

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051922\
 Data File : VU048634.D
 Acq On : 19 May 2022 16:53
 Operator : SY/MD
 Sample : N2848-22ME
 Misc : 4.88g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 EXY12ME

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0 % of largest Peak

Start Thrs: 0.2

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\SFAMULM051622WMA.M
 Title : VOC Analysis

Signal : TIC: VU048634.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.478	36	43	60	rVB	73072	90187	9.76%	1.067%
2	1.597	73	80	103	rVB	52189	93021	10.06%	1.100%
3	2.536	360	372	385	rBV	176506	283454	30.66%	3.352%
4	2.954	492	502	517	rBV3	5464	11695	1.27%	0.138%
5	3.031	517	526	541	rVB4	3790	7136	0.77%	0.084%
6	3.099	544	547	550	rBV	716	527	0.06%	0.006%
7	3.183	559	573	591	rBV	6413	15455	1.67%	0.183%
8	3.295	603	608	610	rBV	271	267	0.03%	0.003%
9	3.565	688	692	694	rBV3	950	780	0.08%	0.009%
10	3.967	810	817	818	rBV	507	552	0.06%	0.007%
11	4.019	831	833	837	rVB2	380	296	0.03%	0.004%
12	4.051	837	843	846	rBV2	778	849	0.09%	0.010%
13	4.109	857	861	863	rBB	265	170	0.02%	0.002%
14	4.253	903	906	910	rBB	337	196	0.02%	0.002%
15	4.459	965	970	975	rBV2	496	437	0.05%	0.005%
16	4.539	990	995	998	rBB	264	168	0.02%	0.002%
17	4.629	1010	1023	1037	rBV	71432	197051	21.32%	2.330%
18	5.060	1140	1157	1185	rBV	187898	445902	48.24%	5.274%
19	5.199	1189	1200	1220	rVV2	26731	62065	6.71%	0.734%
20	5.578	1315	1318	1322	rBB2	318	222	0.02%	0.003%
21	5.716	1341	1361	1393	rBV3	352041	924425	100.00%	10.933%
22	5.825	1393	1395	1398	rVB	481	256	0.03%	0.003%
23	5.880	1409	1412	1415	rBV2	351	279	0.03%	0.003%
24	5.970	1429	1440	1457	rBV5	8471	19142	2.07%	0.226%
25	6.163	1481	1500	1513	rBV3	10613	31106	3.36%	0.368%
26	6.247	1514	1526	1552	rVB	264390	531257	57.47%	6.283%
27	6.398	1560	1573	1592	rBV3	26982	56649	6.13%	0.670%
28	6.597	1625	1635	1646	rBV7	4581	10446	1.13%	0.124%
29	6.690	1650	1664	1681	rVV	199462	402117	43.50%	4.756%
30	6.780	1682	1692	1723	rVV3	17927	59244	6.41%	0.701%
31	6.928	1727	1738	1767	rVB	19675	58270	6.30%	0.689%
32	7.186	1807	1818	1850	rVB4	47702	115298	12.47%	1.364%
33	7.292	1850	1851	1856	rBV2	292	261	0.03%	0.003%
34	7.372	1871	1876	1878	rBV	1100	975	0.11%	0.012%
35	7.497	1913	1915	1919	rBV	282	198	0.02%	0.002%
36	7.565	1923	1936	1957	rVB2	116847	229594	24.84%	2.715%

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Instrument :
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 ClientSampleId :
 EXY12ME

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0 % of largest Peak

Start Thrs: 0.2

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\SFAMULM051622WMA.M
 Title : VOC Analysis

37	7.687	1969	1974	1976	rBV3	777	733	0.08%	0.009%
38	7.793	1996	2007	2018	rBV5	3780	8291	0.90%	0.098%
39	7.896	2025	2039	2054	rBV	422480	762439	82.48%	9.017%
40	8.182	2116	2128	2153	rVB3	101635	218185	23.60%	2.580%
41	8.311	2165	2168	2170	rBV	302	183	0.02%	0.002%
42	8.346	2173	2179	2184	rBV3	487	492	0.05%	0.006%
43	8.440	2203	2208	2211	rBV2	364	374	0.04%	0.004%
44	8.468	2211	2217	2218	rBV2	1376	901	0.10%	0.011%
45	8.597	2252	2257	2258	rBV	248	200	0.02%	0.002%
46	8.652	2258	2274	2307	rVV	242547	553588	59.88%	6.547%
47	8.764	2307	2309	2312	rBV3	663	450	0.05%	0.005%
48	8.854	2334	2337	2341	rVB	288	196	0.02%	0.002%
49	8.874	2341	2343	2346	rVB2	455	265	0.03%	0.003%
50	8.903	2349	2352	2354	rBV	547	319	0.03%	0.004%
51	8.931	2354	2361	2362	rBV	1188	961	0.10%	0.011%
52	9.115	2413	2418	2423	rBB	198	251	0.03%	0.003%
53	9.173	2434	2436	2439	rBB	311	204	0.02%	0.002%
54	9.340	2486	2488	2491	rBV	299	225	0.02%	0.003%
55	9.420	2500	2513	2533	rBV	356897	670497	72.53%	7.930%
56	9.546	2548	2552	2556	rBV3	414	436	0.05%	0.005%
57	9.575	2557	2561	2569	rVB5	859	1071	0.12%	0.013%
58	9.690	2588	2597	2598	rBV4	2315	2309	0.25%	0.027%
59	9.777	2620	2624	2626	rBV	207	200	0.02%	0.002%
60	9.960	2677	2681	2683	rBV	217	188	0.02%	0.002%
61	9.999	2689	2693	2701	rVB2	229	233	0.03%	0.003%
62	10.176	2745	2748	2750	rBV	257	202	0.02%	0.002%
63	10.192	2750	2753	2757	rVB2	273	197	0.02%	0.002%
64	10.269	2774	2777	2781	rVV2	265	259	0.03%	0.003%
65	10.542	2856	2862	2864	rBV	213	280	0.03%	0.003%
66	10.578	2870	2873	2876	rBV	168	176	0.02%	0.002%
67	10.642	2888	2893	2894	rBV	401	309	0.03%	0.004%
68	10.671	2898	2902	2908	rVB4	1627	1874	0.20%	0.022%
69	10.764	2917	2931	2947	rVB	338374	562221	60.82%	6.649%
70	10.848	2952	2957	2959	rBV	208	224	0.02%	0.003%
71	10.893	2961	2971	2983	rVV	4100	6358	0.69%	0.075%
72	10.938	2983	2985	2989	rVB2	582	342	0.04%	0.004%
73	10.967	2990	2994	2996	rBV	188	176	0.02%	0.002%
74	11.057	3013	3022	3030	rBV6	14003	20817	2.25%	0.246%
75	11.115	3030	3040	3058	rVB	168602	253970	27.47%	3.004%
76	11.311	3096	3101	3107	rVB4	1482	1657	0.18%	0.020%

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Instrument :
 MSVOA_U
 ClientSampleId :
 EXY12ME

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0 % of largest Peak

Start Thrs: 0.2

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\SFAMULM051622WMA.M
 Title : VOC Analysis

77	11.459	3141	3147	3149	rBV4	1528	1478	0.16%	0.017%
78	11.578	3176	3184	3187	rBV4	1693	2166	0.23%	0.026%
79	11.636	3199	3202	3205	rBV3	1020	752	0.08%	0.009%
80	11.700	3215	3222	3228	rVB4	1517	1963	0.21%	0.023%
81	11.751	3233	3238	3243	rBV2	781	703	0.08%	0.008%
82	11.816	3243	3258	3271	rBV	494192	792048	85.68%	9.367%
83	12.195	3363	3376	3391	rVV	527167	852115	92.18%	10.078%
84	12.330	3408	3418	3427	rVB3	28587	40391	4.37%	0.478%
85	12.375	3427	3432	3436	rBV2	379	398	0.04%	0.005%
86	12.398	3436	3439	3443	rBV3	398	354	0.04%	0.004%
87	12.645	3513	3516	3520	rBV	306	241	0.03%	0.003%
88	12.690	3525	3530	3534	rVB	363	291	0.03%	0.003%
89	12.774	3546	3556	3564	rVB3	2191	3500	0.38%	0.041%
90	12.832	3572	3574	3579	rBV	270	187	0.02%	0.002%
91	12.899	3587	3595	3603	rBV7	2098	2884	0.31%	0.034%
92	13.021	3631	3633	3636	rBV	297	198	0.02%	0.002%
93	13.221	3688	3695	3699	rBV4	1217	1575	0.17%	0.019%
94	13.391	3740	3748	3759	rBV5	15083	21373	2.31%	0.253%
95	13.590	3807	3810	3812	rBV	315	166	0.02%	0.002%
96	14.034	3945	3948	3949	rBV	332	213	0.02%	0.003%
97	14.070	3956	3959	3960	rBV2	410	250	0.03%	0.003%
98	14.336	4036	4042	4051	rBV4	931	1315	0.14%	0.016%
99	15.150	4293	4295	4300	rBV	513	352	0.04%	0.004%
100	15.372	4356	4364	4371	rBV7	5514	8199	0.89%	0.097%

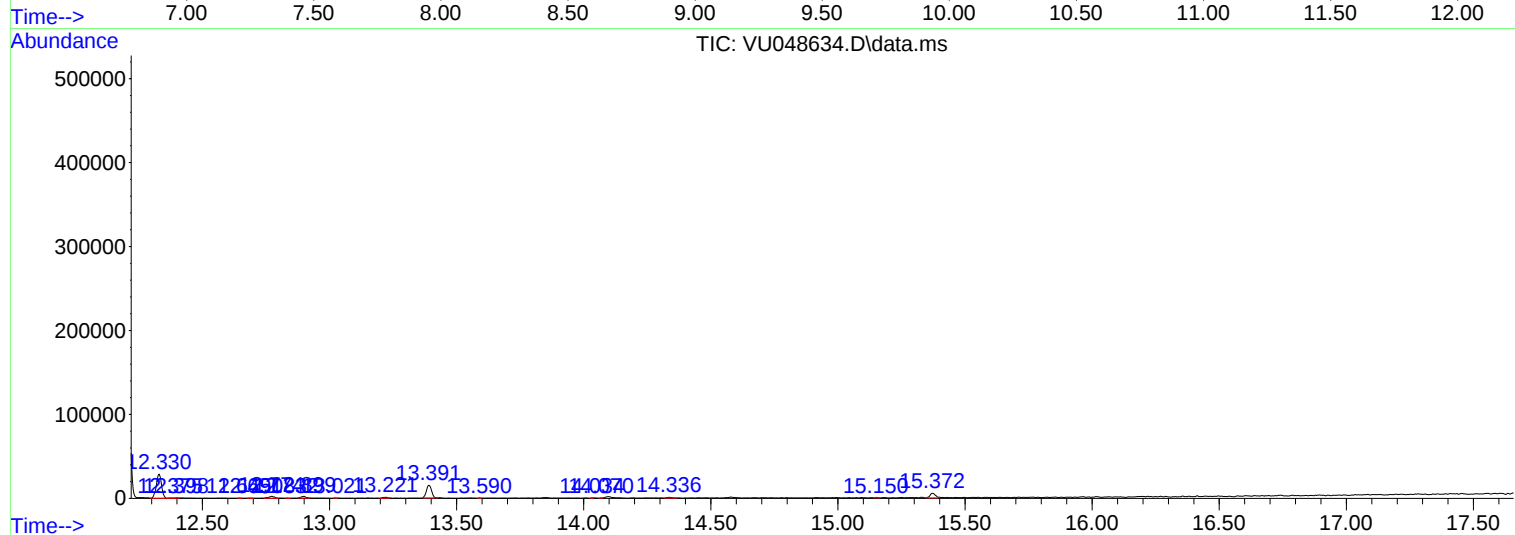
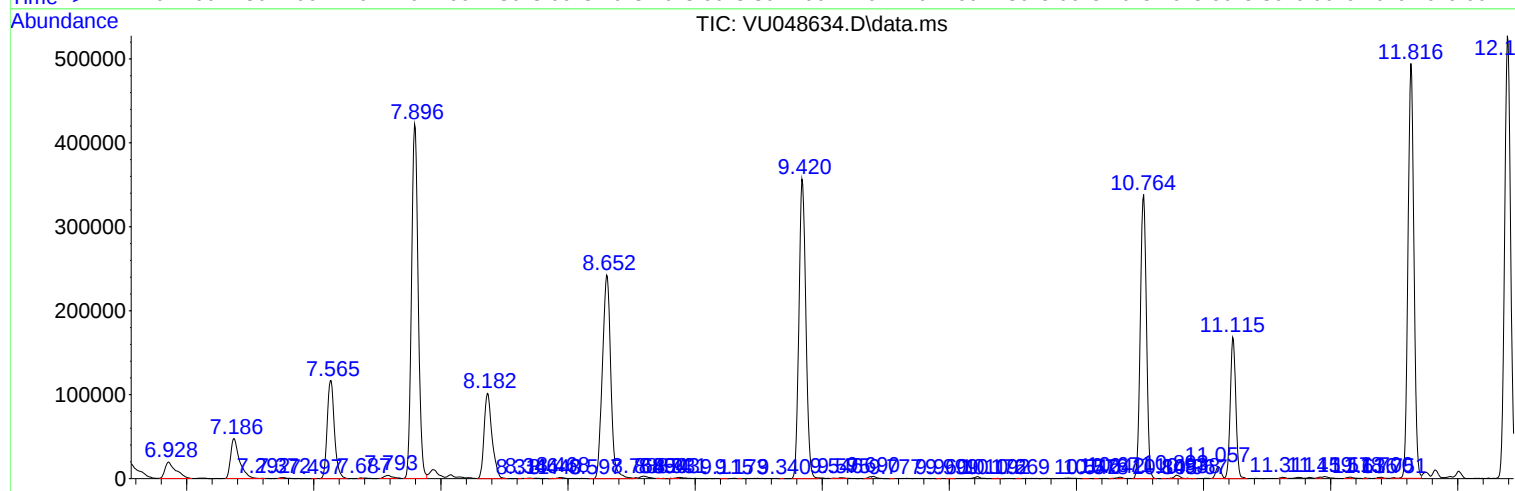
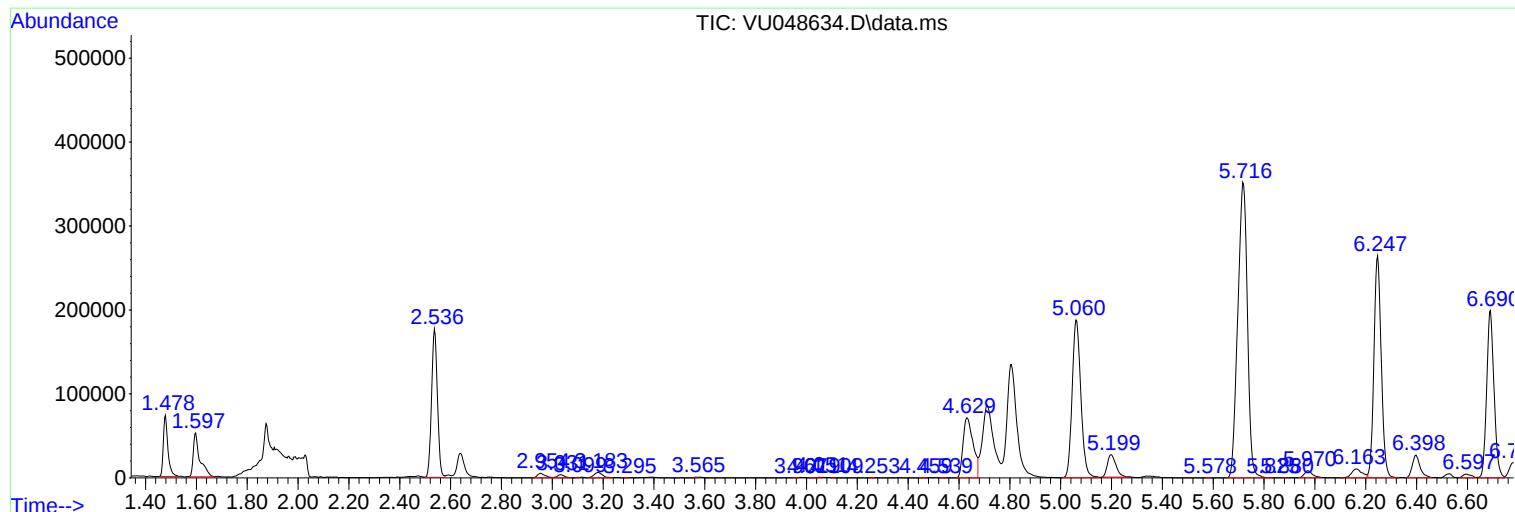
Sum of corrected areas: 8455312

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ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXY12ME

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM051622WMA.M
Quant Title : VOC Analysis

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



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ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXY12ME

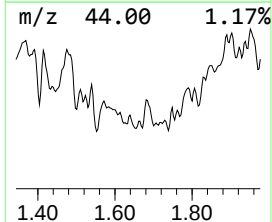
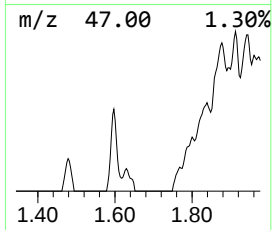
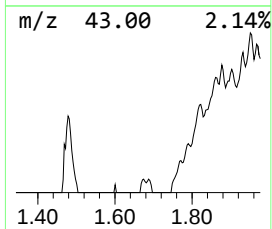
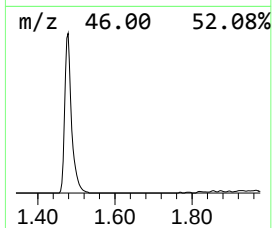
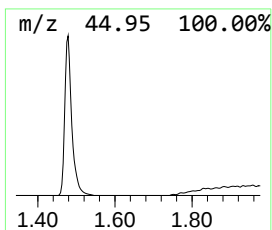
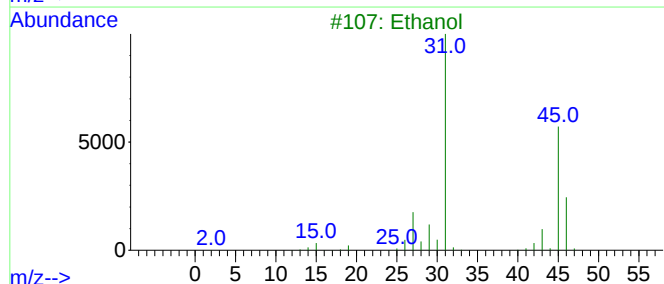
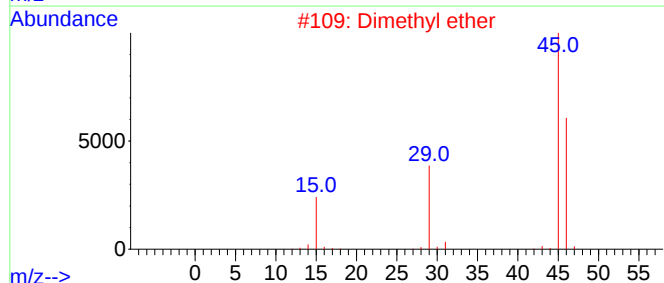
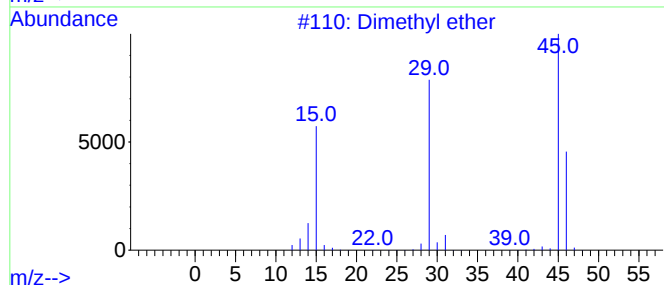
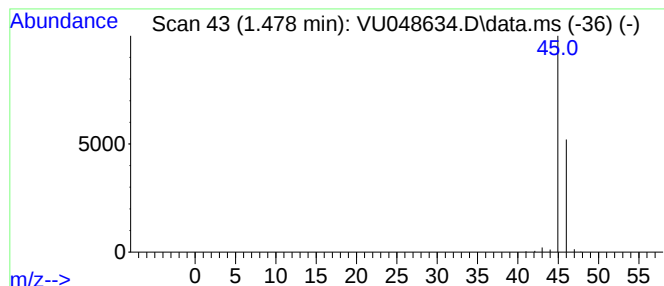
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM051622WMA.M
Quant Title : VOC Analysis

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

Peak Number 1 Dimethyl ether Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.478	8.49 ug/L	90187	1,4-Difluorobenzene	6.247

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dimethyl ether	46	C2H6O	000115-10-6	90
2			Dimethyl ether	46	C2H6O	000115-10-6	9
3			Ethanol	46	C2H6O	000064-17-5	7
4			Formic acid	46	CH2O2	000064-18-6	7
5			Ethanol	46	C2H6O	000064-17-5	7



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EXY12ME

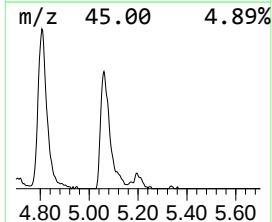
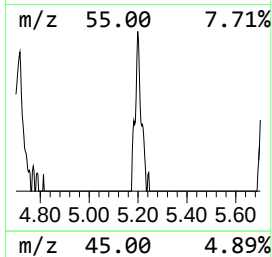
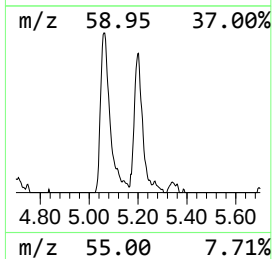
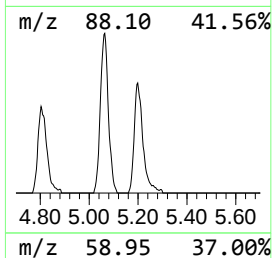
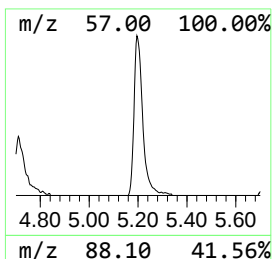
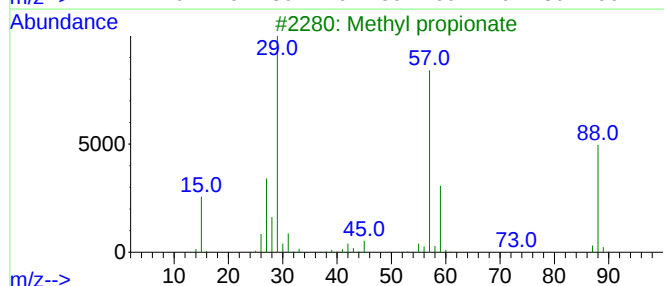
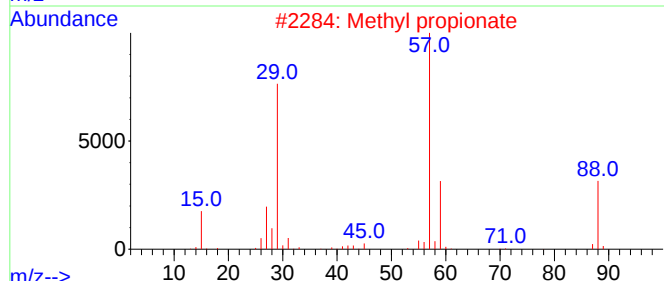
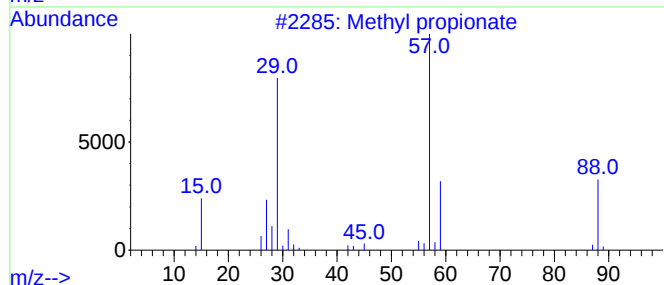
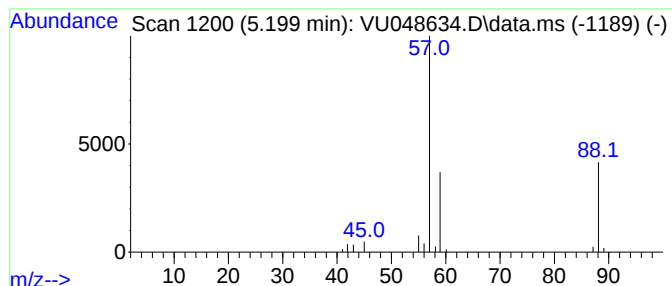
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM051622WMA.M
Quant Title : VOC Analysis

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

Peak Number 2 Methyl propionate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.199	5.84 ug/L	62065	1,4-Difluorobenzene	6.247

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methyl propionate			88	C4H8O2	000554-12-1	90
2	Methyl propionate			88	C4H8O2	000554-12-1	90
3	Methyl propionate			88	C4H8O2	000554-12-1	83
4	Methyl propionate			88	C4H8O2	000554-12-1	83
5	Methyl propionate			88	C4H8O2	000554-12-1	83



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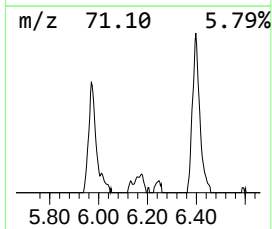
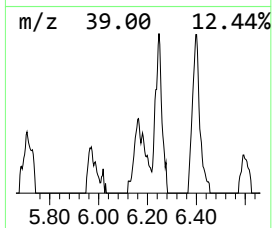
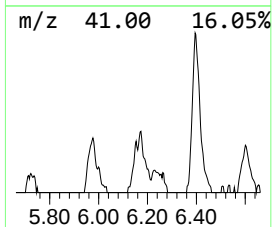
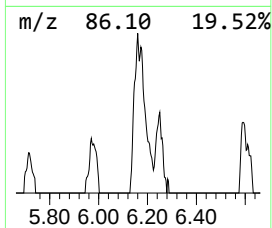
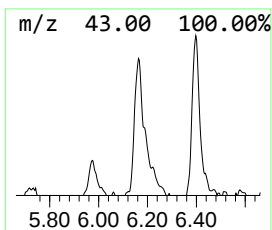
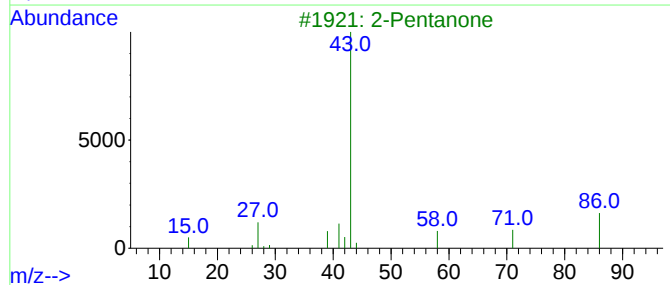
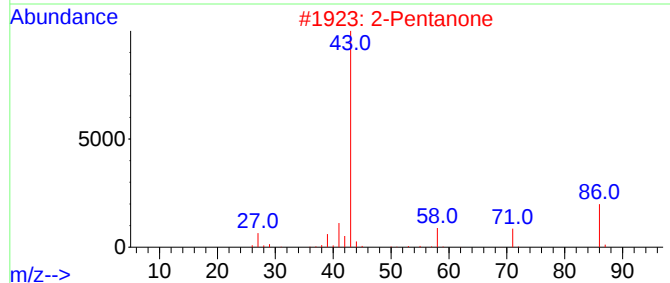
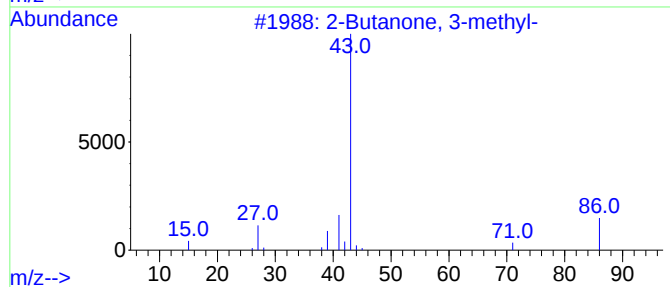
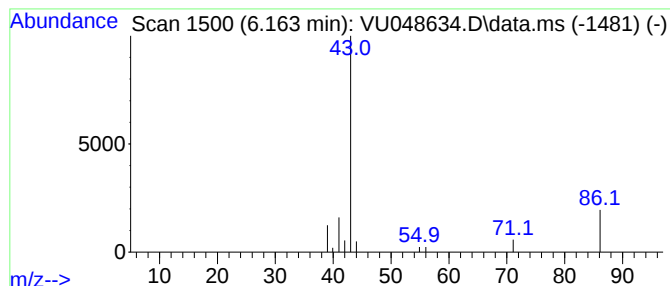
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Quant Title : VOC Analysis

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

Peak Number 3 2-Butanone, 3-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.163	2.93 ug/L	31106	1,4-Difluorobenzene	6.247

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butanone, 3-methyl-			86	C5H10O	000563-80-4	64
2	2-Pentanone			86	C5H10O	000107-87-9	56
3	2-Pentanone			86	C5H10O	000107-87-9	56
4	2-Pentanone			86	C5H10O	000107-87-9	40
5	2-Pentanone			86	C5H10O	000107-87-9	9



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051922\
Data File : VU048634.D
Acq On : 19 May 2022 16:53
Operator : SY/MD
Sample : N2848-22ME
Misc : 4.88g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXY12ME

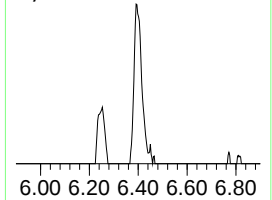
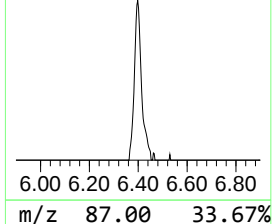
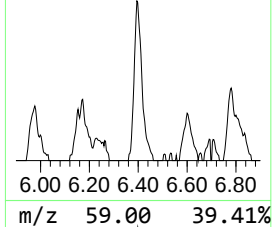
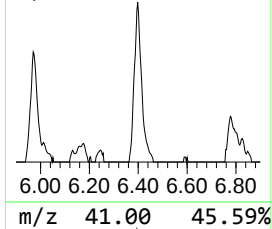
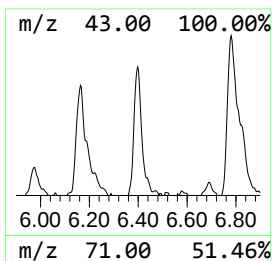
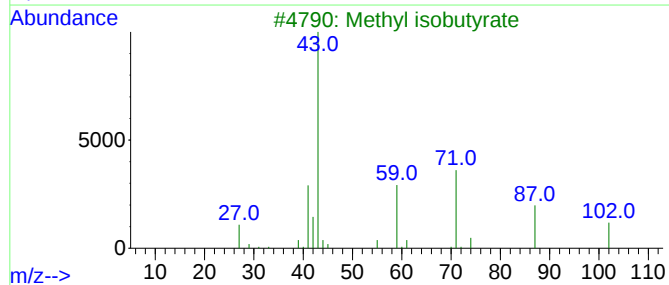
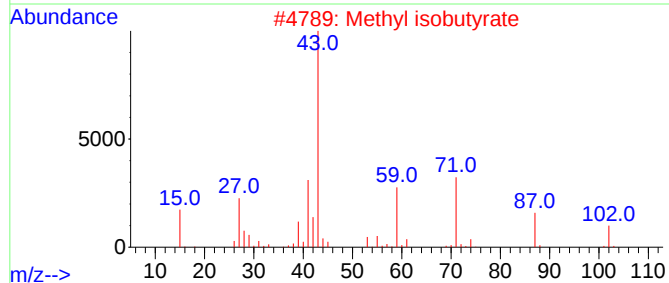
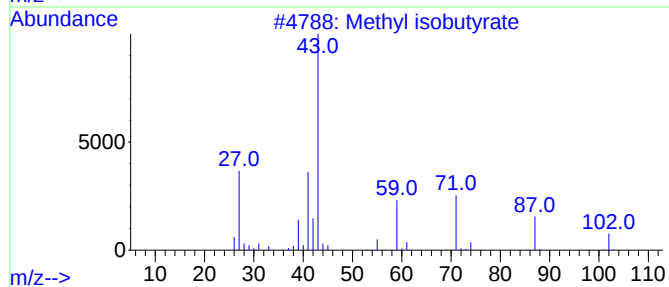
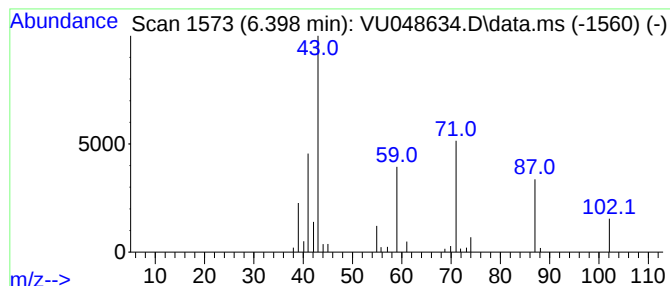
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Quant Title : VOC Analysis

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

Peak Number 4 Methyl isobutyrate Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.398	5.33 ug/L	56649	1,4-Difluorobenzene	6.247

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl isobutyrate	102	C5H10O2	000547-63-7	91
2			Methyl isobutyrate	102	C5H10O2	000547-63-7	91
3			Methyl isobutyrate	102	C5H10O2	000547-63-7	91
4			Methyl isobutyrate	102	C5H10O2	000547-63-7	43
5			Cycloserine	102	C3H6N2O2	000068-41-7	35



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051922\
Data File : VU048634.D
Acq On : 19 May 2022 16:53
Operator : SY/MD
Sample : N2848-22ME
Misc : 4.88g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXY12ME

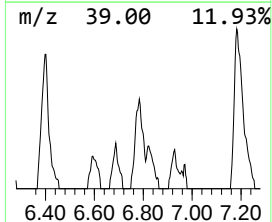
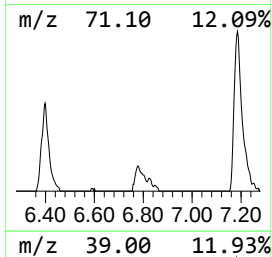
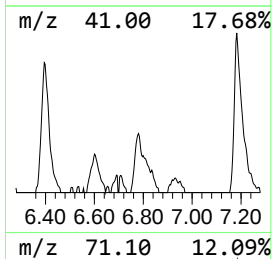
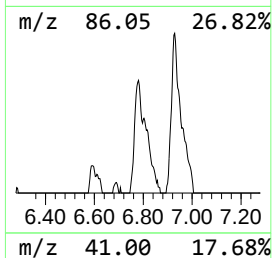
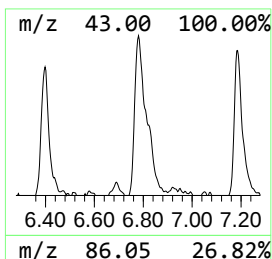
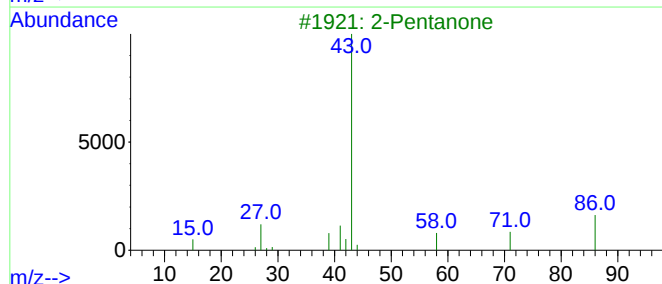
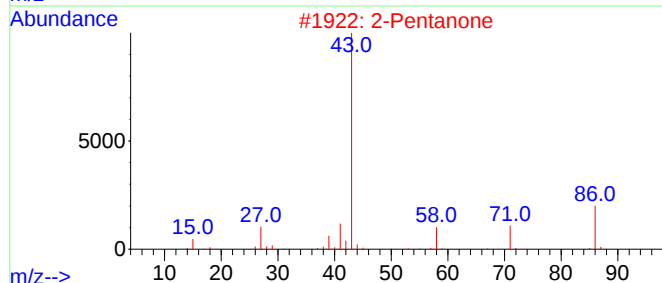
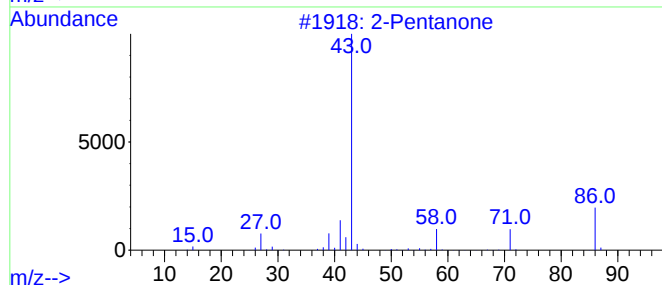
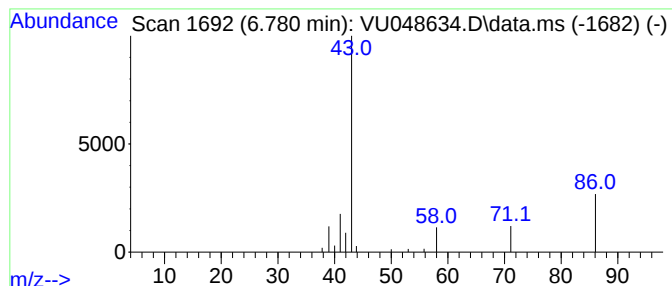
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Quant Title : VOC Analysis

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

Peak Number 5 2-Pentanone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.780	5.58 ug/L	59244	1,4-Difluorobenzene	6.247

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone	86	C5H10O	000107-87-9	86
2			2-Pentanone	86	C5H10O	000107-87-9	78
3			2-Pentanone	86	C5H10O	000107-87-9	78
4			2-Pentanone	86	C5H10O	000107-87-9	78
5			2-Butanone, 3-methyl-	86	C5H10O	000563-80-4	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051922\
Data File : VU048634.D
Acq On : 19 May 2022 16:53
Operator : SY/MD
Sample : N2848-22ME
Misc : 4.88g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXY12ME

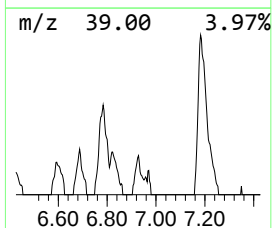
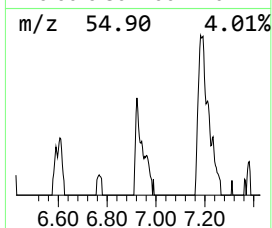
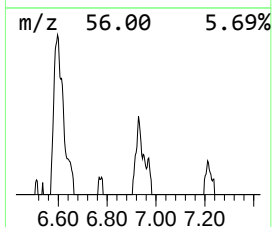
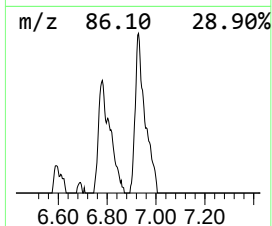
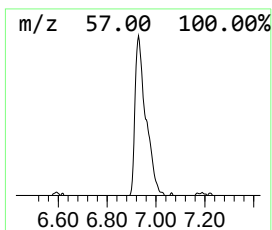
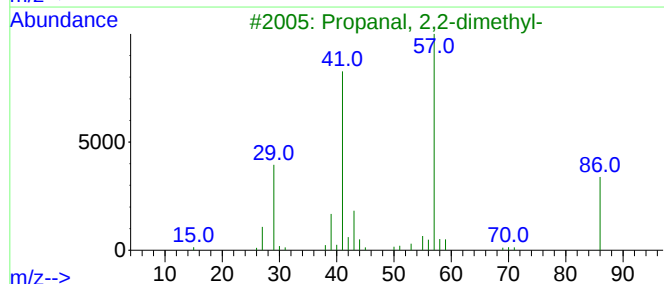
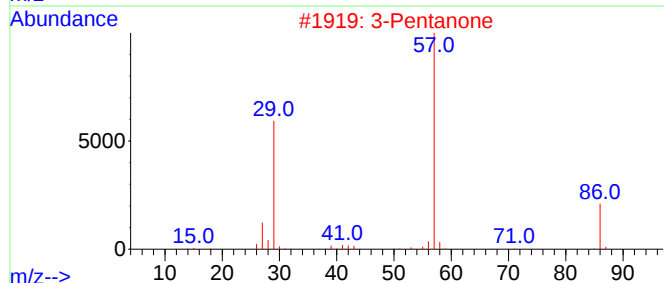
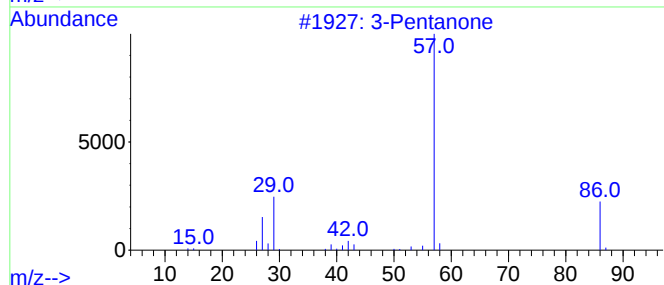
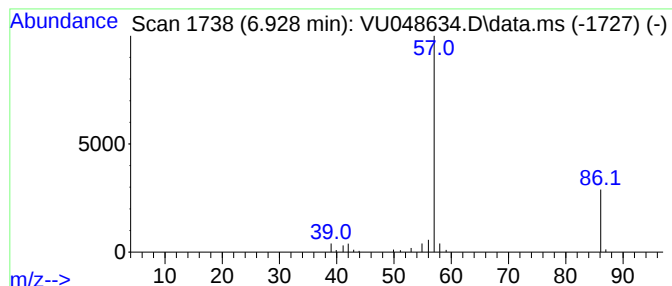
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Quant Title : VOC Analysis

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

Peak Number 6 3-Pentanone Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.928	5.48 ug/L	58270	1,4-Difluorobenzene	6.247

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Pentanone	86	C5H10O	000096-22-0	80
2			3-Pentanone	86	C5H10O	000096-22-0	45
3			Propanal, 2,2-dimethyl-	86	C5H10O	000630-19-3	9
4			3-Pentanone	86	C5H10O	000096-22-0	9
5			Propane, 2-bromo-2-methyl-	136	C4H9Br	000507-19-7	9



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051922\
Data File : VU048634.D
Acq On : 19 May 2022 16:53
Operator : SY/MD
Sample : N2848-22ME
Misc : 4.88g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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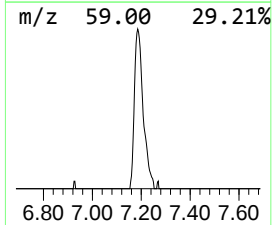
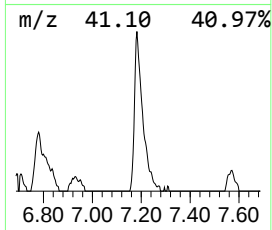
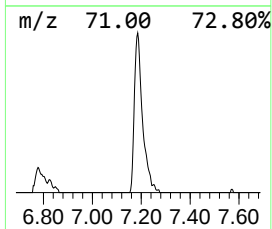
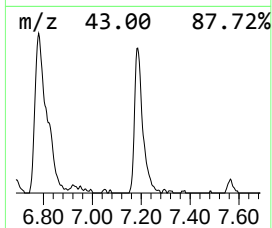
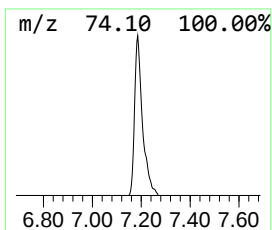
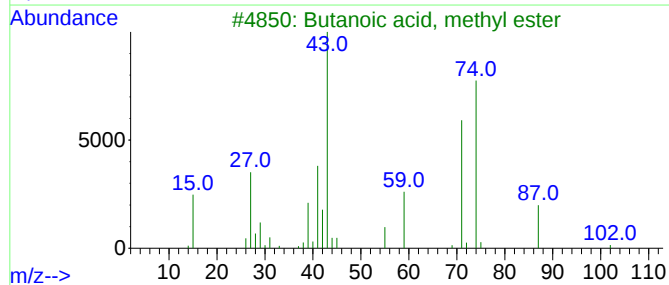
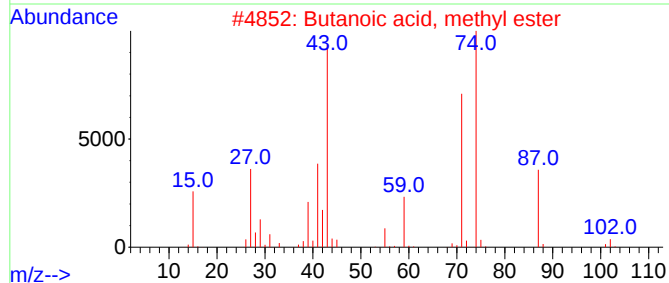
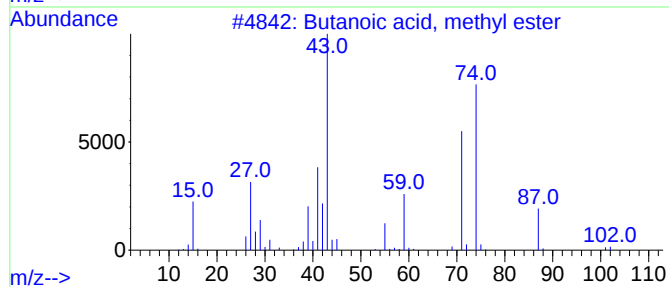
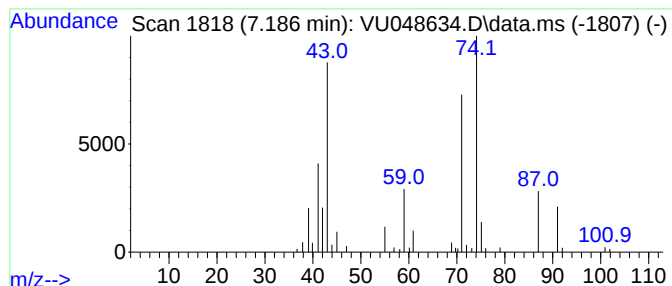
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Quant Title : VOC Analysis

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

Peak Number 7 Butanoic acid, methyl ester Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.186	10.85 ug/L	115298	1,4-Difluorobenzene	6.247

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butanoic acid, methyl ester	102	C5H10O2	000623-42-7	87
2			Butanoic acid, methyl ester	102	C5H10O2	000623-42-7	72
3			Butanoic acid, methyl ester	102	C5H10O2	000623-42-7	72
4			Butanoic acid, methyl ester	102	C5H10O2	000623-42-7	72
5			Methyl isobutyrate	102	C5H10O2	000547-63-7	37



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051922\
Data File : VU048634.D
Acq On : 19 May 2022 16:53
Operator : SY/MD
Sample : N2848-22ME
Misc : 4.88g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXY12ME

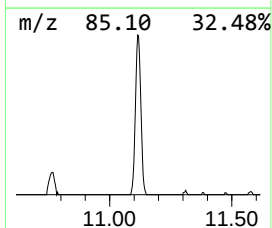
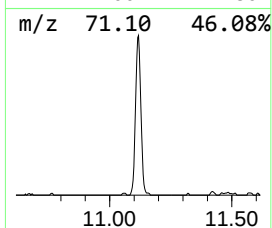
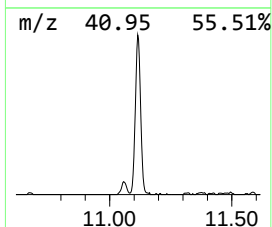
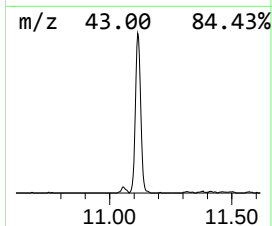
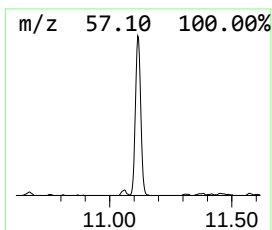
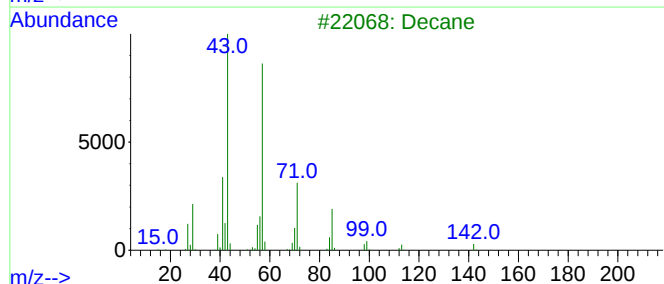
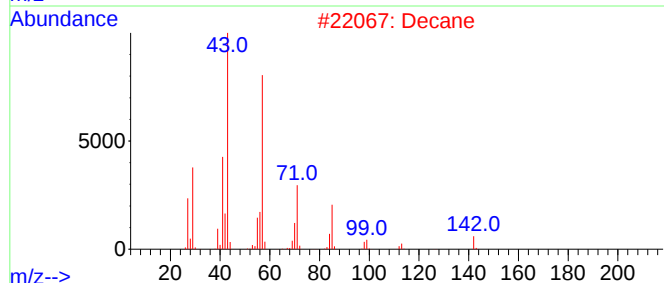
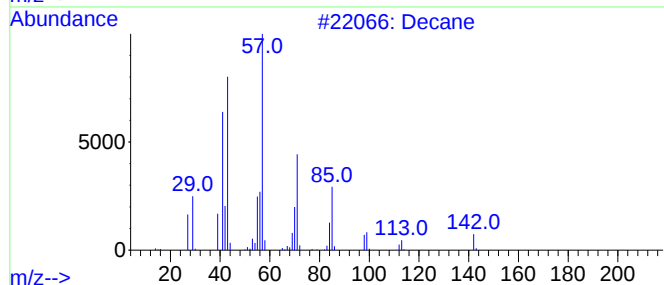
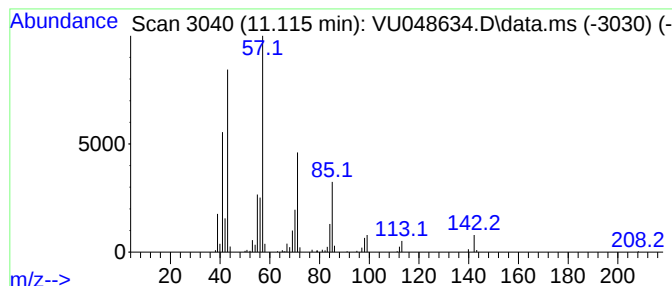
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Quant Title : VOC Analysis

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.115	16.03 ug/L	253970	1,4-Dichlorobenzene-d4	11.816

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane		142	C10H22	000124-18-5	96
2	Decane		142	C10H22	000124-18-5	95
3	Decane		142	C10H22	000124-18-5	90
4	Decane		142	C10H22	000124-18-5	83
5	Undecane		156	C11H24	001120-21-4	78



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051922\
Data File : VU048634.D
Acq On : 19 May 2022 16:53
Operator : SY/MD
Sample : N2848-22ME
Misc : 4.88g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXY12ME

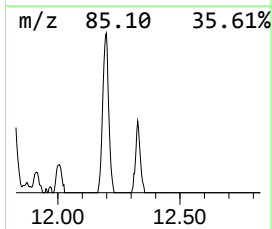
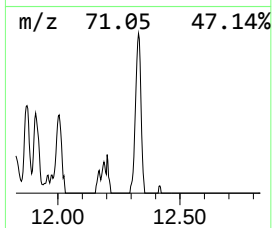
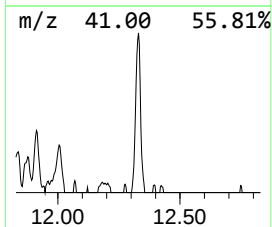
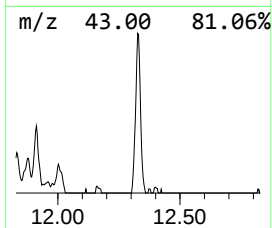
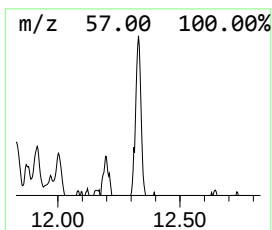
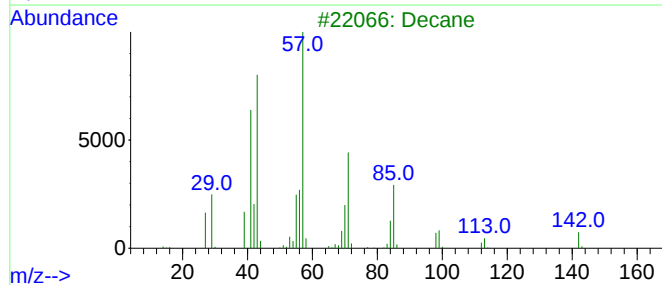
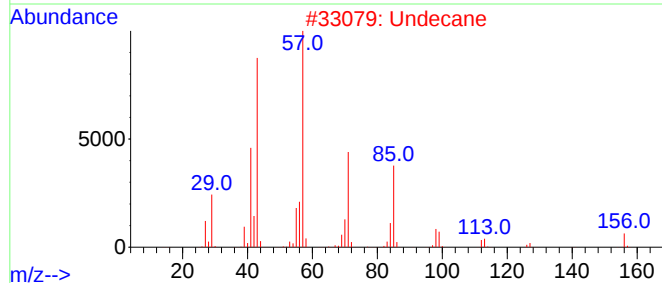
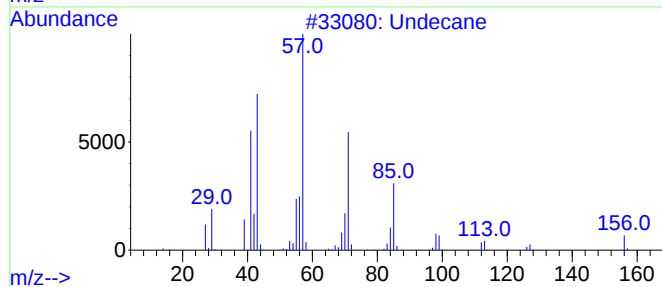
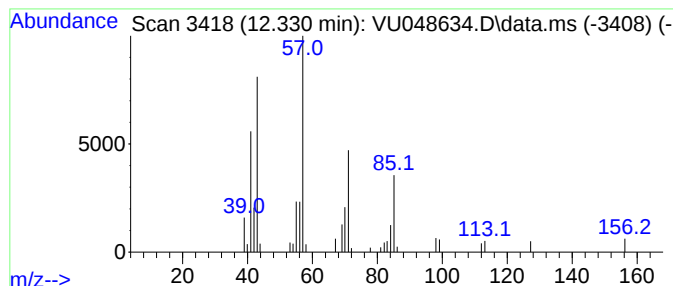
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Quant Title : VOC Analysis

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.330	2.55 ug/L	40391	1,4-Dichlorobenzene-d4	11.816

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane		156	C11H24	001120-21-4	95
2	Undecane		156	C11H24	001120-21-4	87
3	Decane		142	C10H22	000124-18-5	86
4	Undecane		156	C11H24	001120-21-4	74
5	Decane		142	C10H22	000124-18-5	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051922\
 Data File : VU048634.D
 Acq On : 19 May 2022 16:53
 Operator : SY/MD
 Sample : N2848-22ME
 Misc : 4.88g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 EXY12ME

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM051622WMA.M
 Quant Title : VOC Analysis

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Dimethyl ether	1.478	8.5	ug/L	90187	1	6.247	531257	50.0
Methyl propionate	5.199	5.8	ug/L	62065	1	6.247	531257	50.0
2-Butanone, 3-m...	6.163	2.9	ug/L	31106	1	6.247	531257	50.0
Methyl isobutyrate	6.398	5.3	ug/L	56649	1	6.247	531257	50.0
2-Pentanone	6.780	5.6	ug/L	59244	1	6.247	531257	50.0
3-Pentanone	6.928	5.5	ug/L	58270	1	6.247	531257	50.0
Butanoic acid, ...	7.186	10.9	ug/L	115298	1	6.247	531257	50.0
(DEL) Alkane: S...	11.115	16.0	ug/L	253970	3	11.816	792048	50.0
(DEL) Alkane: S...	12.330	2.5	ug/L	40391	3	11.816	792048	50.0