

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU021619\
 Data File : VU029505.D
 Acq On : 15 Feb 2019 18:55
 Operator : JC/SP
 Sample : K1438-16DL 4X
 Misc : 25.0mL/MSVOA U/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 ESXD8DL

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.1

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

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

Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR021219WMA.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	1.183	21	29	49	rBV	1824421	1693196	100.00%	18.307%
2	1.299	54	65	84	rBV	403066	399886	23.62%	4.324%
3	1.399	88	96	107	rBV2	773771	810176	47.85%	8.760%
4	1.679	177	183	198	rVB	92206	121861	7.20%	1.318%
5	1.756	199	207	225	rVB	126318	172282	10.17%	1.863%
6	1.945	258	266	276	rVB	83444	111533	6.59%	1.206%
7	2.273	357	368	382	rBV	188370	290527	17.16%	3.141%
8	2.698	493	500	509	rBV2	17784	31453	1.86%	0.340%
9	2.788	521	528	542	rVB2	28198	51764	3.06%	0.560%
10	2.981	581	588	604	rVB4	20220	42488	2.51%	0.459%
11	4.084	914	931	949	rVB4	19346	50658	2.99%	0.548%
12	4.219	955	973	1009	rBV2	203213	592627	35.00%	6.408%
13	4.646	1090	1106	1127	rBV	98814	241809	14.28%	2.614%
14	4.990	1192	1213	1228	rBV4	26005	63318	3.74%	0.685%
15	5.334	1299	1320	1348	rVB2	220129	595714	35.18%	6.441%
16	5.884	1474	1491	1515	rBV	199308	396016	23.39%	4.282%
17	6.328	1616	1629	1644	rBV	134169	271613	16.04%	2.937%
18	6.408	1644	1654	1668	rVB4	20514	44058	2.60%	0.476%
19	7.228	1896	1909	1933	rBV	77912	142683	8.43%	1.543%
20	7.563	1992	2013	2031	rBV	325137	570971	33.72%	6.173%
21	7.849	2092	2102	2119	rBV	46449	83700	4.94%	0.905%
22	8.315	2234	2247	2290	rBV	373591	687787	40.62%	7.436%
23	9.087	2471	2487	2509	rBV	300552	508340	30.02%	5.496%
24	10.431	2894	2905	2923	rBV	120565	193745	11.44%	2.095%
25	11.482	3220	3232	3268	rBV	317670	527463	31.15%	5.703%
26	11.858	3335	3349	3371	rBV	332110	553279	32.68%	5.982%

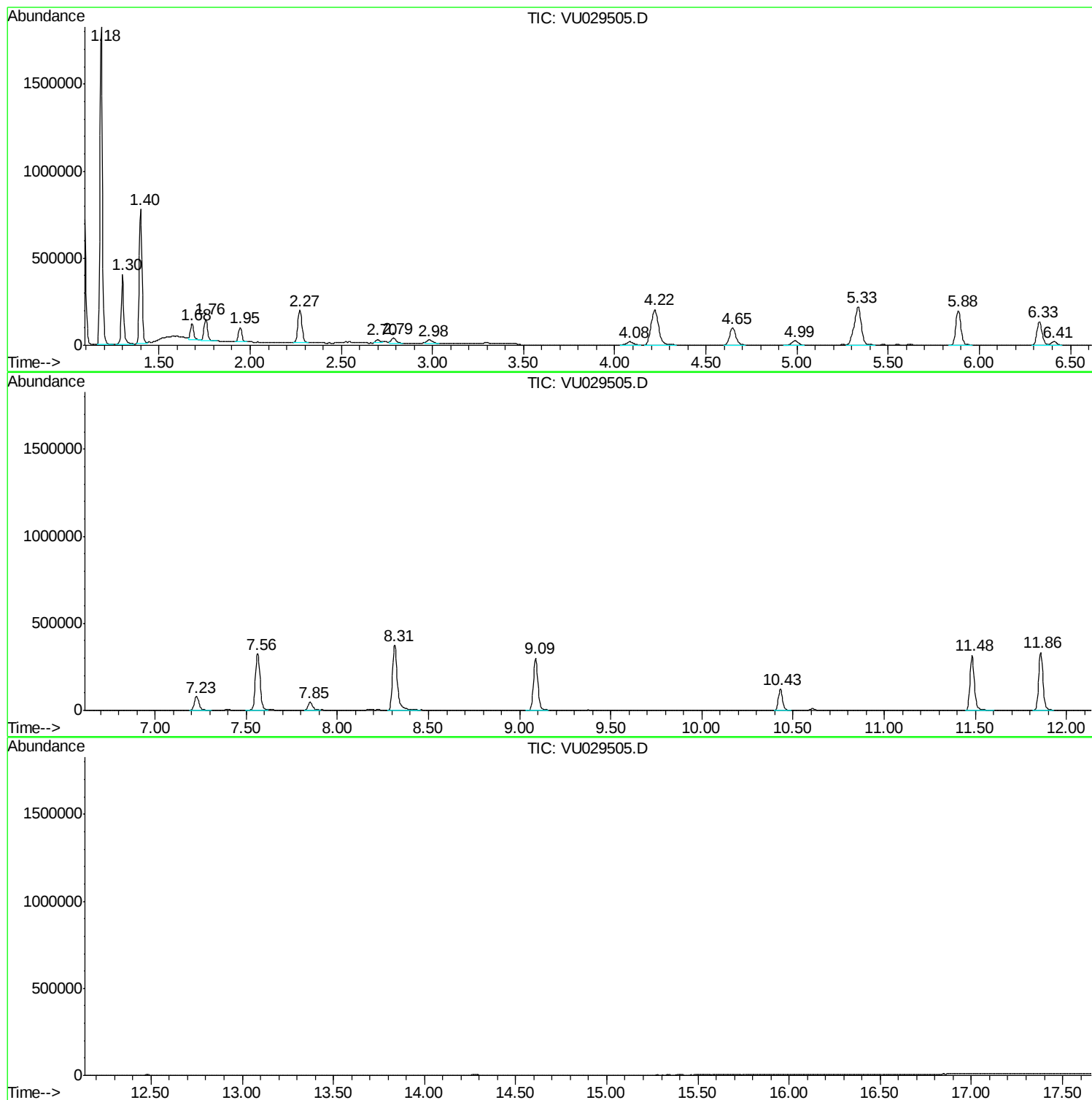
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Quant Title : TRACE VOA SOM01.0

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



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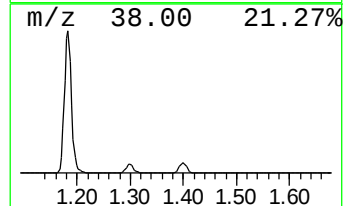
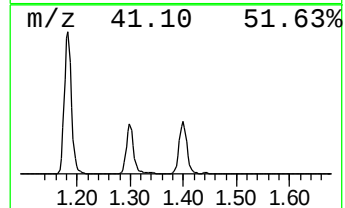
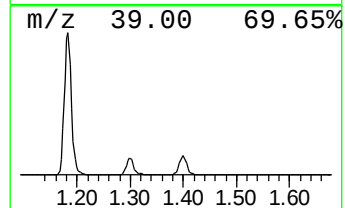
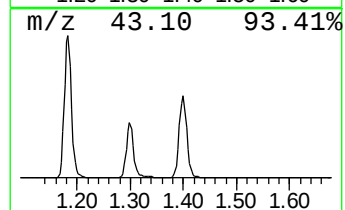
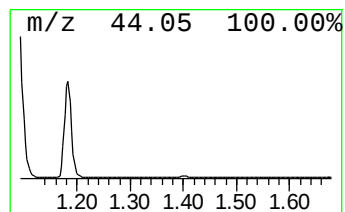
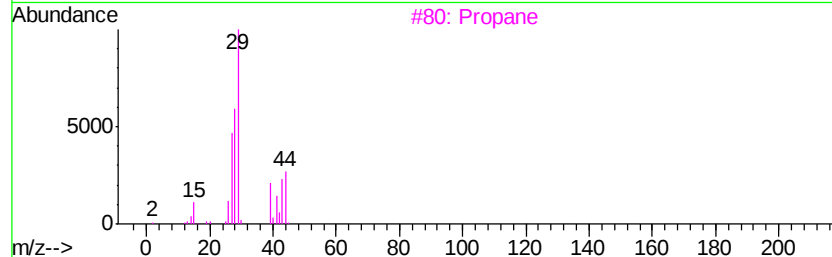
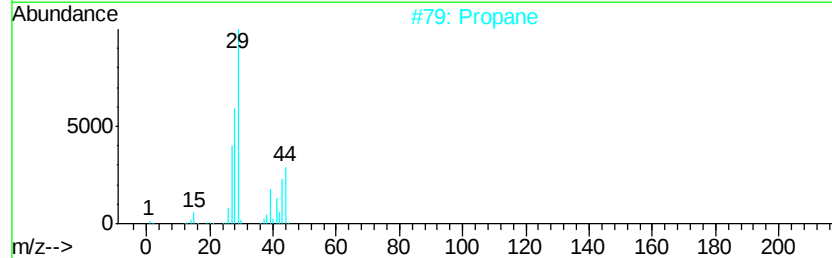
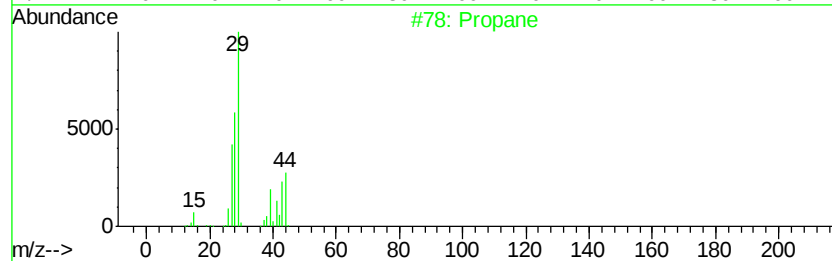
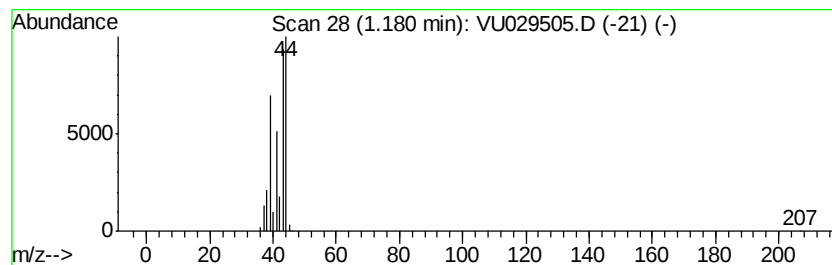
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Quant Title : TRACE VOA SOM01.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.18	21.38 ug/L	1693200	1,4-Difluorobenzene	5.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propane	44	C3H8	000074-98-6	83
2		Propane	44	C3H8	000074-98-6	42
3		Propane	44	C3H8	000074-98-6	9
4		Propene	42	C3H6	000115-07-1	9
5		Acetaldehyde	44	C2H4O	000075-07-0	5



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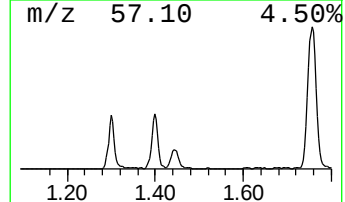
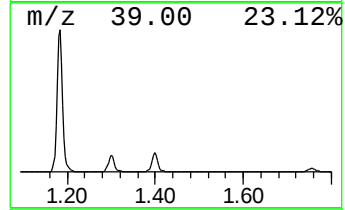
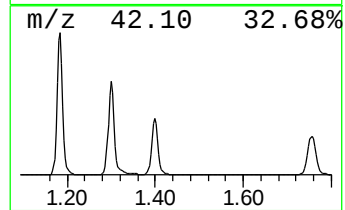
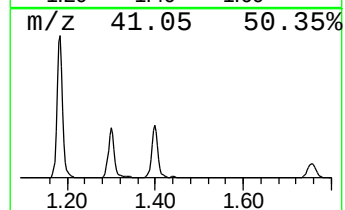
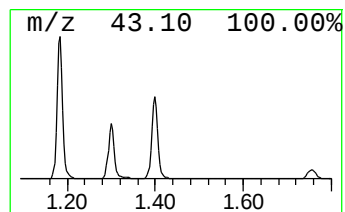
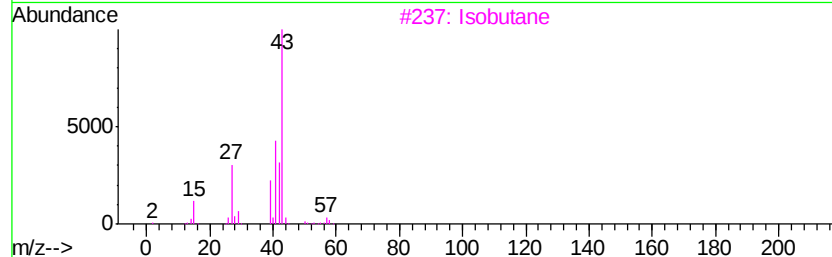
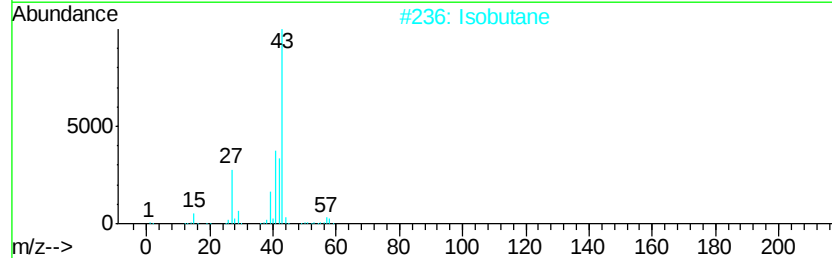
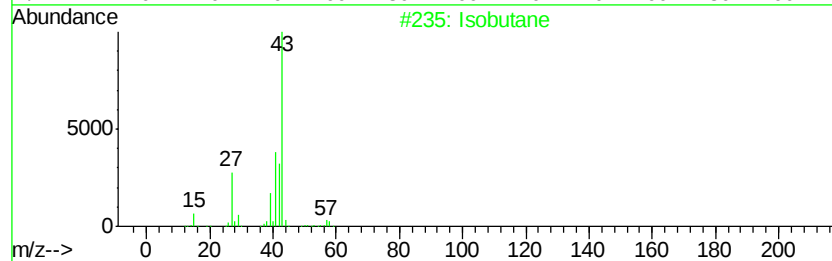
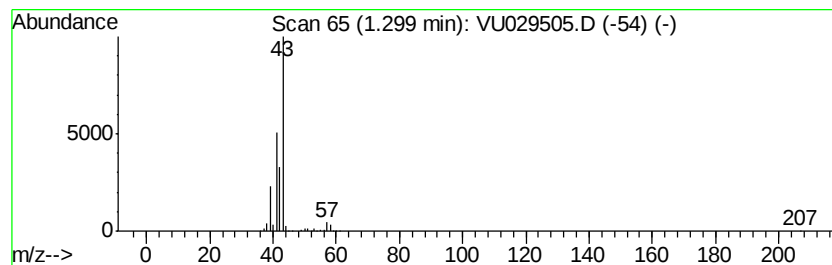
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 Quant Title : TRACE VOA SOM01.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.30	5.05 ug/L	399886	1,4-Difluorobenzene	5.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Isobutane	58	C4H10	000075-28-5	72
2		Isobutane	58	C4H10	000075-28-5	64
3		Isobutane	58	C4H10	000075-28-5	59
4		Isobutane	58	C4H10	000075-28-5	59
5		Propane, 2-nitro-	89	C3H7NO2	000079-46-9	9



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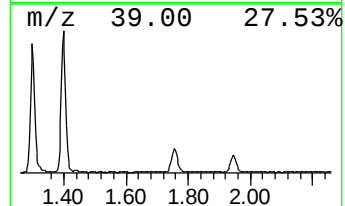
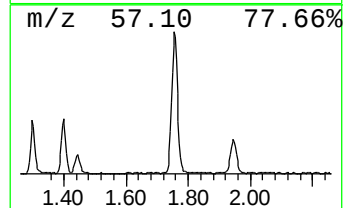
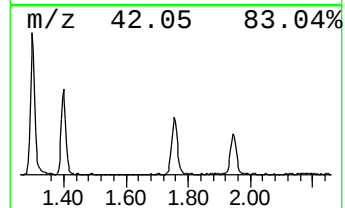
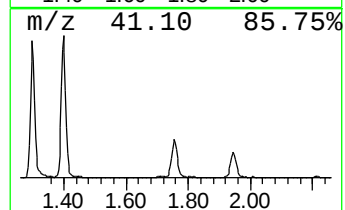
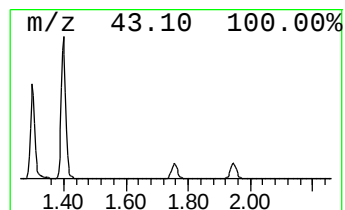
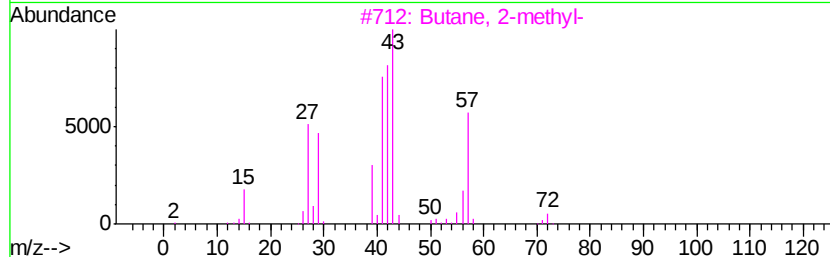
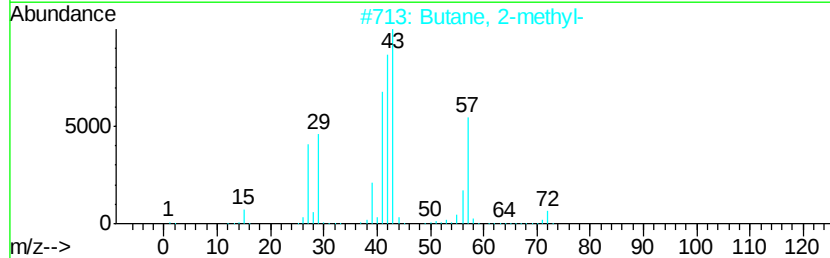
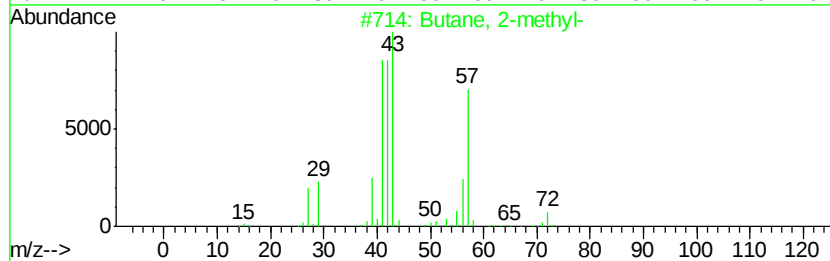
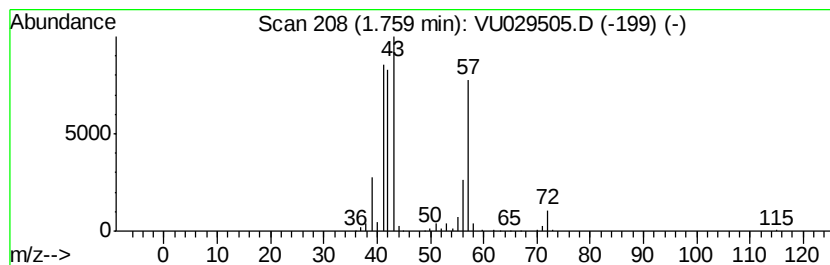
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.76	2.18 ug/L	172282	1,4-Difluorobenzene	5.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2-methyl-	72	C5H12	000078-78-4	94
2		Butane, 2-methyl-	72	C5H12	000078-78-4	91
3		Butane, 2-methyl-	72	C5H12	000078-78-4	86
4		Oxirane, trimethyl-	86	C5H10O	005076-19-7	40
5		Isobutylene epoxide	72	C4H8O	000558-30-5	38



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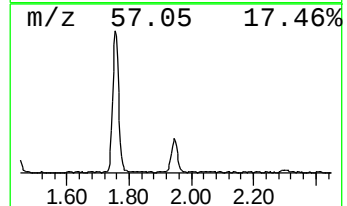
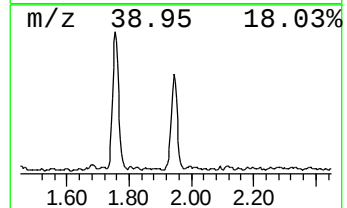
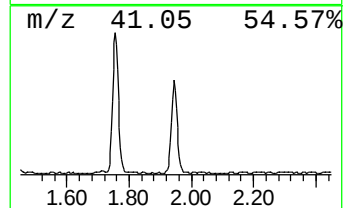
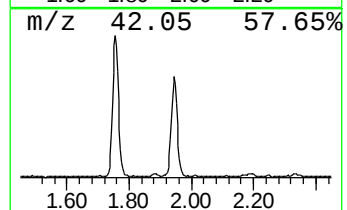
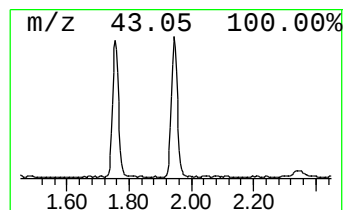
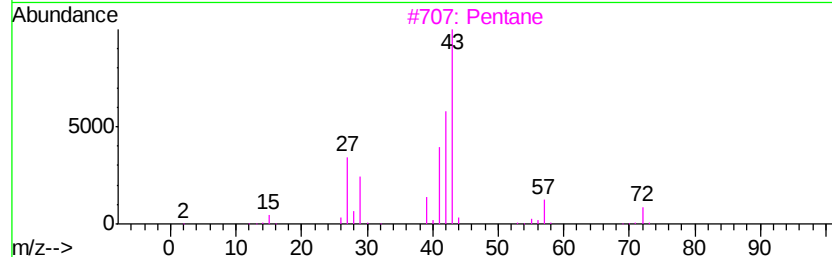
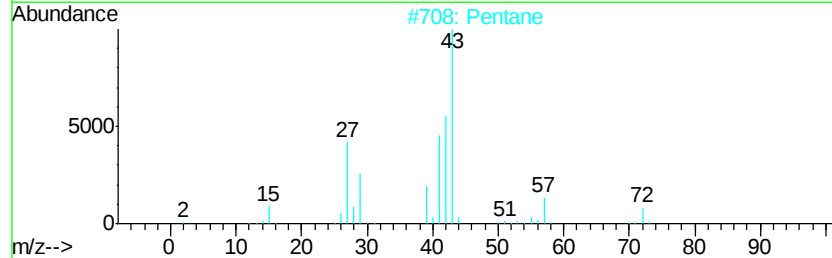
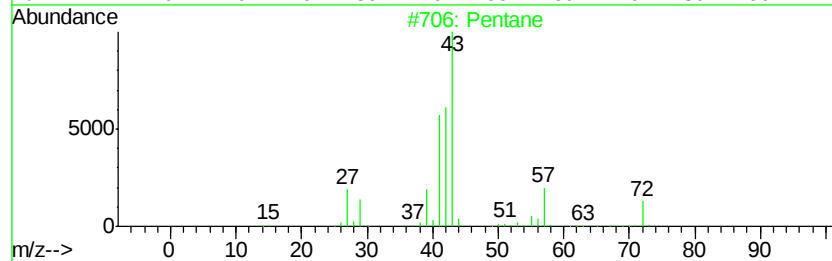
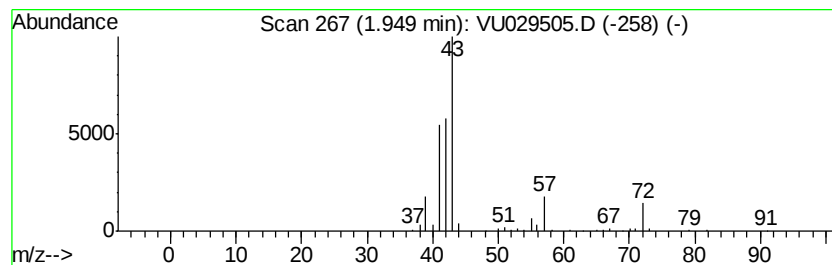
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.95	1.41 ug/L	111533	1,4-Difluorobenzene	5.88

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane		72	C5H12	000109-66-0	91
2	Pentane		72	C5H12	000109-66-0	90
3	Pentane		72	C5H12	000109-66-0	86
4	Pentane		72	C5H12	000109-66-0	86
5	Oxirane, ethyl-		72	C4H8O	000106-88-7	50



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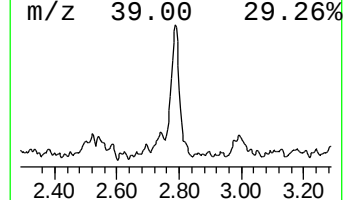
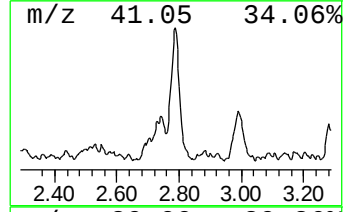
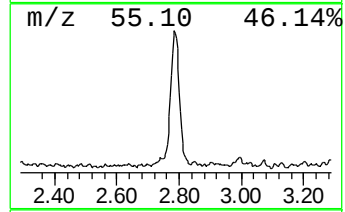
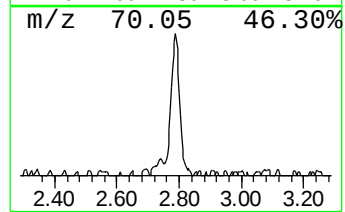
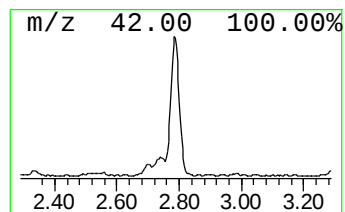
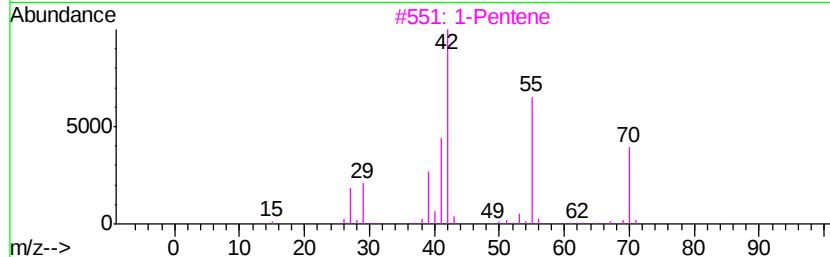
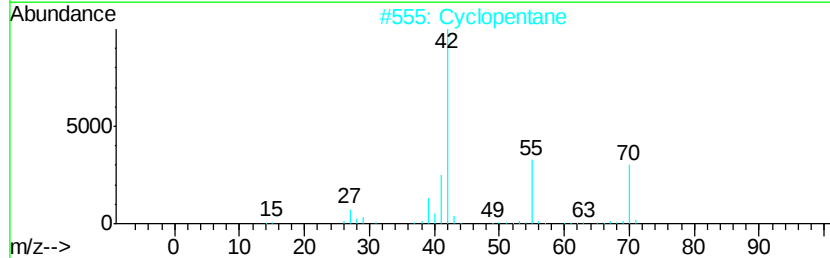
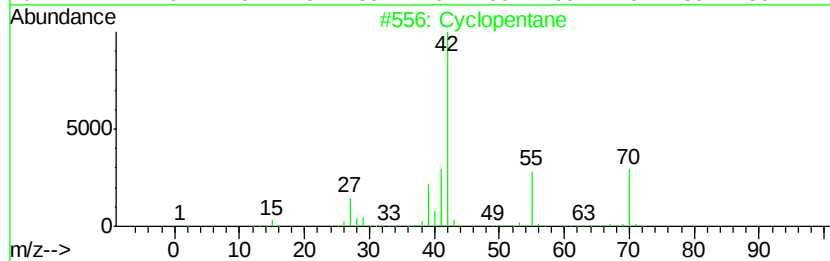
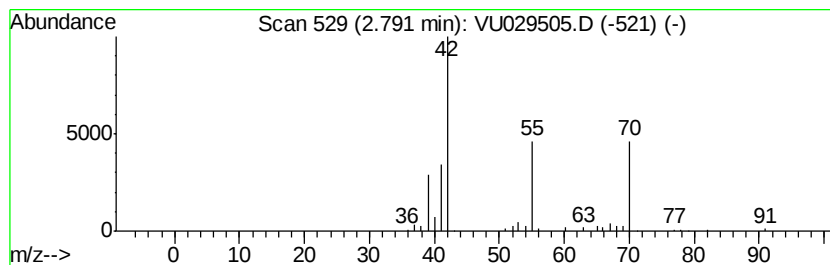
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR021219WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 5 (DEL) Alkane: Cyclic2.79 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.79	0.65 ug/L	51764	1,4-Difluorobenzene	5.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane	70	C5H10	000287-92-3	86
2		Cyclopentane	70	C5H10	000287-92-3	78
3		1-Pentene	70	C5H10	000109-67-1	64
4		1-Pentene	70	C5H10	000109-67-1	64
5		Cyclopropane, ethyl-	70	C5H10	001191-96-4	36



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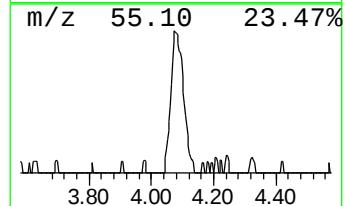
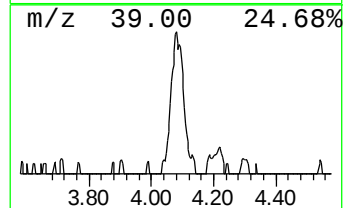
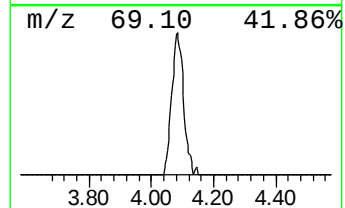
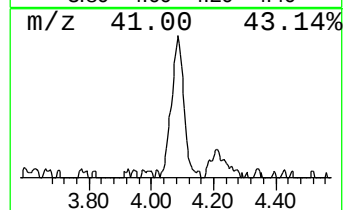
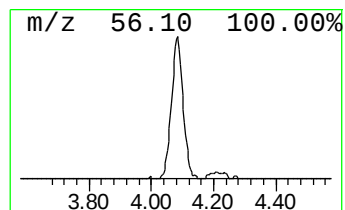
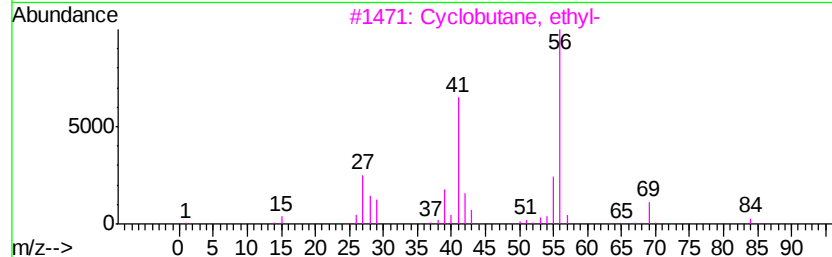
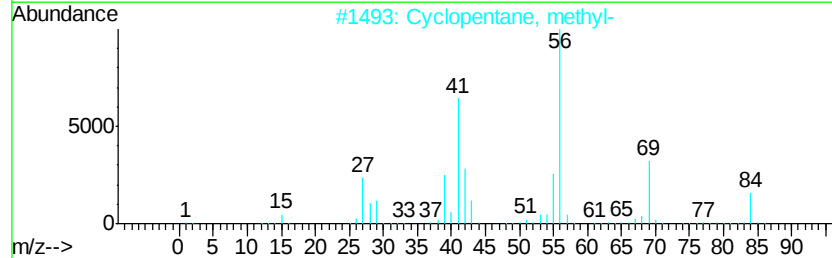
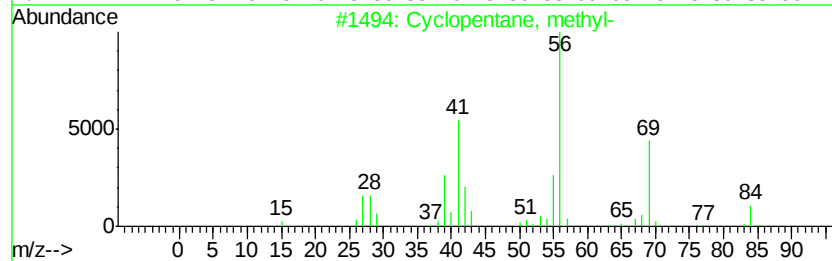
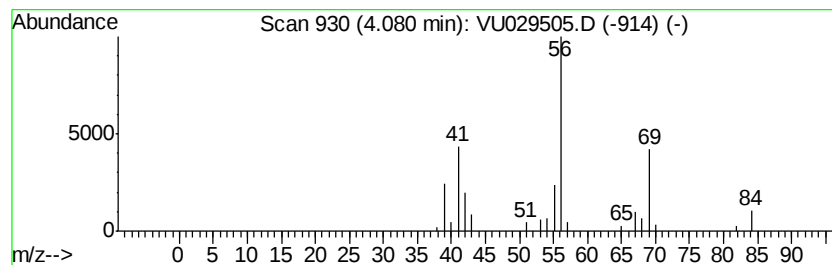
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR021219WMA.M
Quant Title : TRACE VOA SOM01.0

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

Peak Number 6 (DEL) Alkane: Cyclic4.08 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.08	0.64 ug/L	50658	1,4-Difluorobenzene	5.88

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2		Cyclopentane, methyl-	84	C6H12	000096-37-7	86
3		Cyclobutane, ethyl-	84	C6H12	004806-61-5	64
4		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	64
5		Cyclobutane, ethyl-	84	C6H12	004806-61-5	42



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU021619\
Data File : VU029505.D
Acq On : 15 Feb 2019 18:55
Operator : JC/SP
Sample : K1438-16DL 4X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
ESXD8DL

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR021219WMA.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: Str...	1.18	21.4	ug/L	1693200	1	5.88	396016	5.0	
(DEL) Alkane: Str...	1.30	5.0	ug/L	399886	1	5.88	396016	5.0	
(DEL) Alkane: Str...	1.76	2.2	ug/L	172282	1	5.88	396016	5.0	
(DEL) Alkane: Str...	1.95	1.4	ug/L	111533	1	5.88	396016	5.0	
(DEL) Alkane: Cyc...	2.79	0.7	ug/L	51764	1	5.88	396016	5.0	
(DEL) Alkane: Cyc...	4.08	0.6	ug/L	50658	1	5.88	396016	5.0	