

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU011421\
 Data File : VU042035.D
 Acq On : 14 Jan 2021 15:25
 Operator : SY/MD
 Sample : M1129-09
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BFT67

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\SOMUTR011321WMA.M

Title : TRACE VOA SOM01.0

Signal : TIC: VU042035.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.359	3	6	26	rVB	927195	844690	36.23%	8.800%
2	1.491	39	47	53	rBV2	95267	95738	4.11%	0.997%
3	1.597	69	80	89	rBV3	422843	712636	30.56%	7.425%
4	1.736	116	123	139	rVB2	45373	61548	2.64%	0.641%
5	1.916	163	179	192	rVB3	77012	129875	5.57%	1.353%
6	1.999	195	205	215	rBV	90086	119022	5.10%	1.240%
7	2.157	245	254	262	rBV	118985	160674	6.89%	1.674%
8	2.208	262	270	290	rVB3	109540	186563	8.00%	1.944%
9	2.420	319	336	344	rBV3	28462	47453	2.04%	0.494%
10	2.572	369	383	396	rBV2	107203	178705	7.66%	1.862%
11	2.983	492	511	525	rBV9	14613	43812	1.88%	0.456%
12	3.372	613	632	648	rBV4	22385	52905	2.27%	0.551%
13	3.597	687	702	719	rVB4	28353	68357	2.93%	0.712%
14	3.710	723	737	753	rBV2	14268	31086	1.33%	0.324%
15	4.658	1015	1032	1053	rBV2	77786	267426	11.47%	2.786%
16	5.079	1147	1163	1190	rBV3	100830	260951	11.19%	2.719%
17	5.739	1346	1368	1395	rBV2	185313	501246	21.50%	5.222%
18	6.259	1516	1530	1553	rBV	151783	288762	12.38%	3.008%
19	6.552	1599	1621	1652	rBV	1242186	2331762	100.00%	24.294%
20	6.700	1655	1667	1684	rVB	116916	225896	9.69%	2.354%
21	7.115	1782	1796	1806	rBV2	19179	33844	1.45%	0.353%
22	7.571	1925	1938	1951	rBV	62340	107477	4.61%	1.120%
23	7.902	2026	2041	2056	rBV	208806	366220	15.71%	3.815%
24	8.185	2118	2129	2139	rBV2	38854	65043	2.79%	0.678%
25	8.555	2234	2244	2258	rVB2	51021	84476	3.62%	0.880%
26	8.642	2258	2271	2297	rBV	358898	677721	29.06%	7.061%
27	8.793	2308	2318	2335	rVB5	17879	46642	2.00%	0.486%
28	9.420	2499	2513	2531	rBV	203625	336459	14.43%	3.505%
29	9.767	2607	2621	2634	rVB2	19625	34688	1.49%	0.361%
30	10.758	2916	2929	2943	rVB	111331	176748	7.58%	1.841%
31	11.687	3208	3218	3229	rBV3	25125	38972	1.67%	0.406%
32	11.809	3244	3256	3272	rVB	214019	334515	14.35%	3.485%
33	12.188	3362	3374	3390	rBV	219462	352181	15.10%	3.669%
34	12.880	3578	3589	3601	rBV2	62580	93480	4.01%	0.974%
35	12.963	3602	3615	3624	rVV	33288	49336	2.12%	0.514%
36	13.018	3624	3632	3641	rVB	18551	27814	1.19%	0.290%

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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\SOMUTR011321WMA.M
Title : TRACE VOA SOM01.0

37	13.436	3755	3762	3777	rVB4	18378	33443	1.43%	0.348%
38	13.822	3873	3882	3888	rBV3	14777	23739	1.02%	0.247%
39	14.076	3949	3961	3971	rBV	66241	106365	4.56%	1.108%

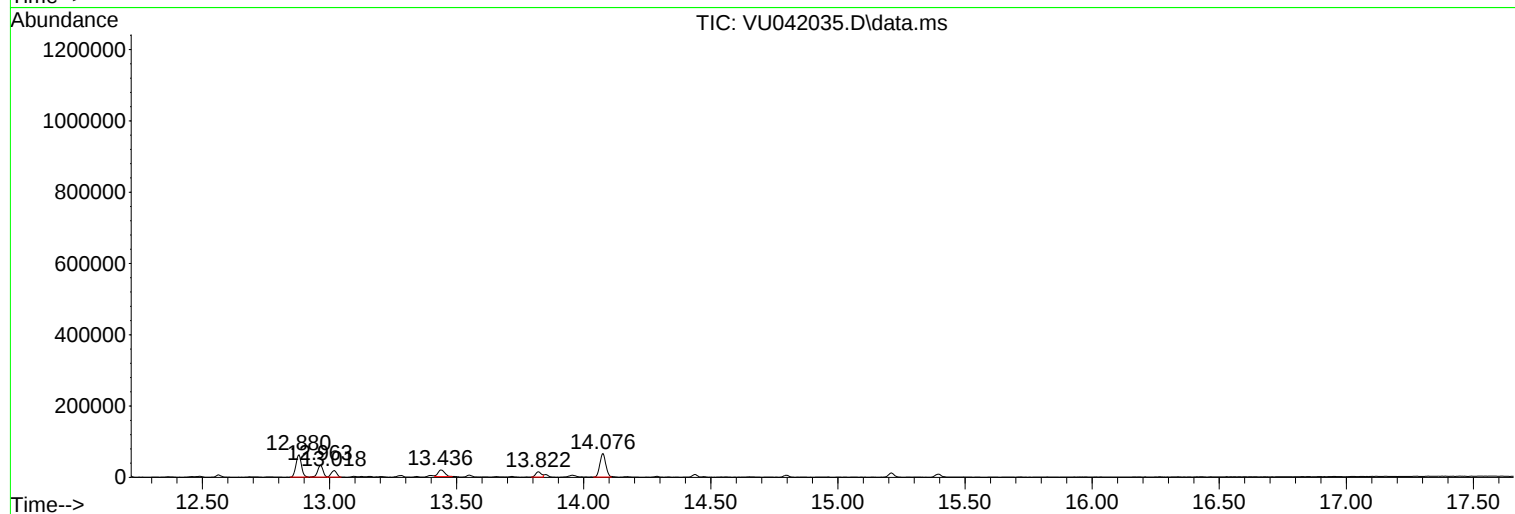
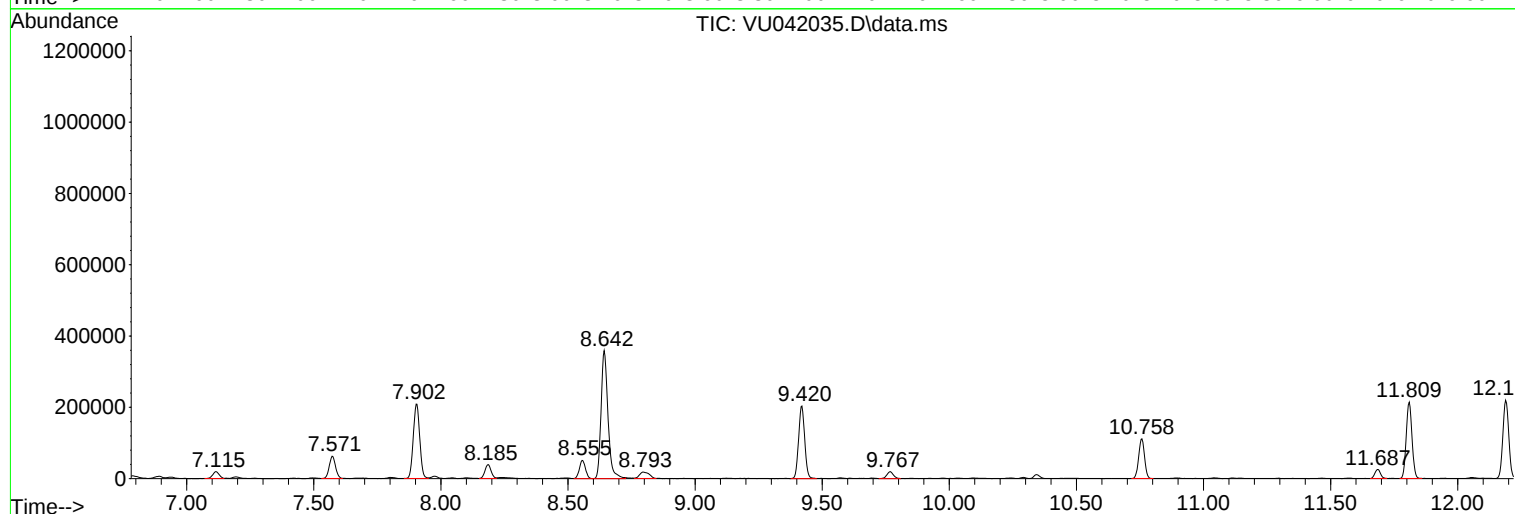
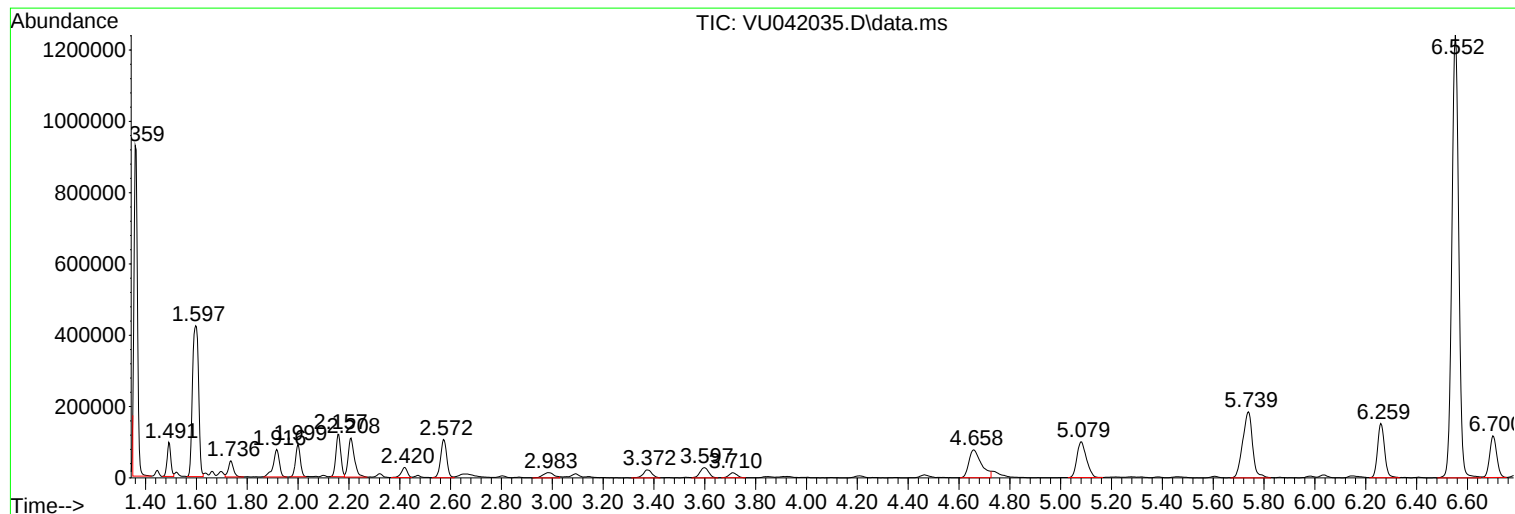
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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR011321WMA.M
Quant Title : TRACE VOA SOM01.0

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



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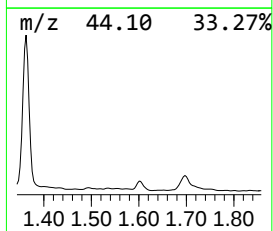
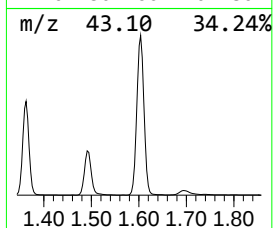
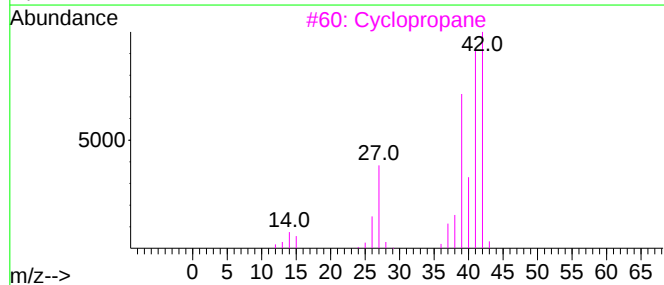
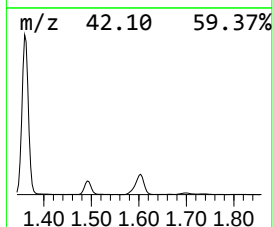
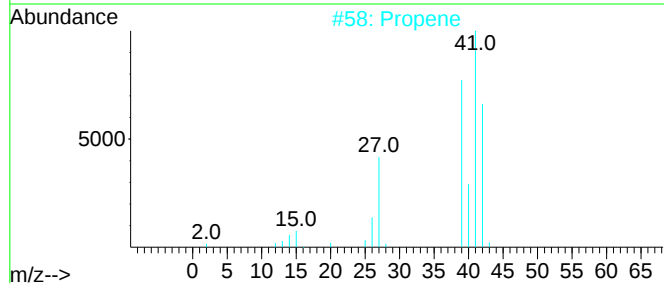
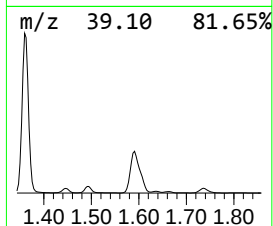
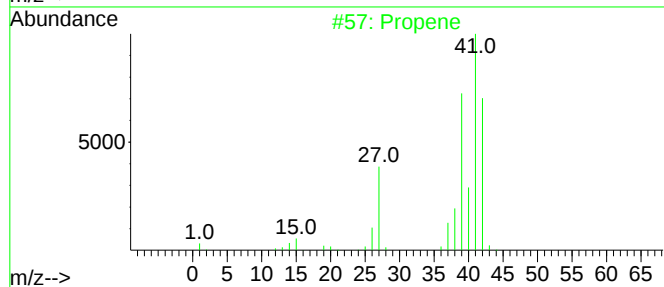
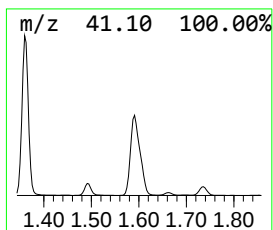
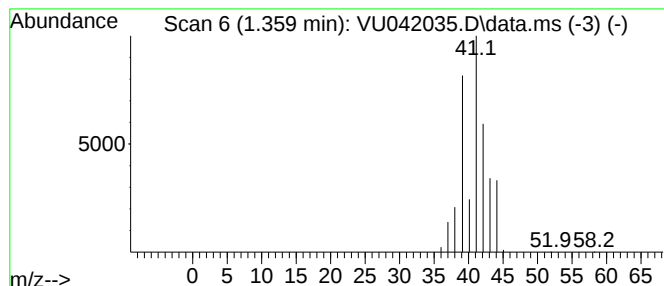
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Quant Title : TRACE VOA SOM01.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.359	14.63 ug/L	844690	1,4-Difluorobenzene	6.259

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			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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ClientSampleId :
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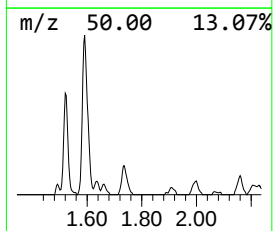
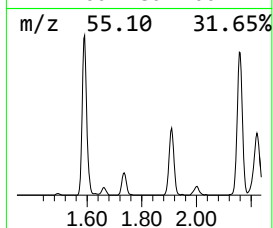
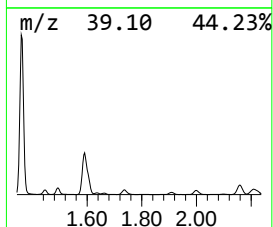
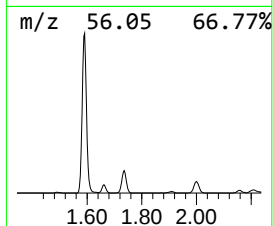
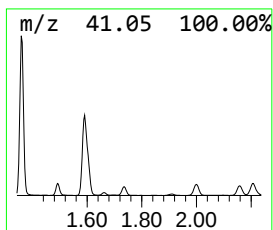
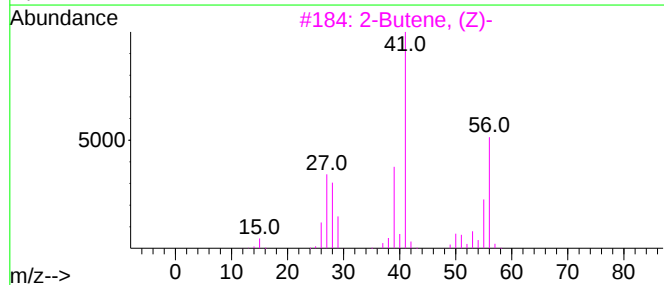
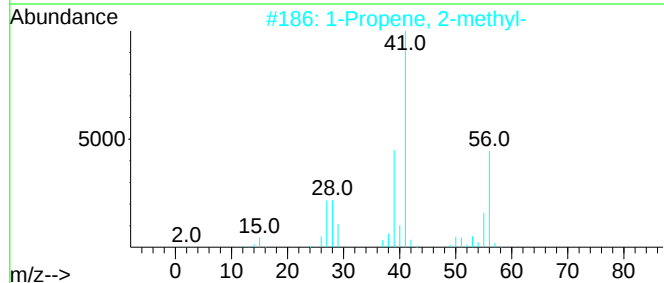
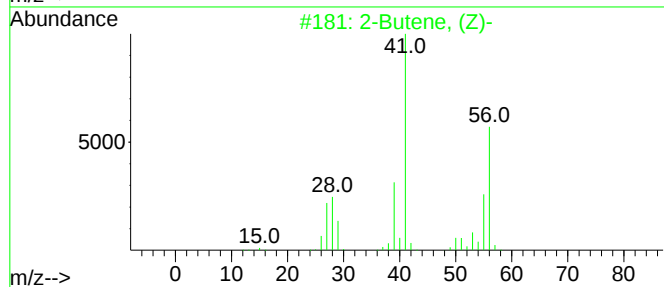
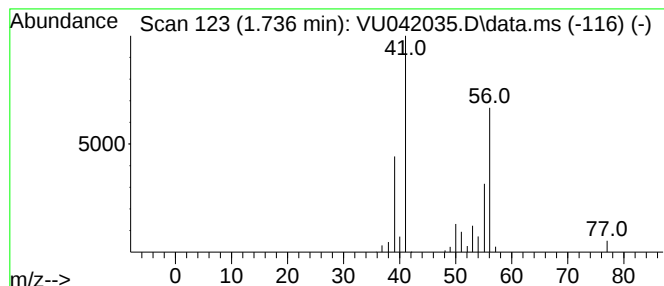
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.736	1.07 ug/L	61548	1,4-Difluorobenzene	6.259

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butene, (Z)-		56	C4H8	000590-18-1	72
2	1-Propene, 2-methyl-		56	C4H8	000115-11-7	72
3	2-Butene, (Z)-		56	C4H8	000590-18-1	68
4	2-Butene, (E)-		56	C4H8	000624-64-6	64
5	2-Butene, (E)-		56	C4H8	000624-64-6	64



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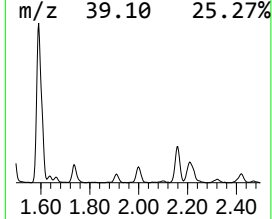
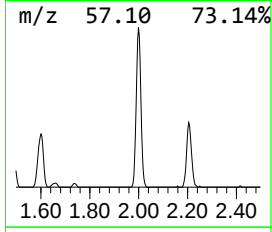
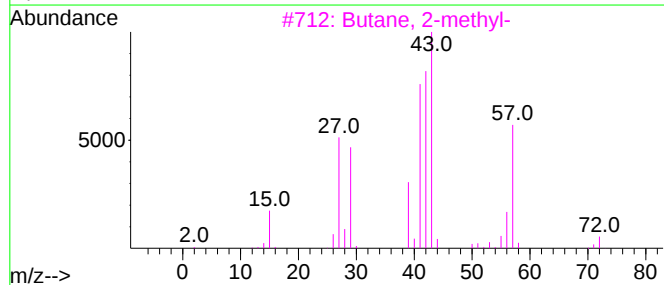
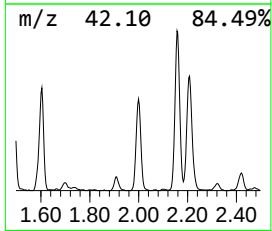
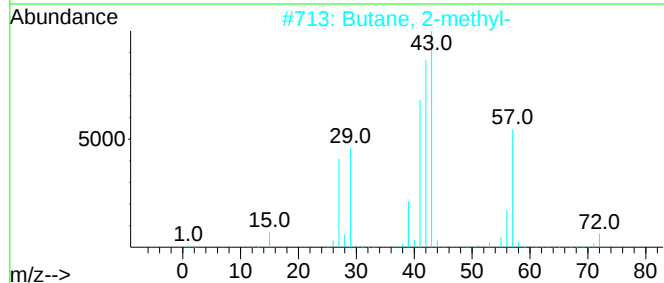
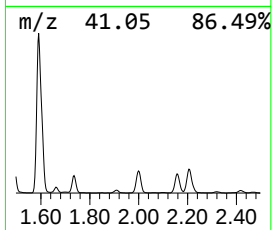
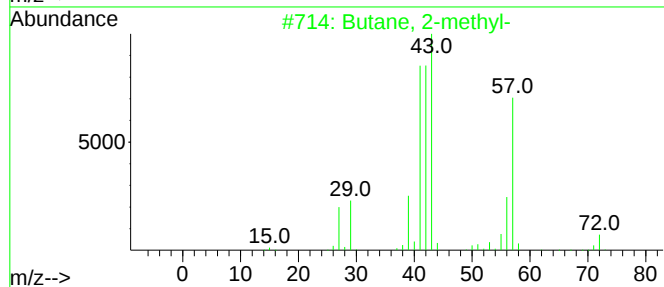
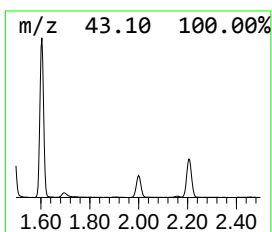
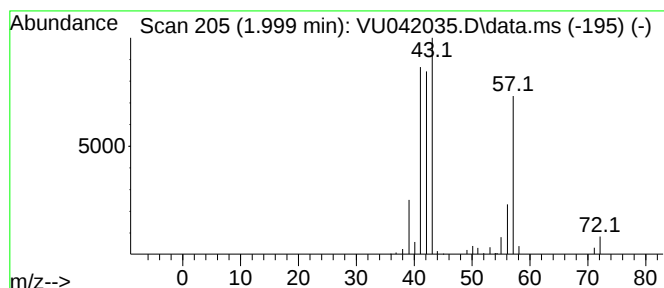
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR011321WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.999	2.06 ug/L	119022	1,4-Difluorobenzene	6.259

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Butane, 2-methyl-	72	C5H12	000078-78-4	91
2		Butane, 2-methyl-	72	C5H12	000078-78-4	91
3		Butane, 2-methyl-	72	C5H12	000078-78-4	90
4		Pentane	72	C5H12	000109-66-0	33
5		3-Buten-1-ol	72	C4H8O	000627-27-0	25



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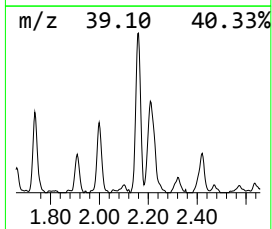
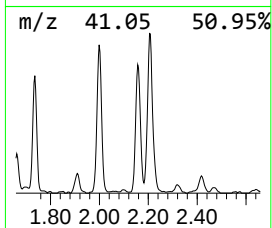
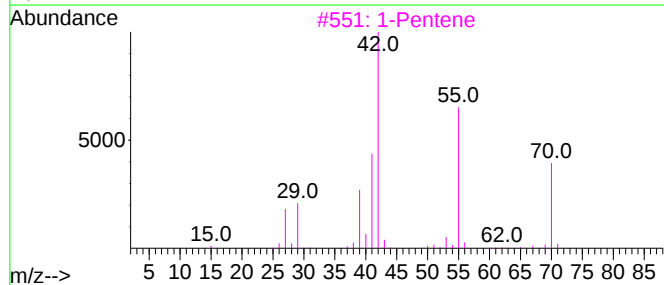
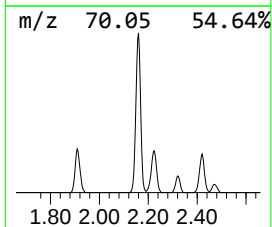
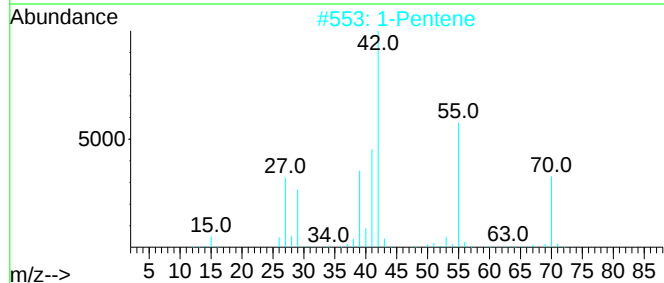
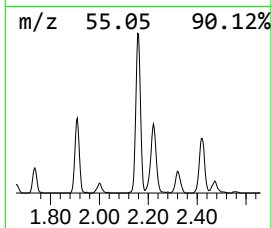
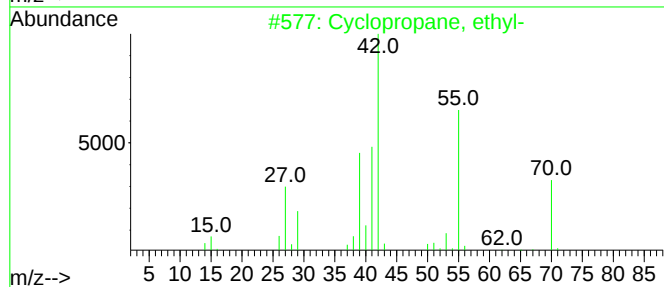
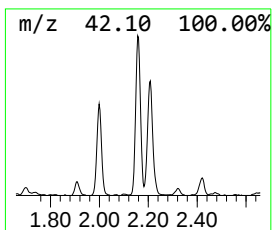
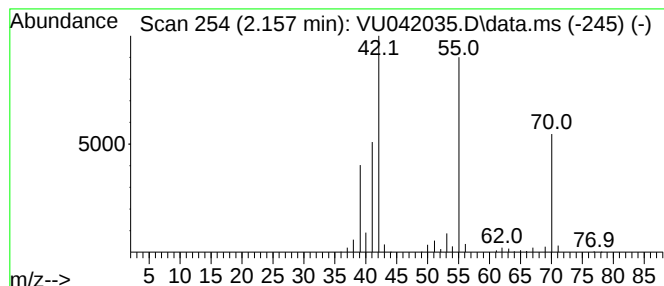
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 (DEL) Alkane: Cyclic2.157 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.157	2.78 ug/L	160674	1,4-Difluorobenzene	6.259

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopropane, ethyl-	70	C5H10	001191-96-4	86
2			1-Pentene	70	C5H10	000109-67-1	70
3			1-Pentene	70	C5H10	000109-67-1	68
4			Cyclobutane, methyl-	70	C5H10	000598-61-8	68
5			Cyclopropane, ethyl-	70	C5H10	001191-96-4	62



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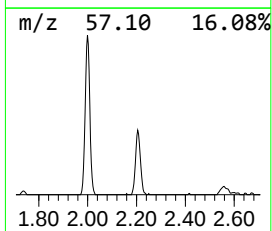
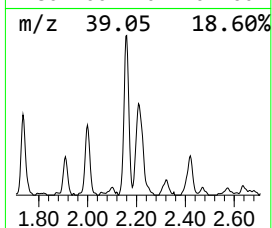
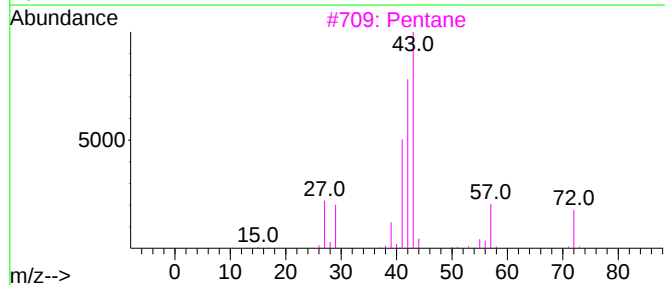
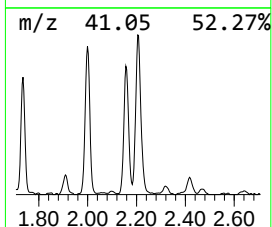
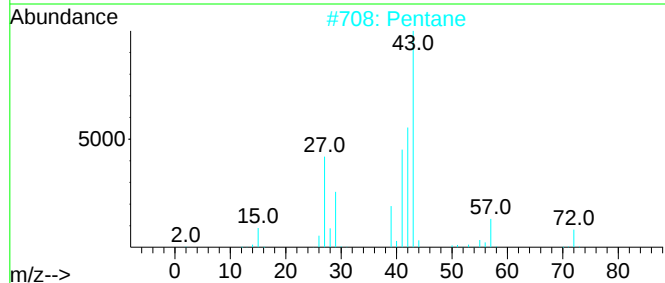
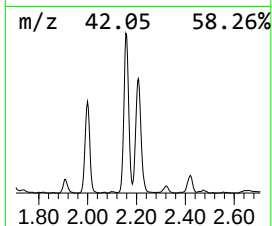
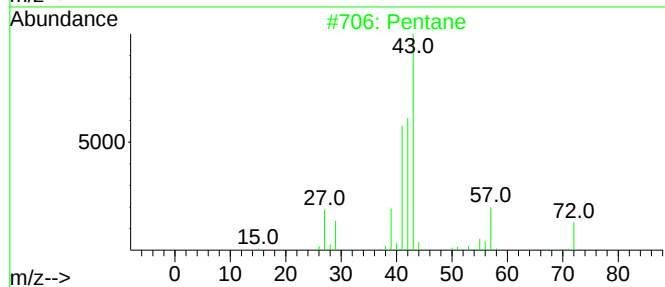
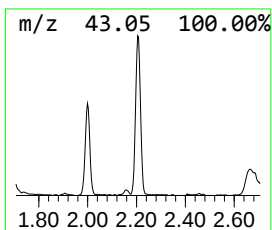
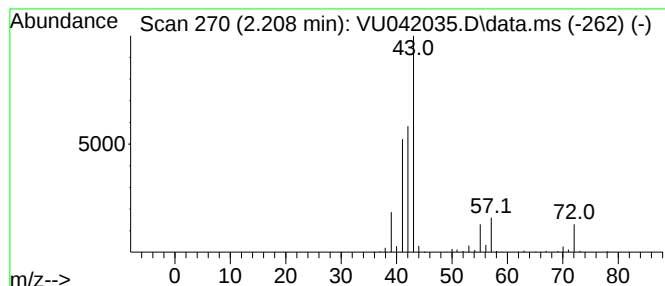
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR011321WMA.M
Quant Title : TRACE VOA SOM01.0

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.
2.208	3.23 ug/L	186563	1,4-Difluorobenzene	6.259

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane	72	C5H12	000109-66-0	90
2	Pentane	72	C5H12	000109-66-0	86
3	Pentane	72	C5H12	000109-66-0	80
4	Pentane	72	C5H12	000109-66-0	80
5	Oxirane, ethyl-	72	C4H8O	000106-88-7	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU011421\
Data File : VU042035.D
Acq On : 14 Jan 2021 15:25
Operator : SY/MD
Sample : M1129-09
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BFT67

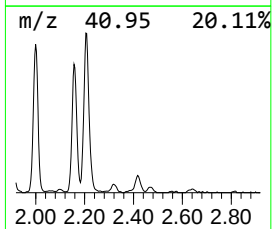
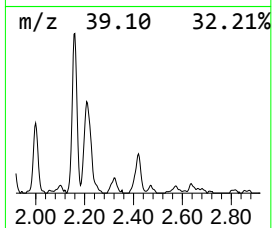
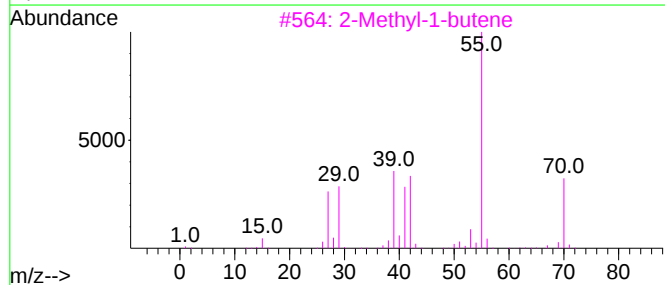
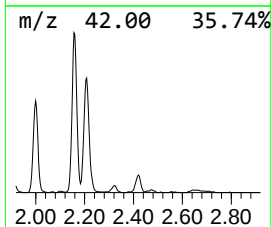
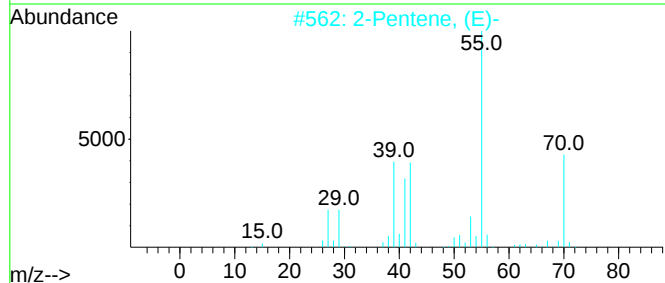
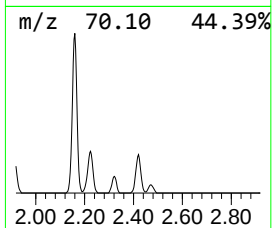
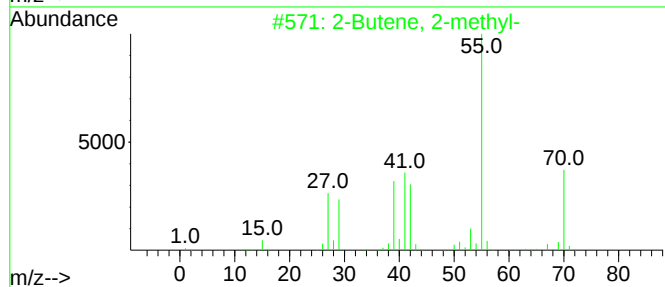
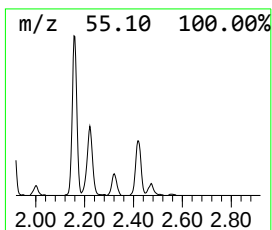
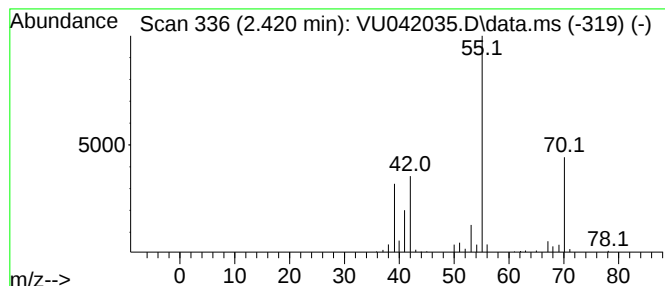
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.420	0.82 ug/L	47453	1,4-Difluorobenzene	6.259

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butene, 2-methyl-	70	C5H10	000513-35-9	90	
2	2-Pentene, (E)-	70	C5H10	000646-04-8	90	
3	2-Methyl-1-butene	70	C5H10	000563-46-2	87	
4	Cyclopropane, 1,2-dimethyl-, cis-	70	C5H10	000930-18-7	87	
5	2-Pentene	70	C5H10	000109-68-2	86	



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU011421\
Data File : VU042035.D
Acq On : 14 Jan 2021 15:25
Operator : SY/MD
Sample : M1129-09
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 15 Sample Multiplier: 1

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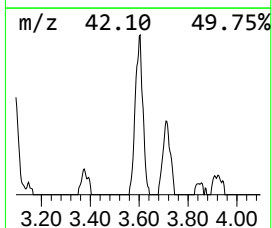
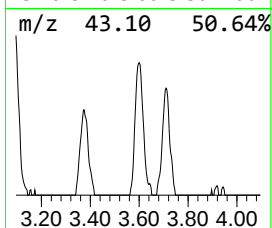
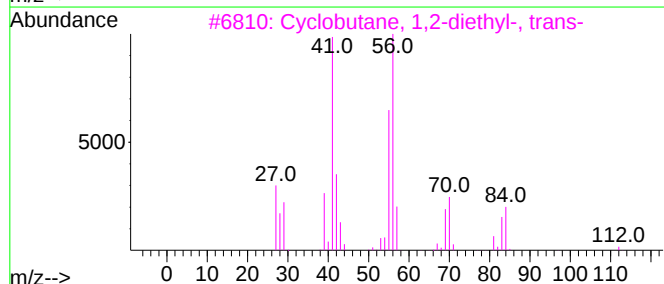
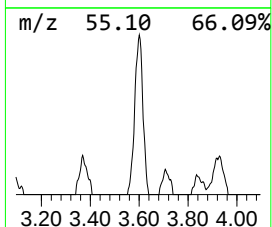
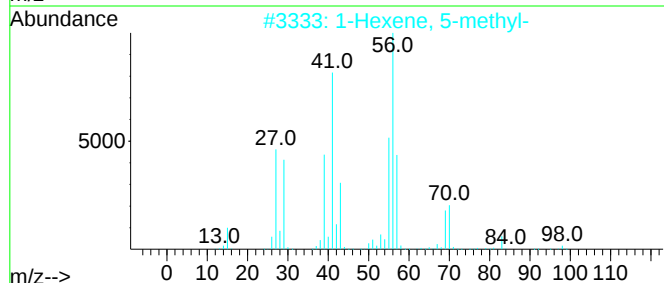
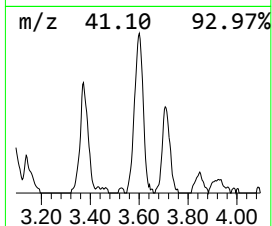
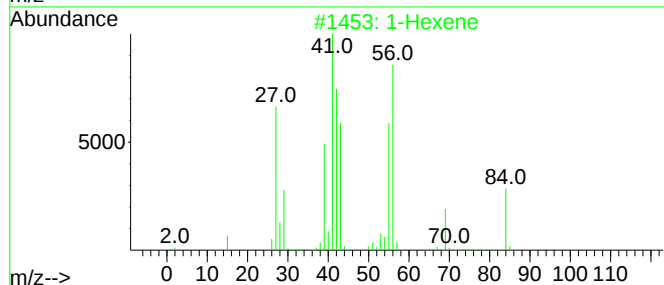
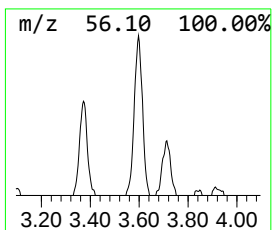
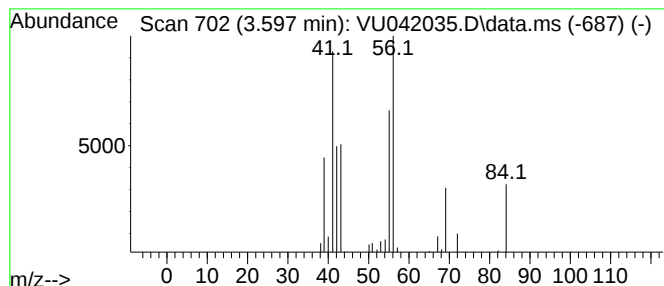
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.597	1.18 ug/L	68357	1,4-Difluorobenzene	6.259

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Hexene	84	C6H12	000592-41-6	76
2			1-Hexene, 5-methyl-	98	C7H14	003524-73-0	53
3			Cyclobutane, 1,2-diethyl-, trans-	112	C8H16	019341-98-1	53
4			Formic acid, hexyl ester	130	C7H14O2	000629-33-4	50
5			1-Pentene, 2-methyl-	84	C6H12	000763-29-1	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU011421\
Data File : VU042035.D
Acq On : 14 Jan 2021 15:25
Operator : SY/MD
Sample : M1129-09
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BFT67

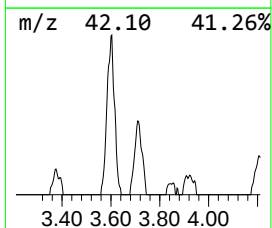
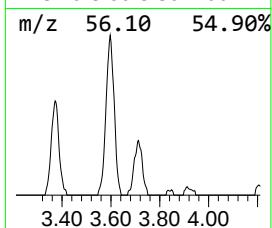
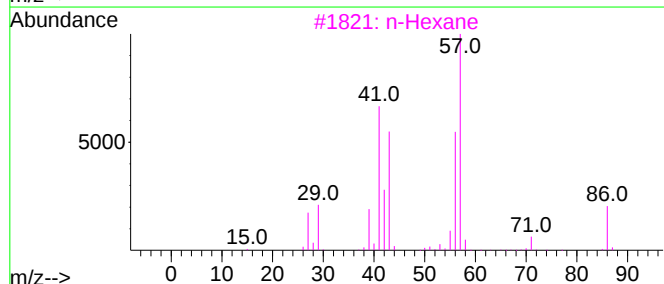
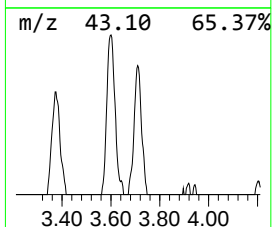
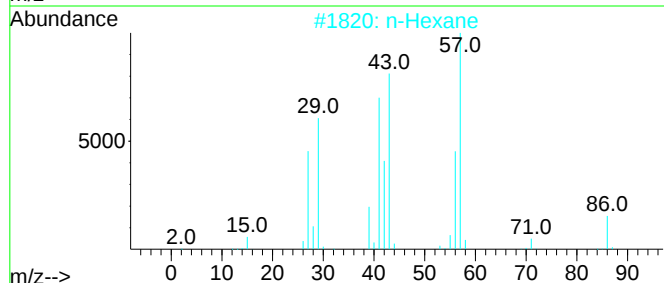
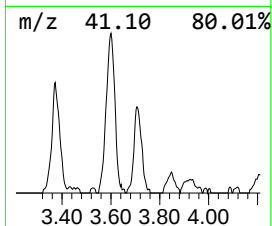
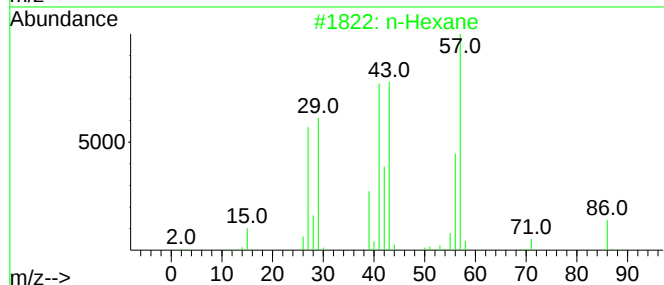
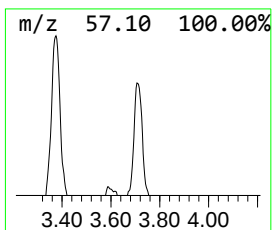
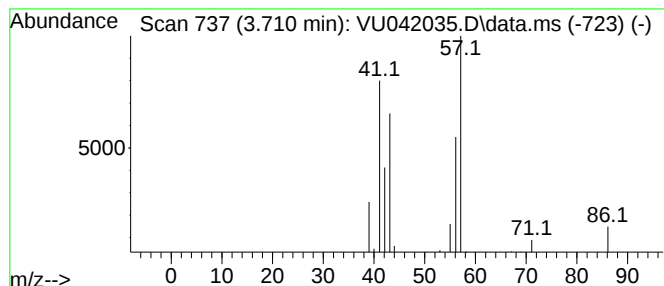
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.710	0.54 ug/L	31086	1,4-Difluorobenzene	6.259

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexane	86	C6H14	000110-54-3	80
2		n-Hexane	86	C6H14	000110-54-3	59
3		n-Hexane	86	C6H14	000110-54-3	53
4		1-Pentene, 4-methyl-	84	C6H12	000691-37-2	42
5		1-Pentene, 4-methyl-	84	C6H12	000691-37-2	36



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU011421\
Data File : VU042035.D
Acq On : 14 Jan 2021 15:25
Operator : SY/MD
Sample : M1129-09
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 15 Sample Multiplier: 1

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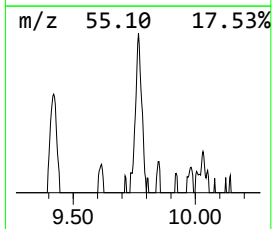
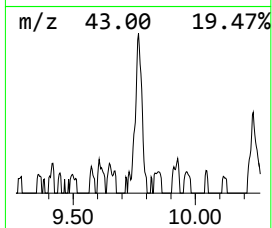
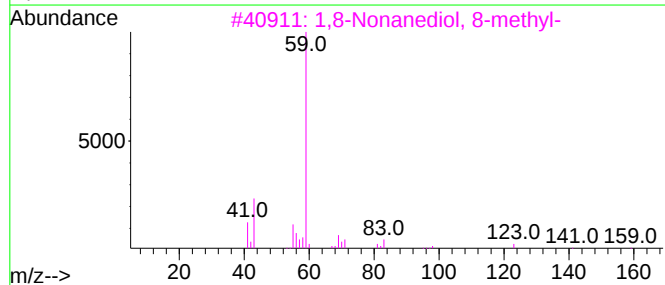
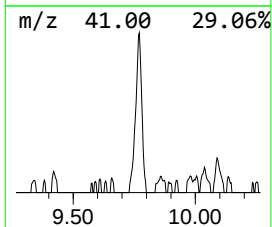
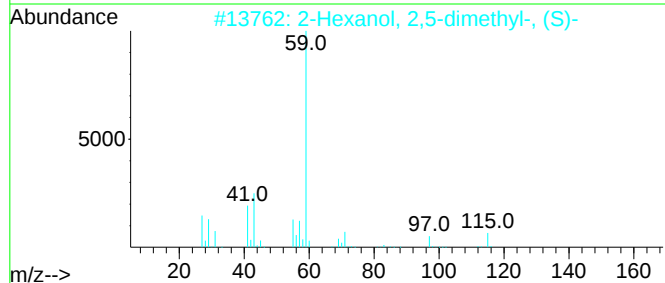
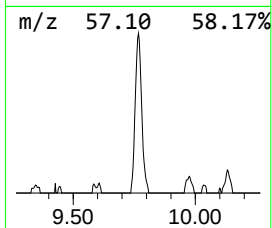
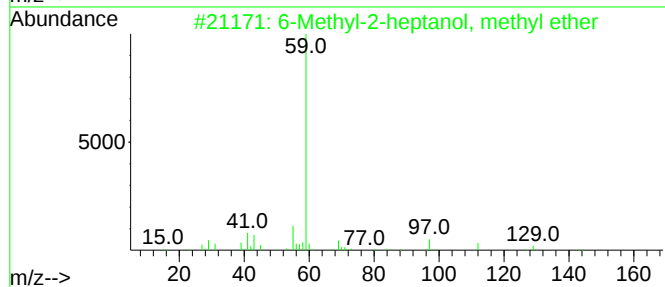
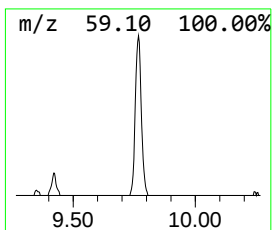
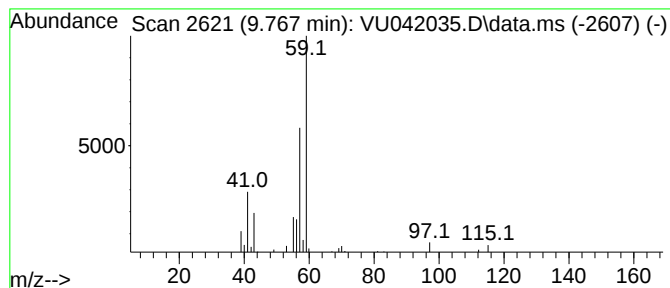
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 10 unknown-01 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.767	0.52 ug/L	34688	Chlorobenzene-d5	9.420

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		6-Methyl-2-heptanol, methyl ether	144	C9H20O	1000365-52-0	45
2		2-Hexanol, 2,5-dimethyl-, (S)-	130	C8H18O	003730-60-7	40
3		1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	39
4		3-Pentanol, 2,2-dimethyl-	116	C7H16O	003970-62-5	39
5		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU011421\
Data File : VU042035.D
Acq On : 14 Jan 2021 15:25
Operator : SY/MD
Sample : M1129-09
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 15 Sample Multiplier: 1

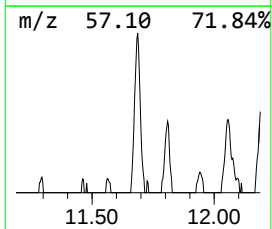
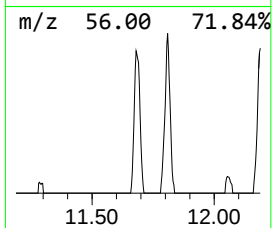
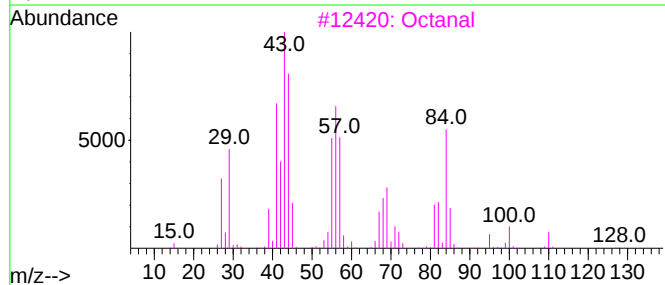
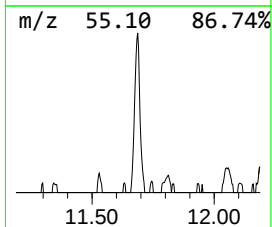
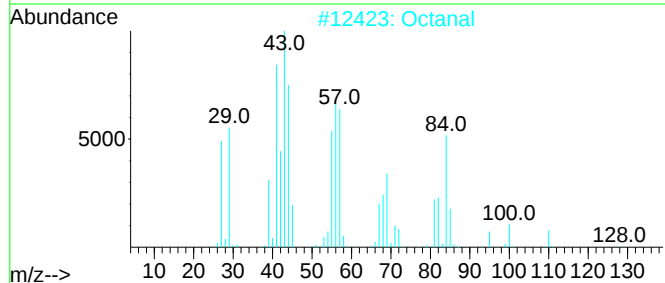
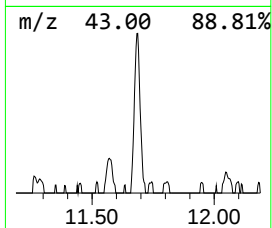
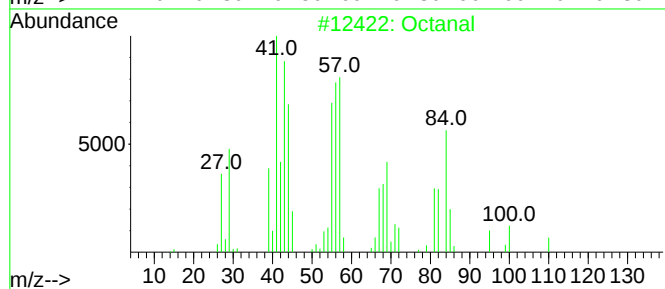
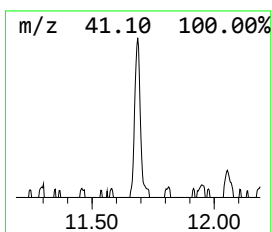
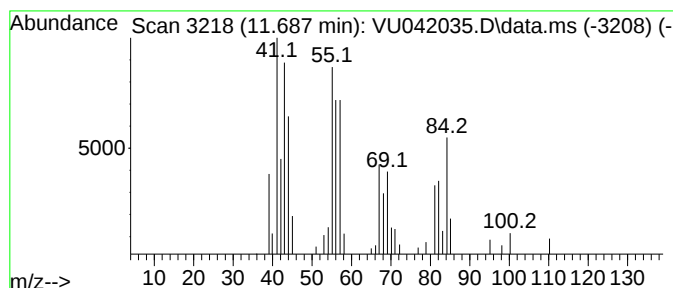
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 11 Octanal Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.687	0.58 ug/L	38972	1,4-Dichlorobenzene-d4		11.809	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Octanal		128	C8H16O	000124-13-0	80
2	Octanal		128	C8H16O	000124-13-0	80
3	Octanal		128	C8H16O	000124-13-0	64
4	Octanal		128	C8H16O	000124-13-0	59
5	Octanal		128	C8H16O	000124-13-0	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU011421\
Data File : VU042035.D
Acq On : 14 Jan 2021 15:25
Operator : SY/MD
Sample : M1129-09
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 15 Sample Multiplier: 1

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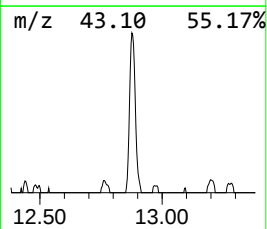
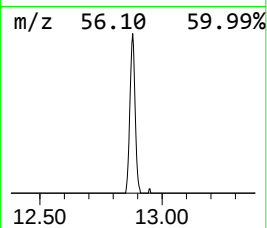
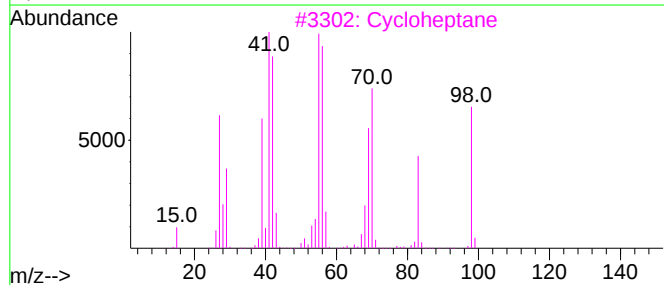
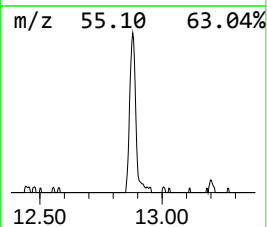
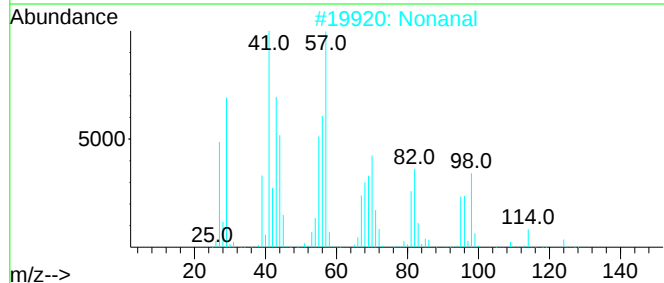
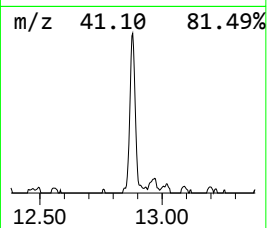
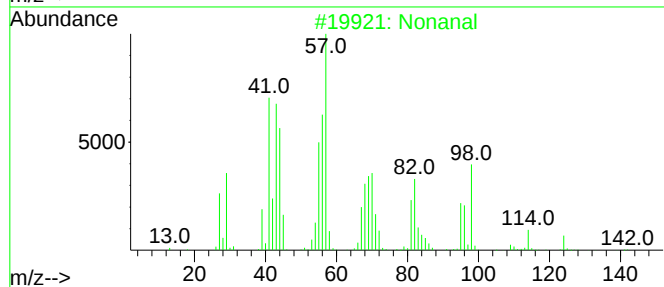
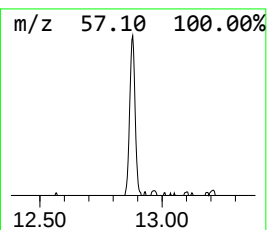
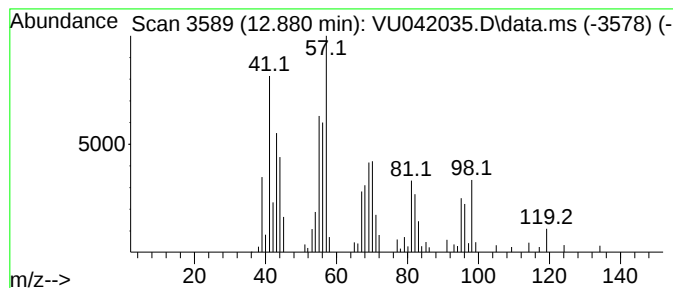
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 12 Nonanal Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.880	1.40 ug/L	93480	1,4-Dichlorobenzene-d4	11.809

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonanal	142	C9H18O	000124-19-6	80
2			Nonanal	142	C9H18O	000124-19-6	59
3			Cycloheptane	98	C7H14	000291-64-5	38
4			2-Propen-1-amine, N-2-propenyl-	97	C6H11N	000124-02-7	25
5			1-Pentanol, 3,4-dimethyl-	116	C7H16O	006570-87-2	22



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU011421\
Data File : VU042035.D
Acq On : 14 Jan 2021 15:25
Operator : SY/MD
Sample : M1129-09
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BFT67

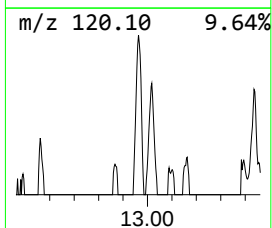
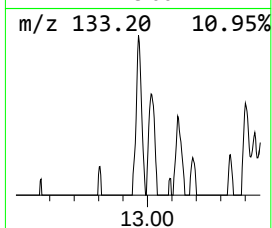
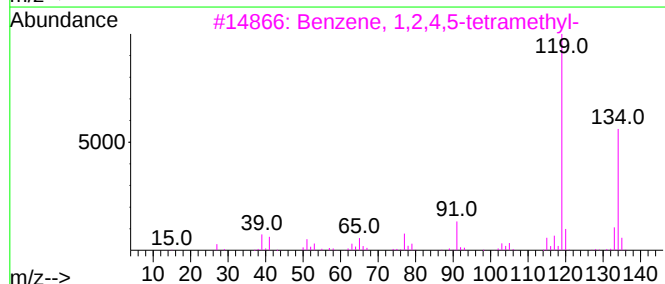
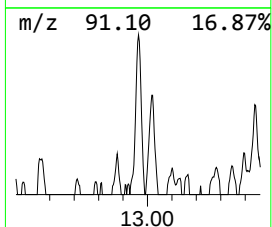
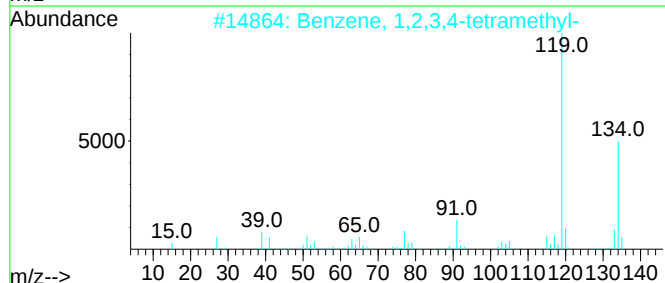
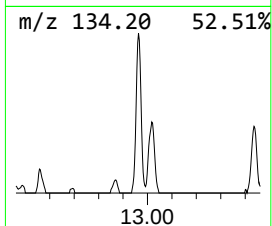
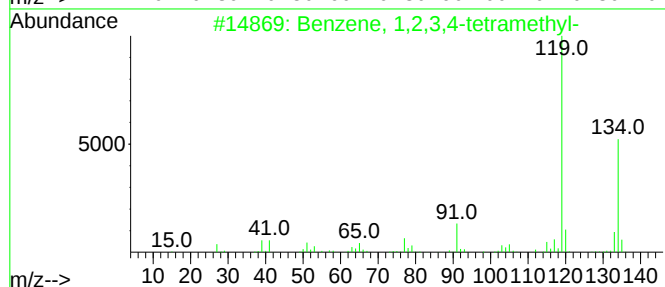
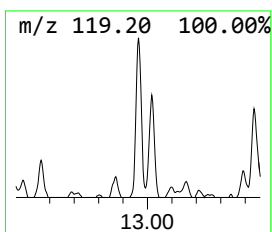
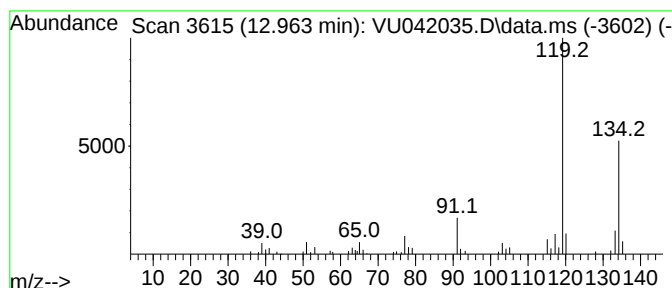
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 13 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.963	0.74 ug/L	49336	1,4-Dichlorobenzene-d4	11.809

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
4			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
5			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU011421\
Data File : VU042035.D
Acq On : 14 Jan 2021 15:25
Operator : SY/MD
Sample : M1129-09
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BFT67

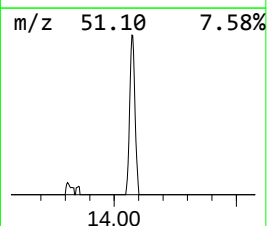
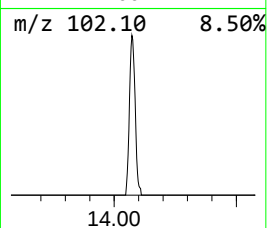
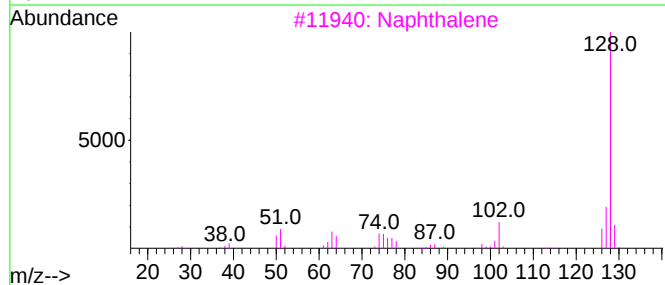
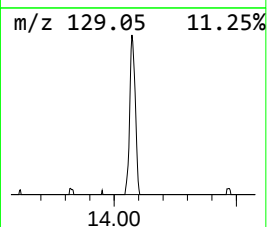
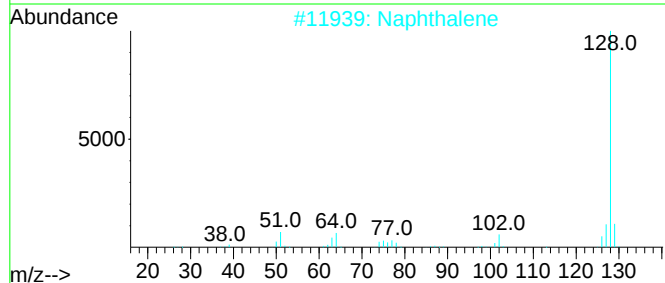
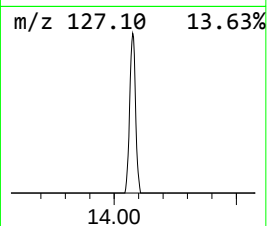
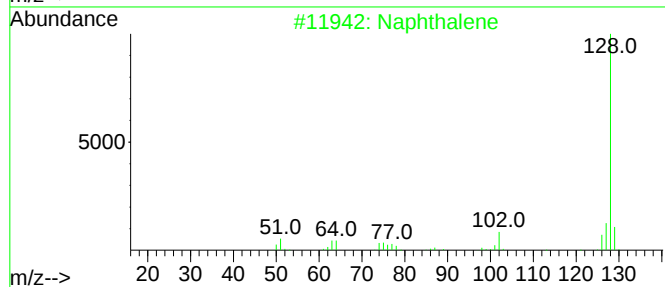
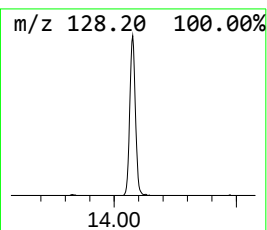
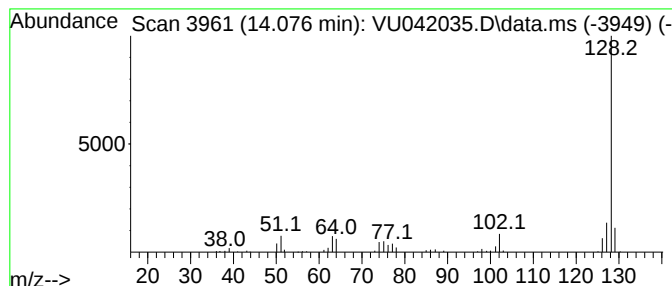
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Quant Title : TRACE VOA SOM01.0

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

Peak Number 14 Azulene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.076	1.59 ug/L	106365	1,4-Dichlorobenzene-d4	11.809

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene	128	C10H8	000091-20-3	95
2			Naphthalene	128	C10H8	000091-20-3	94
3			Naphthalene	128	C10H8	000091-20-3	94
4			Azulene	128	C10H8	000275-51-4	91
5			Azulene	128	C10H8	000275-51-4	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU011421\
Data File : VU042035.D
Acq On : 14 Jan 2021 15:25
Operator : SY/MD
Sample : M1129-09
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BFT67

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR011321WMA.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: S...	1.359	14.6	ug/L	844690	1	6.259	288762	5.0
(DEL) Alkane: S...	1.736	1.1	ug/L	61548	1	6.259	288762	5.0
(DEL) Alkane: S...	1.999	2.1	ug/L	119022	1	6.259	288762	5.0
(DEL) Alkane: C...	2.157	2.8	ug/L	160674	1	6.259	288762	5.0
(DEL) Alkane: S...	2.208	3.2	ug/L	186563	1	6.259	288762	5.0
(DEL) Alkane: S...	2.420	0.8	ug/L	47453	1	6.259	288762	5.0
(DEL) Alkane: S...	3.597	1.2	ug/L	68357	1	6.259	288762	5.0
(DEL) Alkane: S...	3.710	0.5	ug/L	31086	1	6.259	288762	5.0
unknown-01	9.767	0.5	ug/L	34688	2	9.420	336459	5.0
Octanal	11.687	0.6	ug/L	38972	3	11.809	334515	5.0
Nonanal	12.880	1.4	ug/L	93480	3	11.809	334515	5.0
Benzene, 1,2,3,...	12.963	0.7	ug/L	49336	3	11.809	334515	5.0
Azulene	14.076	1.6	ug/L	106365	3	11.809	334515	5.0