

Data Path : Z:\VOASRV\HPCHEM1\MSVOA Y\DATA\VY050820\
 Data File : VY002641.D
 Acq On : 08 May 2020 14:37
 Operator : SY/MD
 Sample : L2554-01
 Misc : 6.41G/5ML/MSVOA Y/SOIL
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 DRUM-1-7

Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 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_Y\METHODS\82Y043020S.M
 Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.871	17	22	35	rVB	287362	393073	30.05%	4.543%
2	2.231	74	81	93	rBV	156748	280305	21.43%	3.240%
3	2.353	96	101	112	rVB	30178	51482	3.94%	0.595%
4	2.469	112	120	128	rBV	16121	30116	2.30%	0.348%
5	2.767	161	169	175	rBV	20340	42908	3.28%	0.496%
6	2.920	188	194	203	rVB5	5746	13555	1.04%	0.157%
7	3.188	229	238	245	rBV	50024	109231	8.35%	1.262%
8	3.292	246	255	267	rVB3	27204	82494	6.31%	0.953%
9	3.462	275	283	293	rVB	16292	37026	2.83%	0.428%
10	3.621	301	309	318	rBV5	7387	19345	1.48%	0.224%
11	3.718	318	325	339	rVB3	10807	29194	2.23%	0.337%
12	3.968	357	366	385	rBV	68877	190940	14.60%	2.207%
13	4.596	457	469	481	rBV6	23030	99511	7.61%	1.150%
14	4.743	485	493	509	rBV4	10461	44664	3.41%	0.516%
15	5.614	621	636	648	rBV3	20050	79364	6.07%	0.917%
16	5.767	652	661	673	rVB3	8968	27949	2.14%	0.323%
17	6.066	700	710	714	rBV4	4563	13242	1.01%	0.153%
18	6.120	714	719	730	rVB3	5712	18150	1.39%	0.210%
19	6.352	750	757	766	rBV3	8293	25697	1.96%	0.297%
20	6.822	824	834	843	rVB6	5024	16256	1.24%	0.188%
21	7.017	854	866	875	rBV	17701	48781	3.73%	0.564%
22	7.553	948	954	961	rVB3	7688	17557	1.34%	0.203%
23	7.742	976	985	990	rBV2	183927	425761	32.55%	4.921%
24	7.815	990	997	1009	rVV	362636	852200	65.14%	9.849%
25	7.901	1009	1011	1021	rVB2	10926	21754	1.66%	0.251%
26	8.163	1044	1054	1065	rBV	162917	363414	27.78%	4.200%
27	8.309	1069	1078	1085	rBV3	15399	34673	2.65%	0.401%
28	8.449	1095	1101	1110	rVB2	15383	33498	2.56%	0.387%
29	8.565	1113	1120	1130	rVB5	5559	14884	1.14%	0.172%
30	8.705	1135	1143	1158	rBV	505001	1029573	78.70%	11.899%
31	8.980	1177	1188	1195	rBV4	8635	19604	1.50%	0.227%
32	9.150	1208	1216	1220	rBV7	7015	16260	1.24%	0.188%
33	9.199	1220	1224	1230	rVB4	8015	16984	1.30%	0.196%
34	9.577	1281	1286	1291	rVV3	12806	29045	2.22%	0.336%

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35	9.754	1307	1315	1322	rBV7	12803	28633	2.19%	0.331%
36	10.065	1362	1366	1373	rVB5	6766	13513	1.03%	0.156%
37	10.193	1380	1387	1393	rBV	773703	1308213	100.00%	15.120%
38	10.254	1393	1397	1403	rVV	42853	79373	6.07%	0.917%
39	10.382	1411	1418	1424	rBV6	6641	13378	1.02%	0.155%
40	10.620	1447	1457	1468	rVB8	5402	15682	1.20%	0.181%
41	11.309	1565	1570	1580	rVB8	9418	29451	2.25%	0.340%
42	11.504	1592	1602	1614	rBV	605509	1005164	76.83%	11.617%
43	11.607	1614	1619	1622	rBV3	7998	15410	1.18%	0.178%
44	11.711	1631	1636	1647	rVB3	26586	66393	5.08%	0.767%
45	12.034	1681	1689	1695	rVB5	8532	20614	1.58%	0.238%
46	12.314	1729	1735	1742	rBV4	18170	32359	2.47%	0.374%
47	12.491	1757	1764	1773	rBV	428210	676913	51.74%	7.824%
48	12.814	1813	1817	1823	rVB5	11623	19737	1.51%	0.228%
49	13.125	1863	1868	1874	rBV	8845	15897	1.22%	0.184%
50	13.278	1887	1893	1899	rBV10	8005	15313	1.17%	0.177%
51	13.436	1912	1919	1928	rVB	444776	676750	51.73%	7.822%
52	13.747	1966	1970	1976	rVB5	9405	19135	1.46%	0.221%
53	13.973	2004	2007	2014	rVB5	9892	16044	1.23%	0.185%
54	14.997	2170	2175	2183	rBV6	14976	28676	2.19%	0.331%
55	15.978	2331	2336	2345	rBV4	14778	27159	2.08%	0.314%

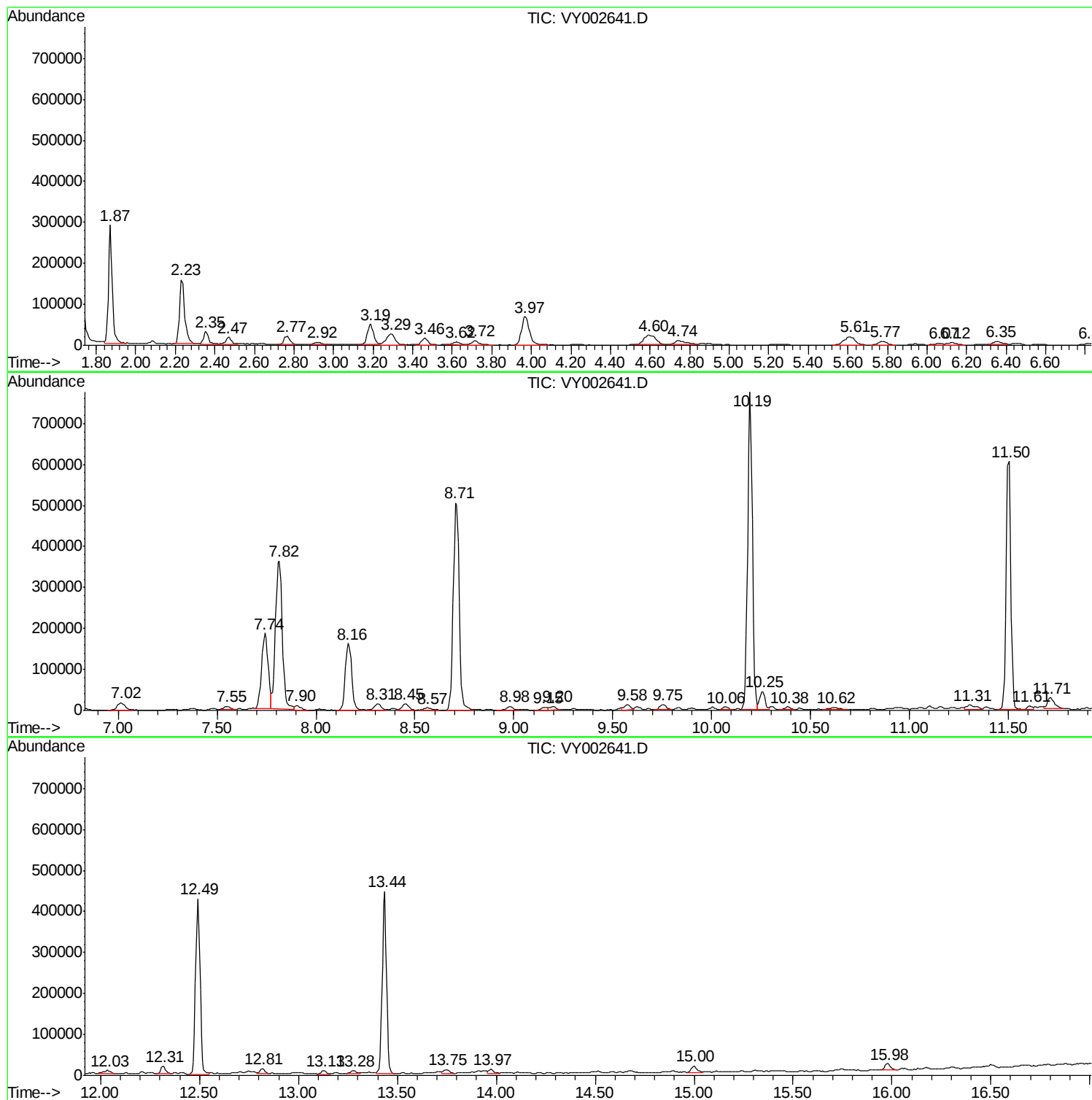
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y043020S.M
Quant Title : SW846 8260

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



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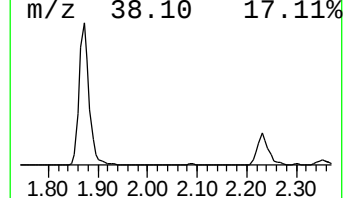
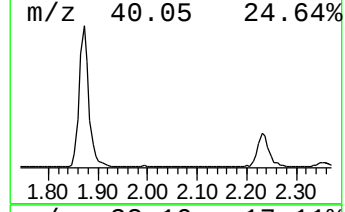
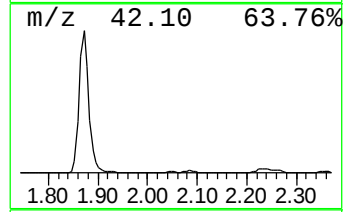
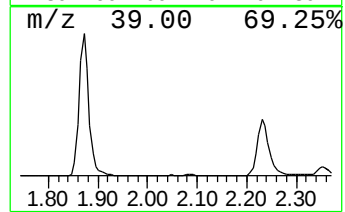
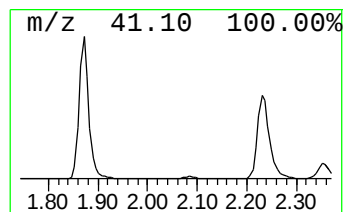
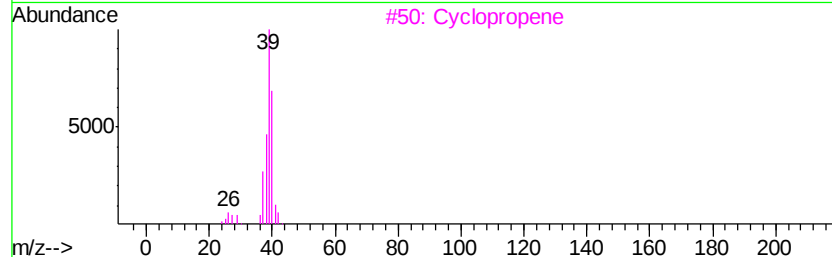
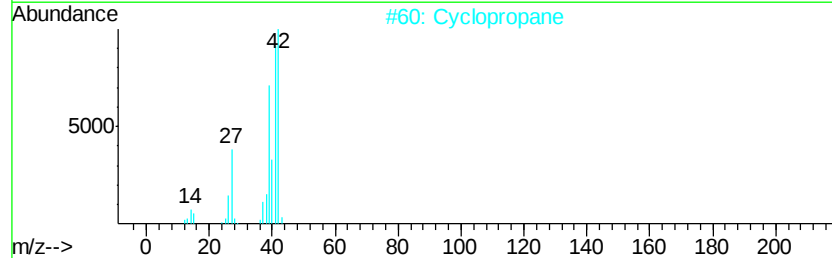
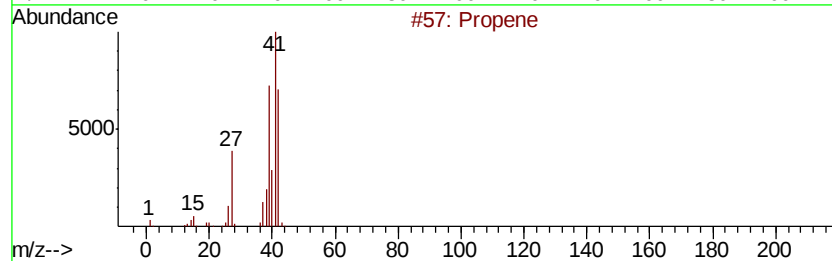
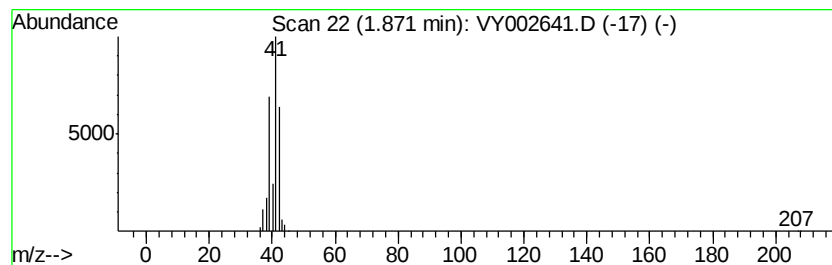
Instrument :
MSVOA_Y
ClientSampleId :
DRUM-1-7

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y043020S.M
Quant Title : SW846 8260

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

Peak Number 1 Propene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.		
1.87	23.06 ug/l	393073	Pentafluorobenzene	7.82		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propene		42	C3H6	000115-07-1	86
2	Cyclopropane		42	C3H6	000075-19-4	78
3	Cyclopropene		40	C3H4	002781-85-3	16
4	Cyanamide		42	CH2N2	000420-04-2	4
5	2-Oxetanone, 4-methyl-		86	C4H6O2	003068-88-0	4



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Misc : 6.41G/5ML/MSVOA Y/SOIL
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_Y
ClientSampleId :
DRUM-1-7

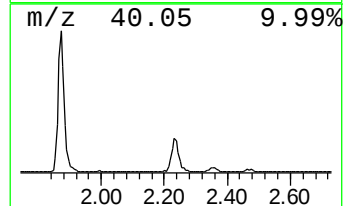
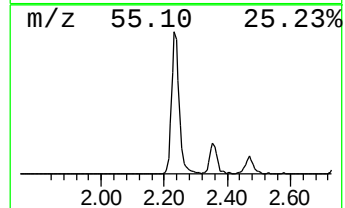
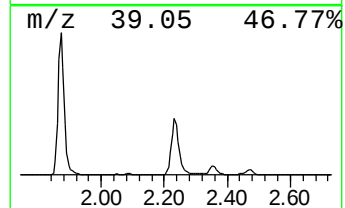
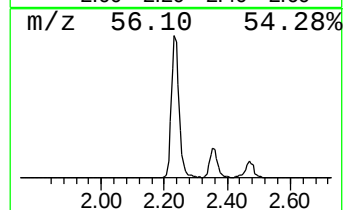
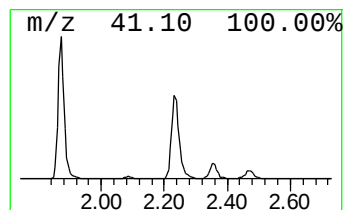
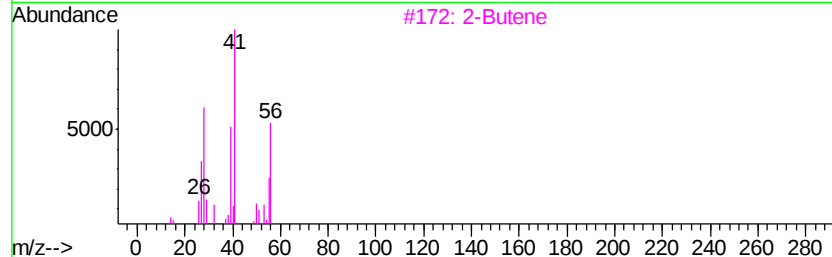
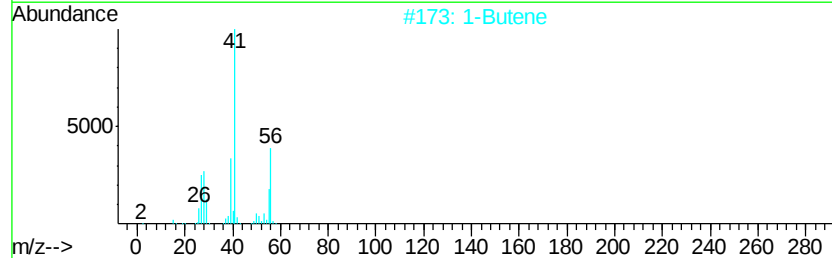
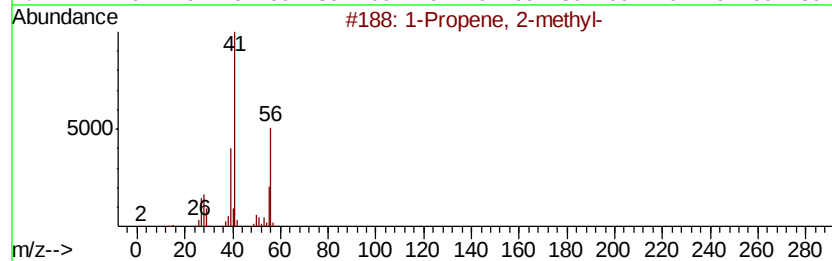
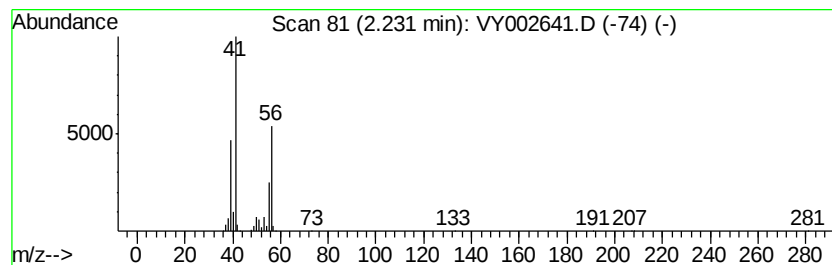
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_Y\METHODS\82Y043020S.M
Quant Title : SW846 8260

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

Peak Number 2 1-Propene, 2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.23	16.45 ug/l	280305	Pentafluorobenzene	7.82

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Propene, 2-methyl-	56	C4H8	000115-11-7	90
2		1-Butene	56	C4H8	000106-98-9	87
3		2-Butene	56	C4H8	000107-01-7	83
4		2-Butene, (E)-	56	C4H8	000624-64-6	80
5		2-Butene, (Z)-	56	C4H8	000590-18-1	49



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Misc : 6.41G/5ML/MSVOA Y/SOIL
ALS Vial : 10 Sample Multiplier: 1

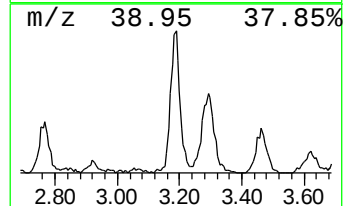
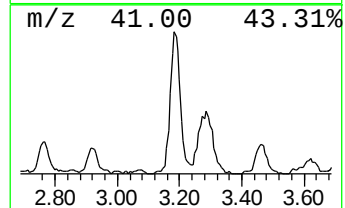
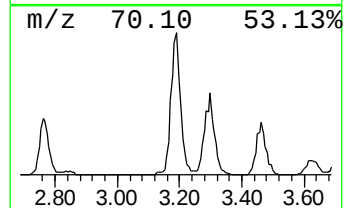
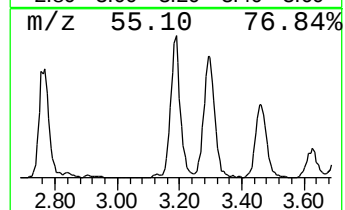
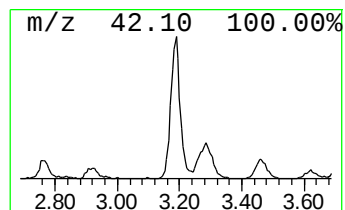
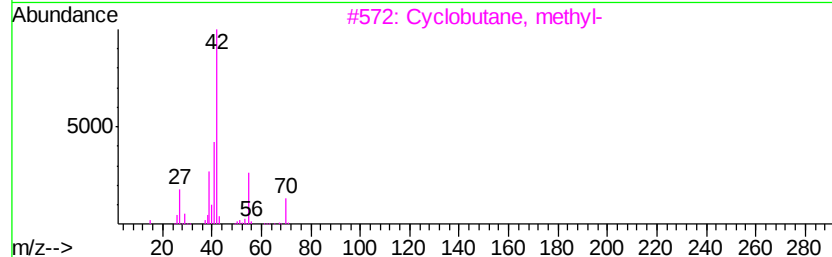
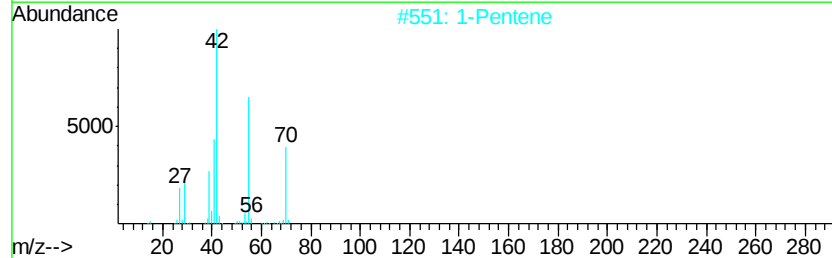
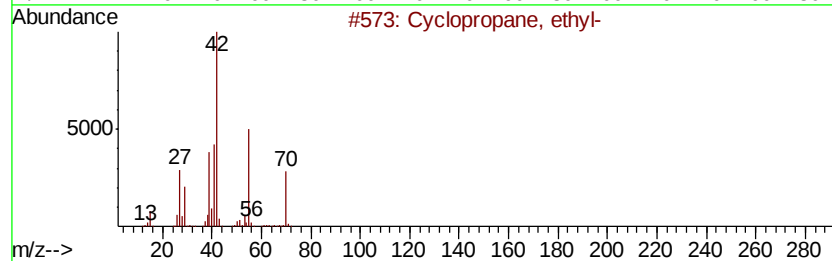
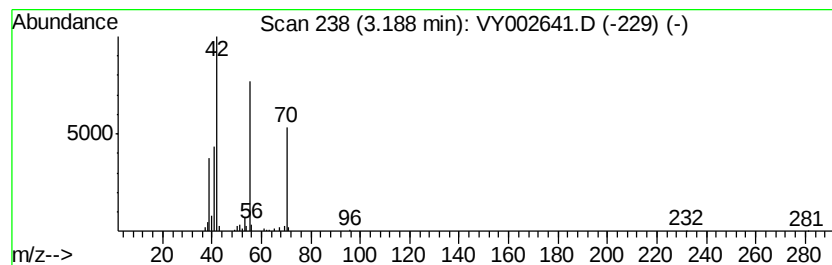
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 3 Cyclopropane, ethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.		
3.19	6.41 ug/l	109231	Pentafluorobenzene	7.82		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cvclopropane, ethvl-	70	C5H10	001191-96-4	91
2		1-Pentene	70	C5H10	000109-67-1	91
3		Cvclobutane, methvl-	70	C5H10	000598-61-8	90
4		Cvclopropane, 1,2-dimethvl-, cis-	70	C5H10	000930-18-7	60
5		2-Pentene	70	C5H10	000109-68-2	47



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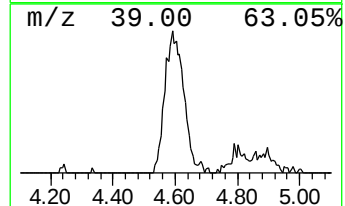
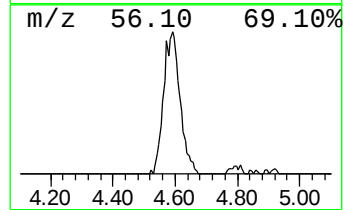
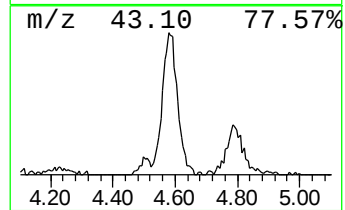
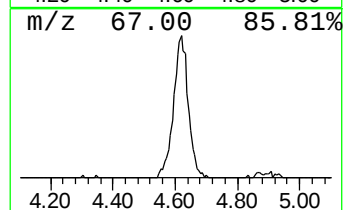
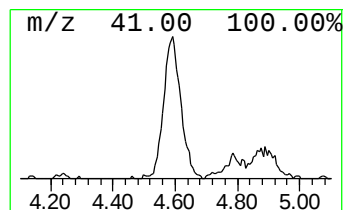
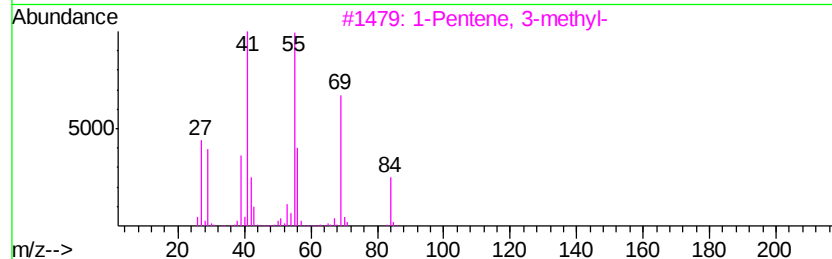
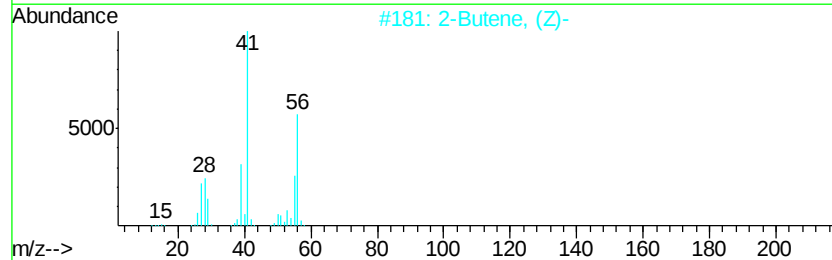
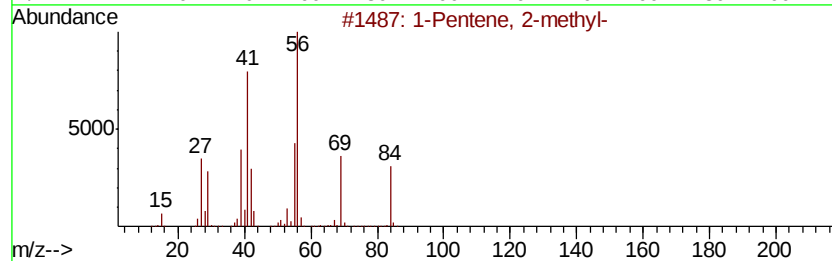
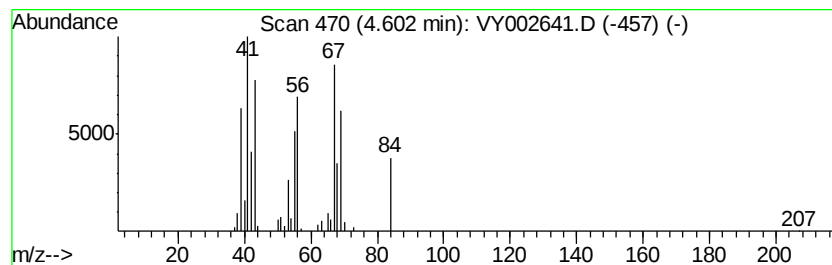
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 4 unknown4.60 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.60	5.84 ug/l	99511	Pentafluorobenzene	7.82

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Pentene, 2-methyl-	84	C6H12	000763-29-1	38
2			2-Butene, (Z)-	56	C4H8	000590-18-1	35
3			1-Pentene, 3-methyl-	84	C6H12	000760-20-3	35
4			Bicyclo[2.1.0]pentane	68	C5H8	000185-94-4	30
5			2-Butene, (E)-	56	C4H8	000624-64-6	30



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propene	1.87	23.1	ug/l	393073	1	7.82	852200	50.0
1-Propene, 2-methyl-	2.23	16.4	ug/l	280305	1	7.82	852200	50.0
Cyclopropane, ethyl-	3.19	6.4	ug/l	109231	1	7.82	852200	50.0
unknown4.60	4.60	5.8	ug/l	99511	1	7.82	852200	50.0