

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX061920\
 Data File : VX016909.D
 Acq On : 20 Jun 2020 05:30
 Operator : JC/SP
 Sample : L3063-08
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TW-12

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_X\METHOD\82X061820W.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	1.160	8	12	22	rVB	143948	138117	8.34%	0.850%
2	1.380	44	48	56	rBV2	72433	124950	7.54%	0.769%
3	1.776	108	113	123	rVB	174715	236025	14.25%	1.453%
4	1.935	135	139	143	rBV	19620	27642	1.67%	0.170%
5	1.983	143	147	157	rVB2	43743	74977	4.53%	0.462%
6	2.252	186	191	201	rVB	40452	72887	4.40%	0.449%
7	2.380	201	212	216	rVV	41119	87941	5.31%	0.541%
8	2.410	216	217	228	rVB	20867	31813	1.92%	0.196%
9	2.575	234	244	253	rBV2	22275	45244	2.73%	0.279%
10	2.825	268	285	289	rBV2	40806	111964	6.76%	0.689%
11	2.867	289	292	295	rVV	33136	61806	3.73%	0.380%
12	2.916	295	300	308	rVV	65937	138611	8.37%	0.853%
13	3.008	308	315	329	rVB	55363	126318	7.62%	0.778%
14	3.160	331	340	351	rBV3	97674	250625	15.13%	1.543%
15	4.282	513	524	531	rBV3	10792	33126	2.00%	0.204%
16	4.373	531	539	552	rVV2	69669	200725	12.12%	1.236%
17	4.501	552	560	571	rVV6	10000	32900	1.99%	0.203%
18	5.275	666	687	702	rVB7	15706	67863	4.10%	0.418%
19	5.464	702	718	727	rBV	178568	534980	32.29%	3.293%
20	5.556	727	733	737	rVV2	18322	40384	2.44%	0.249%
21	5.629	737	745	755	rVB	295155	755803	45.62%	4.652%
22	5.903	780	790	801	rBV6	9177	30061	1.81%	0.185%
23	6.031	802	811	818	rVV	175851	468501	28.28%	2.884%
24	6.111	818	824	837	rVV	209291	563484	34.01%	3.469%
25	6.233	837	844	849	rVV6	10910	33984	2.05%	0.209%
26	6.318	850	858	869	rVV3	33766	111967	6.76%	0.689%
27	6.434	869	877	888	rVB7	11445	33109	2.00%	0.204%
28	6.830	931	942	953	rBV	418654	1020288	61.58%	6.280%
29	7.159	986	996	1002	rBV6	12245	38297	2.31%	0.236%
30	7.232	1003	1008	1016	rVB5	16002	38059	2.30%	0.234%
31	7.440	1031	1042	1051	rVB4	26295	74151	4.48%	0.456%
32	8.037	1132	1140	1149	rVB	29745	62610	3.78%	0.385%
33	8.159	1151	1160	1164	rBV	63791	140230	8.46%	0.863%
34	8.244	1170	1174	1179	rVB4	11106	19204	1.16%	0.118%

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35	8.555	1218	1225	1232	rVB6	12508	27201	1.64%	0.167%
36	8.696	1232	1248	1255	rBV	919359	1460432	88.15%	8.990%
37	8.769	1255	1260	1265	rVB	25216	42744	2.58%	0.263%
38	9.208	1327	1332	1337	rBV	12211	19521	1.18%	0.120%
39	9.287	1341	1345	1352	rVB2	10589	18130	1.09%	0.112%
40	9.500	1371	1380	1384	rBV	32428	54219	3.27%	0.334%
41	9.561	1384	1390	1394	rVB4	9703	16898	1.02%	0.104%
42	10.098	1472	1478	1488	rBV	1145063	1566313	94.54%	9.642%
43	10.348	1511	1519	1524	rVB	76028	107745	6.50%	0.663%
44	11.000	1620	1626	1632	rBV	487622	646691	39.03%	3.981%
45	11.122	1640	1646	1656	rBV	1125587	1406942	84.92%	8.661%
46	11.341	1677	1682	1692	rVV	399339	503501	30.39%	3.099%
47	11.439	1692	1698	1703	rVV	33788	45716	2.76%	0.281%
48	11.494	1703	1707	1714	rVB	19230	26312	1.59%	0.162%
49	11.652	1728	1733	1739	rBV	112531	138896	8.38%	0.855%
50	11.793	1751	1756	1762	rVB	247426	294035	17.75%	1.810%
51	11.902	1768	1774	1776	rBV	19608	25498	1.54%	0.157%
52	11.927	1776	1778	1782	rVB	23774	28691	1.73%	0.177%
53	12.061	1794	1800	1806	rBV	1338640	1656775	100.00%	10.198%
54	12.128	1806	1811	1817	rVB	115667	144074	8.70%	0.887%
55	12.286	1828	1837	1843	rBV	588514	806446	48.68%	4.964%
56	12.353	1843	1848	1858	rVB2	59951	102202	6.17%	0.629%
57	12.445	1858	1863	1867	rBV	22962	28308	1.71%	0.174%
58	12.500	1867	1872	1877	rVB	47454	61995	3.74%	0.382%
59	12.573	1877	1884	1887	rBV	37776	53510	3.23%	0.329%
60	12.652	1891	1897	1900	rVV	150988	180702	10.91%	1.112%
61	12.683	1900	1902	1906	rVB	22386	26878	1.62%	0.165%
62	12.737	1906	1911	1919	rBV2	69920	103705	6.26%	0.638%
63	12.890	1931	1936	1941	rBV4	21855	29710	1.79%	0.183%
64	12.963	1941	1948	1951	rVV	95285	124183	7.50%	0.764%
65	13.000	1951	1954	1960	rVB	70012	87338	5.27%	0.538%
66	13.152	1971	1979	1984	rBV6	10432	25859	1.56%	0.159%
67	13.207	1984	1988	1998	rVB	57921	74300	4.48%	0.457%
68	13.329	2004	2008	2014	rVB2	140389	192167	11.60%	1.183%
69	13.463	2025	2030	2036	rBV2	17119	22640	1.37%	0.139%
70	13.621	2052	2056	2062	rVB4	17139	31554	1.90%	0.194%
71	13.701	2067	2069	2078	rVB4	12929	27016	1.63%	0.166%

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72	13.817	2082	2088	2094	rBV	131123	168866	10.19%	1.039%
73	14.255	2153	2160	2165	rBV3	10217	20542	1.24%	0.126%
74	14.822	2248	2253	2258	rVB	14033	18596	1.12%	0.114%

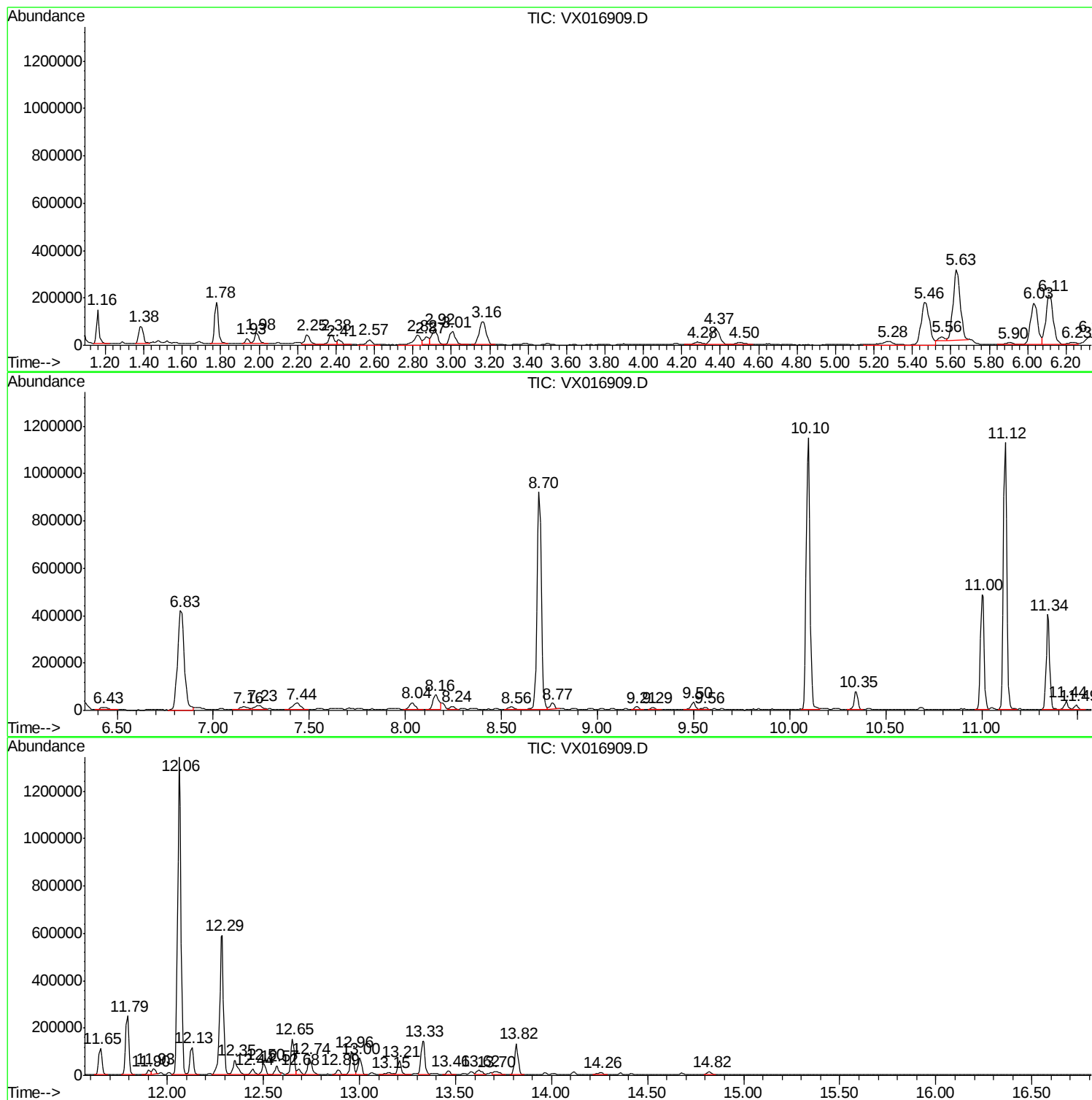
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061820W.M
Quant Title : SW846 8260

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



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Instrument :
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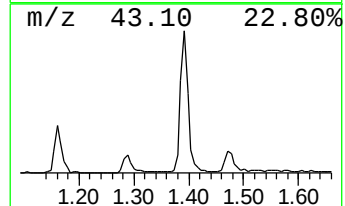
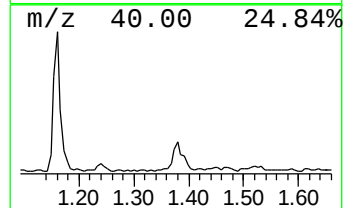
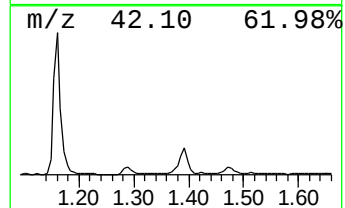
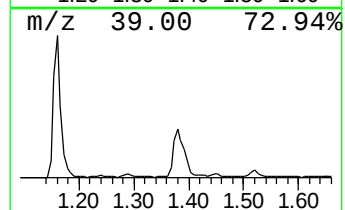
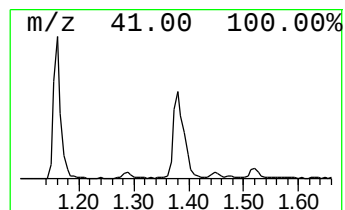
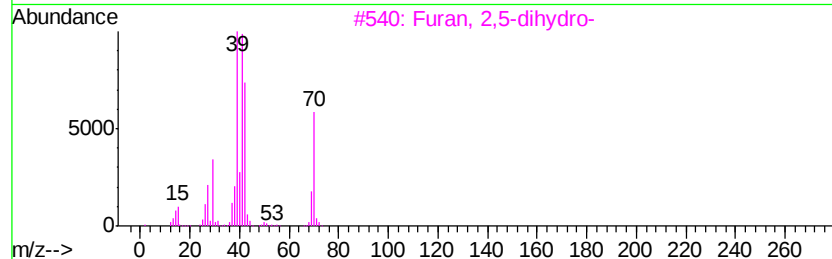
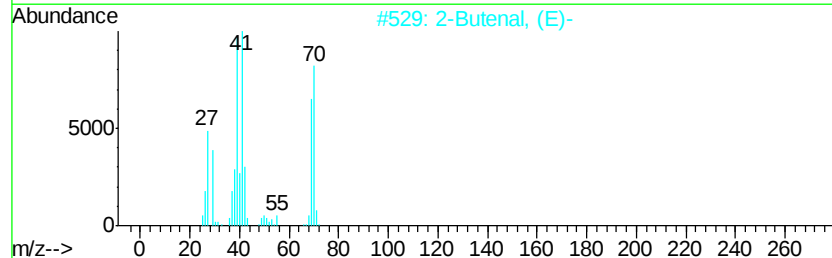
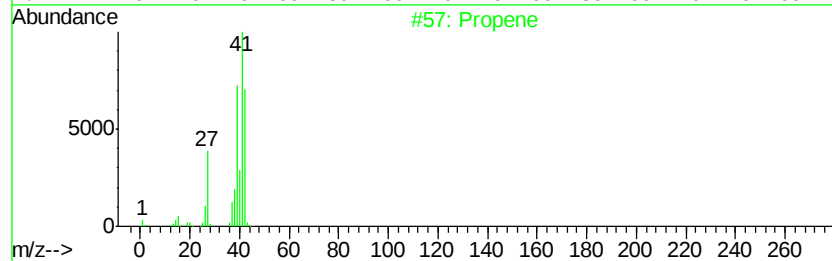
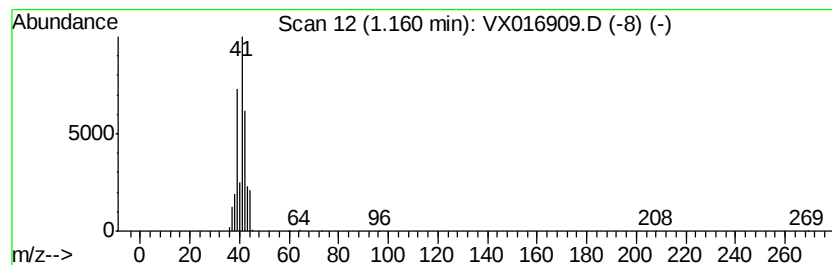
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061820W.M
Quant Title : SW846 8260

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

Peak Number 1 Propene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.16	9.14 ug/l	138117	Pentafluorobenzene	5.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Propene	42	C3H6	000115-07-1	80
2		2-Butenal, (E)-	70	C4H6O	000123-73-9	64
3		Furan, 2,5-dihydro-	70	C4H6O	001708-29-8	56
4		Cyclopropane	42	C3H6	000075-19-4	50
5		Dimethylketene	70	C4H6O	000598-26-5	39



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Instrument :
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ClientSampled :
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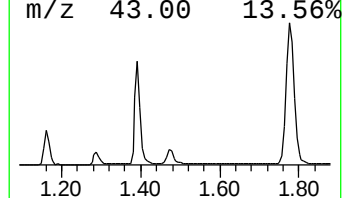
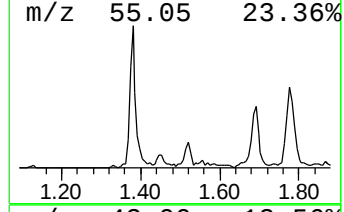
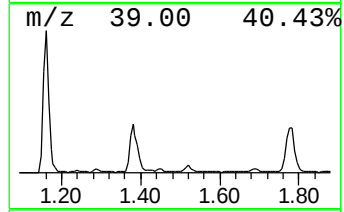
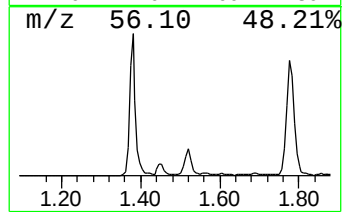
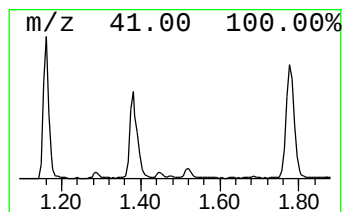
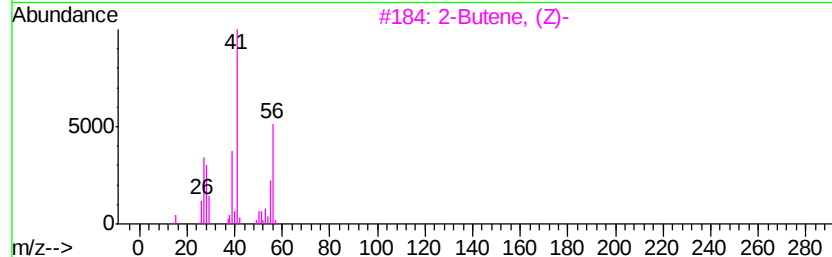
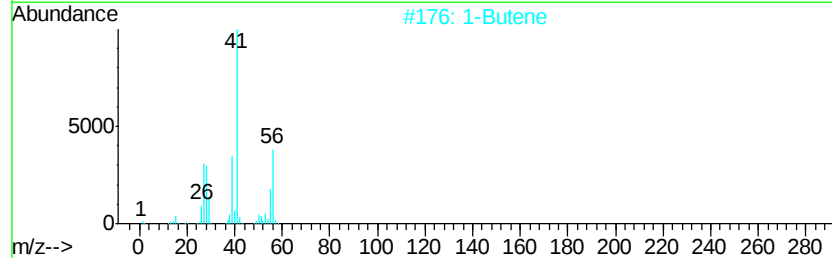
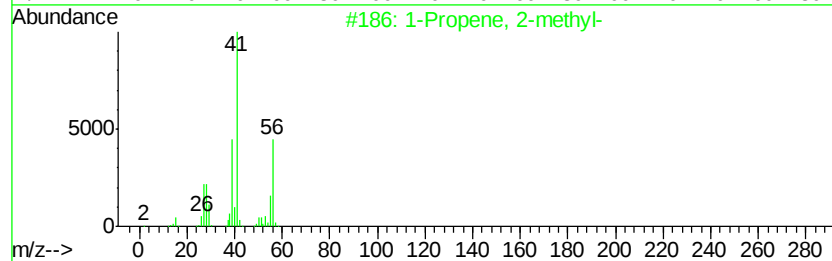
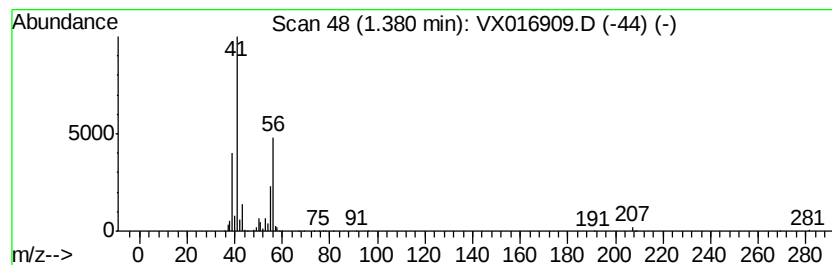
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061820W.M
Quant Title : SW846 8260

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

Peak Number 2 1-Propene, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.38	8.27 ug/l	124950	Pentafluorobenzene	5.63

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, (Z)-	56	C4H8	000590-18-1	87
4		2-Butene, (E)-	56	C4H8	000624-64-6	86
5		2-Butene	56	C4H8	000107-01-7	78



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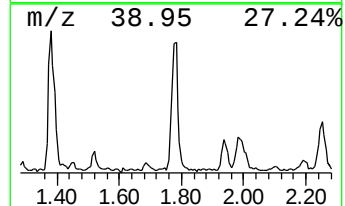
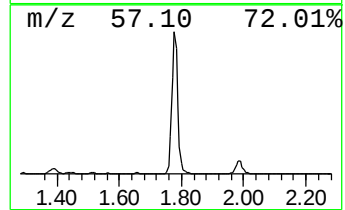
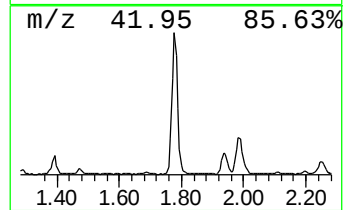
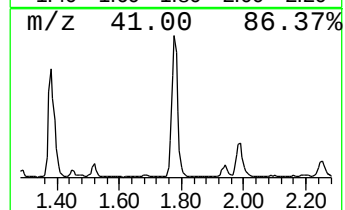
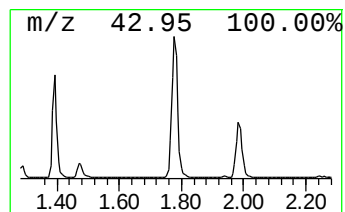
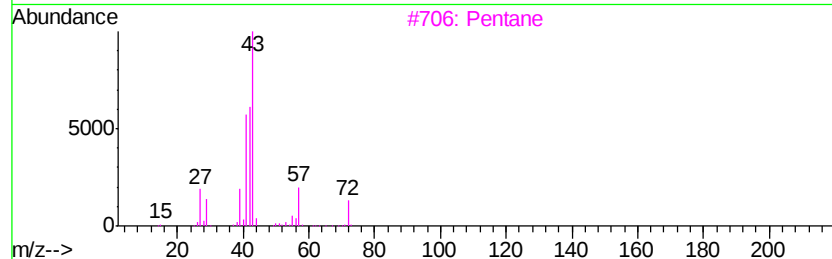
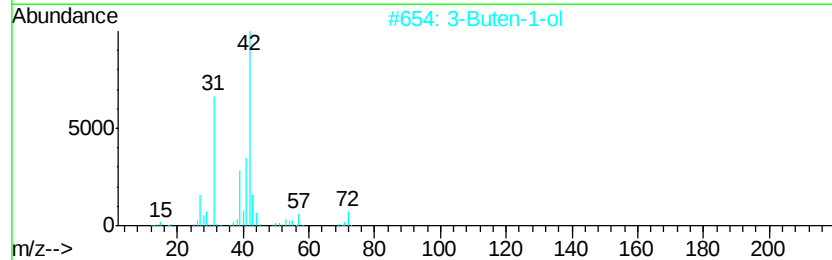
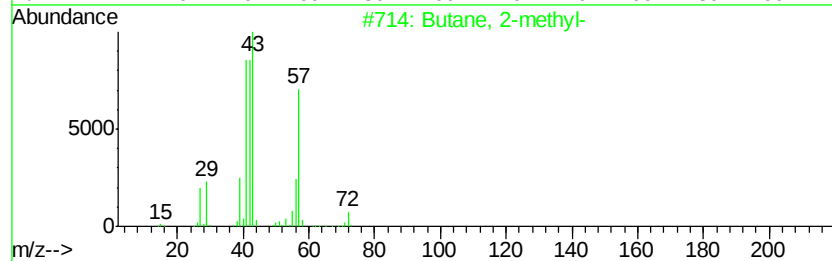
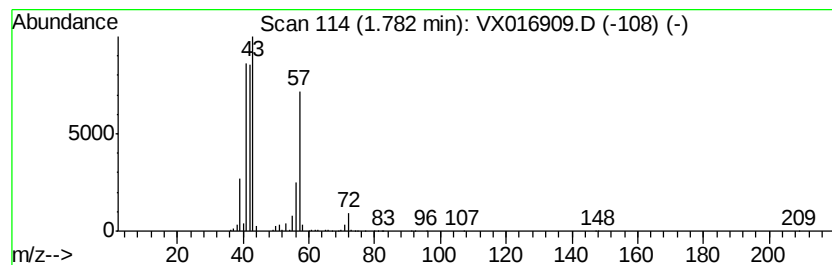
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Quant Title : SW846 8260

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

Peak Number 3 Butane, 2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.78	15.61 ug/l	236025	Pentafluorobenzene	5.63

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Butane, 2-methyl-	72	C5H12	000078-78-4	91
2		3-Buten-1-ol	72	C4H8O	000627-27-0	40
3		Pentane	72	C5H12	000109-66-0	39
4		Cyclobutane, methyl-	70	C5H10	000598-61-8	25
5		1-Propene, 2-methyl-	56	C4H8	000115-11-7	10



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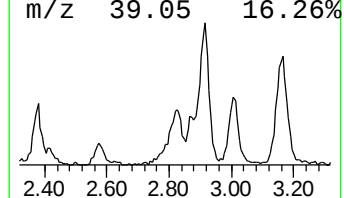
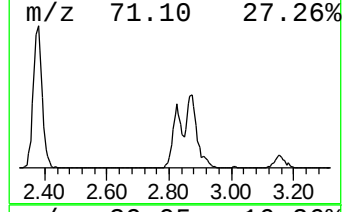
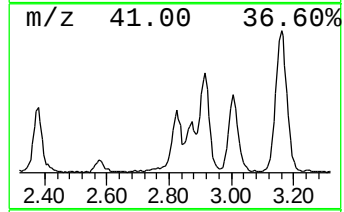
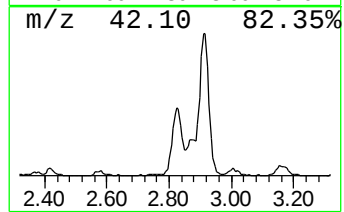
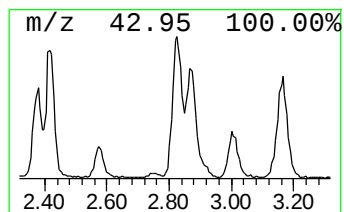
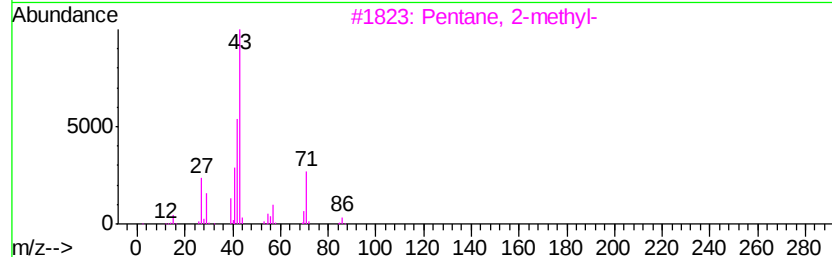
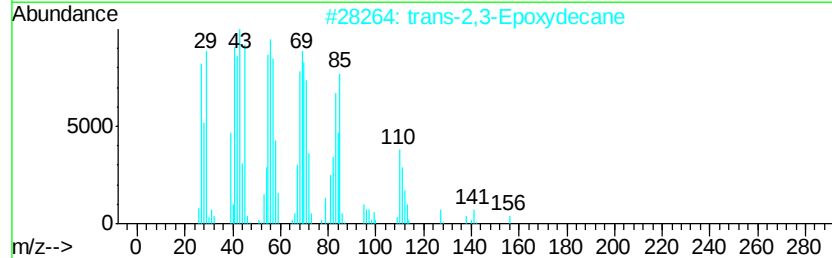
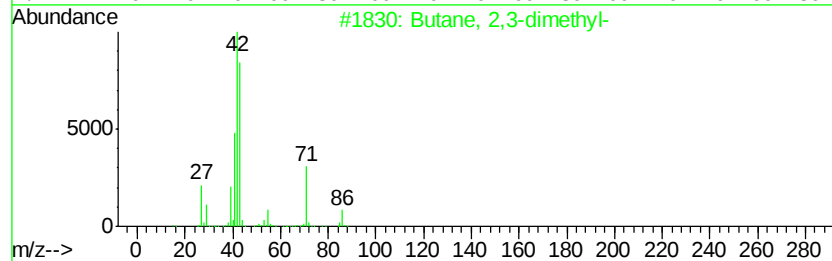
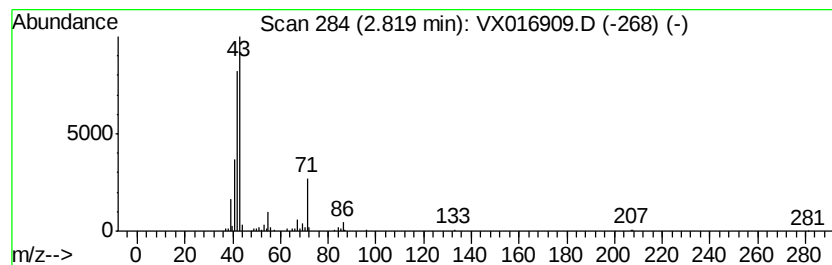
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Butane, 2,3-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.82	7.41 ug/l	111964	Pentafluorobenzene	5.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	87
2		trans-2,3-Epoxydecane	156	C10H20O	054125-39-2	72
3		Pentane, 2-methyl-	86	C6H14	000107-83-5	47
4		Pentane	72	C5H12	000109-66-0	36
5		Pyrrolidine	71	C4H9N	000123-75-1	22



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX061920\
Data File : VX016909.D
Acq On : 20 Jun 2020 05:30
Operator : JC/SP
Sample : L3063-08
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 42 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TW-12

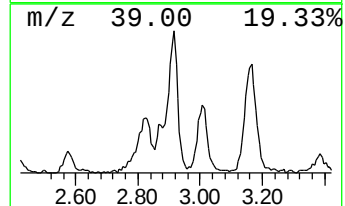
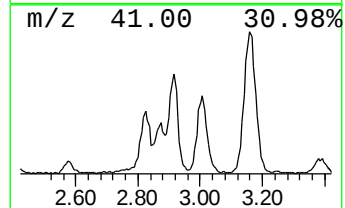
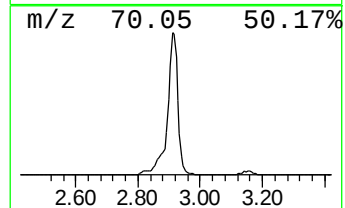
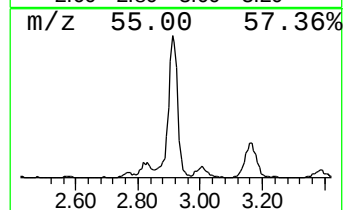
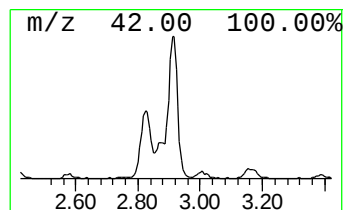
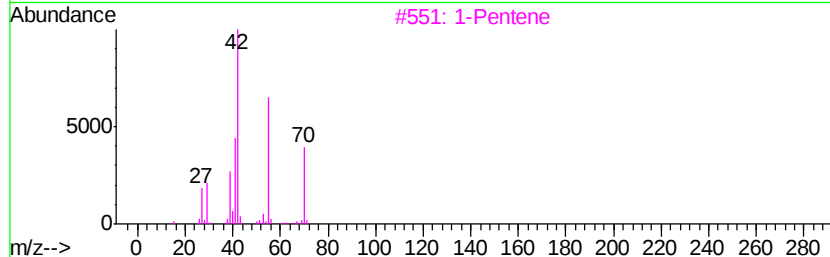
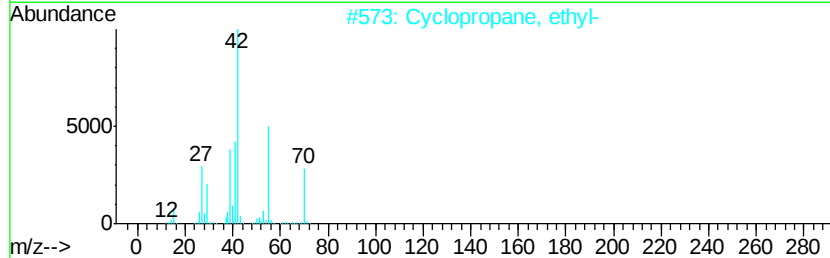
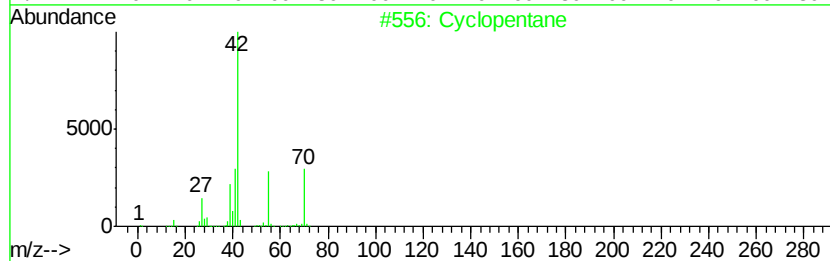
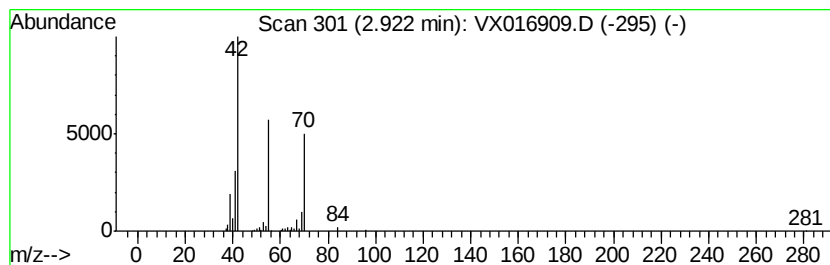
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Quant Title : SW846 8260

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

Peak Number 5 Cyclopentane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.92	9.17 ug/l	138611	Pentafluorobenzene	5.63

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopentane	70	C5H10	000287-92-3	72
2		Cyclopropane, ethyl-	70	C5H10	001191-96-4	56
3		1-Pentene	70	C5H10	000109-67-1	42
4		2-Butene-1,4-diol	88	C4H8O2	000110-64-5	39
5		2-Pentene, (E)-	70	C5H10	000646-04-8	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX061920\
Data File : VX016909.D
Acq On : 20 Jun 2020 05:30
Operator : JC/SP
Sample : L3063-08
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 42 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TW-12

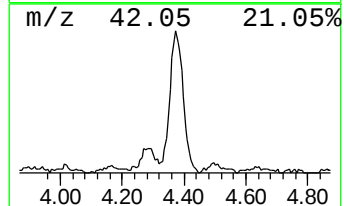
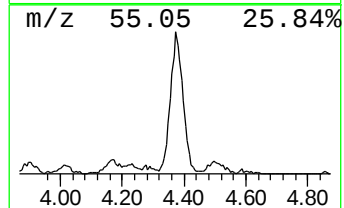
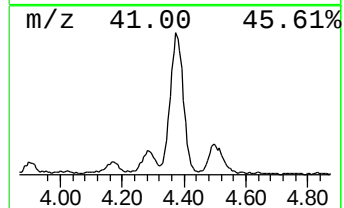
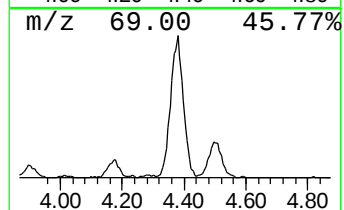
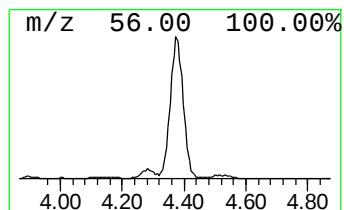
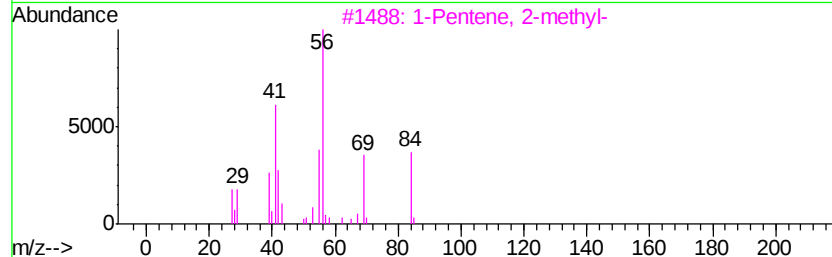
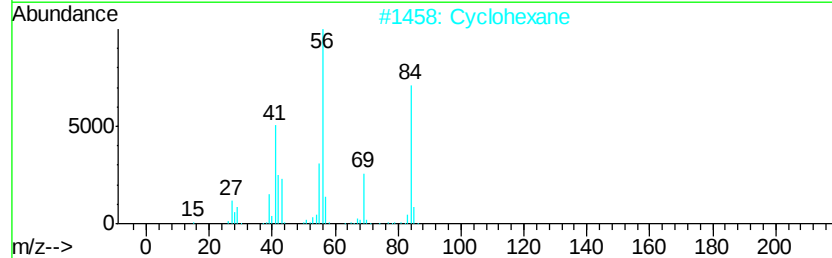
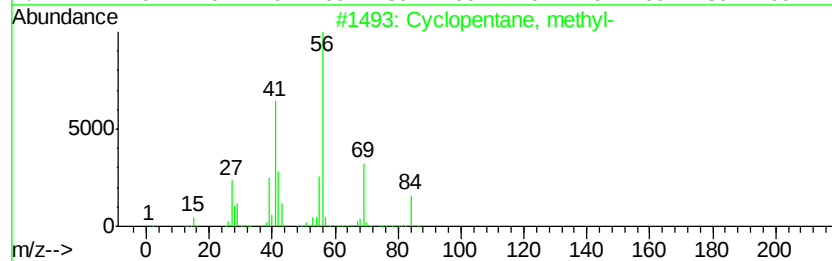
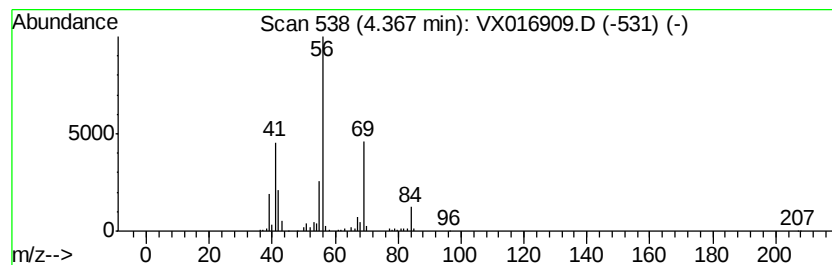
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Quant Title : SW846 8260

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

Peak Number 6 Cyclopentane, methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.37	13.28 ug/l	200725	Pentafluorobenzene	5.63

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2		Cyclohexane	84	C6H12	000110-82-7	83
3		1-Pentene, 2-methyl-	84	C6H12	000763-29-1	72
4		Cyclobutane, ethyl-	84	C6H12	004806-61-5	64
5		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX061920\
Data File : VX016909.D
Acq On : 20 Jun 2020 05:30
Operator : JC/SP
Sample : L3063-08
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 42 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TW-12

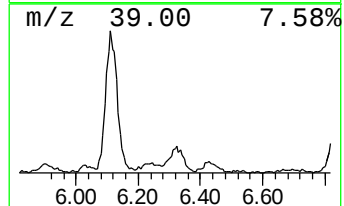
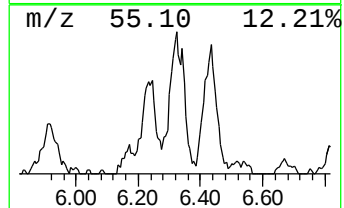
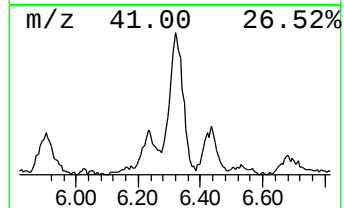
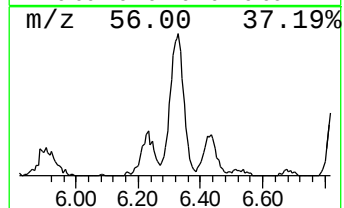
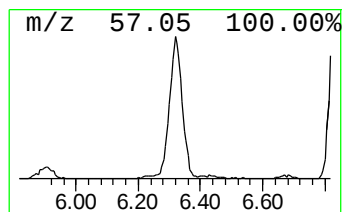
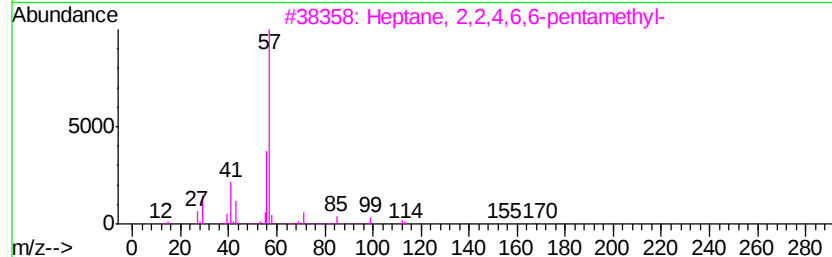
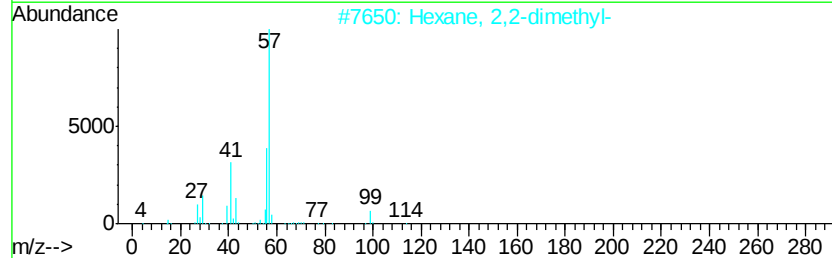
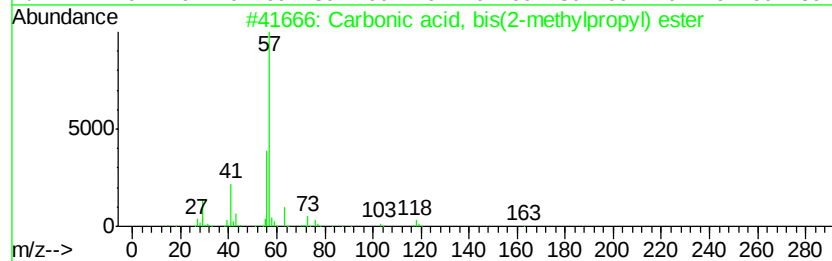
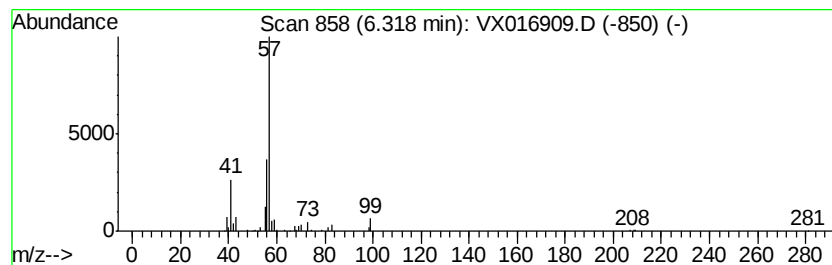
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Quant Title : SW846 8260

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

Peak Number 7 Carbonic acid, bis(2-methyl... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.32	5.49 ug/l	111967	1,4-Difluorobenzene	6.83

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Carbonic acid, bis(2-methylpropyl-...	174	C9H18O3	000539-92-4	59
2		Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	53
3		Heptane, 2,2,4,6,6-pentamethyl-	170	C12H26	013475-82-6	53
4		Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	53
5		Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX061920\
Data File : VX016909.D
Acq On : 20 Jun 2020 05:30
Operator : JC/SP
Sample : L3063-08
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 42 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TW-12

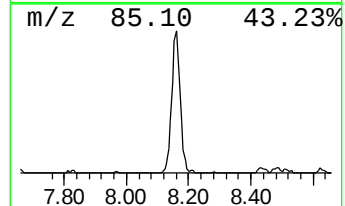
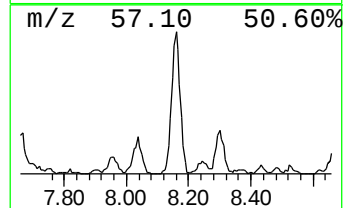
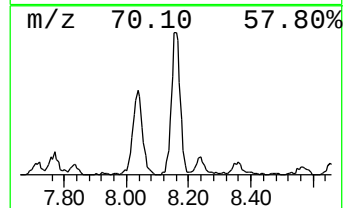
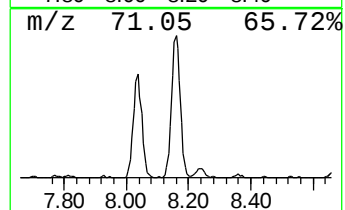
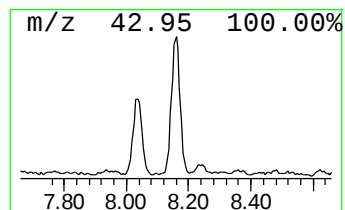
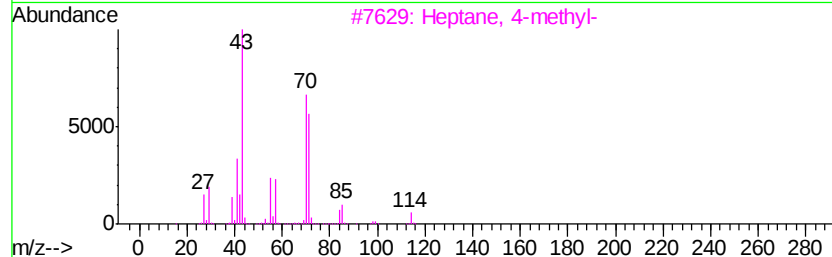
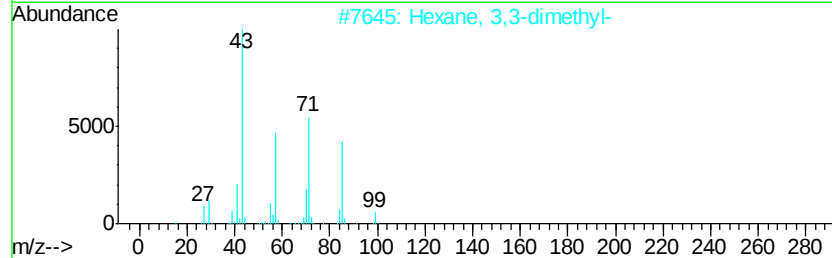
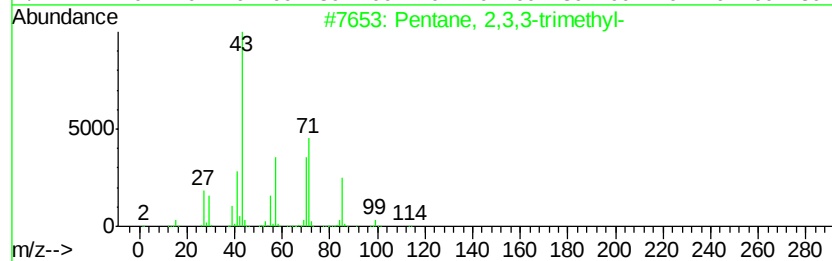
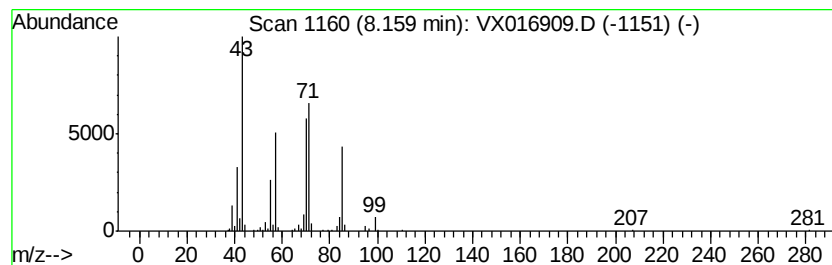
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Quant Title : SW846 8260

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

Peak Number 8 Pentane, 2,3,3-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.16	6.87 ug/l	140230	1,4-Difluorobenzene	6.83

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	90
2			Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	72
3			Heptane, 4-methyl-	114	C8H18	000589-53-7	53
4			Undecane, 5,7-dimethyl-	184	C13H28	017312-83-3	53
5			Undecane, 2,7-dimethyl-	184	C13H28	017301-24-5	53



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX061920\
Data File : VX016909.D
Acq On : 20 Jun 2020 05:30
Operator : JC/SP
Sample : L3063-08
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 42 Sample Multiplier: 1

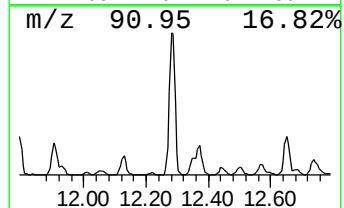
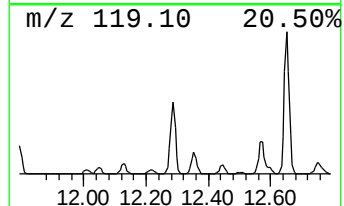
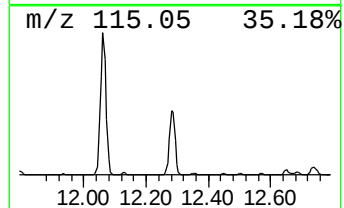
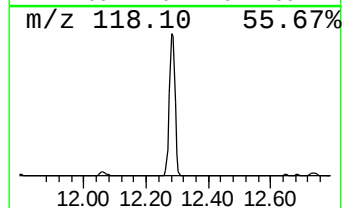
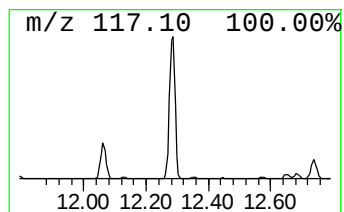
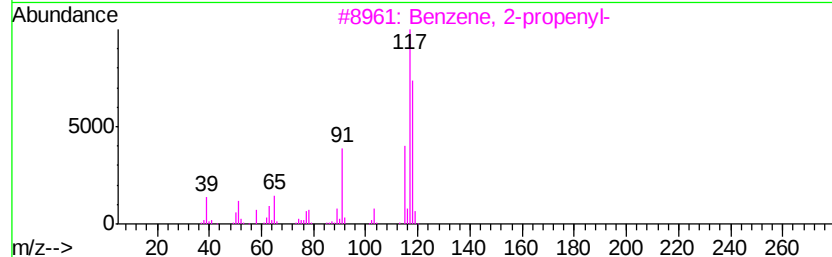
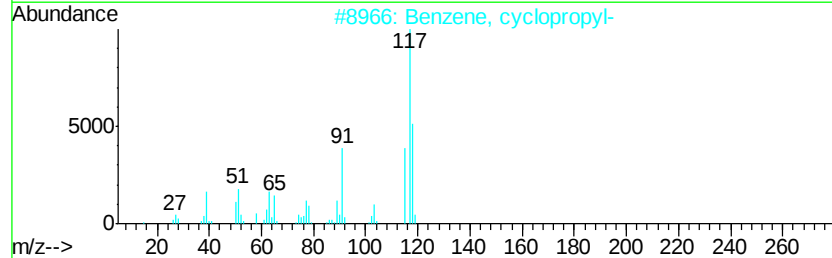
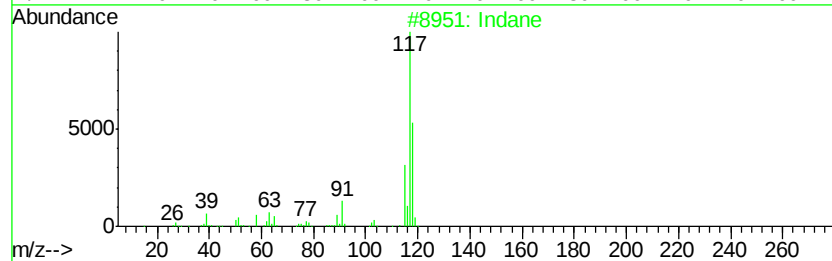
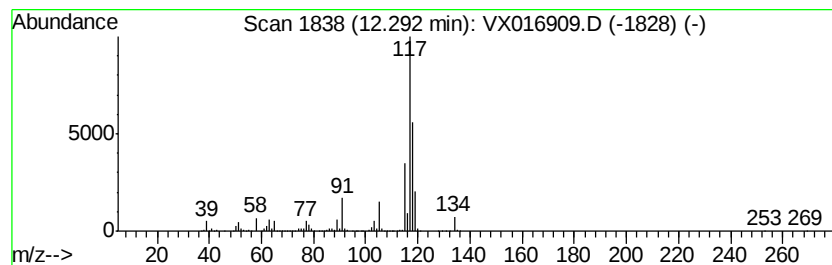
Instrument :
MSVOA_X
ClientSampleId :
TW-12

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X061820W.M
Quant Title : SW846 8260

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

Peak Number 9 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.29	24.34 ug/l	806446	1,4-Dichlorobenzene-d4	12.06
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Indane		118 C9H10	000496-11-7 76
2	Benzene, cyclopropyl-		118 C9H10	000873-49-4 76
3	Benzene, 2-propenyl-		118 C9H10	000300-57-2 76
4	Deltacyclene		118 C9H10	007785-10-6 58
5	Benzene, 1-ethenyl-2-methyl-		118 C9H10	000611-15-4 58



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA X\DATA\VX061920\
Data File : VX016909.D
Acq On : 20 Jun 2020 05:30
Operator : JC/SP
Sample : L3063-08
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 42 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
TW-12

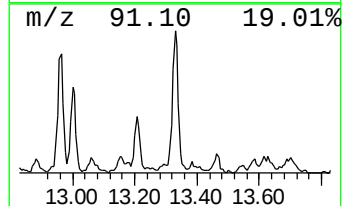
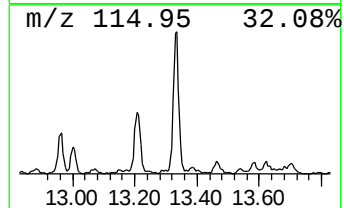
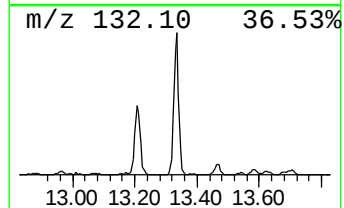
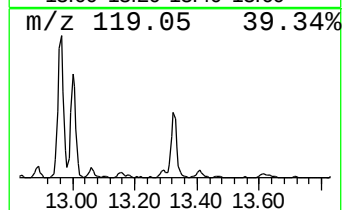
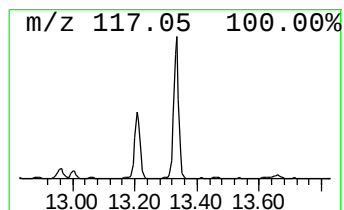
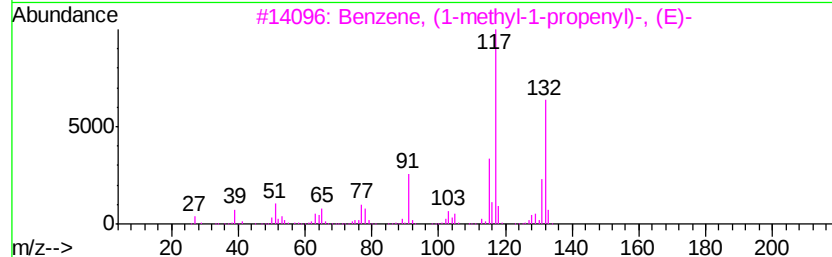
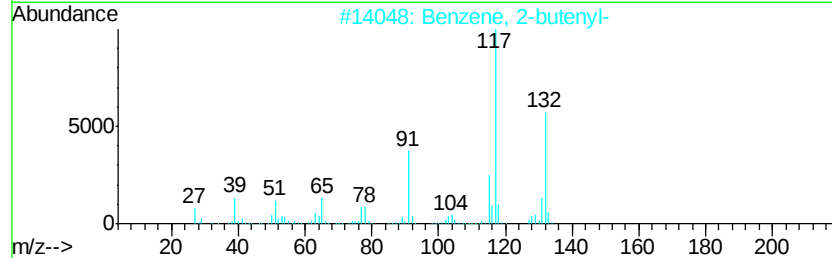
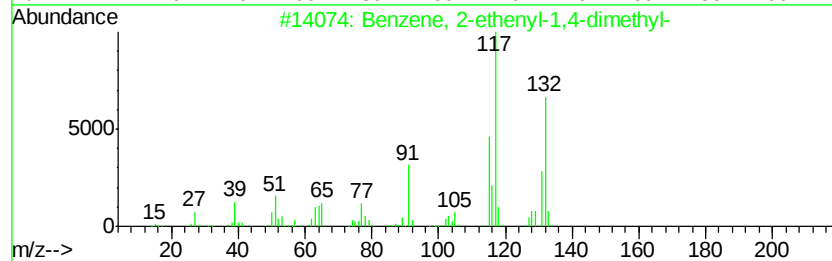
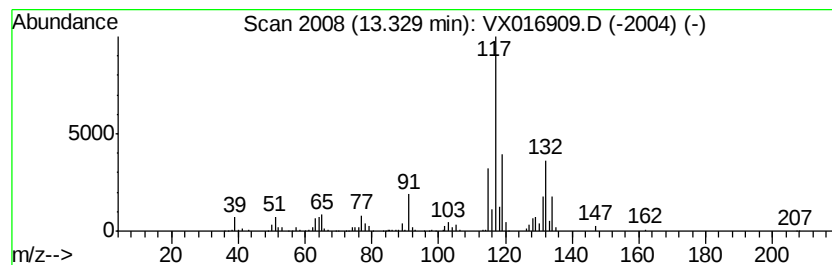
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Quant Title : SW846 8260

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

Peak Number 10 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.33	5.80 ug/l	192167	1,4-Dichlorobenzene-d4	12.06

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	90
2		Benzene, 2-butenyl-	132	C10H12	001560-06-1	76
3		Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	76
4		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	70
5		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	70



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_X\DATA\VX061920\
Data File : VX016909.D
Acq On : 20 Jun 2020 05:30
Operator : JC/SP
Sample : L3063-08
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 42 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TW-12

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X061820W.M
Quant Title : SW846 8260

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Propene	1.16	9.1	ug/l	138117	1	5.63	755803	50.0
1-Propene, 2-methyl-	1.38	8.3	ug/l	124950	1	5.63	755803	50.0
Butane, 2-methyl-	1.78	15.6	ug/l	236025	1	5.63	755803	50.0
Butane, 2,3-dimet...	2.82	7.4	ug/l	111964	1	5.63	755803	50.0
Cyclopentane	2.92	9.2	ug/l	138611	1	5.63	755803	50.0
Cyclopentane, met...	4.37	13.3	ug/l	200725	1	5.63	755803	50.0
Carbonic acid, bi...	6.32	5.5	ug/l	111967	2	6.83	1020290	50.0
Pentane, 2,3,3-tr...	8.16	6.9	ug/l	140230	2	6.83	1020290	50.0
Indane	12.29	24.3	ug/l	806446	4	12.06	1656780	50.0
Benzene, 2-etheny...	13.33	5.8	ug/l	192167	4	12.06	1656780	50.0