LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV102122\
Data File : VV028670.D
Acq On : 21 Oct 2022 18:43
Operator : SY/MD
Sample : N5216-16
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BHA75

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\SFAMVTR102022WMA.M
Title : TRACE VOA SFAM1.0

Signal : TIC: VV028670.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.947	271	282	298	rBV	6162372	8372169	11.80%	4.334%
2	3.291	677	700	729	rBV	7355992	17105643	24.11%	8.854%
3	3.735	822	838	857	rVB	540533	1310252	1.85%	0.678%
4	5.082	1229	1257	1281	rBV2	24665803	52595211	74.14%	27.225%
5	8.886	2412	2440	2470	rBV2	44732278	70940925	100.00%	36.721%
6	9.133	2508	2517	2534	rVB	456839	749687	1.06%	0.388%
7	9.538	2632	2643	2651	rVV	945363	1540058	2.17%	0.797%
8	9.924	2751	2763	2780	rBV	4026585	6137973	8.65%	3.177%
9	10.262	2859	2868	2881	rVB	1501863	2203689	3.11%	1.141%
10	10.349	2882	2895	2908	rVV2	2283412	3928296	5.54%	2.033%
11	10.416	2908	2916	2940	rVB	689366	1200443	1.69%	0.621%
12	10.731	3001	3014	3027	rBV	1548320	2324644	3.28%	1.203%
13	11.265	3164	3180	3191	rBV3	758014	1733310	2.44%	0.897%
14	11.345	3191	3205	3228	rVB2	1343249	3323925	4.69%	1.721%
15	11.519	3250	3259	3278	rVB	3593033	5956642	8.40%	3.083%
16	11.622	3279	3291	3317	rVB6	693751	1782651	2.51%	0.923%
17	12.008	3395	3411	3426	rBV	1510156	2454985	3.46%	1.271%
18	12.236	3471	3482	3493	rBV4	802822	1564584	2.21%	0.810%
19	12.403	3519	3534	3543	rBV	1095852	1699869	2.40%	0.880%
20	12.715	3622	3631	3643	rVB	478701	733826	1.03%	0.380%
21	12.869	3662	3679	3689	rBV2	1043863	1753870	2.47%	0.908%
22	13.371	3824	3835	3843	rBV4	420257	915479	1.29%	0.474%
23	13.419	3843	3850	3866	rVB	588495	1105988	1.56%	0.572%
24	14.490	4174	4183	4230	rVB2	614051	1755816	2.48%	0.909%

Sum of corrected areas: 193189935

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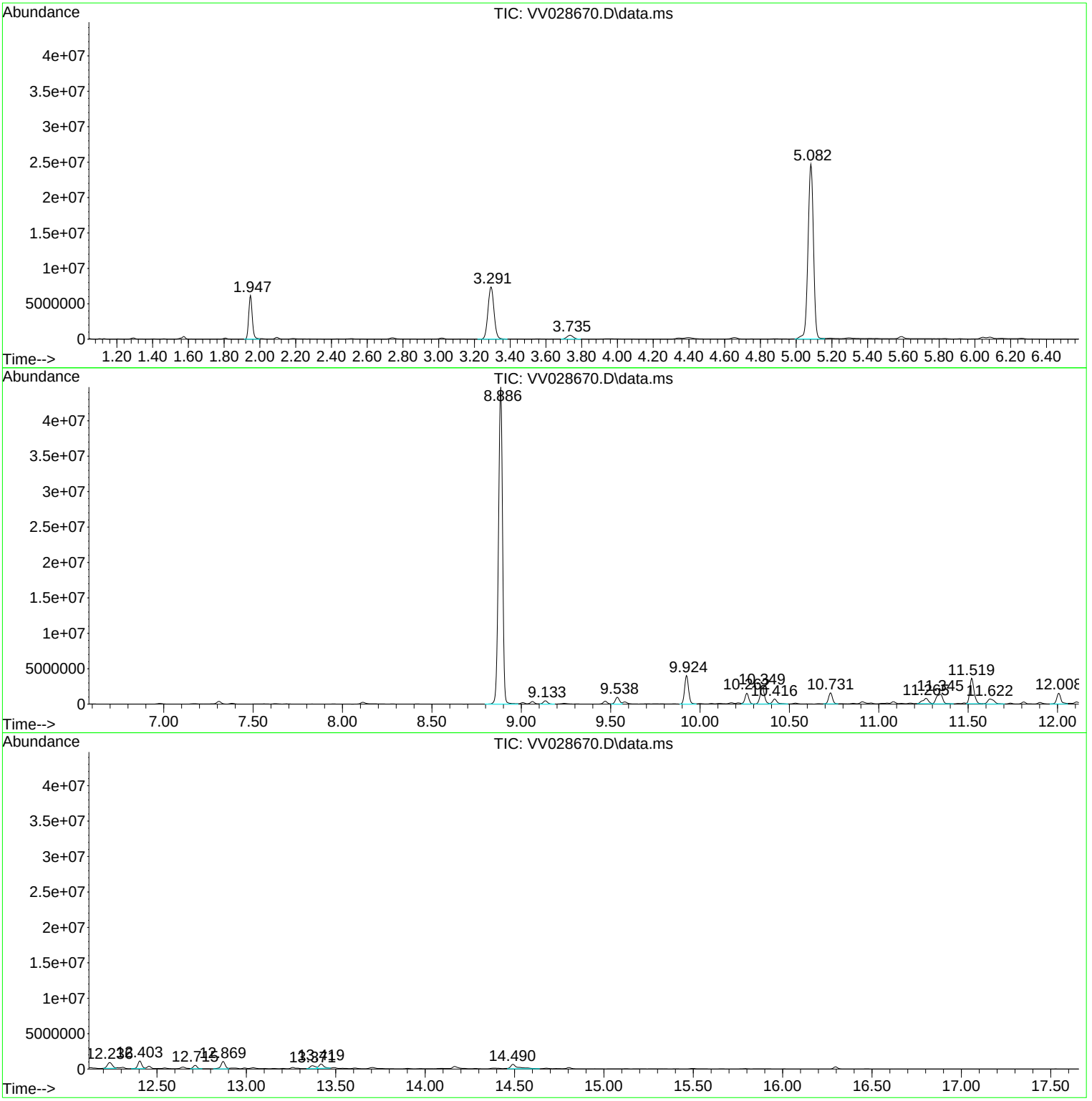
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BHA75

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR102022WMA.M
Quant Title : TRACE VOA SFAM1.0

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV102122\
Data File : VV028670.D
Acq On : 21 Oct 2022 18:43
Operator : SY/MD
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Instrument :
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ClientSampleId :
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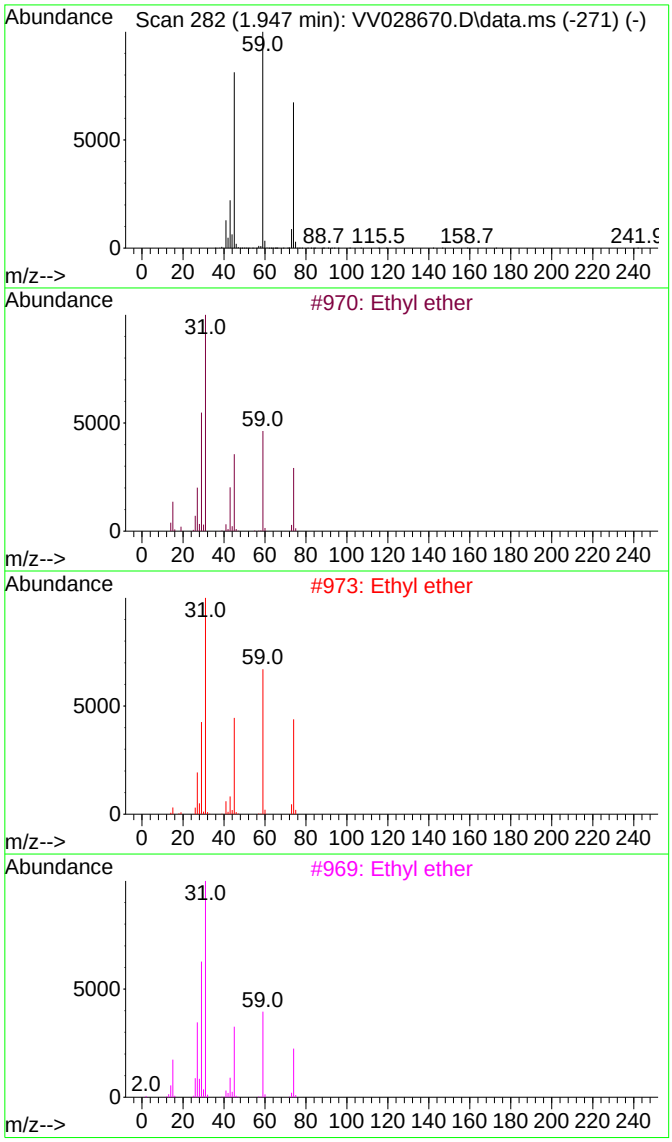
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR102022WMA.M
Quant Title : TRACE VOA SFAM1.0

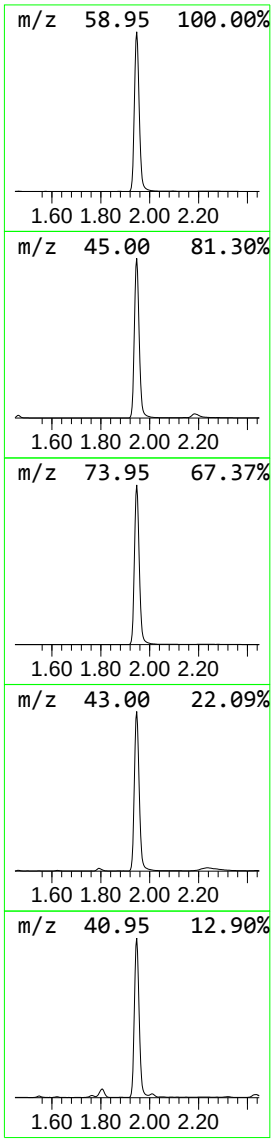
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Ethyl ether Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.947	0.80 ug/L	8372170	1,4-Difluorobenzene	5.590

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethyl ether	74	C4H10O	000060-29-7	90
2	Ethyl ether	74	C4H10O	000060-29-7	90
3	Ethyl ether	74	C4H10O	000060-29-7	90
4	Ethyl ether	74	C4H10O	000060-29-7	72
5	Ethyl ether	74	C4H10O	000060-29-7	59





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV102122\
Data File : VV028670.D
Acq On : 21 Oct 2022 18:43
Operator : SY/MD
Sample : N5216-16
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BHA75

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR102022WMA.M
Quant Title : TRACE VOA SFAM1.0

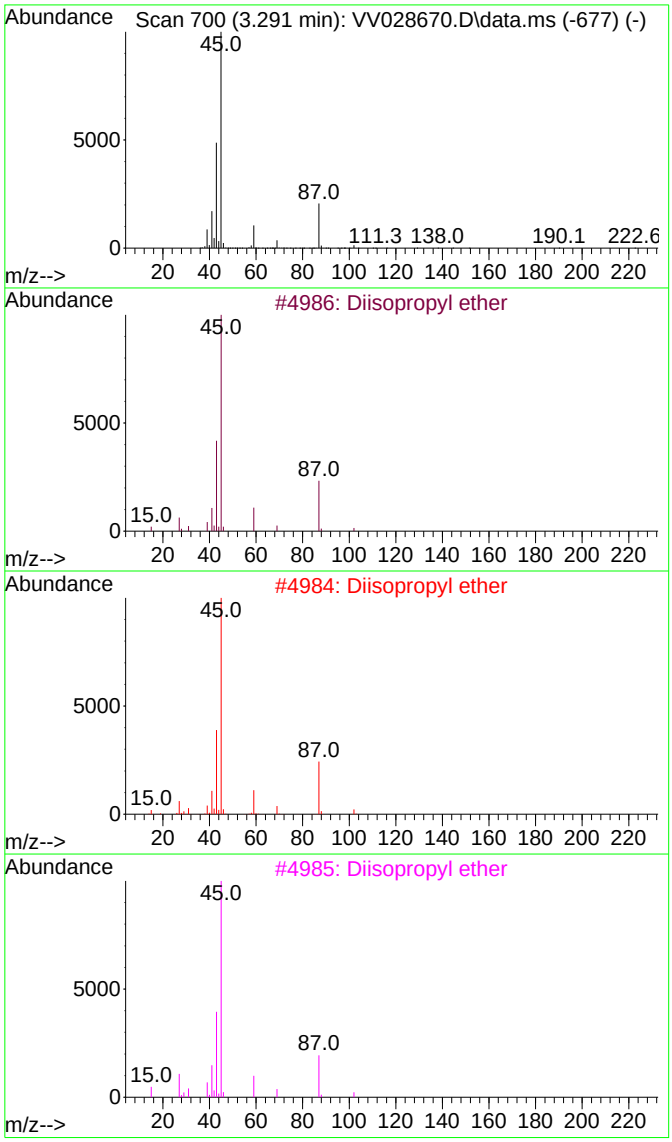
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

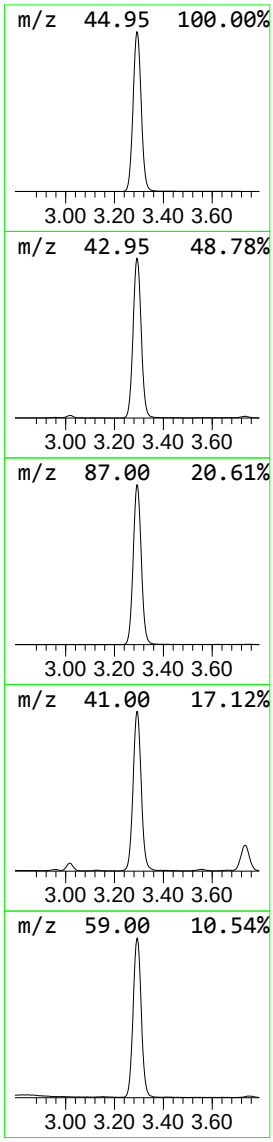
****

Peak Number 2 Diisopropyl ether Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.291	1.63 ug/L	17105600	1,4-Difluorobenzene	5.590

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diisopropyl ether	102	C6H14O	000108-20-3	91
2	Diisopropyl ether	102	C6H14O	000108-20-3	86
3	Diisopropyl ether	102	C6H14O	000108-20-3	83
4	Diisopropyl ether	102	C6H14O	000108-20-3	83
5	Diisopropyl ether	102	C6H14O	000108-20-3	83





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Data File : VV028670.D
Acq On : 21 Oct 2022 18:43
Operator : SY/MD
Sample : N5216-16
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BHA75

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR102022WMA.M
Quant Title : TRACE VOA SFAM1.0

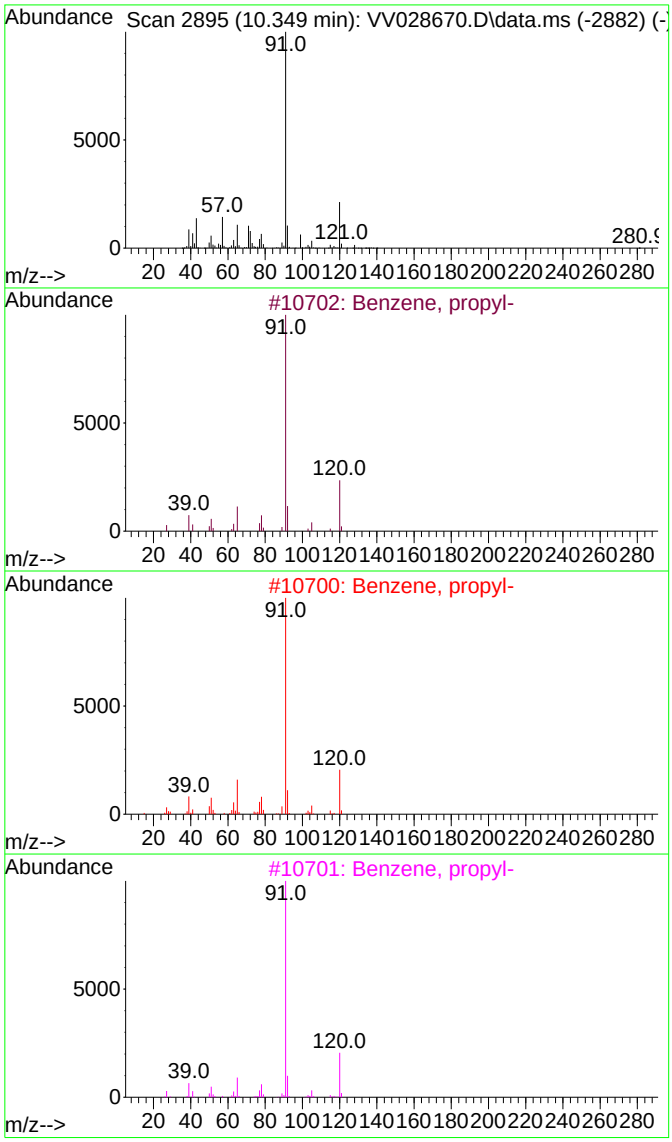
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

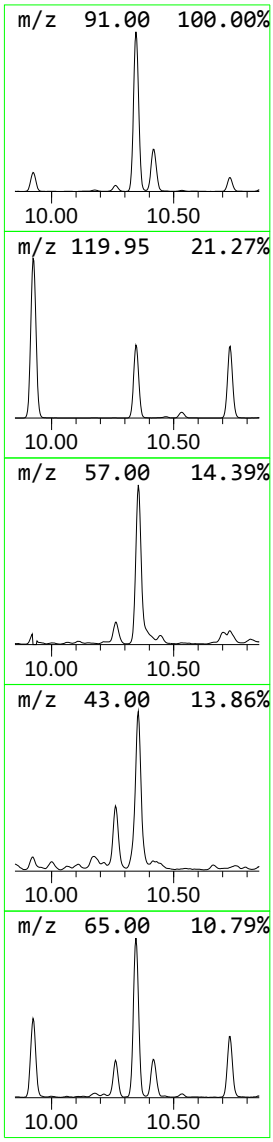
****

Peak Number 3 Benzene, propyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.349	11.33 ug/L	3928300	1,4-Dichlorobenzene-d4	11.242

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, propyl-	120	C9H12	000103-65-1	91
2	Benzene, propyl-	120	C9H12	000103-65-1	64
3	Benzene, propyl-	120	C9H12	000103-65-1	64
4	Benzene, propyl-	120	C9H12	000103-65-1	64
5	1,2-Ethanediamine, N,N'-bis(phen...	240	C16H20N2	000140-28-3	50





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV102122\
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Acq On : 21 Oct 2022 18:43
Operator : SY/MD
Sample : N5216-16
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ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BHA75

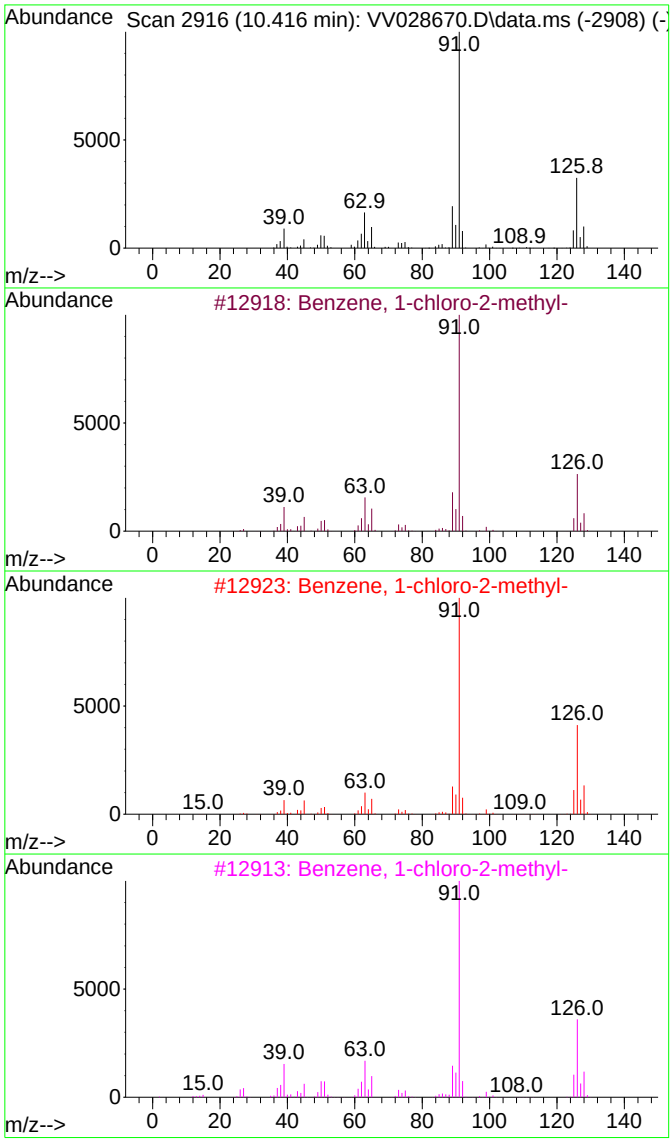
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR102022WMA.M
Quant Title : TRACE VOA SFAM1.0

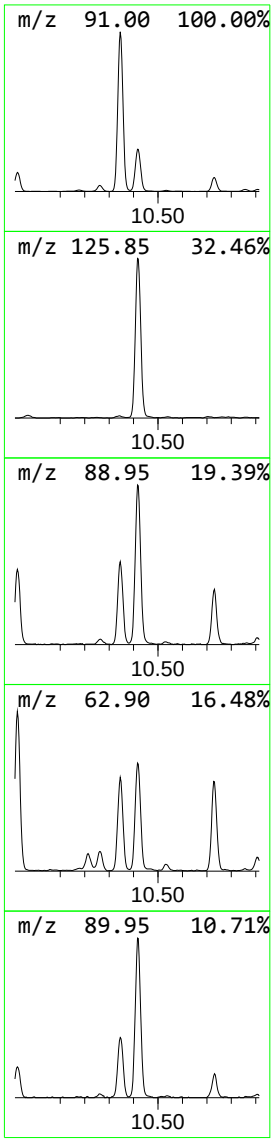
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Benzene, 1-chloro-2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.416	3.46 ug/L	1200440	1,4-Dichlorobenzene-d4	11.242

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-chloro-2-methyl-	126	C7H7Cl	000095-49-8	96
2	Benzene, 1-chloro-2-methyl-	126	C7H7Cl	000095-49-8	93
3	Benzene, 1-chloro-2-methyl-	126	C7H7Cl	000095-49-8	93
4	Benzene, 1-chloro-2-methyl-	126	C7H7Cl	000095-49-8	90
5	Benzene, 1-chloro-2-methyl-	126	C7H7Cl	000095-49-8	90





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV102122\
Data File : VV028670.D
Acq On : 21 Oct 2022 18:43
Operator : SY/MD
Sample : N5216-16
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BHA75

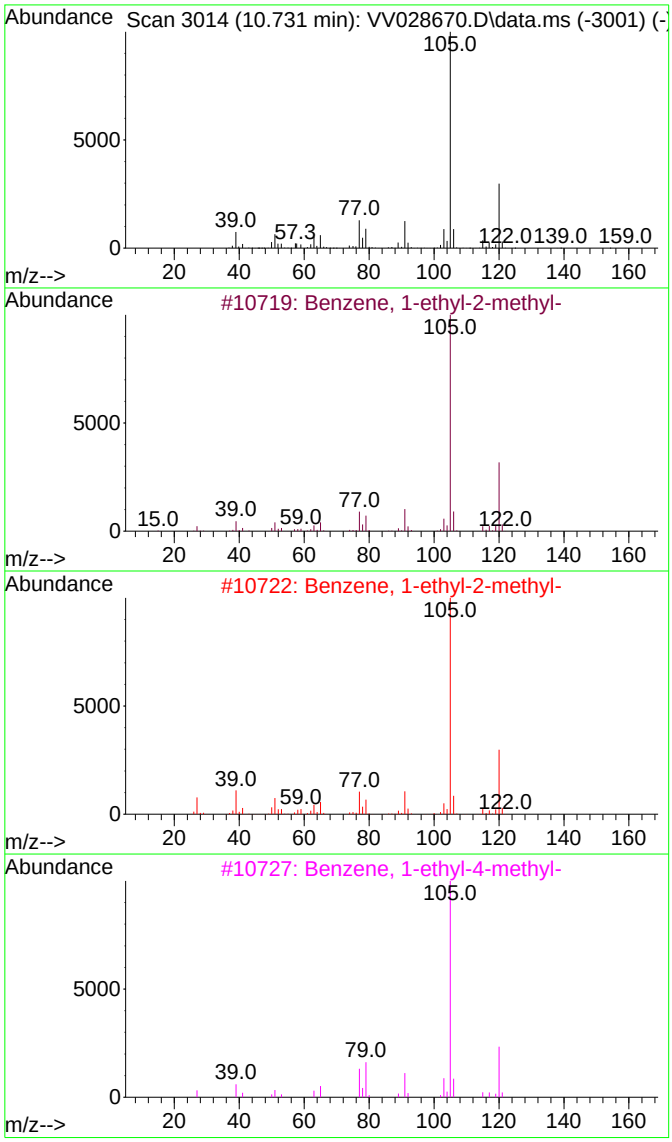
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR102022WMA.M
Quant Title : TRACE VOA SFAM1.0

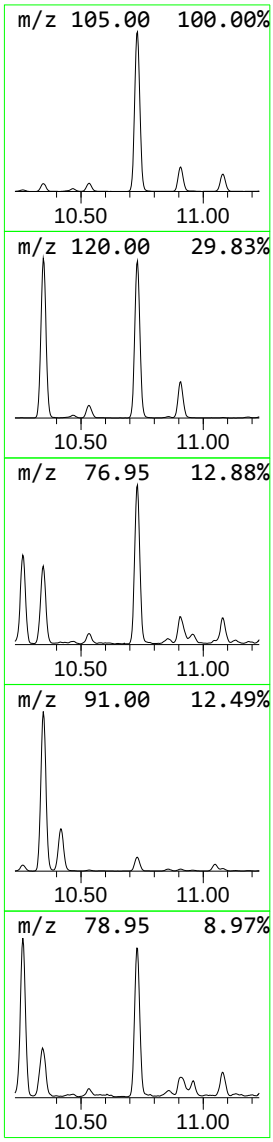
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzene, 1-ethyl-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.731	6.71 ug/L	2324640	1,4-Dichlorobenzene-d4	11.242

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
2	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
3	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
4	Mesitylene	120	C9H12	000108-67-8	90
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV102122\
Data File : VV028670.D
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Sample : N5216-16
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ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BHA75

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR102022WMA.M
Quant Title : TRACE VOA SFAM1.0

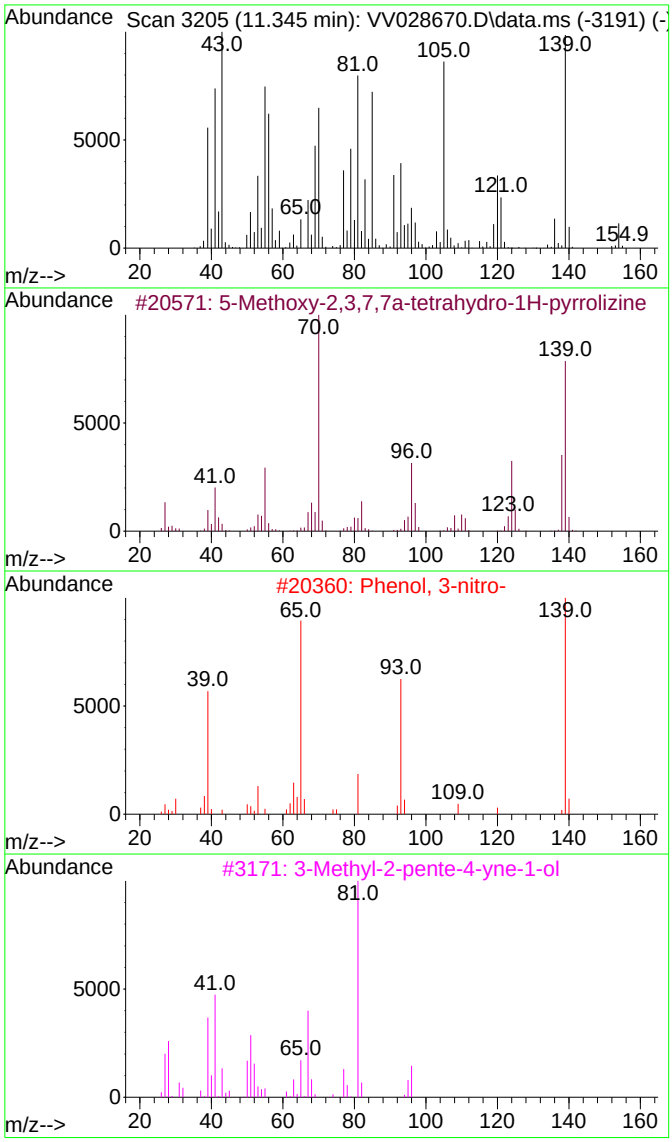
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

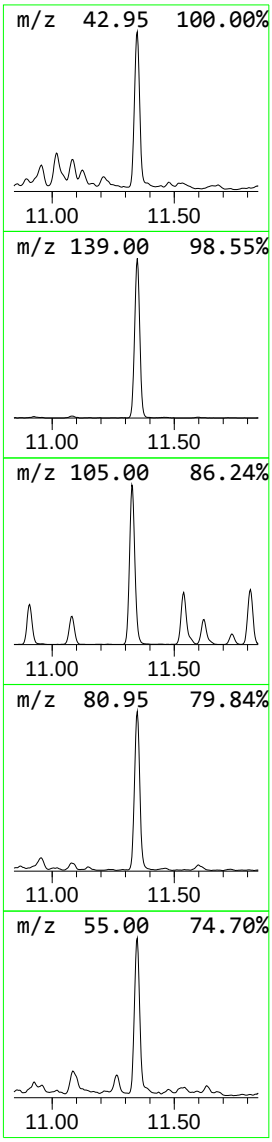
****

Peak Number 6 unknown-01 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.345	9.59 ug/L	3323930	1,4-Dichlorobenzene-d4	11.242

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Methoxy-2,3,7,7a-tetrahydro-1H...	139	C8H13NO	1000193-86-7	18
2	Phenol, 3-nitro-	139	C6H5NO3	000554-84-7	14
3	3-Methyl-2-pente-4-yne-1-ol	96	C6H8O	1000432-32-6	11
4	Cyclobutanone, 2,3,3-trimethyl-	112	C7H12O	028290-01-9	11
5	Pyridine-2,3,5,6-tetraamine	139	C5H9N5	038926-45-3	11





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Instrument :
MSVOA_V
ClientSampleId :
BHA75

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR102022WMA.M
Quant Title : TRACE VOA SFAM1.0

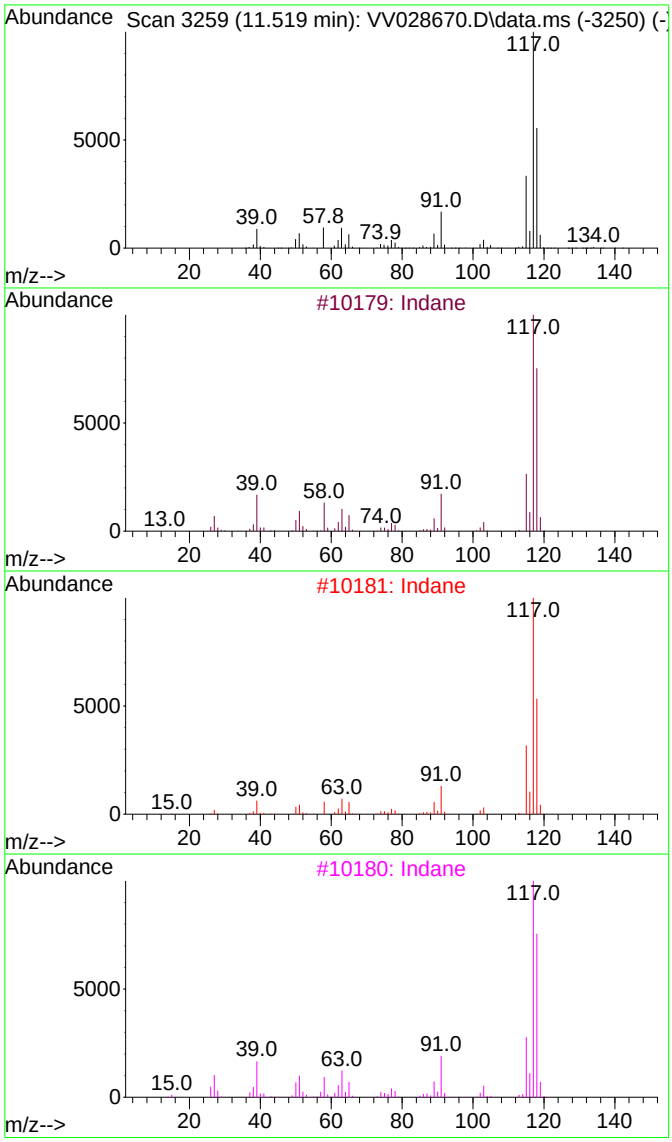
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

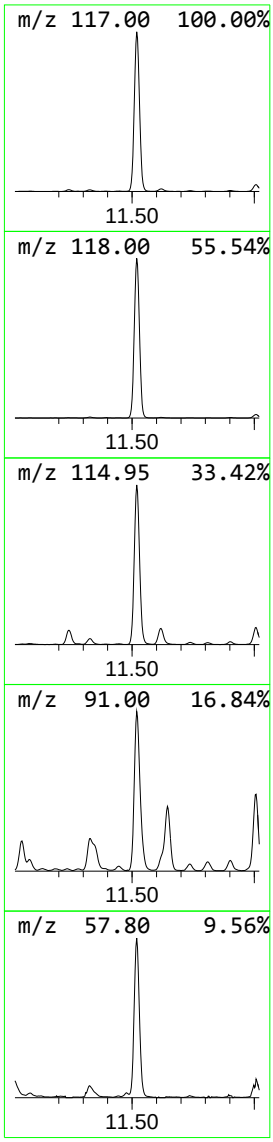
****

Peak Number 7 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.519	17.18 ug/L	5956640	1,4-Dichlorobenzene-d4	11.242

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	96
2	Indane	118	C9H10	000496-11-7	93
3	Indane	118	C9H10	000496-11-7	90
4	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	81
5	Indane	118	C9H10	000496-11-7	81





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Instrument :
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ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR102022WMA.M
Quant Title : TRACE VOA SFAM1.0

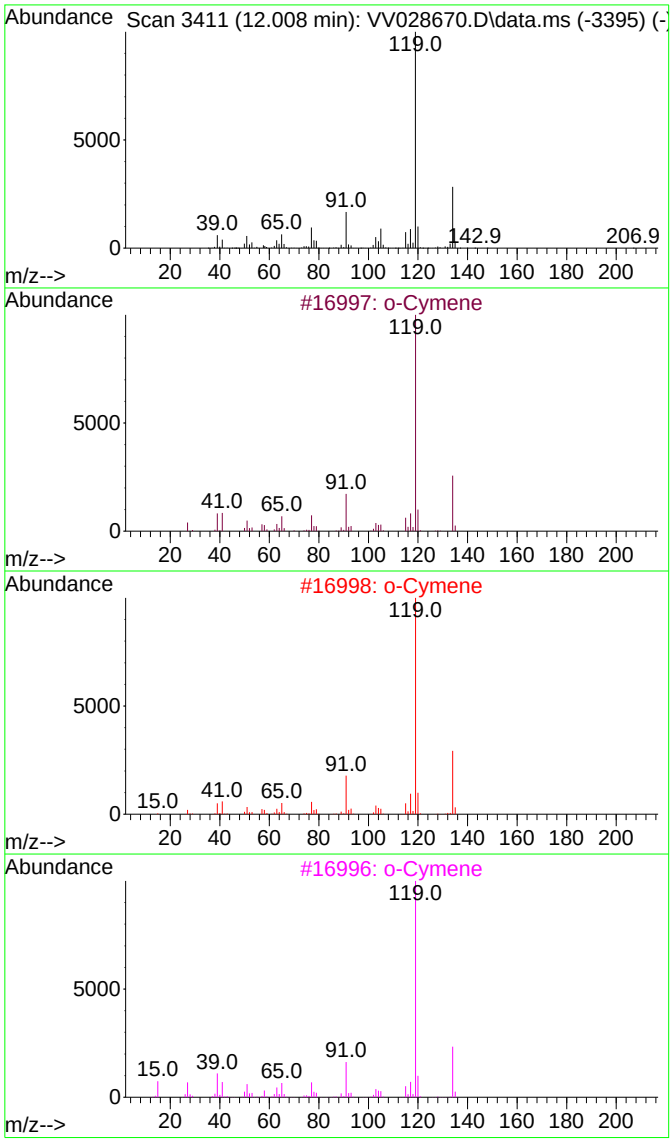
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

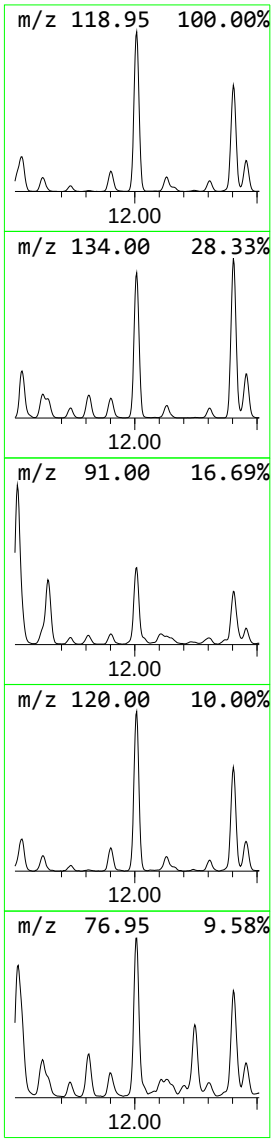
****

Peak Number 8 o-Cymene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.008	7.08 ug/L	2454990	1,4-Dichlorobenzene-d4	11.242

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	95
2	o-Cymene	134	C10H14	000527-84-4	95
3	o-Cymene	134	C10H14	000527-84-4	95
4	p-Cymene	134	C10H14	000099-87-6	95
5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95





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Instrument :
MSVOA_V
ClientSampleId :
BHA75

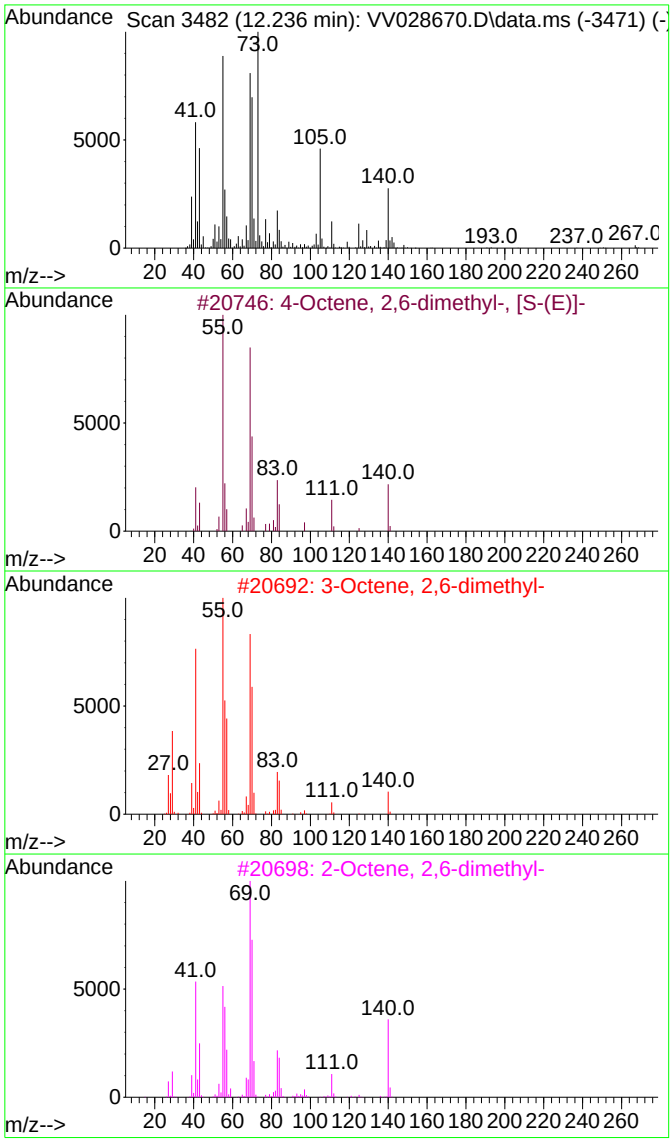
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR102022WMA.M
Quant Title : TRACE VOA SFAM1.0

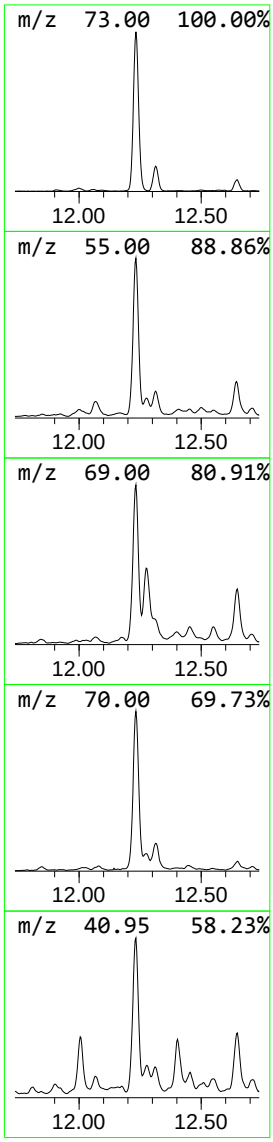
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.236	4.51 ug/L	1564580	1,4-Dichlorobenzene-d4	11.242

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Octene, 2,6-dimethyl-, [S-(E)]-	140	C10H20	062960-76-3	68
2	3-Octene, 2,6-dimethyl-	140	C10H20	006874-28-8	59
3	2-Octene, 2,6-dimethyl-	140	C10H20	004057-42-5	58
4	Cyclohexane, 1,2,3-trimethyl-, (...	126	C9H18	001839-88-9	58
5	trans-4-Decene	140	C10H20	019398-89-1	53





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Data File : VV028670.D
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ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BHA75

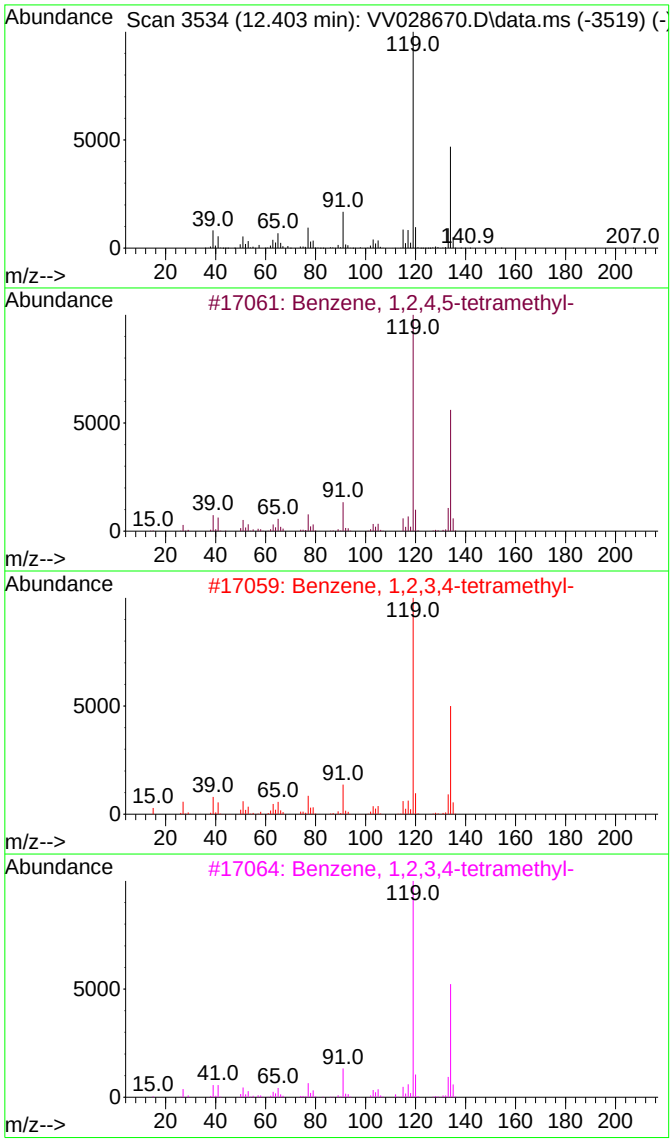
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR102022WMA.M
Quant Title : TRACE VOA SFAM1.0

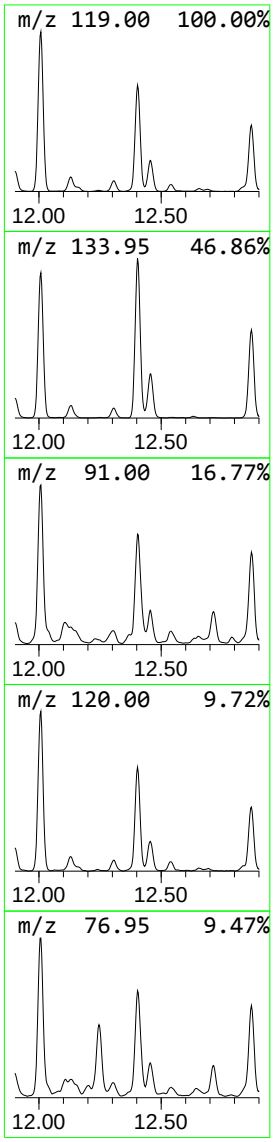
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.403	4.90 ug/L	1699870	1,4-Dichlorobenzene-d4	11.242

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
4	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96





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Data File : VV028670.D
Acq On : 21 Oct 2022 18:43
Operator : SY/MD
Sample : N5216-16
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BHA75

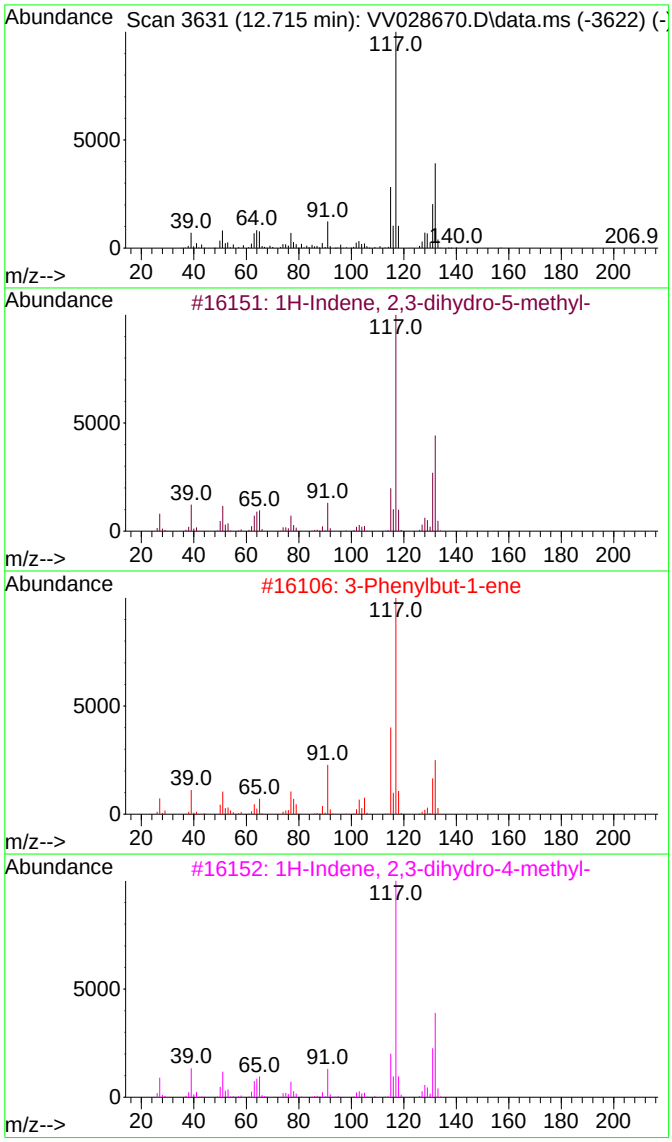
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR102022WMA.M
Quant Title : TRACE VOA SFAM1.0

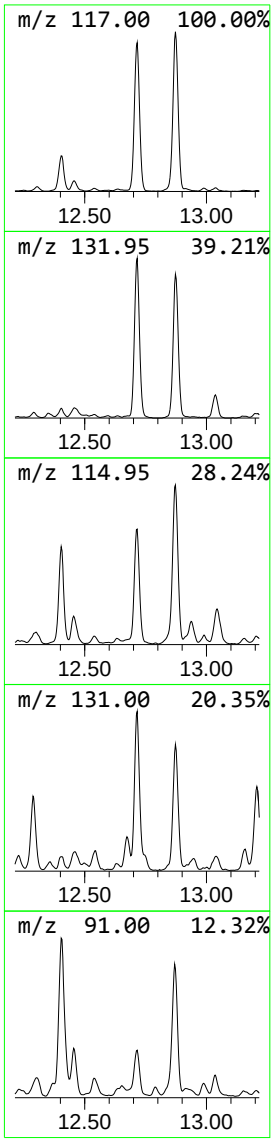
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.715	2.12 ug/L	733826	1,4-Dichlorobenzene-d4	11.242

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	96
2	3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
3	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
4	Indan, 1-methyl-	132	C10H12	000767-58-8	90
5	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87





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Data File : VV028670.D
Acq On : 21 Oct 2022 18:43
Operator : SY/MD
Sample : N5216-16
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BHA75

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR102022WMA.M
Quant Title : TRACE VOA SFAM1.0

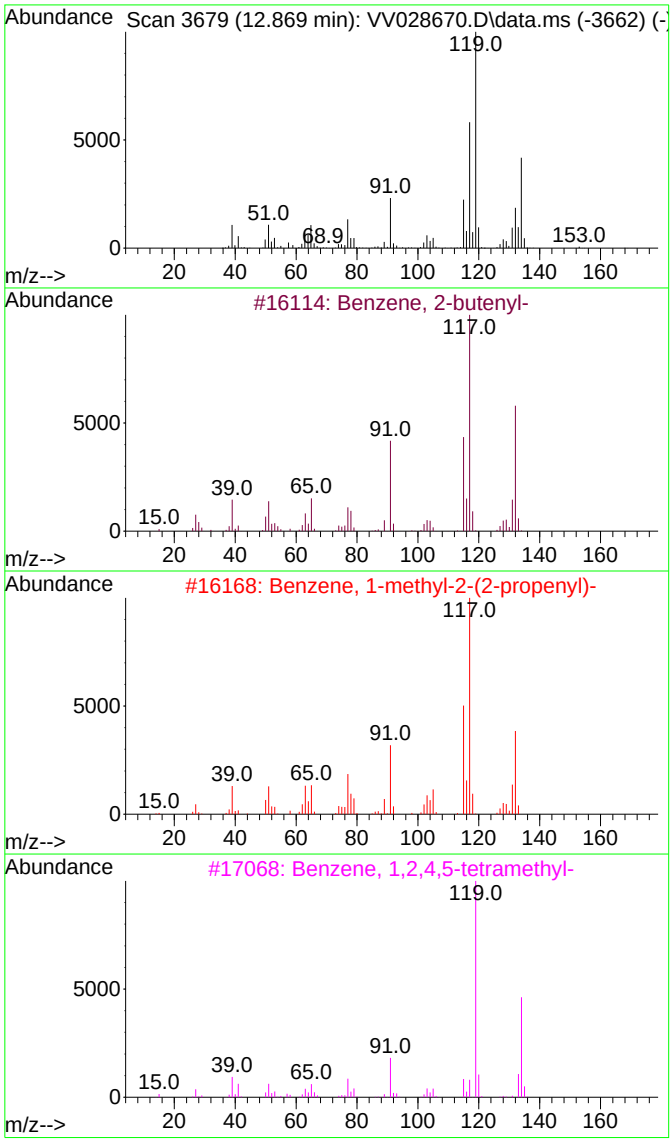
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

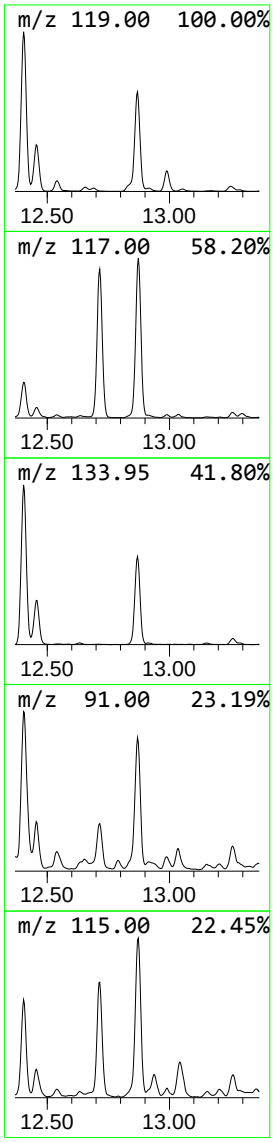
****

Peak Number 12 Benzene, 2-butenyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.869	5.06 ug/L	1753870	1,4-Dichlorobenzene-d4	11.242

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-butenyl-	132	C10H12	001560-06-1	74
2	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	70
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	64
4	p-Cymene	134	C10H14	000099-87-6	64
5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	64





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Data File : VV028670.D
Acq On : 21 Oct 2022 18:43
Operator : SY/MD
Sample : N5216-16
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BHA75

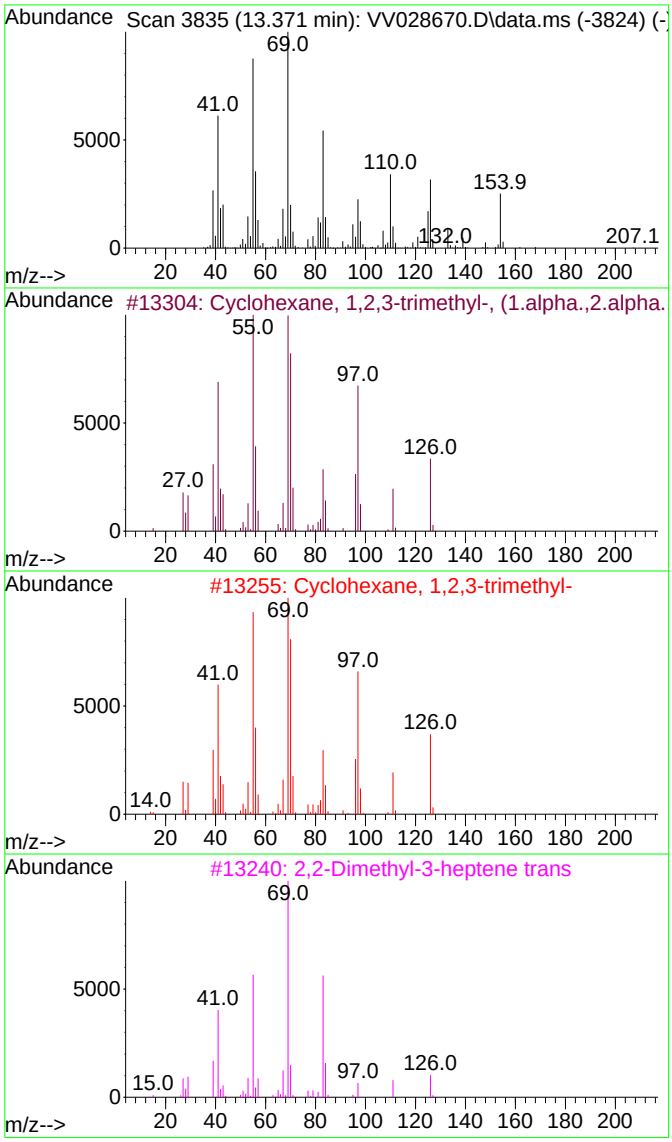
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR102022WMA.M
Quant Title : TRACE VOA SFAM1.0

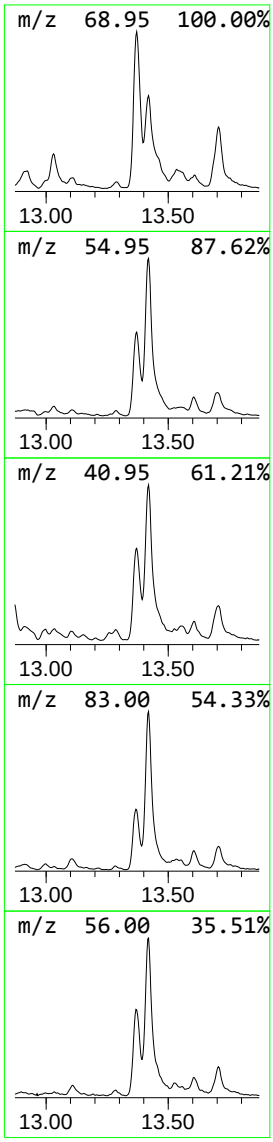
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 (DEL) Alkane: Cyclic13.371 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.371	2.64 ug/L	915479	1,4-Dichlorobenzene-d4	11.242

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,2,3-trimethyl-, (...	126	C9H18	001839-88-9	55
2	Cyclohexane, 1,2,3-trimethyl-	126	C9H18	001678-97-3	55
3	2,2-Dimethyl-3-heptene trans	126	C9H18	019550-75-5	53
4	1-Decene, 5-methyl-	154	C11H22	054244-79-0	52
5	5-Undecene, (E)-	154	C11H22	000764-97-6	52





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Data File : VV028670.D
Acq On : 21 Oct 2022 18:43
Operator : SY/MD
Sample : N5216-16
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BHA75

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR102022WMA.M
Quant Title : TRACE VOA SFAM1.0

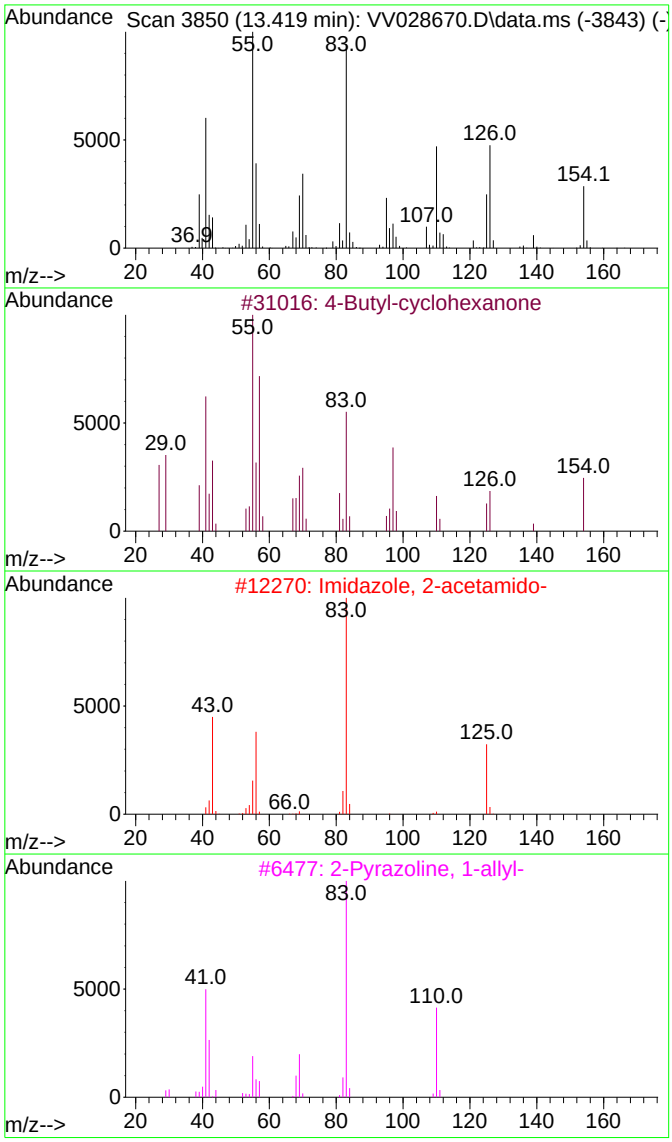
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

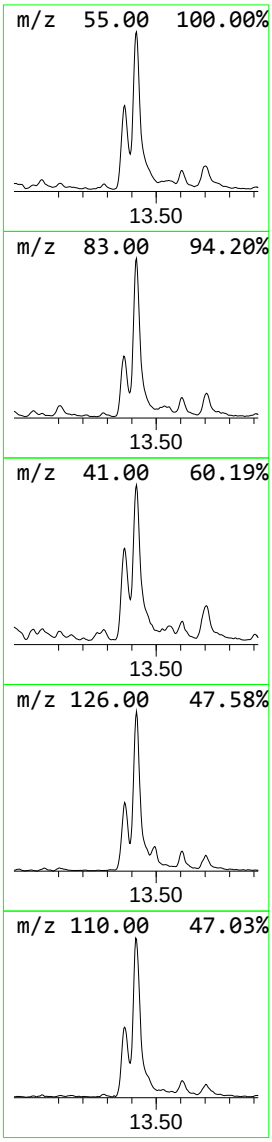
****

Peak Number 14 unknown-02 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.419	3.19 ug/L	1105990	1,4-Dichlorobenzene-d4	11.242

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Butyl-cyclohexanone	154	C10H18O	061203-82-5	43
2	Imidazole, 2-acetamido-	125	C5H7N3O	052737-49-2	38
3	2-Pyrazoline, 1-allyl-	110	C6H10N2	019804-38-7	35
4	Cyclopentane, 1-ethyl-3-methyl-	112	C8H16	003726-47-4	30
5	2-Octene, 2,3,7-trimethyl-	154	C11H22	033933-75-4	30





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV102122\
Data File : VV028670.D
Acq On : 21 Oct 2022 18:43
Operator : SY/MD
Sample : N5216-16
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BHA75

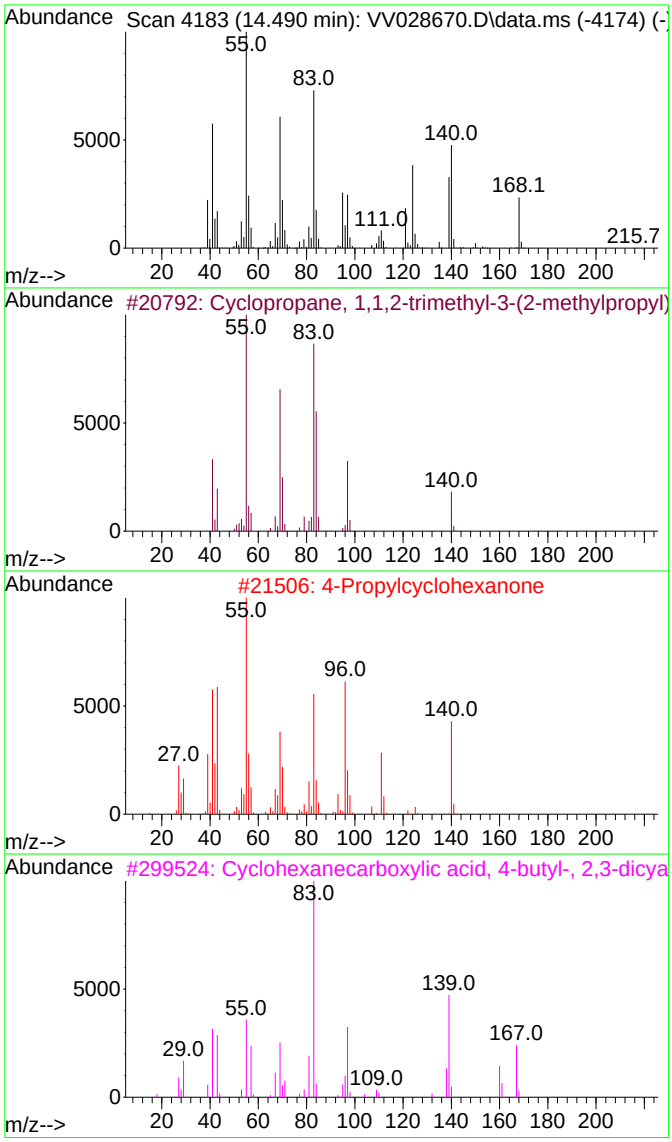
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR102022WMA.M
Quant Title : TRACE VOA SFAM1.0

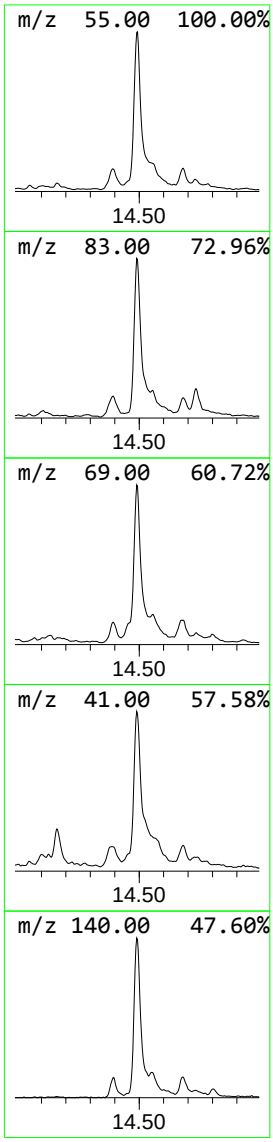
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 (DEL) Alkane: Cyclic14.490 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.490	5.06 ug/L	1755820	1,4-Dichlorobenzene-d4	11.242

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopropane, 1,1,2-trimethyl-3-...	140	C10H20	041977-43-9	38
2	4-Propylcyclohexanone	140	C9H16O	040649-36-3	38
3	Cyclohexanecarboxylic acid, 4-bu...	396	C24H32N2O3	075941-51-4	27
4	1,3-Cyclohexanedione, 5,5-dimethyl-	140	C8H12O2	000126-81-8	27
5	2,4,6-Trimethyl-3-heptene	140	C10H20	1000113-56-9	27





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV102122\
Data File : VV028670.D
Acq On : 21 Oct 2022 18:43
Operator : SY/MD
Sample : N5216-16
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BHA75

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR102022WMA.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
Ethyl ether	1.947	0.8	ug/L	8372170	1	5.590	52595200	5.0
Diisopropyl ether	3.291	1.6	ug/L	17105600	1	5.590	52595200	5.0
Benzene, propyl-	10.349	11.3	ug/L	3928300	3	11.242	1733310	5.0
Benzene, 1-chlo...	10.416	3.5	ug/L	1200440	3	11.242	1733310	5.0
Benzene, 1-ethy...	10.731	6.7	ug/L	2324640	3	11.242	1733310	5.0
unknown-01	11.345	9.6	ug/L	3323930	3	11.242	1733310	5.0
Indane	11.519	17.2	ug/L	5956640	3	11.242	1733310	5.0
o-Cymene	12.008	7.1	ug/L	2454990	3	11.242	1733310	5.0
(DEL) Alkane: S...	12.236	4.5	ug/L	1564580	3	11.242	1733310	5.0
Benzene, 1,2,4,...	12.403	4.9	ug/L	1699870	3	11.242	1733310	5.0
1H-Indene, 2,3-...	12.715	2.1	ug/L	733826	3	11.242	1733310	5.0
Benzene, 2-bute...	12.869	5.1	ug/L	1753870	3	11.242	1733310	5.0
(DEL) Alkane: C...	13.371	2.6	ug/L	915479	3	11.242	1733310	5.0
unknown-02	13.419	3.2	ug/L	1105990	3	11.242	1733310	5.0
(DEL) Alkane: C...	14.490	5.1	ug/L	1755820	3	11.242	1733310	5.0