LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050222\
Data File : VU048376.D
Acq On : 02 May 2022 19:30
Operator : SY/MD
Sample : N2639-09
Misc : 5.86g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXL19

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\SFAMULM042522WMA.M
Title : VOC Analysis

Signal : TIC: VU048376.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	35	43	68	rBV	217292	365114	12.43%	1.635%
2	1.600	73	81	117	rVB	118697	347159	11.82%	1.555%
3	1.771	121	134	135	rBV6	14091	23380	0.80%	0.105%
4	2.536	360	372	387	rBV	410594	734537	25.01%	3.289%
5	2.954	493	502	518	rVV2	10020	22438	0.76%	0.100%
6	3.034	519	527	530	rBV7	1877	2609	0.09%	0.012%
7	3.179	557	572	588	rBV2	31945	81781	2.78%	0.366%
8	3.369	628	631	633	rBV3	792	565	0.02%	0.003%
9	3.494	666	670	676	rBV	713	667	0.02%	0.003%
10	3.575	688	695	707	rVB6	3491	7787	0.27%	0.035%
11	3.626	709	711	716	rVB3	887	663	0.02%	0.003%
12	3.655	716	720	722	rBV2	521	449	0.02%	0.002%
13	3.671	722	725	733	rVB3	995	904	0.03%	0.004%
14	3.755	746	751	754	rBV2	720	647	0.02%	0.003%
15	3.784	757	760	764	rBV	584	573	0.02%	0.003%
16	4.054	837	844	847	rBV3	1172	1316	0.04%	0.006%
17	4.243	899	903	908	rBV2	505	679	0.02%	0.003%
18	4.301	916	921	923	rBV2	685	692	0.02%	0.003%
19	4.420	954	958	963	rBV2	613	600	0.02%	0.003%
20	4.462	963	971	976	rBV5	1069	1769	0.06%	0.008%
21	4.636	1011	1025	1039	rBV	233727	694471	23.64%	3.110%
22	5.063	1139	1158	1188	rBV	526419	1266255	43.11%	5.670%
23	5.202	1190	1201	1231	rVB	59503	142944	4.87%	0.640%
24	5.719	1341	1362	1390	rBV2	1156936	2937528	100.00%	13.154%
25	5.973	1430	1441	1455	rBV3	16805	39547	1.35%	0.177%
26	6.099	1478	1480	1483	rBV2	711	497	0.02%	0.002%
27	6.166	1483	1501	1514	rVV	21139	59775	2.03%	0.268%
28	6.247	1514	1526	1555	rVB	349736	726109	24.72%	3.251%
29	6.398	1561	1573	1604	rVB	52315	118541	4.04%	0.531%
30	6.526	1604	1613	1626	rBV5	5746	10949	0.37%	0.049%
31	6.600	1626	1636	1649	rVV2	10321	22011	0.75%	0.099%
32	6.690	1649	1664	1683	rVV	646074	1360966	46.33%	6.094%
33	6.783	1684	1693	1726	rVV2	32650	120839	4.11%	0.541%
34	6.935	1728	1740	1769	rVB	36247	118571	4.04%	0.531%
35	7.189	1807	1819	1849	rBV3	92792	230681	7.85%	1.033%
36	7.369	1871	1875	1876	rBV3	950	619	0.02%	0.003%

LSC Area Percent Report

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Acq On : 02 May 2022 19:30
Operator : SY/MD
Sample : N2639-09
Misc : 5.86g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXL19

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\SFAMULM042522WMA.M
Title : VOC Analysis

37	7.439	1893	1897	1902	rVB4	584	583	0.02%	0.003%
38	7.568	1923	1937	1960	rBV	352235	692673	23.58%	3.102%
39	7.687	1967	1974	1975	rBV3	2045	1755	0.06%	0.008%
40	7.790	1997	2006	2018	rBV3	5047	9793	0.33%	0.044%
41	7.899	2026	2040	2057	rBV	1338861	2459472	83.73%	11.013%
42	8.185	2113	2129	2154	rBV	258727	579837	19.74%	2.596%
43	8.356	2176	2182	2191	rVB6	2045	3302	0.11%	0.015%
44	8.555	2236	2244	2253	rVB4	3685	5565	0.19%	0.025%
45	8.661	2253	2277	2307	rBV	827752	1923822	65.49%	8.615%
46	8.931	2355	2361	2363	rBV3	1022	1181	0.04%	0.005%
47	9.134	2421	2424	2428	rVB2	803	575	0.02%	0.003%
48	9.172	2432	2436	2439	rVB3	726	496	0.02%	0.002%
49	9.314	2477	2480	2484	rVV	797	532	0.02%	0.002%
50	9.423	2500	2514	2540	rBV	459481	929914	31.66%	4.164%
51	9.555	2550	2555	2557	rBV4	1286	1373	0.05%	0.006%
52	9.578	2557	2562	2581	rVB8	2806	5417	0.18%	0.024%
53	9.700	2589	2600	2609	rVV4	6244	11066	0.38%	0.050%
54	9.944	2669	2676	2678	rBV	399	459	0.02%	0.002%
55	10.108	2718	2727	2736	rBV5	3755	6999	0.24%	0.031%
56	10.233	2760	2766	2771	rBV	528	551	0.02%	0.002%
57	10.272	2771	2778	2780	rBV2	542	627	0.02%	0.003%
58	10.439	2826	2830	2840	rBV	624	986	0.03%	0.004%
59	10.484	2840	2844	2846	rBV	907	720	0.02%	0.003%
60	10.539	2856	2861	2867	rBV3	976	974	0.03%	0.004%
61	10.606	2879	2882	2886	rVB2	761	604	0.02%	0.003%
62	10.674	2886	2903	2910	rBV5	4430	9262	0.32%	0.041%
63	10.767	2915	2932	2951	rBV	1189187	1935145	65.88%	8.665%
64	10.896	2967	2972	2980	rVB3	3407	4297	0.15%	0.019%
65	11.060	3014	3023	3030	rVV3	25715	38195	1.30%	0.171%
66	11.118	3030	3041	3058	rVB	353817	516503	17.58%	2.313%
67	11.211	3065	3070	3076	rVB4	1628	1959	0.07%	0.009%
68	11.314	3095	3102	3110	rVV9	2495	3573	0.12%	0.016%
69	11.417	3128	3134	3141	rBV6	3945	5416	0.18%	0.024%
70	11.471	3141	3151	3159	rBV5	4525	9455	0.32%	0.042%
71	11.577	3172	3184	3196	rBV7	3768	7848	0.27%	0.035%
72	11.642	3198	3204	3205	rBV4	1587	1483	0.05%	0.007%
73	11.696	3215	3221	3231	rBV7	2625	3820	0.13%	0.017%
74	11.819	3247	3259	3272	rBV	610401	980479	33.38%	4.390%
75	11.877	3273	3277	3282	rVV2	12520	14899	0.51%	0.067%
76	11.912	3283	3288	3296	rVB3	17131	22999	0.78%	0.103%

LSC Area Percent Report

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Operator : SY/MD
Sample : N2639-09
Misc : 5.86g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXL19

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\SFAMULM042522WMA.M
Title : VOC Analysis

77	12.005	3311	3317	3332	rVB	18399	26864	0.91%	0.120%
78	12.198	3358	3377	3394	rBV	1540058	2486711	84.65%	11.135%
79	12.330	3408	3418	3430	rVB2	70459	99992	3.40%	0.448%
80	12.439	3450	3452	3455	rVB2	905	459	0.02%	0.002%
81	12.523	3475	3478	3483	rBV4	875	754	0.03%	0.003%
82	12.761	3548	3552	3553	rBV	864	645	0.02%	0.003%
83	13.339	3729	3732	3735	rBV2	588	513	0.02%	0.002%
84	13.394	3737	3749	3758	rBV2	31843	47770	1.63%	0.214%
85	13.597	3807	3812	3816	rBV3	842	985	0.03%	0.004%
86	13.754	3855	3861	3865	rBV3	583	714	0.02%	0.003%
87	13.844	3887	3889	3899	rVB3	671	926	0.03%	0.004%
88	14.063	3954	3957	3962	rVB	739	493	0.02%	0.002%
89	14.156	3983	3986	3992	rVB2	997	809	0.03%	0.004%
90	14.188	3992	3996	3998	rBV4	674	569	0.02%	0.003%
91	14.455	4072	4079	4084	rVB5	2120	2714	0.09%	0.012%
92	14.648	4133	4139	4147	rVV2	2559	3705	0.13%	0.017%
93	14.693	4149	4153	4156	rBV3	1171	819	0.03%	0.004%
94	14.728	4161	4164	4166	rVB	1221	642	0.02%	0.003%
95	14.741	4166	4168	4171	rBV	762	446	0.02%	0.002%
96	14.841	4195	4199	4203	rBV	2145	1764	0.06%	0.008%
97	14.941	4226	4230	4234	rBV3	980	1018	0.03%	0.005%
98	15.146	4292	4294	4297	rBV2	755	493	0.02%	0.002%
99	15.182	4302	4305	4308	rBV4	887	706	0.02%	0.003%
100	15.375	4356	4365	4372	rBV5	8531	12372	0.42%	0.055%

Sum of corrected areas: 22332124

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050222\
Data File : VU048376.D
Acq On : 02 May 2022 19:30
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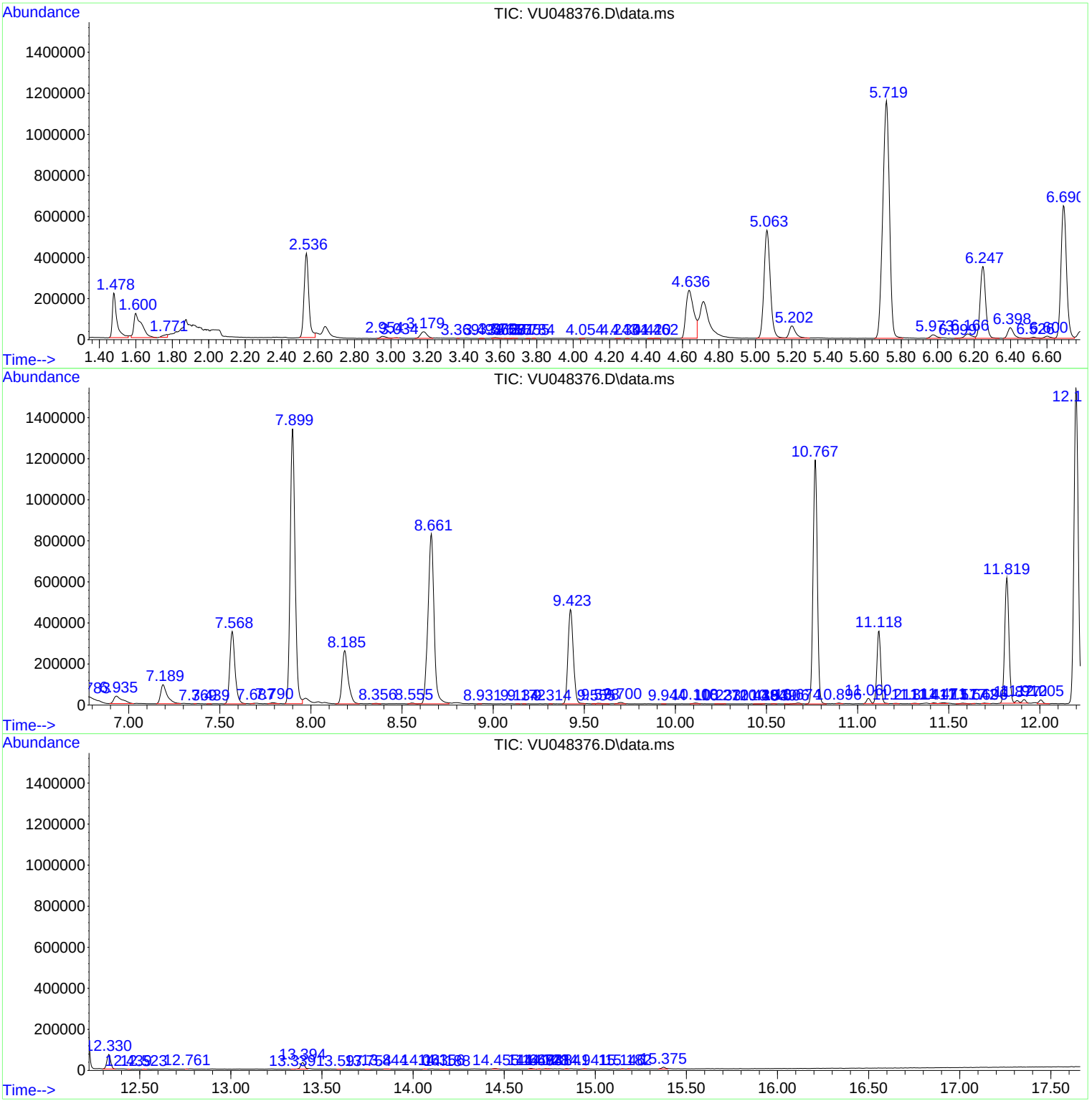
Misc : 5.86g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXL19

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM042522WMA.M
Quant Title : VOC Analysis

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050222\
Data File : VU048376.D
Acq On : 02 May 2022 19:30
Operator : SY/MD
Sample : N2639-09
Misc : 5.86g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXL19

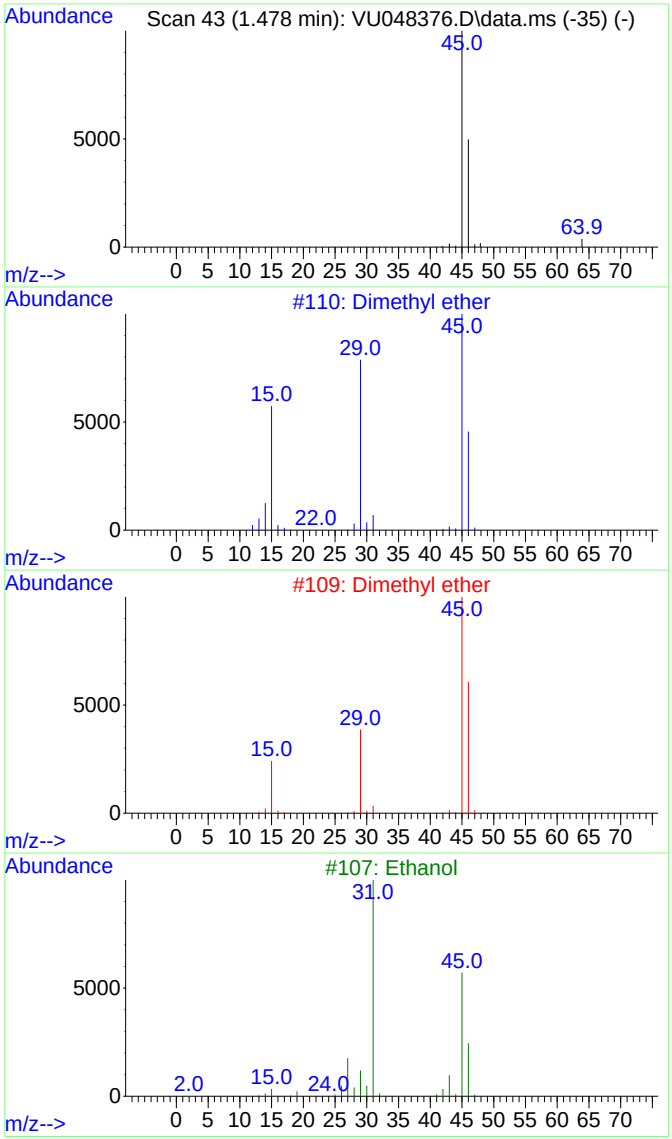
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM042522WMA.M
Quant Title : VOC Analysis

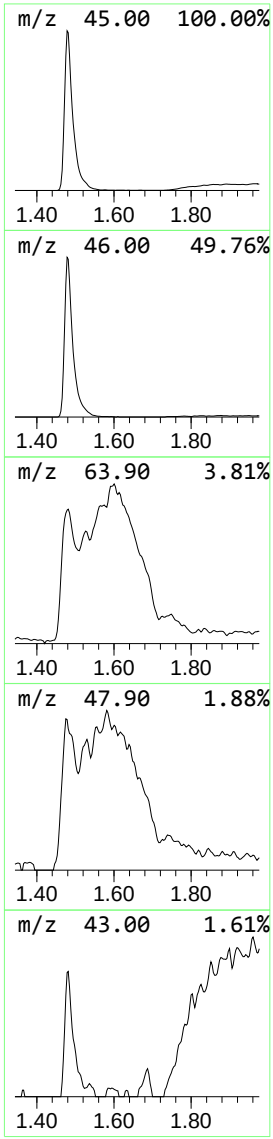
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Dimethyl ether Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.478	25.14 ug/L	365114	1,4-Difluorobenzene	6.250

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dimethyl ether	46	C2H6O	000115-10-6	86
2	Dimethyl ether	46	C2H6O	000115-10-6	64
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





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050222\
Data File : VU048376.D
Acq On : 02 May 2022 19:30
Operator : SY/MD
Sample : N2639-09
Misc : 5.86g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXL19

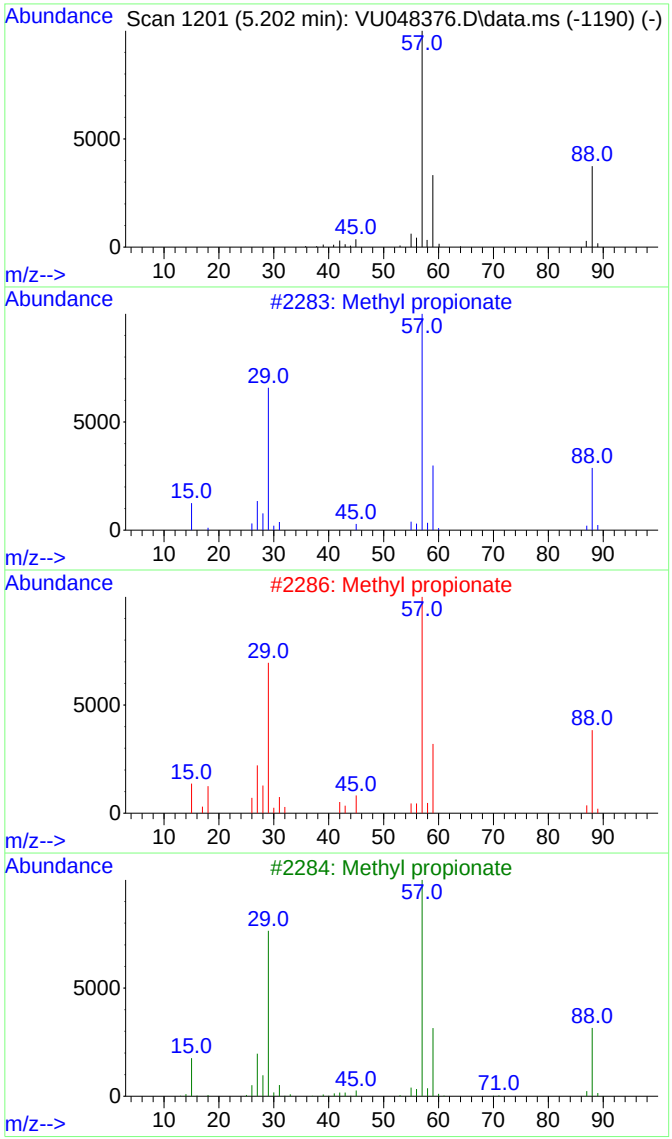
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM042522WMA.M
Quant Title : VOC Analysis

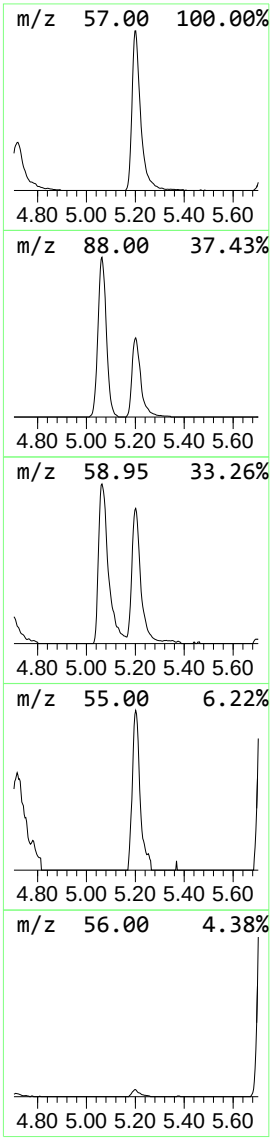
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Methyl propionate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.202	9.84 ug/L	142944	1,4-Difluorobenzene	6.250

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	80
3	Methyl propionate	88	C4H8O2	000554-12-1	80
4	Methyl propionate	88	C4H8O2	000554-12-1	80
5	Methyl propionate	88	C4H8O2	000554-12-1	78





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050222\
Data File : VU048376.D
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ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXL19

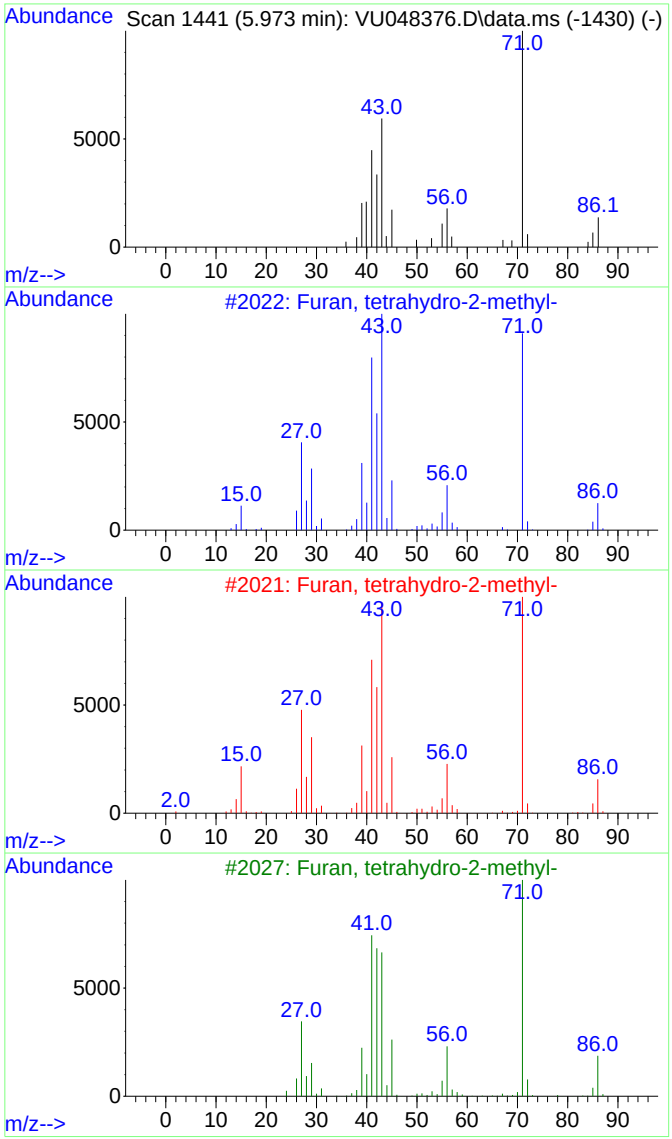
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM042522WMA.M
Quant Title : VOC Analysis

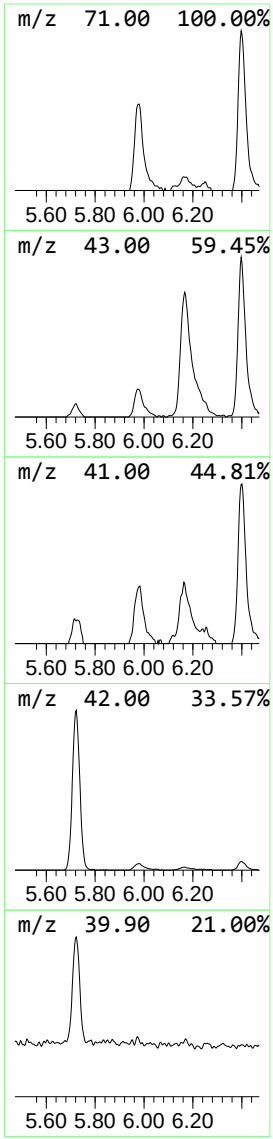
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Furan, tetrahydro-2-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.973	2.72 ug/L	39547	1,4-Difluorobenzene	6.250

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Furan, tetrahydro-2-methyl-	86	C5H10O	000096-47-9	91
2	Furan, tetrahydro-2-methyl-	86	C5H10O	000096-47-9	91
3	Furan, tetrahydro-2-methyl-	86	C5H10O	000096-47-9	78
4	trans-3-Penten-2-ol	86	C5H10O	003899-34-1	53
5	Pentane, 2-methyl-	86	C6H14	000107-83-5	53





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050222\
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ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXL19

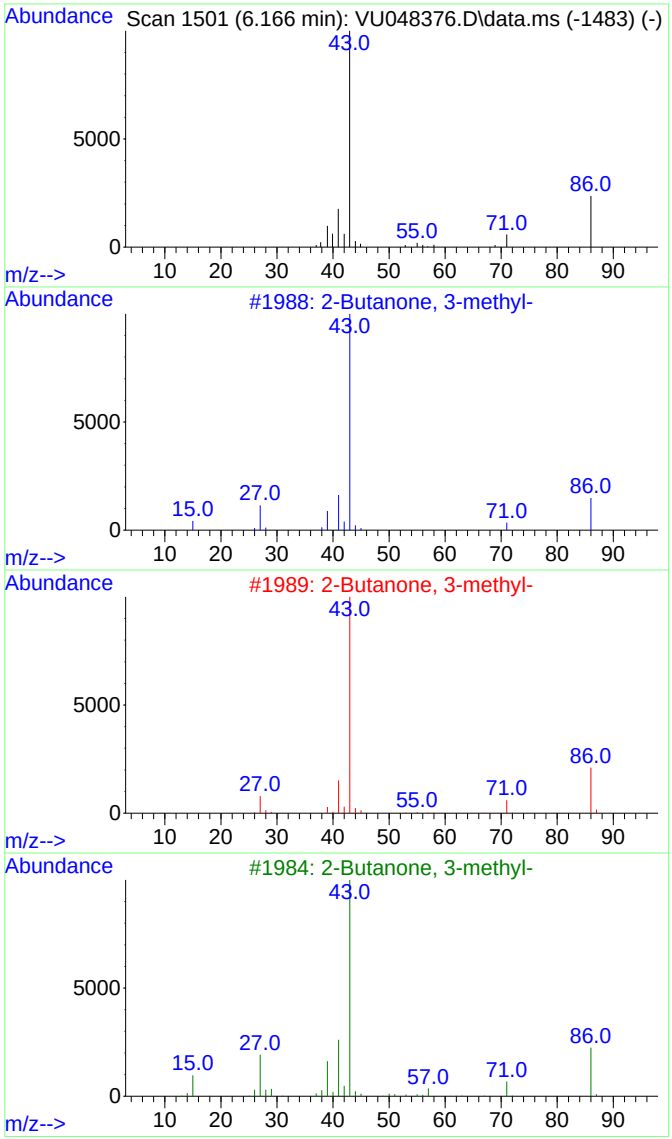
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM042522WMA.M
Quant Title : VOC Analysis

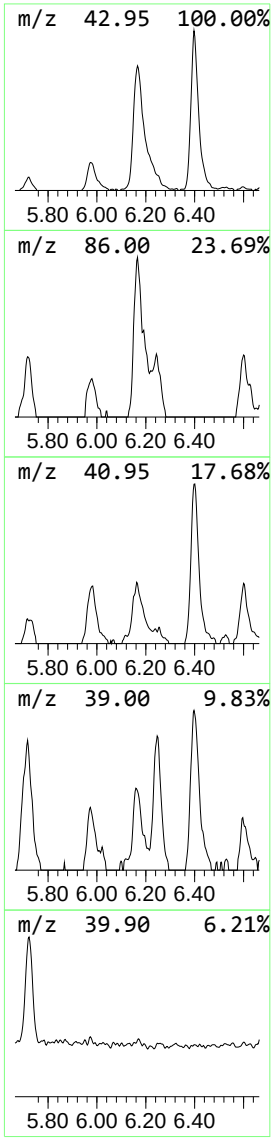
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 2-Butanone, 3-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.166	4.12 ug/L	59775	1,4-Difluorobenzene	6.250

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Butanone, 3-methyl-	86	C5H10O	000563-80-4	72
2	2-Butanone, 3-methyl-	86	C5H10O	000563-80-4	59
3	2-Butanone, 3-methyl-	86	C5H10O	000563-80-4	50
4	2-Pentanone	86	C5H10O	000107-87-9	45
5	2-Pentanone	86	C5H10O	000107-87-9	42





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050222\
Data File : VU048376.D
Acq On : 02 May 2022 19:30
Operator : SY/MD
Sample : N2639-09
Misc : 5.86g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXL19

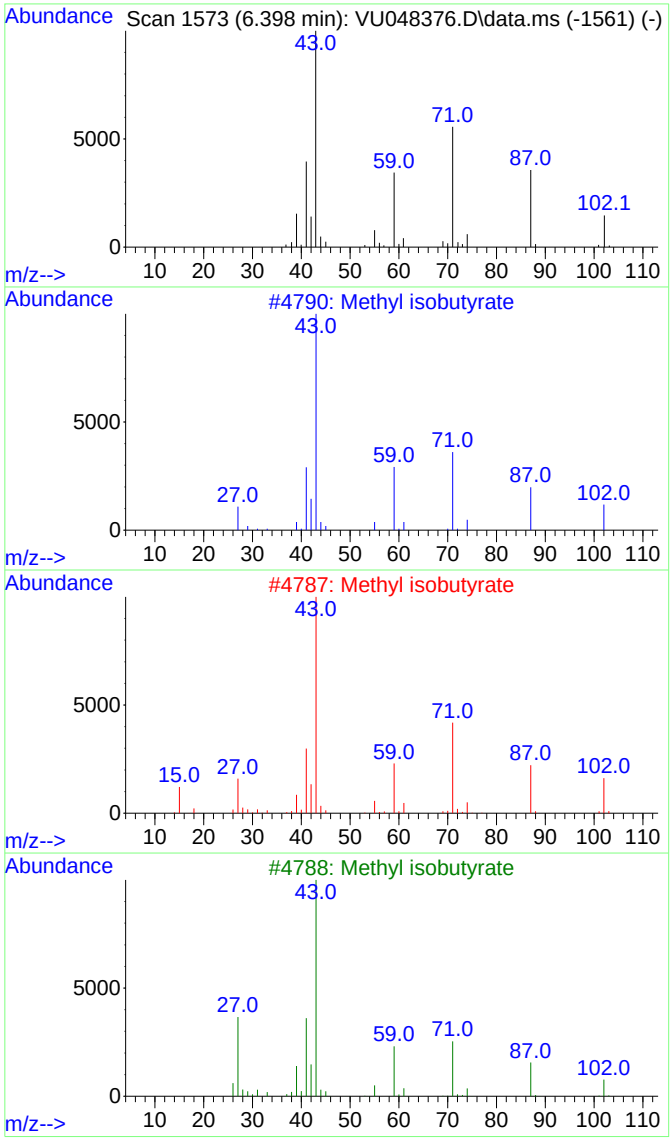
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM042522WMA.M
Quant Title : VOC Analysis

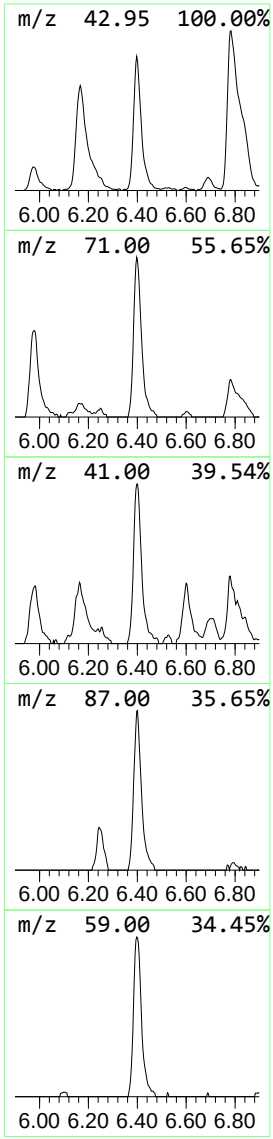
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Methyl isobutyrate Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.398	8.16 ug/L	118541	1,4-Difluorobenzene	6.250

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	72
3	Methyl isobutyrate	102	C5H10O2	000547-63-7	64
4	Methyl isobutyrate	102	C5H10O2	000547-63-7	64
5	Butyric acid hydrazide	102	C4H10N2O	003538-65-6	50





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050222\
Data File : VU048376.D
Acq On : 02 May 2022 19:30
Operator : SY/MD
Sample : N2639-09
Misc : 5.86g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXL19

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM042522WMA.M
Quant Title : VOC Analysis

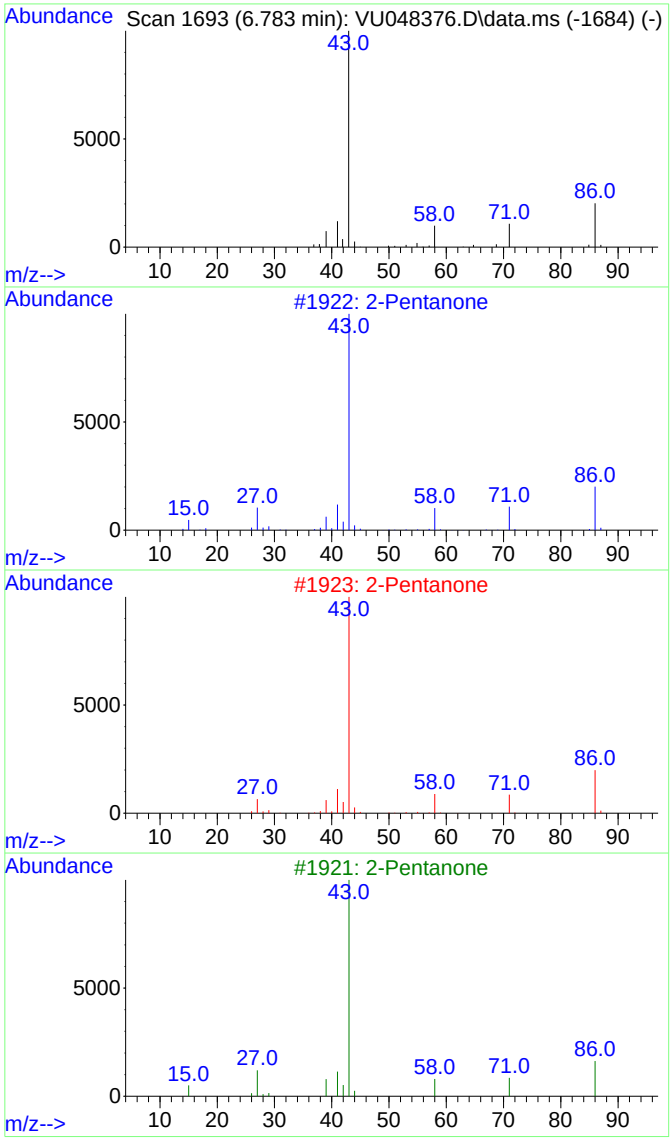
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

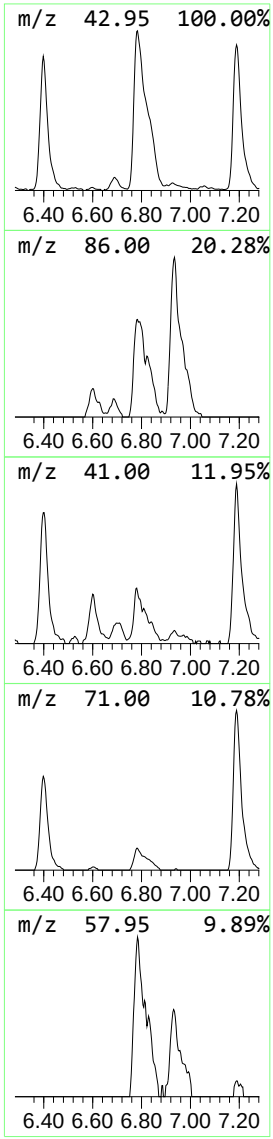
****

Peak Number 7 2-Pentanone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.783	8.32 ug/L	120839	1,4-Difluorobenzene	6.250

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone	86	C5H10O	000107-87-9	80
2	2-Pentanone	86	C5H10O	000107-87-9	80
3	2-Pentanone	86	C5H10O	000107-87-9	59
4	2-Pentanone	86	C5H10O	000107-87-9	59
5	Propane, 2-(ethenyloxy)-	86	C5H10O	000926-65-8	42





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Data File : VU048376.D
Acq On : 02 May 2022 19:30
Operator : SY/MD
Sample : N2639-09
Misc : 5.86g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXL19

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM042522WMA.M
Quant Title : VOC Analysis

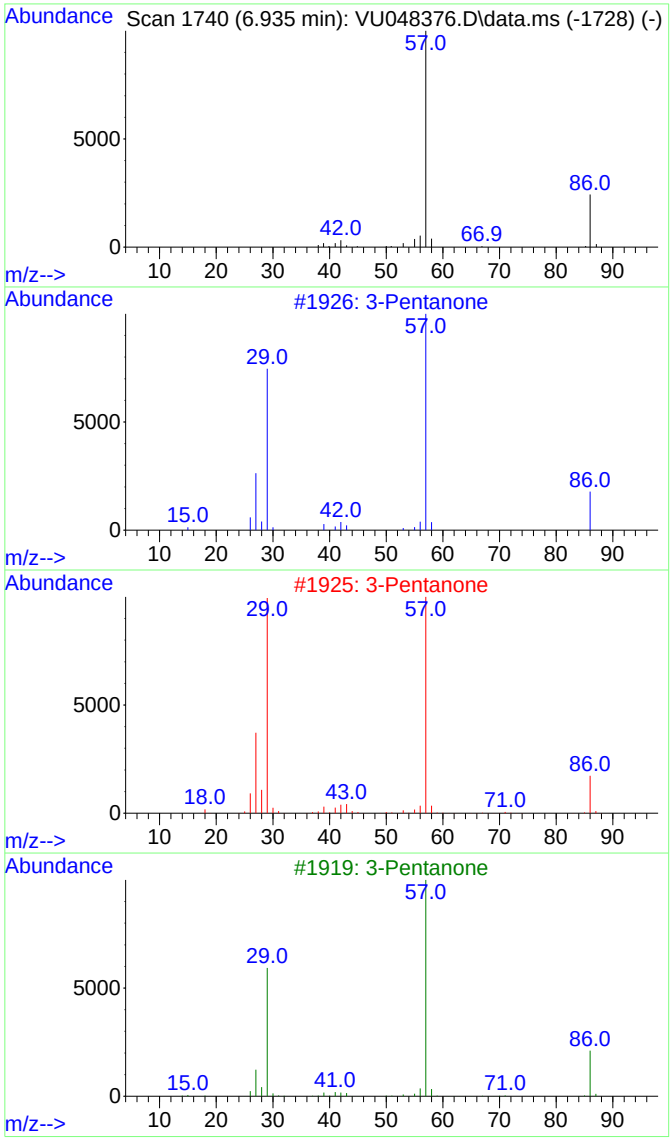
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

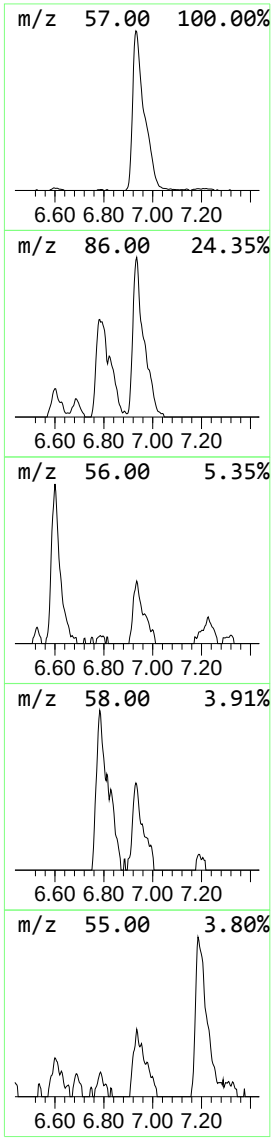
****

Peak Number 8 3-Pentanone Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.935	8.16 ug/L	118571	1,4-Difluorobenzene	6.250

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	72
3	3-Pentanone	86	C5H10O	000096-22-0	72
4	3-Pentanone	86	C5H10O	000096-22-0	72
5	Propanal, 2,2-dimethyl-	86	C5H10O	000630-19-3	9





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Data File : VU048376.D
Acq On : 02 May 2022 19:30
Operator : SY/MD
Sample : N2639-09
Misc : 5.86g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXL19

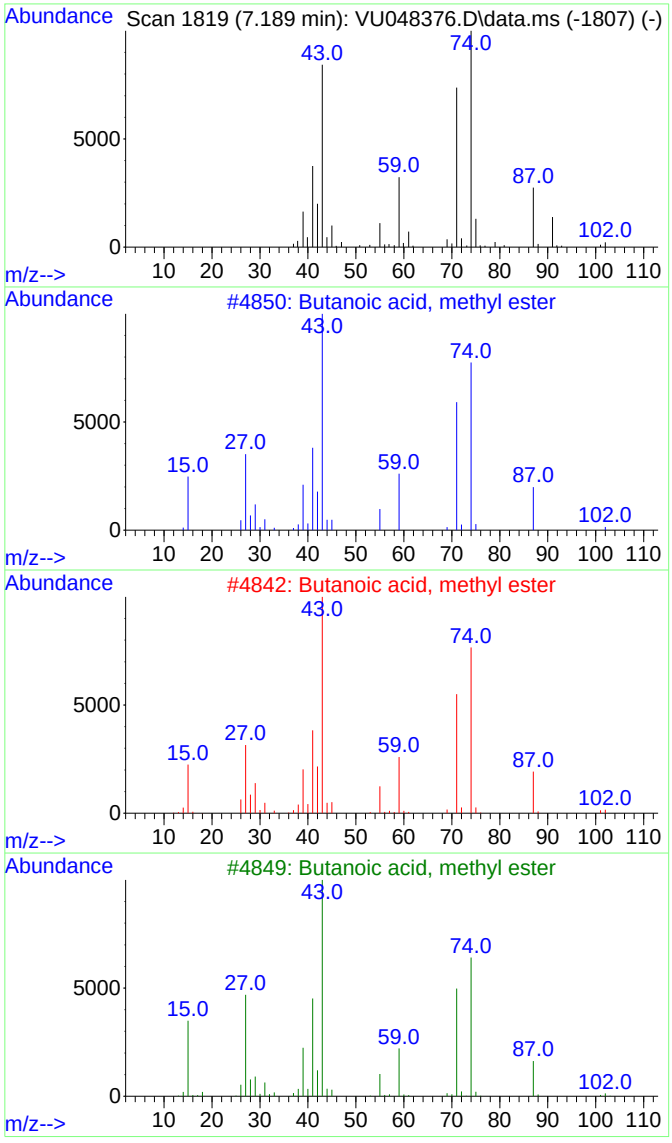
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM042522WMA.M
Quant Title : VOC Analysis

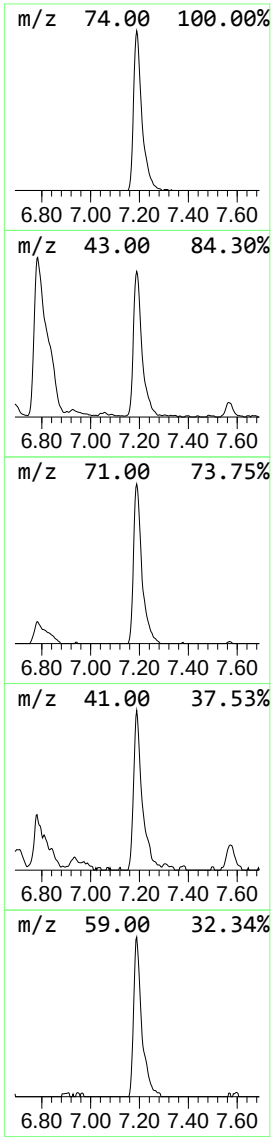
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Butanoic acid, methyl ester Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.189	15.88 ug/L	230681	1,4-Difluorobenzene	6.250

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid, methyl ester	102	C5H10O2	000623-42-7	90
2	Butanoic acid, methyl ester	102	C5H10O2	000623-42-7	90
3	Butanoic acid, methyl ester	102	C5H10O2	000623-42-7	86
4	Butanoic acid, methyl ester	102	C5H10O2	000623-42-7	64
5	Methyl isobutyrate	102	C5H10O2	000547-63-7	37





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Data File : VU048376.D
Acq On : 02 May 2022 19:30
Operator : SY/MD
Sample : N2639-09
Misc : 5.86g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXL19

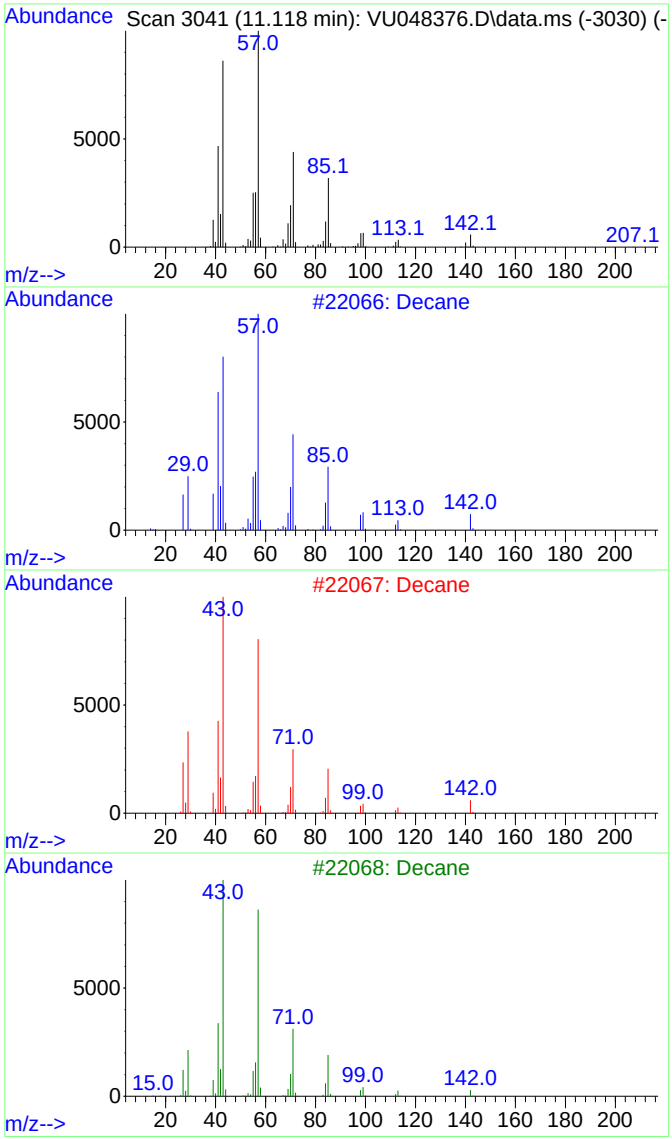
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM042522WMA.M
Quant Title : VOC Analysis

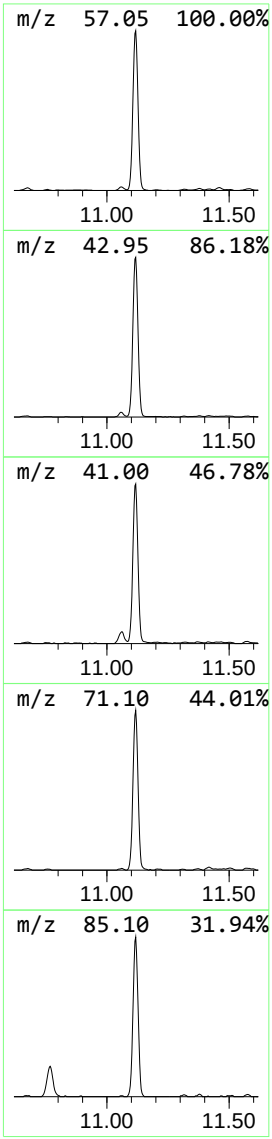
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.118	26.34 ug/L	516503	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane	142	C10H22	000124-18-5	96
2	Decane	142	C10H22	000124-18-5	93
3	Decane	142	C10H22	000124-18-5	87
4	Decane	142	C10H22	000124-18-5	83
5	Octane, 4-ethyl-	142	C10H22	015869-86-0	72





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Data File : VU048376.D
Acq On : 02 May 2022 19:30
Operator : SY/MD
Sample : N2639-09
Misc : 5.86g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXL19

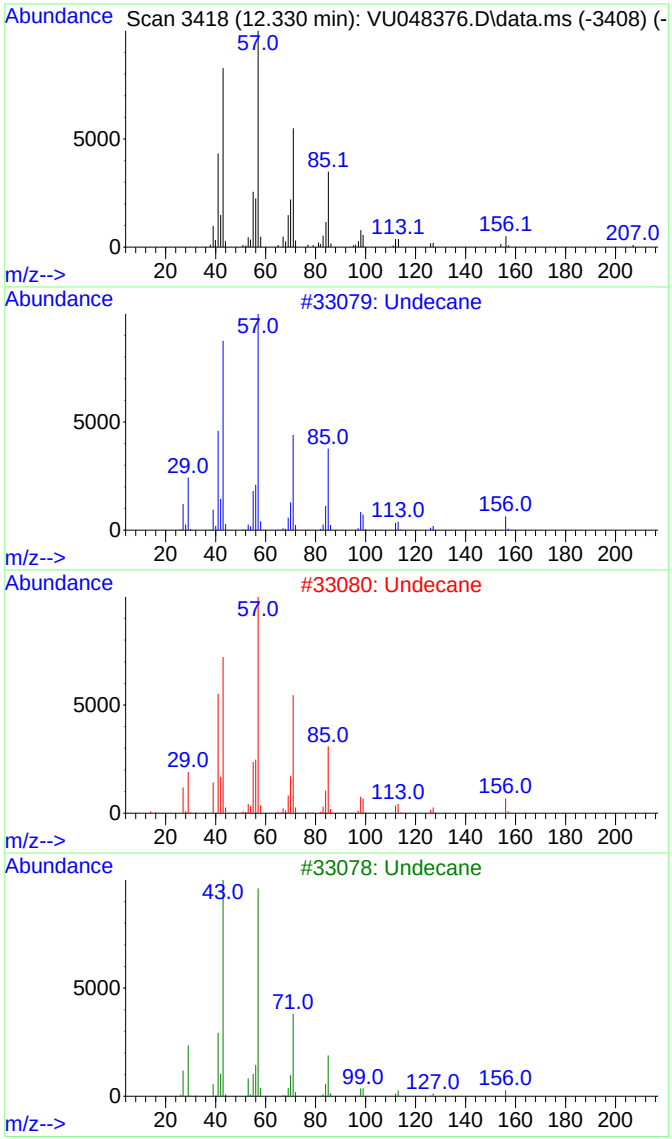
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM042522WMA.M
Quant Title : VOC Analysis

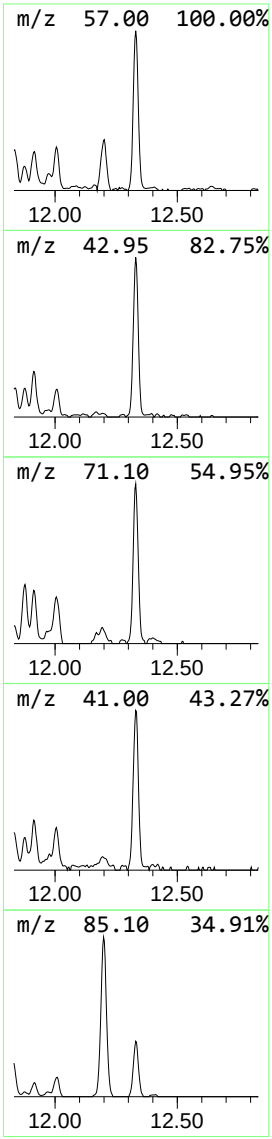
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.330	5.10 ug/L	99992	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane	156	C11H24	001120-21-4	93
2	Undecane	156	C11H24	001120-21-4	90
3	Undecane	156	C11H24	001120-21-4	87
4	Undecane	156	C11H24	001120-21-4	81
5	Carbonic acid, dodecyl prop-1-en...	270	C16H30O3	1000382-90-7	78





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU050222\
Data File : VU048376.D
Acq On : 02 May 2022 19:30
Operator : SY/MD
Sample : N2639-09
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ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXL19

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM042522WMA.M
Quant Title : VOC Analysis

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Dimethyl ether	1.478	25.1	ug/L	365114	1	6.250	726109	50.0
Methyl propionate	5.202	9.8	ug/L	142944	1	6.250	726109	50.0
Furan, tetrahyd...	5.973	2.7	ug/L	39547	1	6.250	726109	50.0
2-Butanone, 3-m...	6.166	4.1	ug/L	59775	1	6.250	726109	50.0
Methyl isobutyrate	6.398	8.2	ug/L	118541	1	6.250	726109	50.0
2-Pentanone	6.783	8.3	ug/L	120839	1	6.250	726109	50.0
3-Pentanone	6.935	8.2	ug/L	118571	1	6.250	726109	50.0
Butanoic acid, ...	7.189	15.9	ug/L	230681	1	6.250	726109	50.0
(DEL) Alkane: S...	11.118	26.3	ug/L	516503	3	11.819	980479	50.0
(DEL) Alkane: S...	12.330	5.1	ug/L	99992	3	11.819	980479	50.0