LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV111122\
Data File : VV028910.D
Acq On : 12 Nov 2022 05:51
Operator : SY/MD
Sample : N5590-09
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GBMY0

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\SFAMVTR111122WMA.M
Title : TRACE VOA SFAM1.0

Signal : TIC: VV028910.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.214	47	54	72	rVB	1612395	1386922	1.95%	1.024%
2	1.307	72	83	93	rBV	7484655	7339805	10.34%	5.418%
3	1.632	174	184	212	rVB	11776753	14481377	20.39%	10.690%
4	1.806	228	238	255	rVB2	1224074	1666717	2.35%	1.230%
5	2.024	297	306	322	rVB	882740	1236484	1.74%	0.913%
6	2.497	423	453	458	rBV3	1644158	3650916	5.14%	2.695%
7	2.574	469	477	501	rVB	1846450	3493972	4.92%	2.579%
8	2.757	518	534	560	rBV	1284990	2527693	3.56%	1.866%
9	3.757	829	845	866	rVB	1812753	4452288	6.27%	3.287%
10	3.902	875	890	935	rVB2	557209	1584800	2.23%	1.170%
11	4.674	1111	1130	1145	rBV	1630477	4034146	5.68%	2.978%
12	4.760	1145	1157	1185	rVB	518244	1406498	1.98%	1.038%
13	5.098	1231	1262	1279	rBV2	30393753	71018271	100.00%	52.427%
14	5.227	1293	1302	1320	rVB5	672940	1919736	2.70%	1.417%
15	5.320	1320	1331	1370	rVB3	300271	805559	1.13%	0.595%
16	5.612	1411	1422	1433	rBV	341207	742136	1.04%	0.548%
17	6.648	1730	1744	1764	rVB	390896	820771	1.16%	0.606%
18	6.773	1766	1783	1800	rBV2	663162	1610993	2.27%	1.189%
19	7.310	1941	1950	1966	rVB	506052	861397	1.21%	0.636%
20	8.085	2181	2191	2207	rBV	428079	835441	1.18%	0.617%
21	8.847	2415	2428	2440	rBV	486910	788841	1.11%	0.582%
22	11.242	3162	3173	3191	rBV	514234	836803	1.18%	0.618%
23	11.519	3246	3259	3279	rBV	2903676	4880874	6.87%	3.603%
24	11.619	3279	3290	3316	rVB4	696655	1246928	1.76%	0.921%
25	12.107	3431	3442	3463	rBV	531958	876126	1.23%	0.647%
26	12.873	3659	3680	3694	rBV2	610740	956032	1.35%	0.706%

Sum of corrected areas: 135461526

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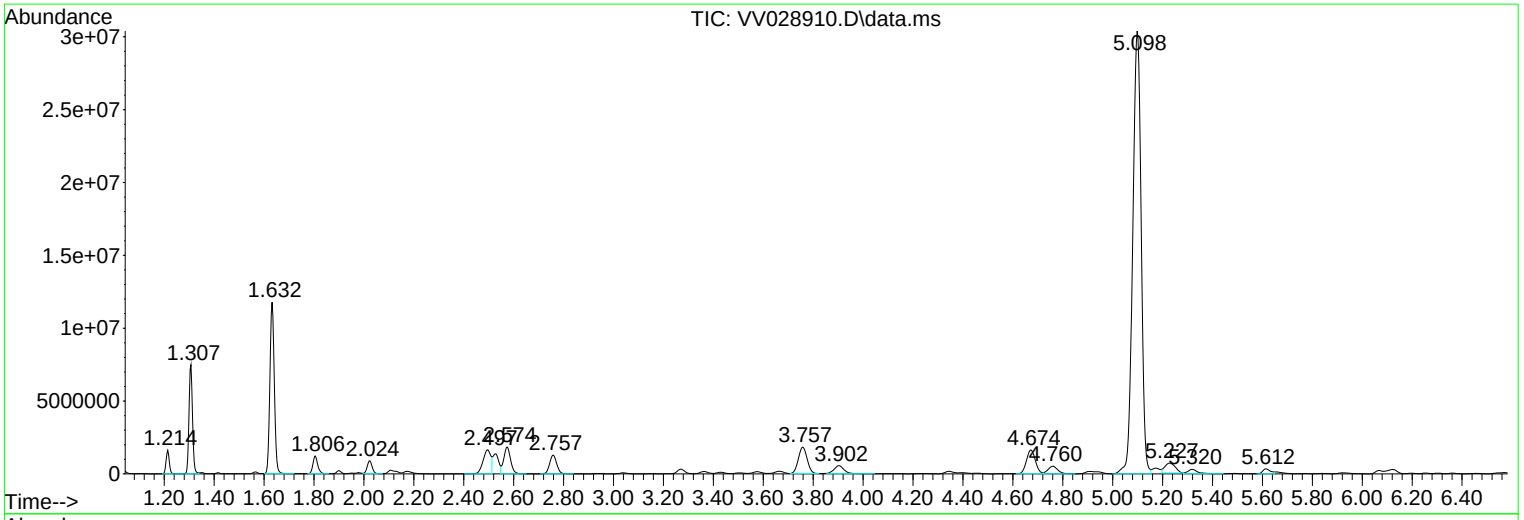
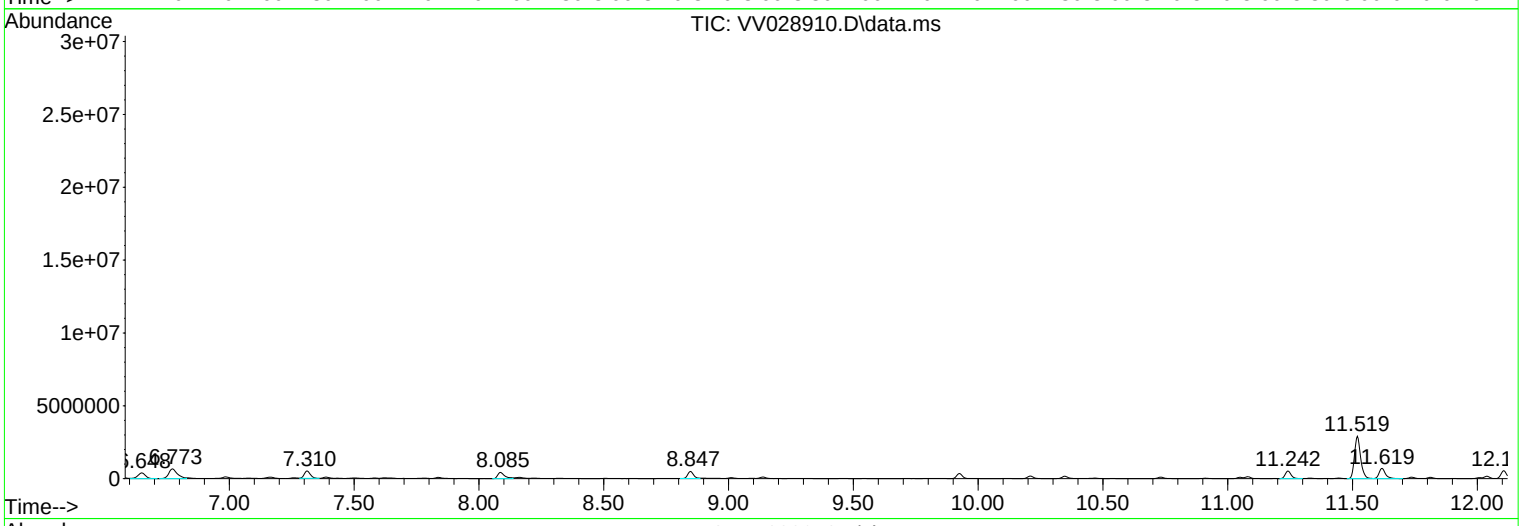
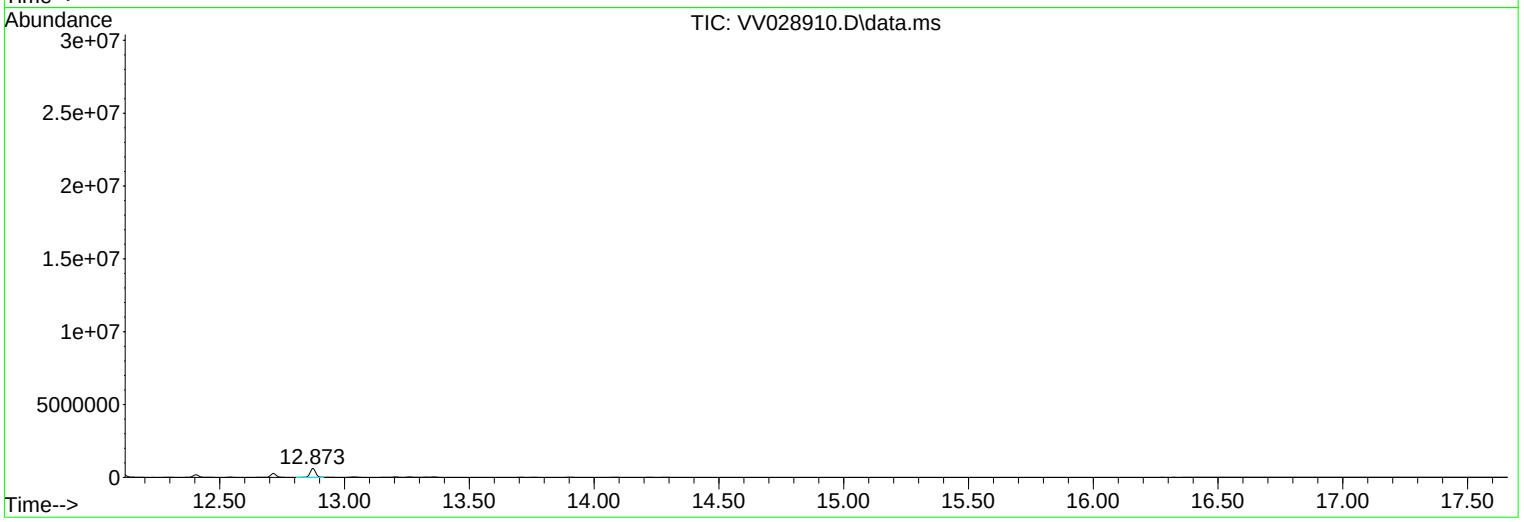
ClientSampleId :
GBMY0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR111122WMA.M
Quant Title : TRACE VOA SFAM1.0

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P







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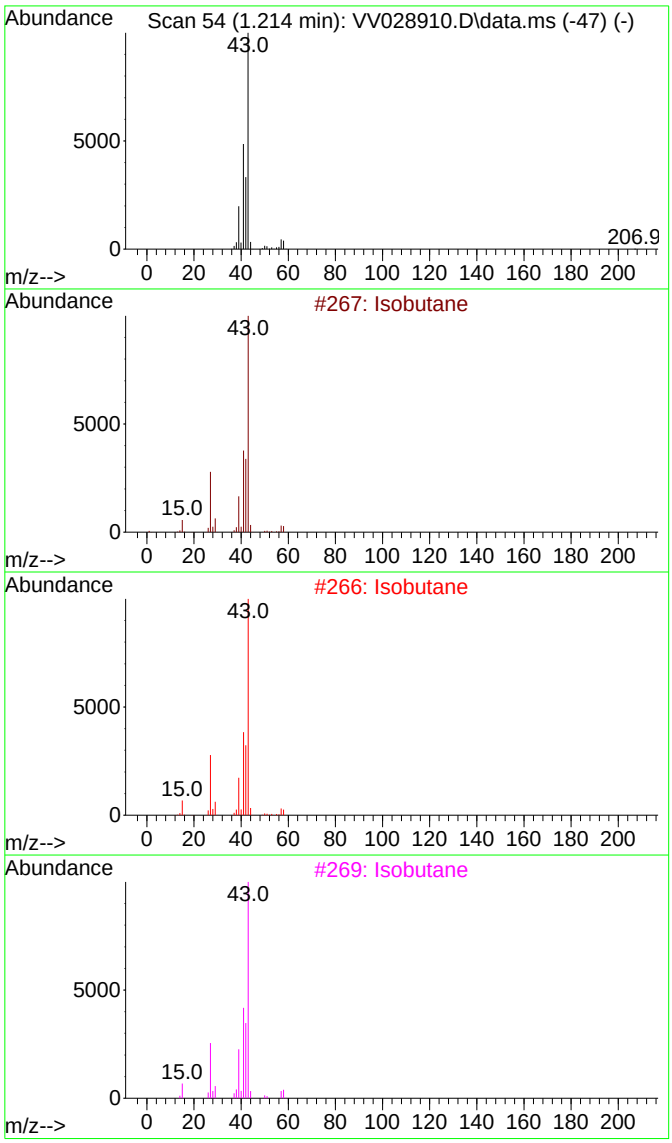
Quant Title : TRACE VOA SFAM1.0

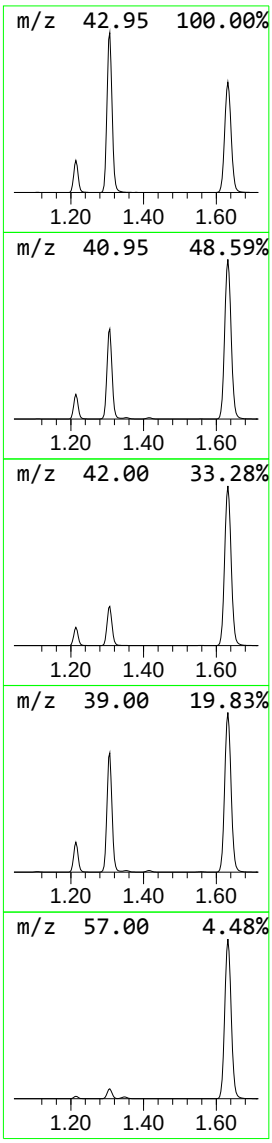
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.214	9.34 ug/L	1386920	1,4-Difluorobenzene	5.612

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Isobutane	58	C4H10	000075-28-5	80
2	Isobutane	58	C4H10	000075-28-5	72
3	Isobutane	58	C4H10	000075-28-5	72
4	Isobutane	58	C4H10	000075-28-5	64
5	Isobutane	58	C4H10	000075-28-5	59





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Operator : SY/MD
Sample : N5590-09
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ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GBMY0

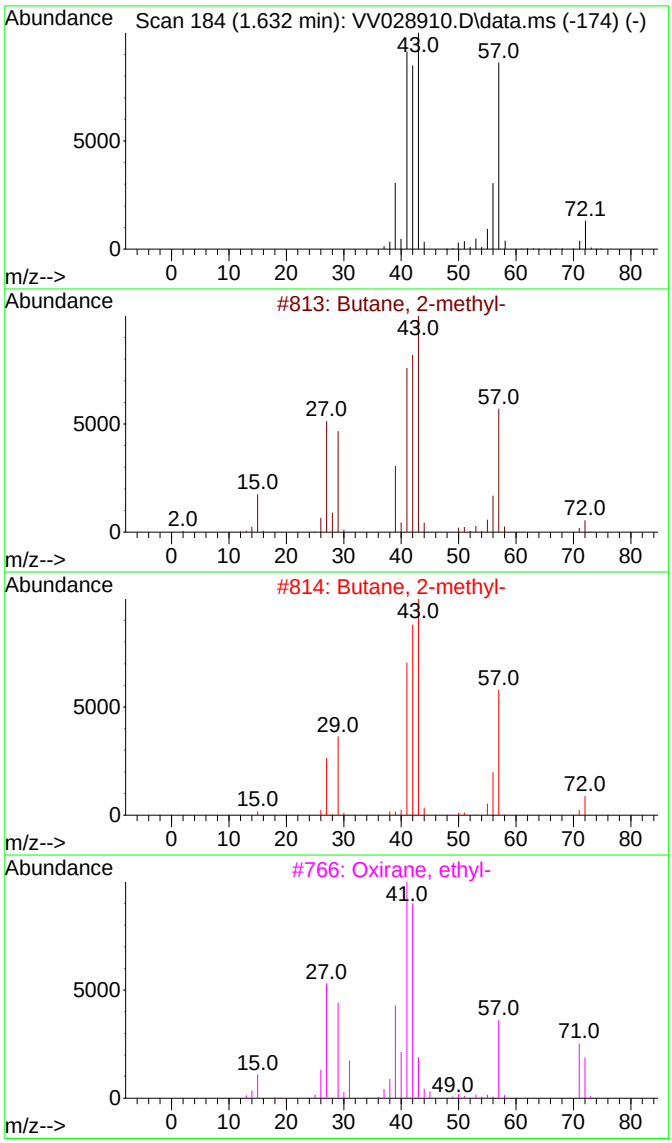
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR111122WMA.M
Quant Title : TRACE VOA SFAM1.0

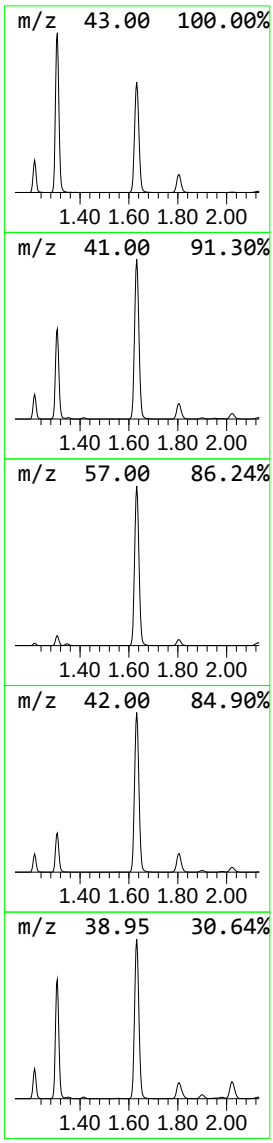
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.632	97.57 ug/L	14481400	1,4-Difluorobenzene	5.612

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-	72	C5H12	000078-78-4	83
2	Butane, 2-methyl-	72	C5H12	000078-78-4	80
3	Oxirane, ethyl-	72	C4H8O	000106-88-7	47
4	Pentane	72	C5H12	000109-66-0	42
5	Pentane	72	C5H12	000109-66-0	40





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Operator : SY/MD
Sample : N5590-09
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GBMY0

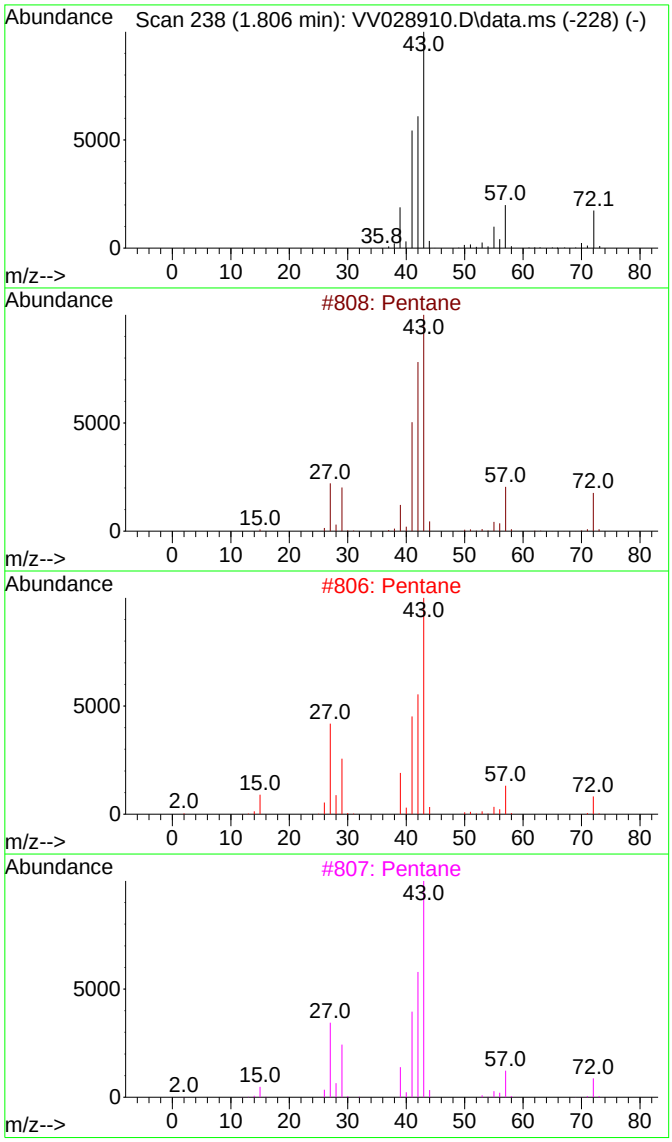
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR111122WMA.M
Quant Title : TRACE VOA SFAM1.0

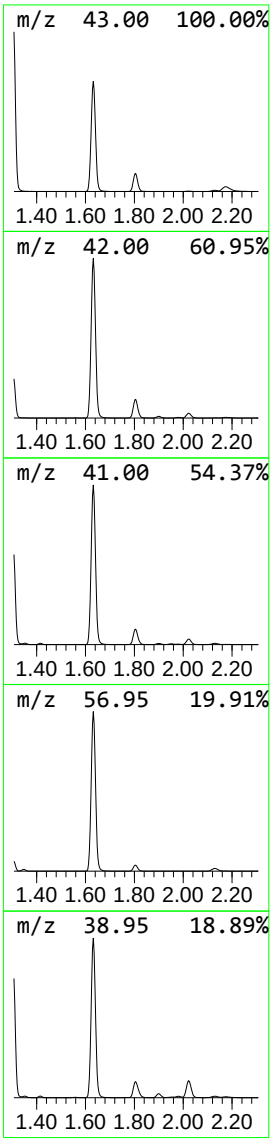
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.806	11.23 ug/L	1666720	1,4-Difluorobenzene	5.612

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane	72	C5H12	000109-66-0	86
2	Pentane	72	C5H12	000109-66-0	86
3	Pentane	72	C5H12	000109-66-0	80
4	Pentane	72	C5H12	000109-66-0	72
5	Pentane	72	C5H12	000109-66-0	56





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Operator : SY/MD
Sample : N5590-09
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ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GBMY0

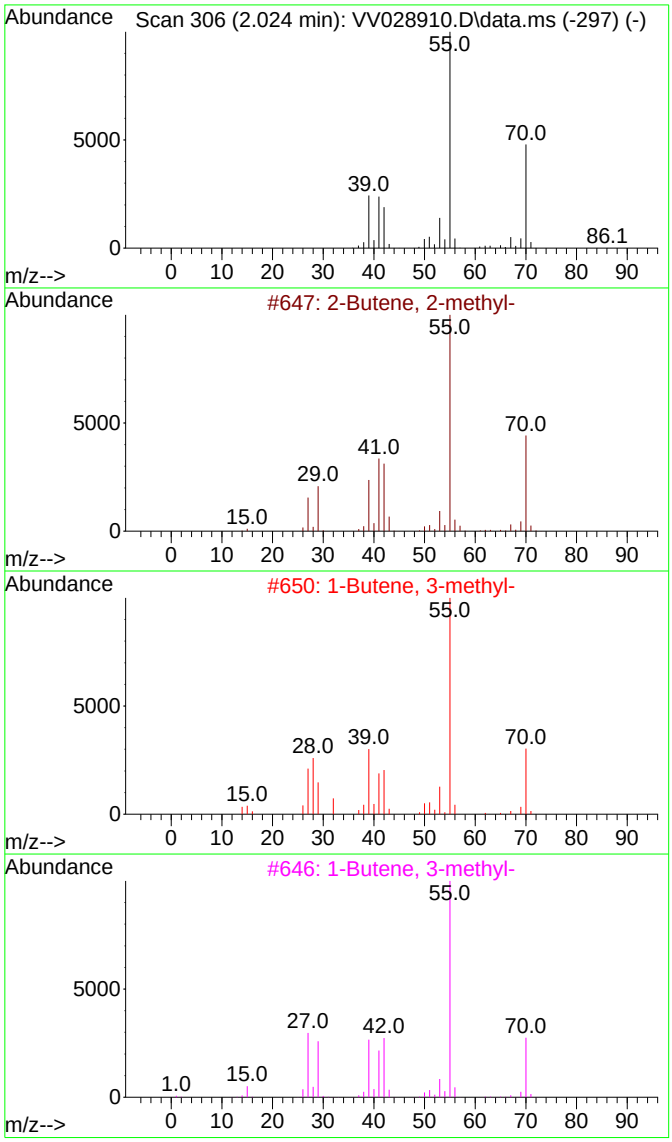
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR111122WMA.M
Quant Title : TRACE VOA SFAM1.0

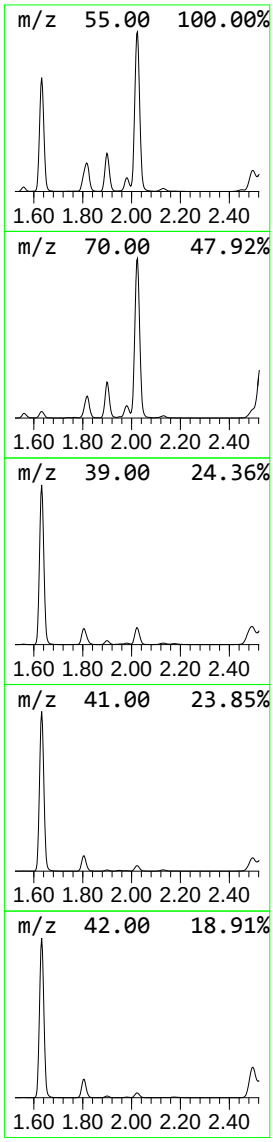
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.024	8.33 ug/L	1236480	1,4-Difluorobenzene	5.612

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butene, 2-methyl-	70	C5H10	000513-35-9	90
2	1-Butene, 3-methyl-	70	C5H10	000563-45-1	86
3	1-Butene, 3-methyl-	70	C5H10	000563-45-1	86
4	Cyclopropane, 1,2-dimethyl-, cis-	70	C5H10	000930-18-7	83
5	2-Butene, 2-methyl-	70	C5H10	000513-35-9	80





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Sample : N5590-09
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ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GBMY0

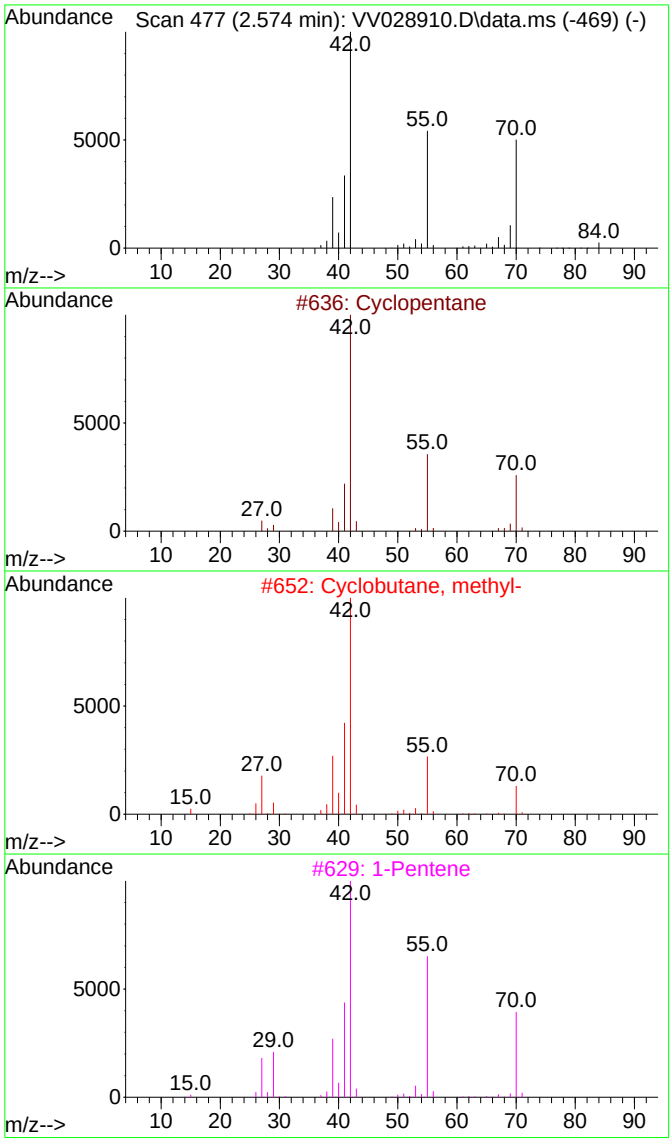
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR111122WMA.M
Quant Title : TRACE VOA SFAM1.0

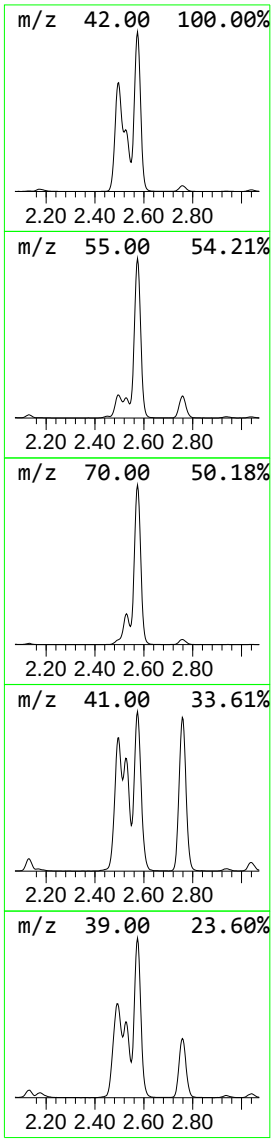
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 (DEL) Alkane: Cyclic2.574 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.574	23.54 ug/L	3493970	1,4-Difluorobenzene	5.612

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane	70	C5H10	000287-92-3	86
2	Cyclobutane, methyl-	70	C5H10	000598-61-8	80
3	1-Pentene	70	C5H10	000109-67-1	78
4	Cyclobutane, methyl-	70	C5H10	000598-61-8	72
5	1-Pentene	70	C5H10	000109-67-1	50





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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR111122WMA.M
Quant Title : TRACE VOA SFAM1.0

TIC Library : C:\Database\NIST20.L
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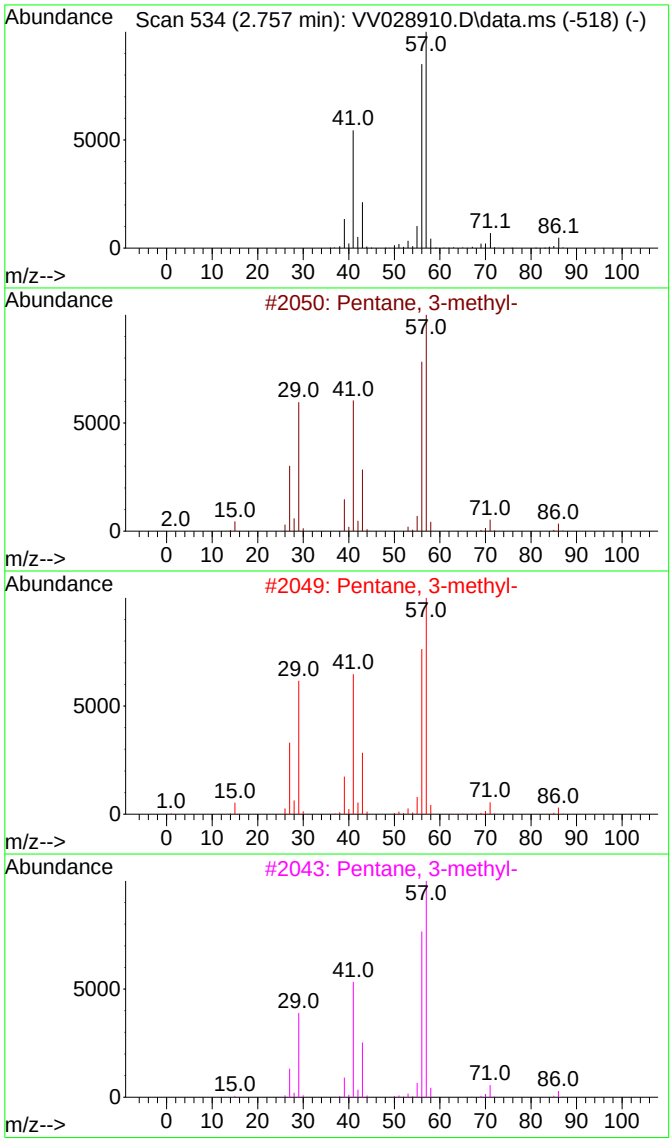
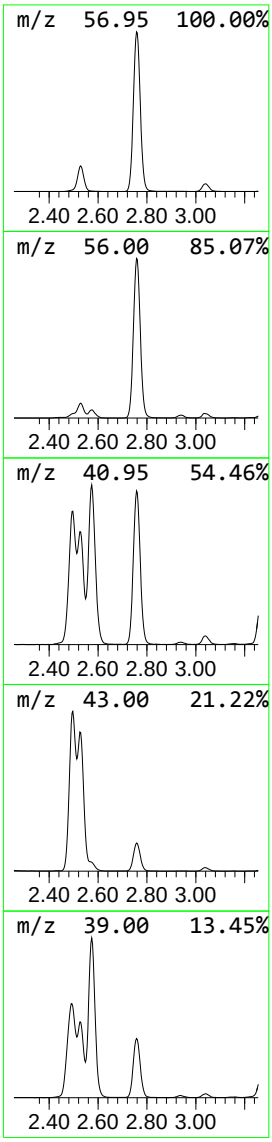
****
Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.757	17.03 ug/L	2527690	1,4-Difluorobenzene	5.612

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 3-methyl-	86	C6H14	000096-14-0	91
2	Pentane, 3-methyl-	86	C6H14	000096-14-0	91
3	Pentane, 3-methyl-	86	C6H14	000096-14-0	91
4	Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	78
5	Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	64





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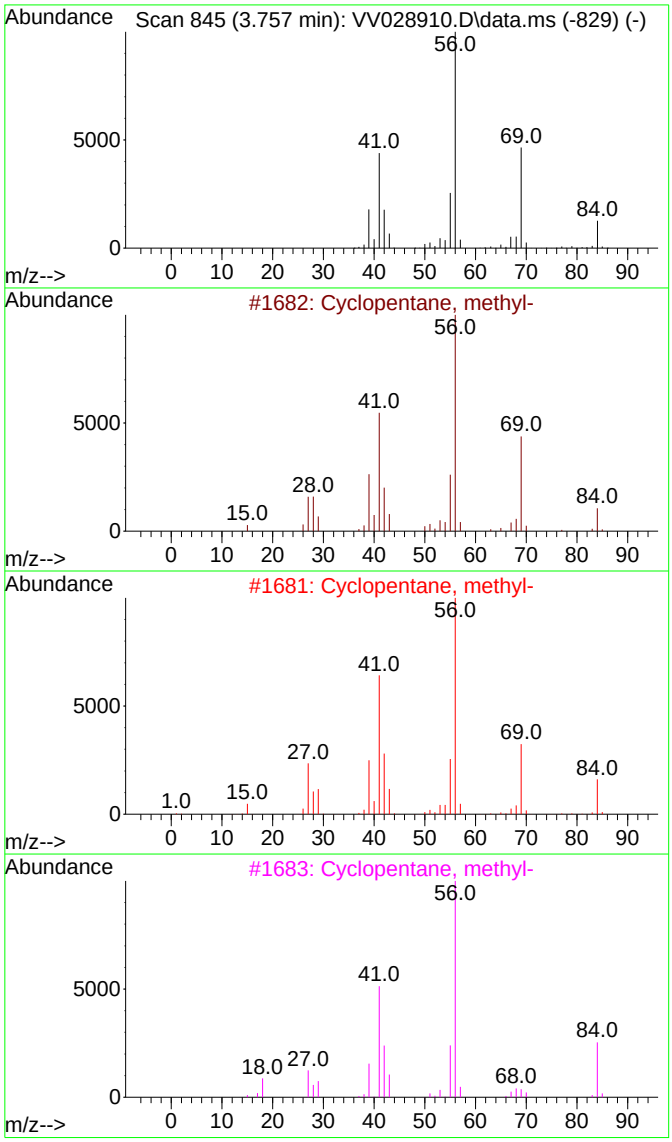
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR111122WMA.M
Quant Title : TRACE VOA SFAM1.0

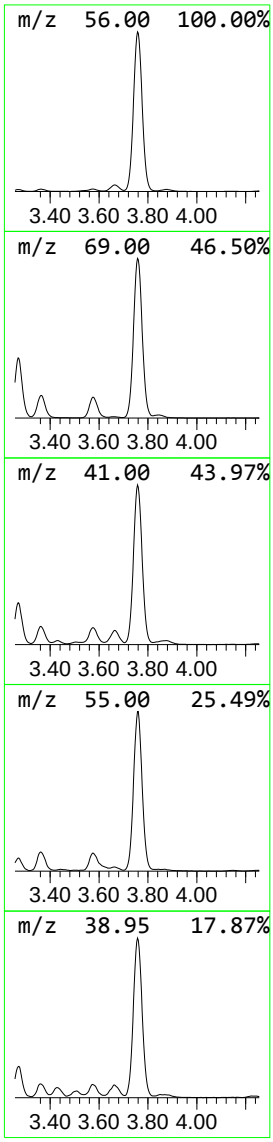
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 (DEL) Alkane: Cyclic3.757 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.757	30.00 ug/L	4452290	1,4-Difluorobenzene	5.612

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	94
2	Cyclopentane, methyl-	84	C6H12	000096-37-7	90
3	Cyclopentane, methyl-	84	C6H12	000096-37-7	86
4	Cyclobutane, ethyl-	84	C6H12	004806-61-5	78
5	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	72





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TIC Library : C:\Database\NIST20.L
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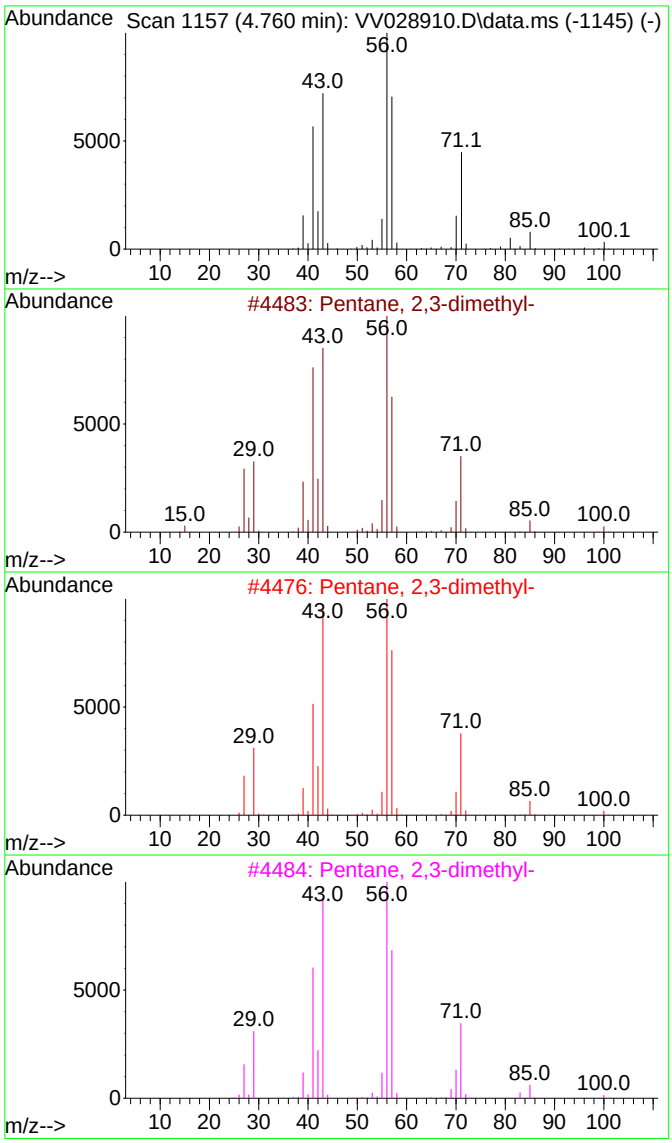
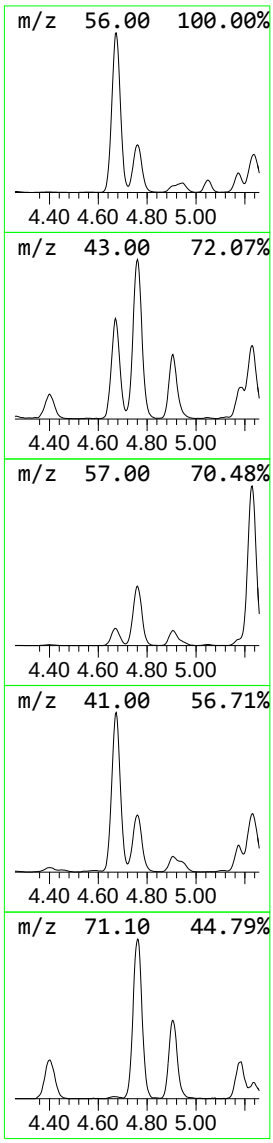
****
Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.760	9.48 ug/L	1406500	1,4-Difluorobenzene	5.612

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	91
2	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	91
3	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	91
4	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	64
5	Heptane	100	C7H16	000142-82-5	43





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ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GBMY0

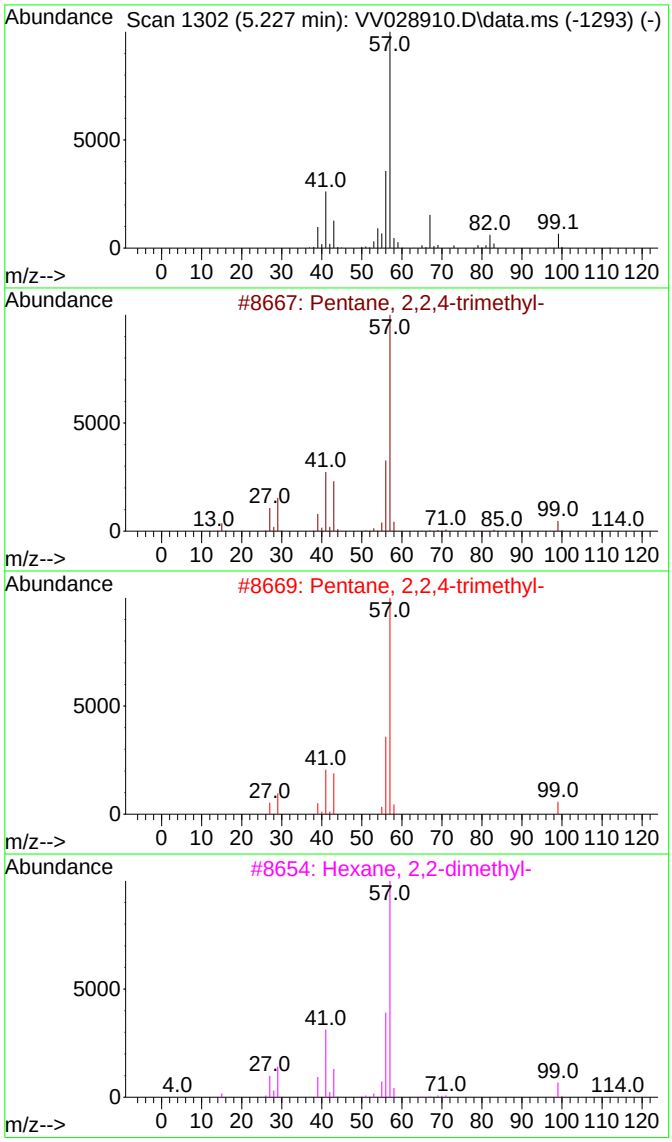
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR111122WMA.M
Quant Title : TRACE VOA SFAM1.0

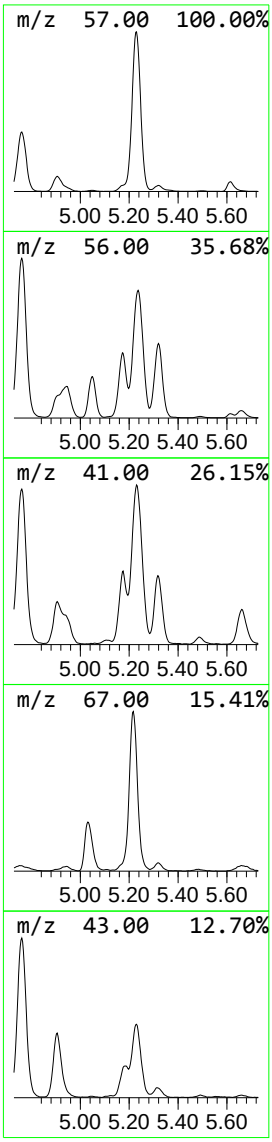
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.227	12.93 ug/L	1919740	1,4-Difluorobenzene	5.612

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	50
2	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	50
3	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	50
4	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	50
5	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	50





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV111122\
Data File : VV028910.D
Acq On : 12 Nov 2022 05:51
Operator : SY/MD
Sample : N5590-09
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GBMY0

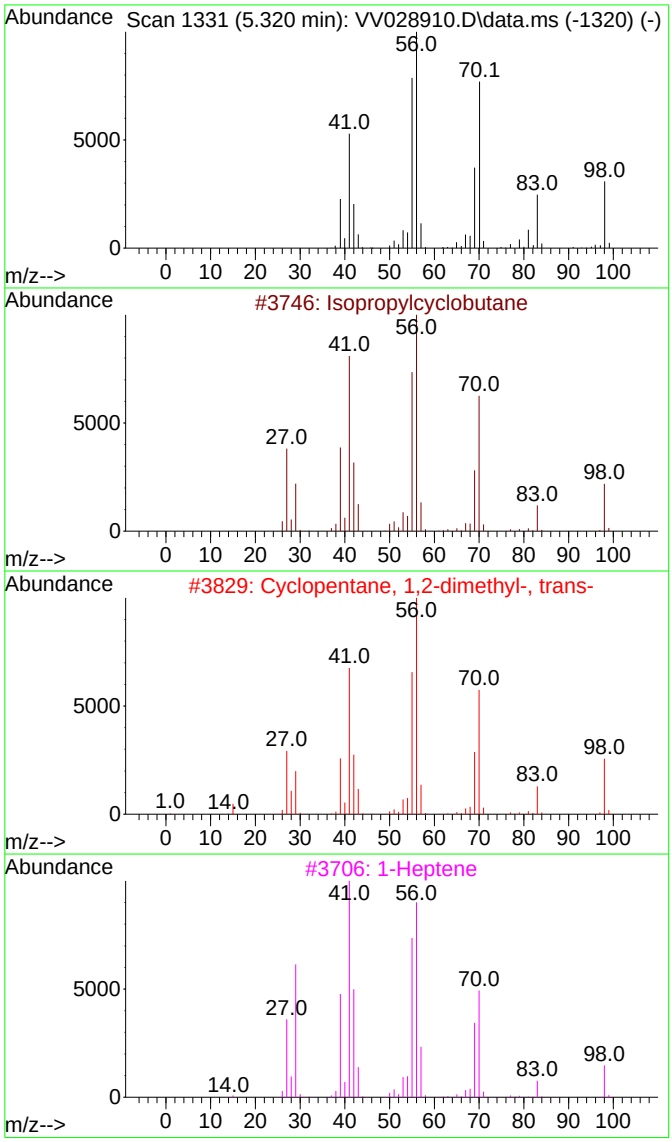
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR111122WMA.M
Quant Title : TRACE VOA SFAM1.0

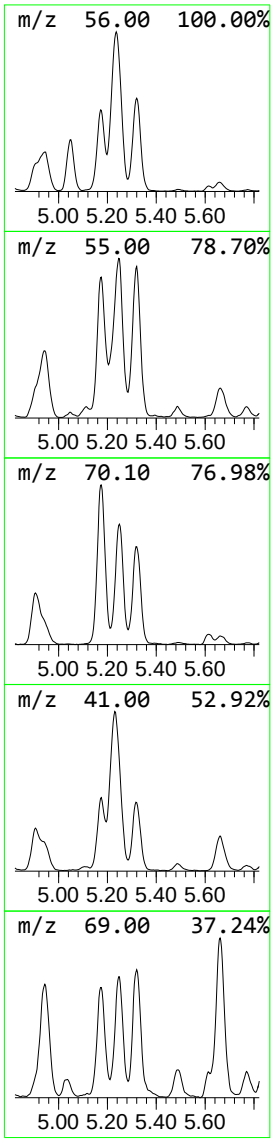
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 (DEL) Alkane: Cyclic5.320 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.320	5.43 ug/L	805559	1,4-Difluorobenzene	5.612

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Isopropylcyclobutane	98	C7H14	000872-56-0	91
2	Cyclopentane, 1,2-dimethyl-, trans-	98	C7H14	000822-50-4	91
3	1-Heptene	98	C7H14	000592-76-7	90
4	Cyclopentane, 1,2-dimethyl-	98	C7H14	002452-99-5	80
5	Cyclobutanone, 3-ethyl-	98	C6H10O	056335-73-0	59





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Data File : VV028910.D
Acq On : 12 Nov 2022 05:51
Operator : SY/MD
Sample : N5590-09
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GBMY0

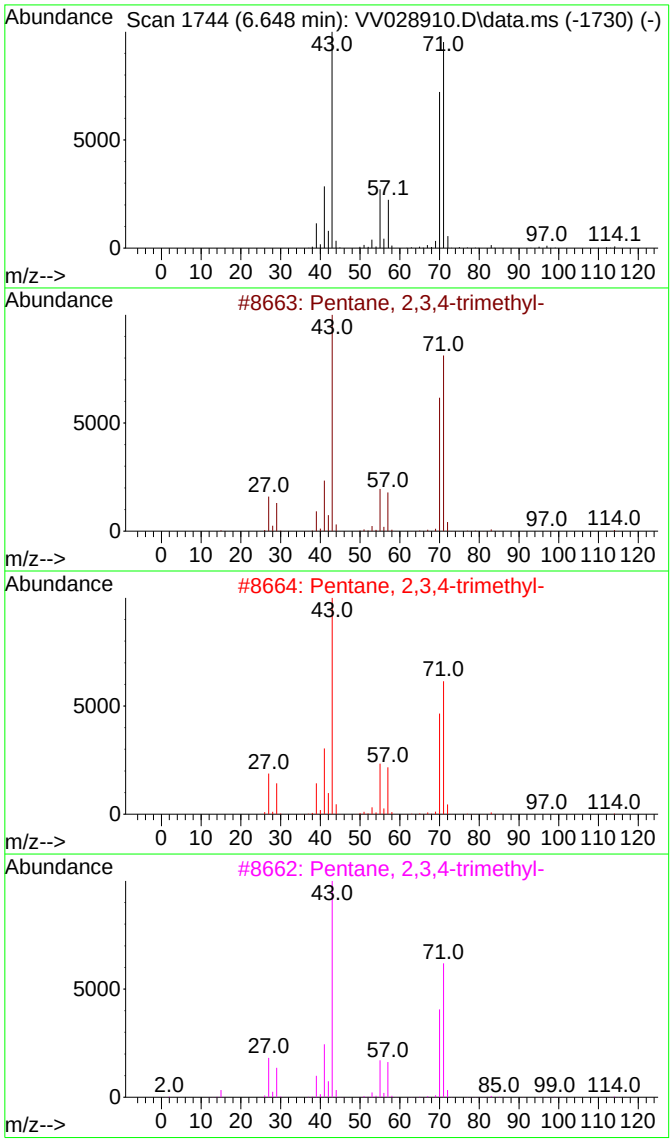
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR111122WMA.M
Quant Title : TRACE VOA SFAM1.0

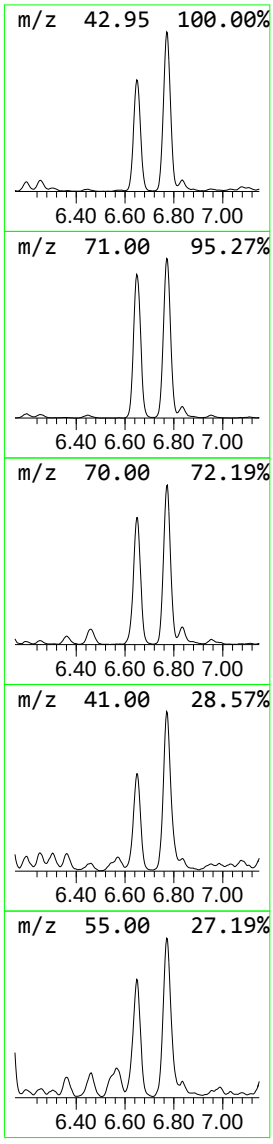
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.648	5.53 ug/L	820771	1,4-Difluorobenzene	5.612

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	91
2	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	83
3	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	80
4	Heptane, 4-methyl-	114	C8H18	000589-53-7	74
5	Pentane, 3-ethyl-	100	C7H16	000617-78-7	74





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Data File : VV028910.D
Acq On : 12 Nov 2022 05:51
Operator : SY/MD
Sample : N5590-09
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GBMY0

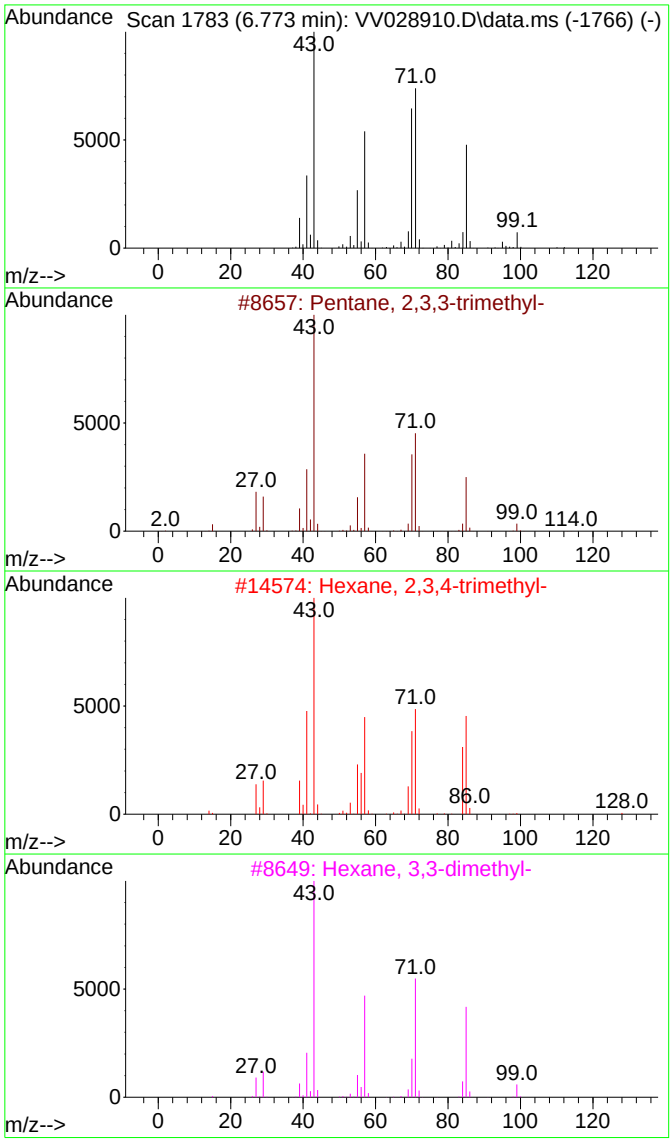
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR111122WMA.M
Quant Title : TRACE VOA SFAM1.0

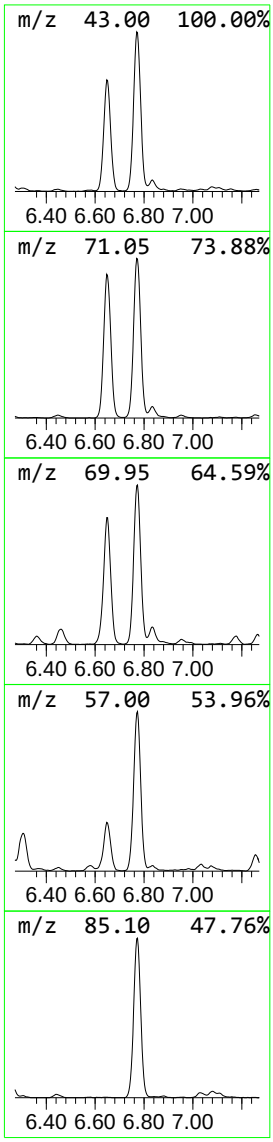
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.773	10.85 ug/L	1610990	1,4-Difluorobenzene	5.612

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	90
2	Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	78
3	Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	64
4	Hexane, 3-methyl-	100	C7H16	000589-34-4	59
5	Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	59





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Data File : VV028910.D
Acq On : 12 Nov 2022 05:51
Operator : SY/MD
Sample : N5590-09
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GBMY0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR111122WMA.M
Quant Title : TRACE VOA SFAM1.0

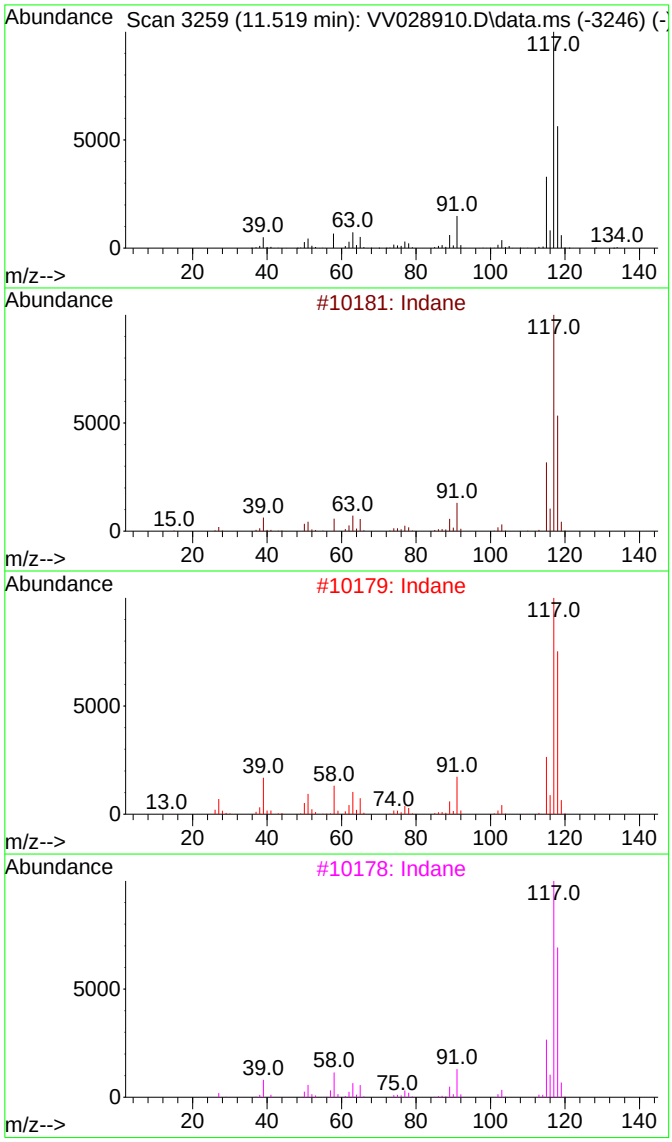
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

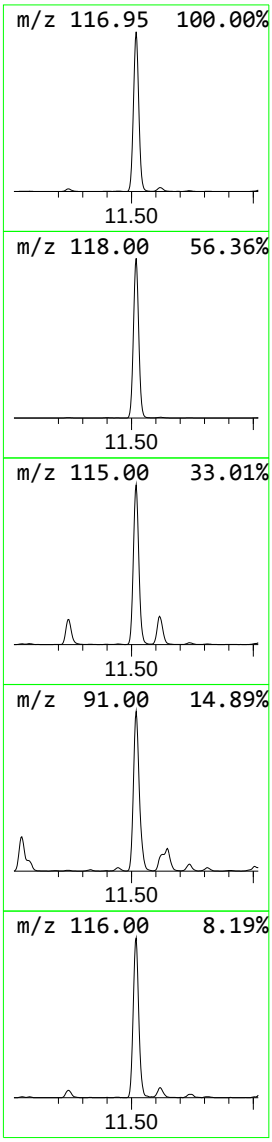
****

Peak Number 13 Indane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.519	29.16 ug/L	4880870	1,4-Dichlorobenzene-d4	11.243

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	95
2	Indane	118	C9H10	000496-11-7	91
3	Indane	118	C9H10	000496-11-7	87
4	Benzene, cyclopropyl-	118	C9H10	000873-49-4	83
5	Benzene, 2-propenyl-	118	C9H10	000300-57-2	83





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Data File : VV028910.D
Acq On : 12 Nov 2022 05:51
Operator : SY/MD
Sample : N5590-09
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GBMY0

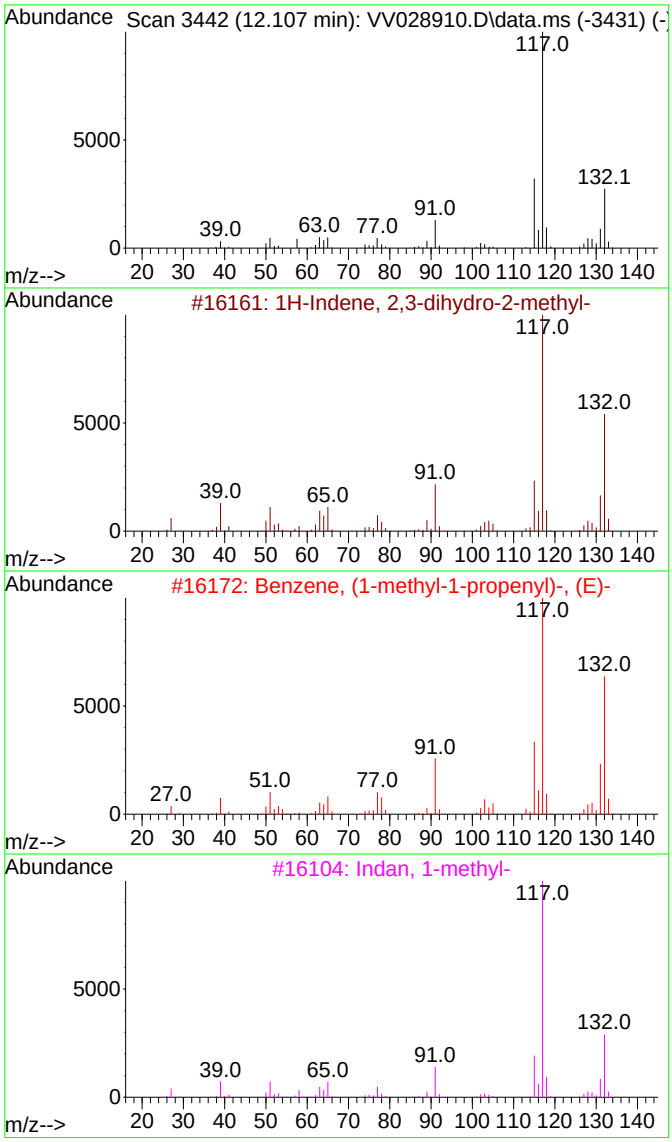
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR111122WMA.M
Quant Title : TRACE VOA SFAM1.0

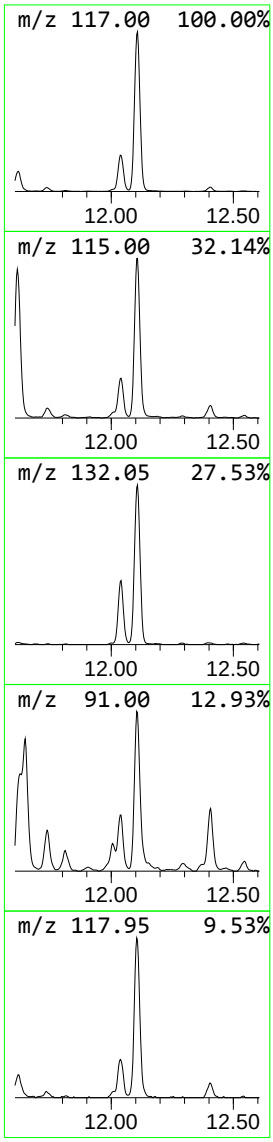
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 1H-Indene, 2,3-dihydro-2-me... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.107	5.23 ug/L	876126	1,4-Dichlorobenzene-d4	11.243

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	87
2	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	87
3	Indan, 1-methyl-	132	C10H12	000767-58-8	87
4	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	86





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MSVOA_V
ClientSampleId :
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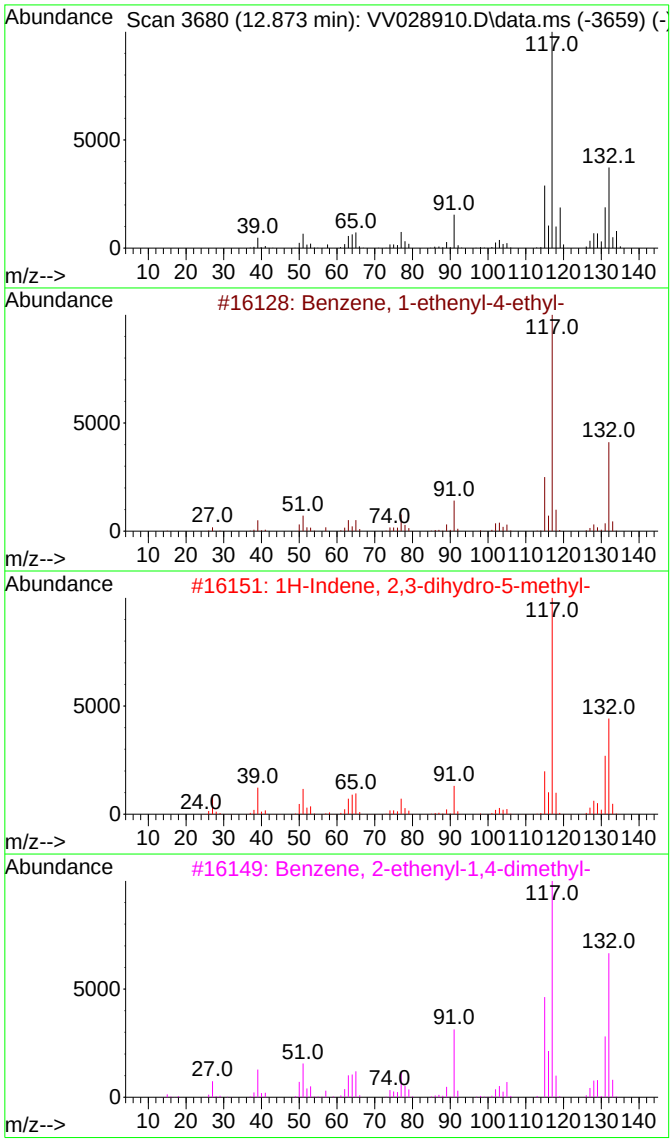
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR111122WMA.M
Quant Title : TRACE VOA SFAM1.0

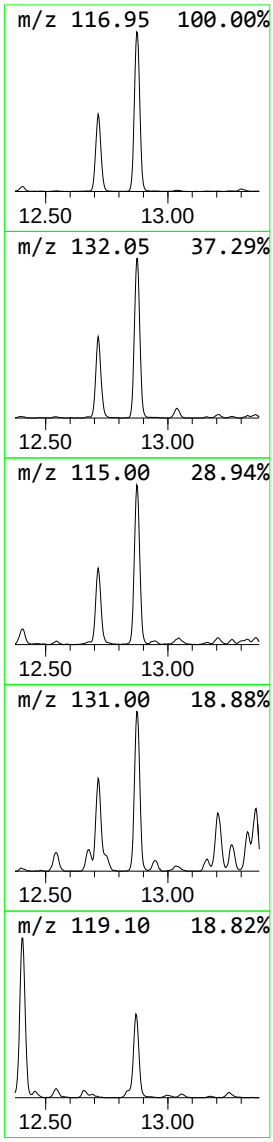
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.873	5.71 ug/L	956032	1,4-Dichlorobenzene-d4	11.243

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	91
2	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	91
3	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	91
4	3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
5	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	90





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV111122\
Data File : VV028910.D
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Operator : SY/MD
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ALS Vial : 44 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
GBMY0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR111122WMA.M
Quant Title : TRACE VOA SFAM1.0

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
(DEL) Alkane: S...	1.214	9.3	ug/L	1386920	1	5.612	742136	5.0
(DEL) Alkane: S...	1.632	97.6	ug/L	14481400	1	5.612	742136	5.0
(DEL) Alkane: S...	1.806	11.2	ug/L	1666720	1	5.612	742136	5.0
(DEL) Alkane: S...	2.024	8.3	ug/L	1236480	1	5.612	742136	5.0
(DEL) Alkane: C...	2.574	23.5	ug/L	3493970	1	5.612	742136	5.0
(DEL) Alkane: S...	2.757	17.0	ug/L	2527690	1	5.612	742136	5.0
(DEL) Alkane: C...	3.757	30.0	ug/L	4452290	1	5.612	742136	5.0
(DEL) Alkane: S...	4.760	9.5	ug/L	1406500	1	5.612	742136	5.0
(DEL) Alkane: S...	5.227	12.9	ug/L	1919740	1	5.612	742136	5.0
(DEL) Alkane: C...	5.320	5.4	ug/L	805559	1	5.612	742136	5.0
(DEL) Alkane: S...	6.648	5.5	ug/L	820771	1	5.612	742136	5.0
(DEL) Alkane: S...	6.773	10.9	ug/L	1610990	1	5.612	742136	5.0
Indane	11.519	29.2	ug/L	4880870	3	11.243	836803	5.0
1H-Indene, 2,3-...	12.107	5.2	ug/L	876126	3	11.243	836803	5.0
Benzene, 1-ethe...	12.873	5.7	ug/L	956032	3	11.243	836803	5.0