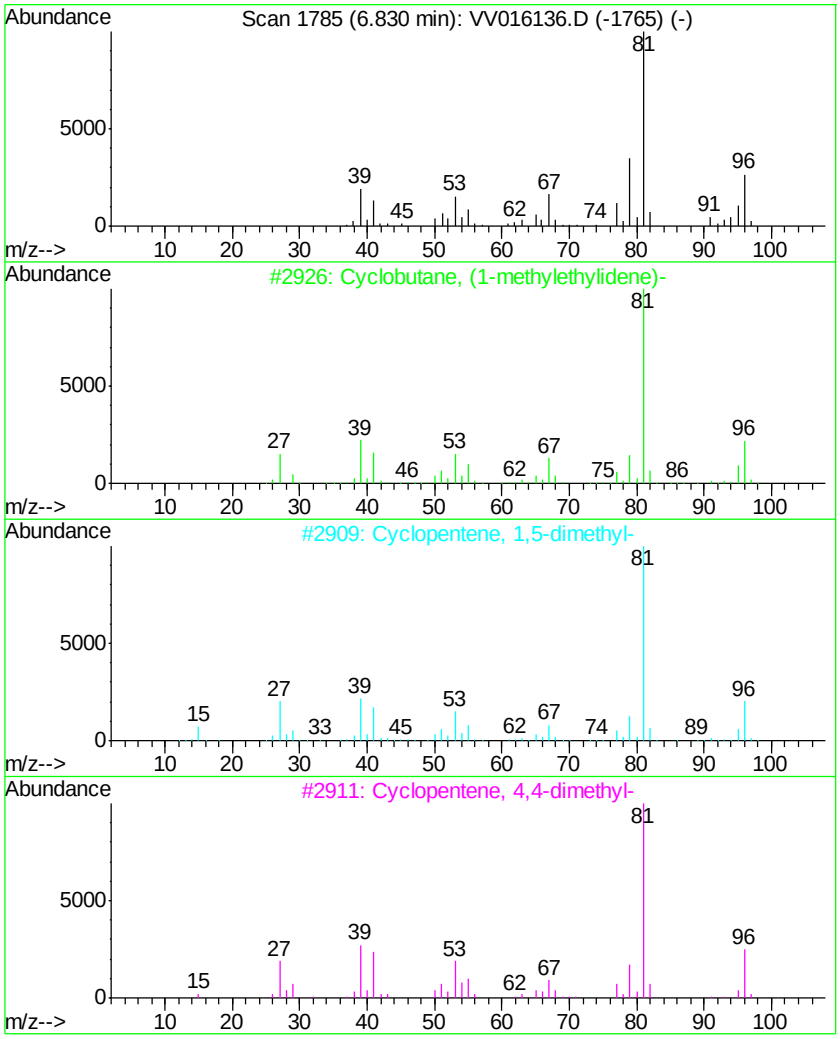
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

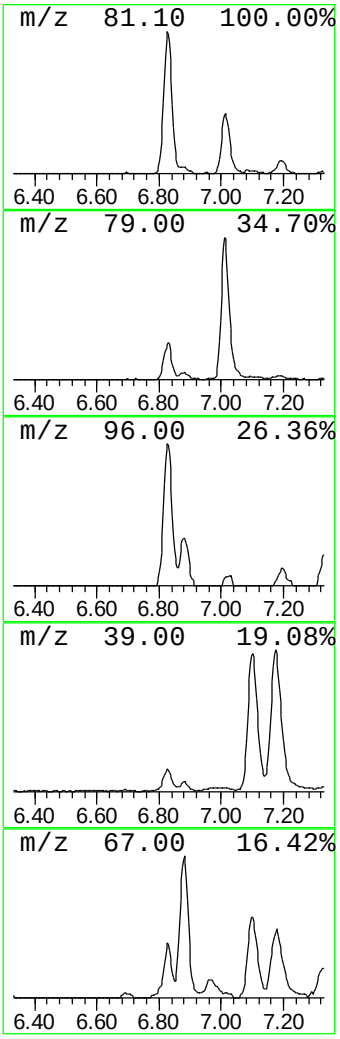
****

Peak Number 6 Cyclobutane, (1-methylethyl... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.83	1.01 ug/L	91874	1,4-Difluorobenzene	5.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclobutane, (1-methylethylidene)-	96	C7H12	001528-22-9	90
2	Cyclopentene, 1,5-dimethyl-	96	C7H12	016491-15-9	87
3	Cyclopentene, 4,4-dimethyl-	96	C7H12	019037-72-0	83
4	3,5-Dimethylcyclopentene	96	C7H12	007459-71-4	72
5	2-Heptyne	96	C7H12	001119-65-9	59





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ALS Vial : 15 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : F9B17

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR050720WMA.M
Quant Title : TRACE VOA SOM01.0

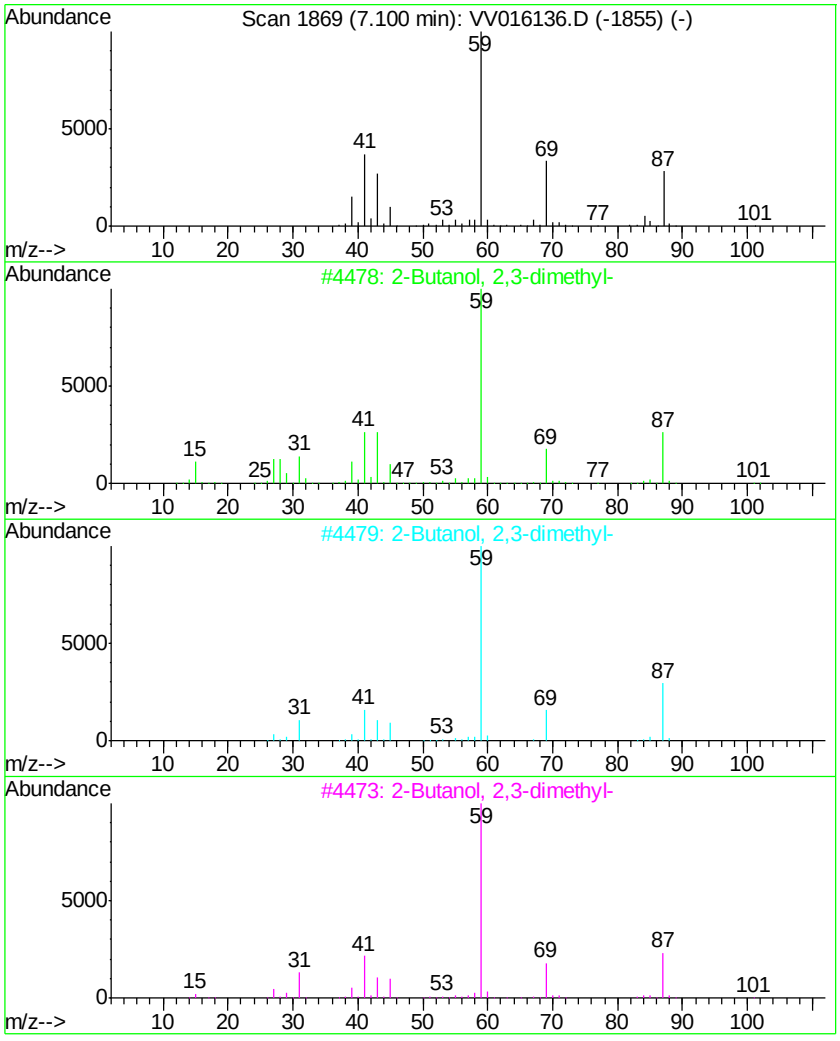
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

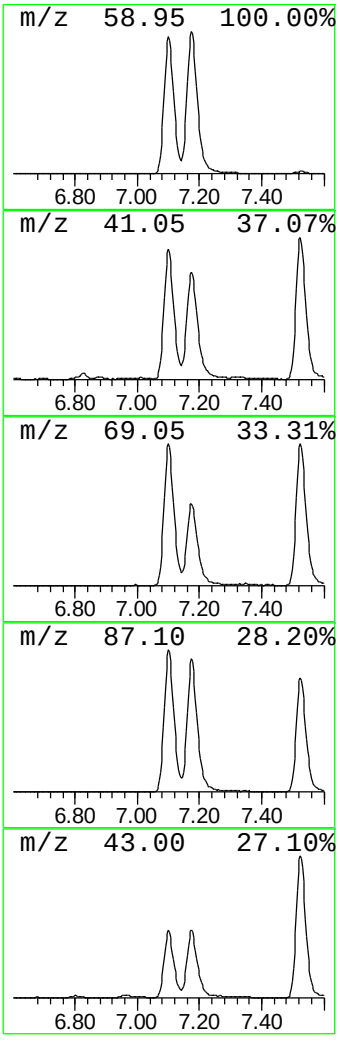
****

Peak Number 8 2-Butanol, 2,3-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.10	7.47 ug/L	680401	1,4-Difluorobenzene	5.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butanol, 2,3-dimethyl-	102	C6H14O	000594-60-5	83
2	2-Butanol, 2,3-dimethyl-	102	C6H14O	000594-60-5	83
3	2-Butanol, 2,3-dimethyl-	102	C6H14O	000594-60-5	74
4	2-Pentanol, 2-methyl-	102	C6H14O	000590-36-3	72
5	Propane, 2-ethoxy-2-methyl-	102	C6H14O	000637-92-3	50





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Instrument : MSVOA_V
ClientSampleId : F9B17

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Quant Title : TRACE VOA SOM01.0

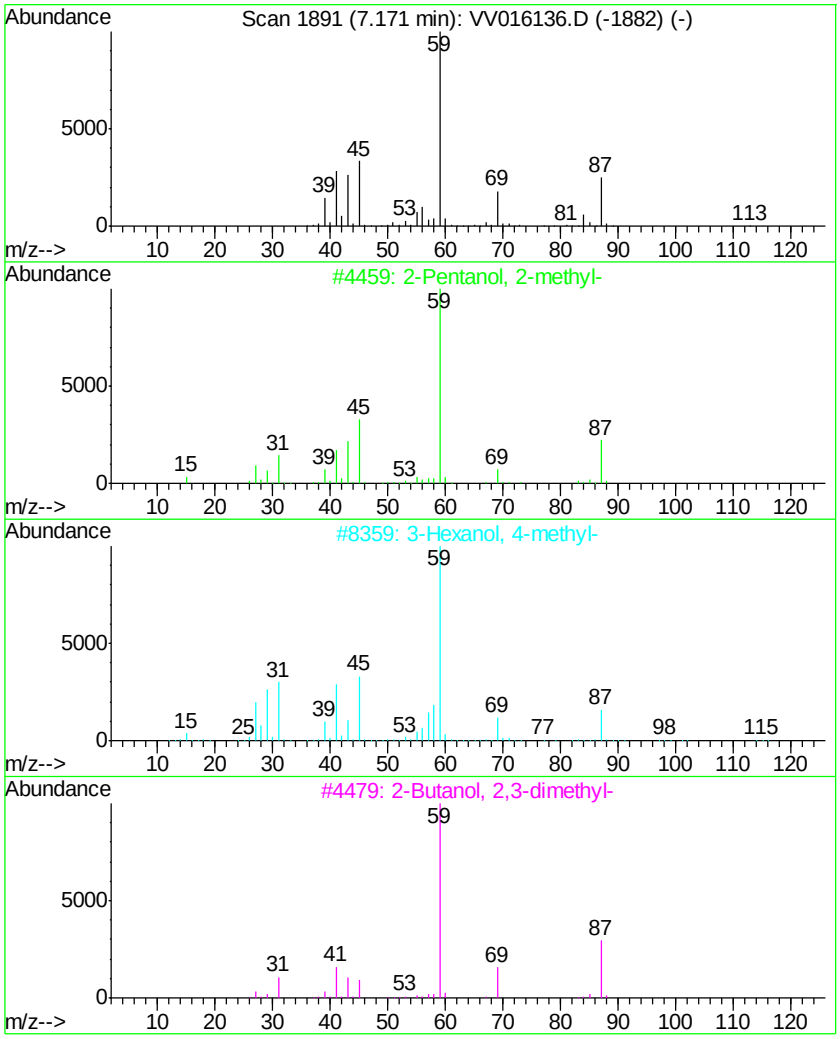
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

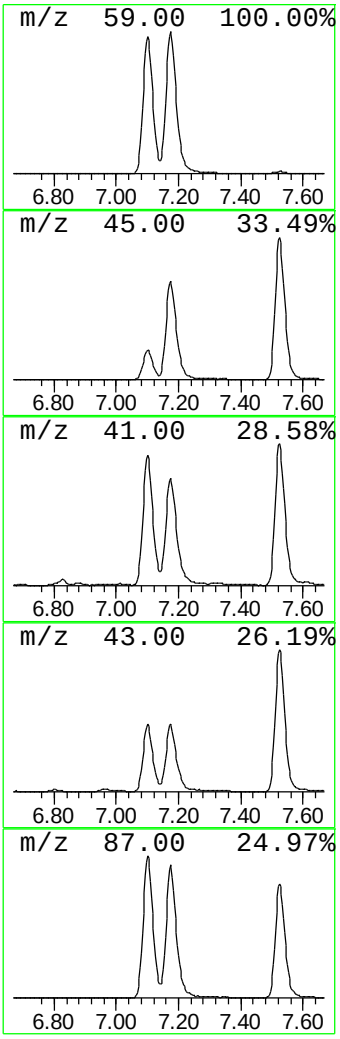
****

Peak Number 9 2-Pentanol, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.17	8.69 ug/L	792060	1,4-Difluorobenzene	5.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanol, 2-methyl-	102	C6H14O	000590-36-3	78
2	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	78
3	2-Butanol, 2,3-dimethyl-	102	C6H14O	000594-60-5	64
4	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	56
5	2-Butanol, 2,3-dimethyl-	102	C6H14O	000594-60-5	53





Library Search Compound Report

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ALS Vial : 15 Sample Multiplier: 1

Instrument : MSVOA_V
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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR050720WMA.M
Quant Title : TRACE VOA SOM01.0

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

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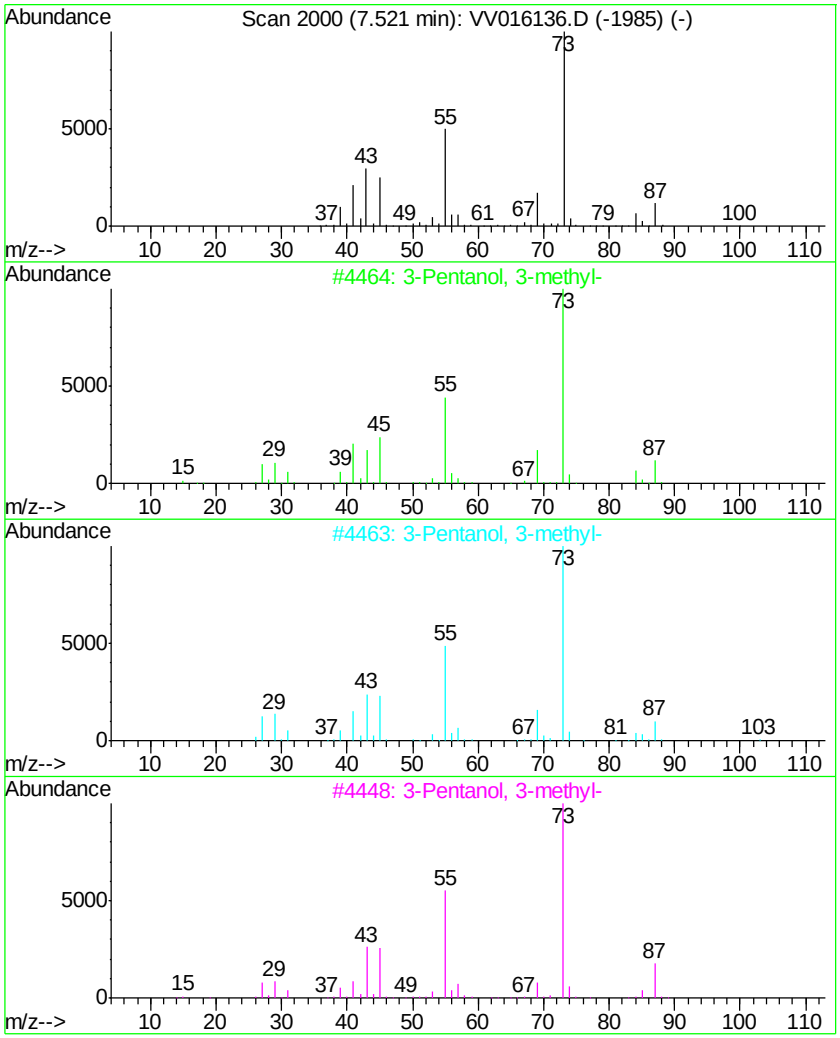
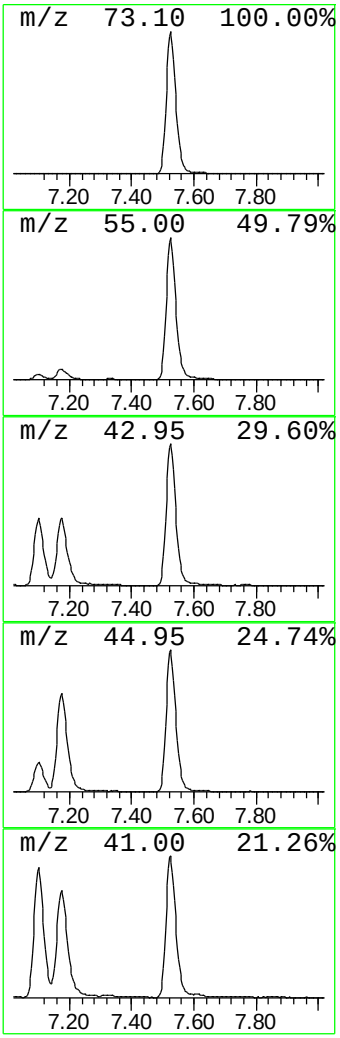
Peak Number 10 3-Pentanol, 3-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.52	14.14 ug/L	1375370	Chlorobenzene-d5	8.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	90
2	3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	64
3	3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	45
4	3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	43
5	4-Heptanol, 3-methyl-	130	C8H18O	001838-73-9	36





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV052020\
Data File : VV016136.D
Acq On : 20 May 2020 17:02
Operator : SY/MD
Sample : L2725-07
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : F9B17

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR050720WMA.M
Quant Title : TRACE VOA SOM01.0

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

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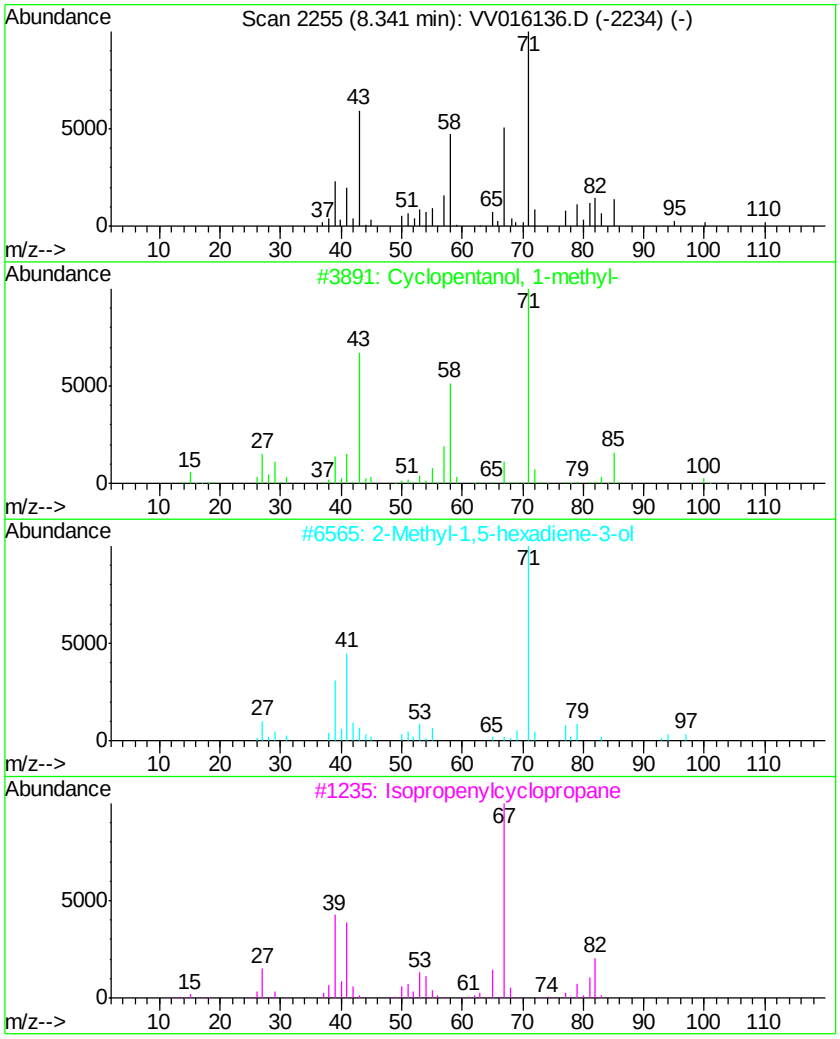
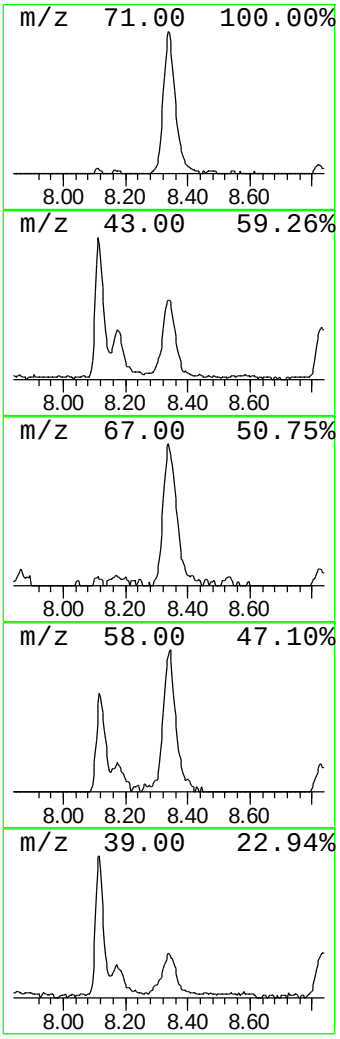
Peak Number 11 Cyclopentanol, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.34	1.07 ug/L	104411	Chlorobenzene-d5	8.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentanol, 1-methyl-	100	C6H12O	001462-03-9	52
2	2-Methyl-1,5-hexadiene-3-ol	112	C7H12O	017123-60-3	27
3	Isopropenylcyclopropane	82	C6H10	004663-22-3	25
4	1,4-Pentadiene, 3-methyl-	82	C6H10	001115-08-8	14
5	1,3-Pentadiene, 3-methyl-, (E)-	82	C6H10	002787-43-1	11





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV052020\
Data File : VV016136.D
Acq On : 20 May 2020 17:02
Operator : SY/MD
Sample : L2725-07
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : F9B17

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR050720WMA.M
Quant Title : TRACE VOA SOM01.0

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

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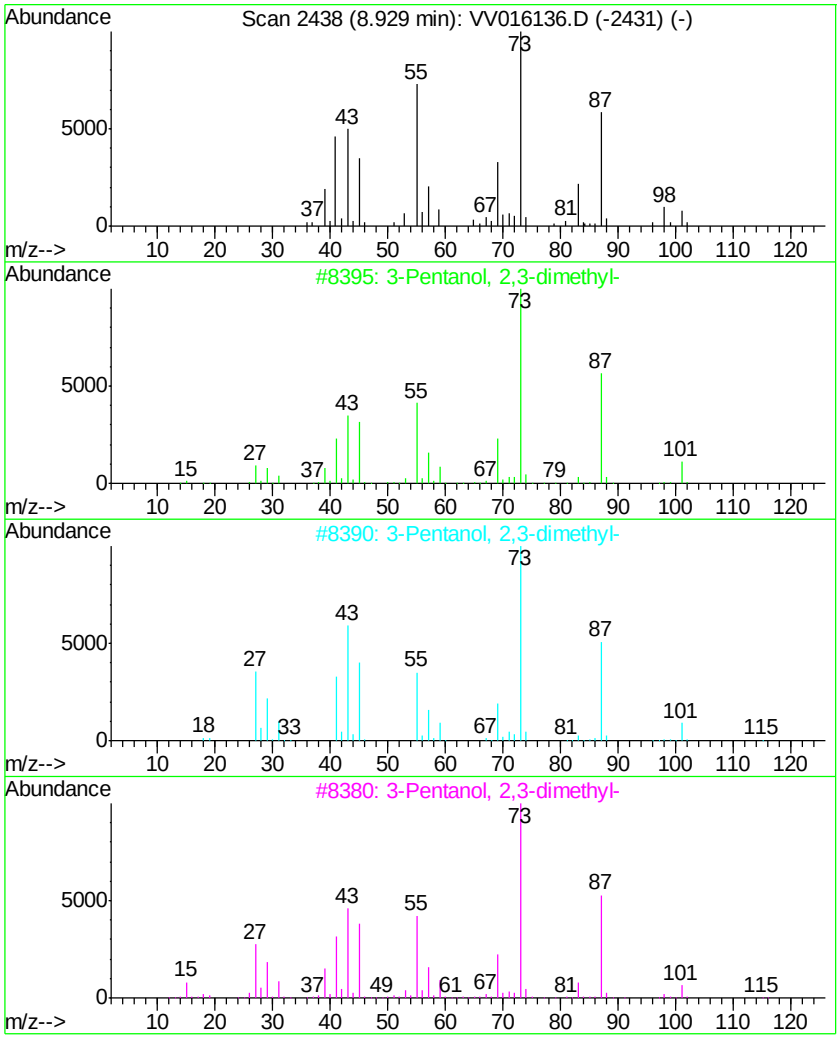
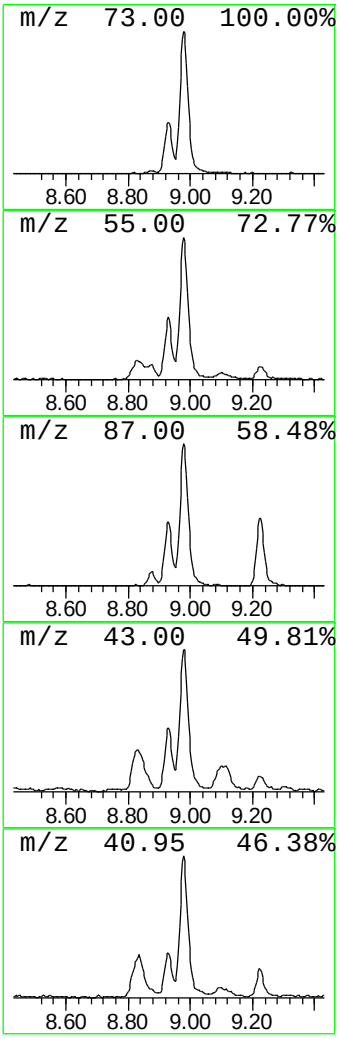
Peak Number 12 3-Pentanol, 2,3-dimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.93	0.76 ug/L	74137	Chlorobenzene-d5	8.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Pentanol, 2,3-dimethyl-	116	C7H16O	000595-41-5	64
2	3-Pentanol, 2,3-dimethyl-	116	C7H16O	000595-41-5	59
3	3-Pentanol, 2,3-dimethyl-	116	C7H16O	000595-41-5	59
4	3-Pentanol, 2,3-dimethyl-	116	C7H16O	000595-41-5	59
5	3-Hexanol, 2,4-dimethyl-	130	C8H18O	013432-25-2	40





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV052020\
Data File : VV016136.D
Acq On : 20 May 2020 17:02
Operator : SY/MD
Sample : L2725-07
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : F9B17

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR050720WMA.M
Quant Title : TRACE VOA SOM01.0

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

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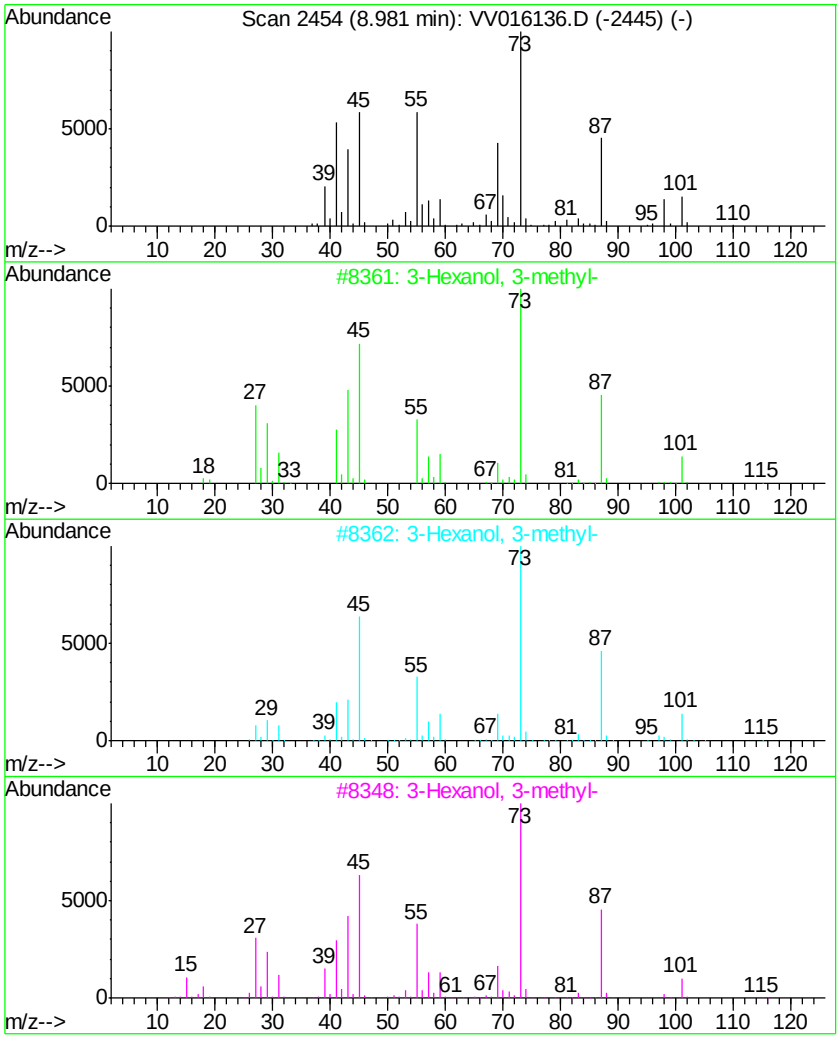
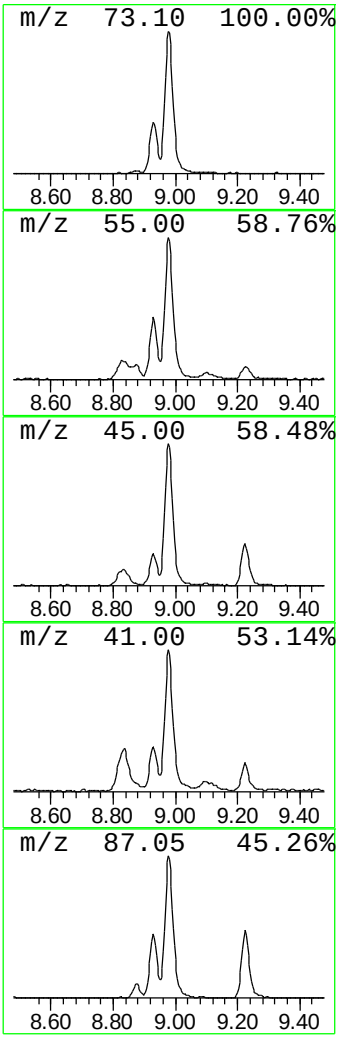
Peak Number 13 3-Hexanol, 3-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.98	2.64 ug/L	257147	Chlorobenzene-d5	8.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Hexanol, 3-methyl-	116	C7H16O	000597-96-6	72
2	3-Hexanol, 3-methyl-	116	C7H16O	000597-96-6	64
3	3-Hexanol, 3-methyl-	116	C7H16O	000597-96-6	64
4	3-Pentanol, 2,3-dimethyl-	116	C7H16O	000595-41-5	59
5	3-Pentanol, 2,3-dimethyl-	116	C7H16O	000595-41-5	50





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV052020\
Data File : VV016136.D
Acq On : 20 May 2020 17:02
Operator : SY/MD
Sample : L2725-07
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : F9B17

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR050720WMA.M
Quant Title : TRACE VOA SOM01.0

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
(DEL) Alkane: Cyc...	2.04	2.4	ug/L	215817	1	5.64	455491	5.0
(DEL) Alkane: Str...	3.30	0.6	ug/L	52893	1	5.64	455491	5.0
(DEL) Alkane: Str...	3.61	0.9	ug/L	79544	1	5.64	455491	5.0
Cyclopentene, 4-m...	4.49	2.9	ug/L	265048	1	5.64	455491	5.0
Amylene hydrate	5.26	24.3	ug/L	2210190	1	5.64	455491	5.0
Cyclobutane, (1-m...	6.83	1.0	ug/L	91874	1	5.64	455491	5.0
2-Butanol, 2,3-di...	7.10	7.5	ug/L	680401	1	5.64	455491	5.0
2-Pentanol, 2-met...	7.17	8.7	ug/L	792060	1	5.64	455491	5.0
3-Pentanol, 3-met...	7.52	14.1	ug/L	1375370	2	8.87	486367	5.0
Cyclopentanol, 1-...	8.34	1.1	ug/L	104411	2	8.87	486367	5.0
3-Pentanol, 2,3-d...	8.93	0.8	ug/L	74137	2	8.87	486367	5.0
3-Hexanol, 3-methyl-	8.98	2.6	ug/L	257147	2	8.87	486367	5.0