LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX050924\
Data File : VX041437.D
Acq On : 09 May 2024 16:20
Operator : JC/MD
Sample : PB160794TB 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
VLEB794

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_X\Method\SFAMXLM050924WMA.M
Title : VOC Analysis

Signal : TIC: VX041437.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.142	7	9	12	rBV2	315	314	0.02%	0.003%
2	1.240	21	25	33	rBV2	15038	24296	1.48%	0.195%
3	1.368	43	46	57	rBV	72464	107635	6.58%	0.863%
4	1.453	57	60	67	rVB	4633	6070	0.37%	0.049%
5	1.569	70	79	81	rBV4	6053	15153	0.93%	0.122%
6	1.648	88	92	108	rVB	68800	121963	7.45%	0.978%
7	2.294	191	198	209	rBV	185952	284099	17.36%	2.279%
8	2.392	209	214	229	rVB	20517	39664	2.42%	0.318%
9	2.569	241	243	246	rBV3	293	316	0.02%	0.003%
10	2.709	262	266	273	rVB2	1294	2518	0.15%	0.020%
11	2.788	273	279	296	rVB	11153	24992	1.53%	0.200%
12	2.947	299	305	311	rBV3	1480	3449	0.21%	0.028%
13	3.008	311	315	317	rBV3	700	1109	0.07%	0.009%
14	3.581	403	409	420	rVV2	7141	18322	1.12%	0.147%
15	3.660	420	422	423	rVV2	350	288	0.02%	0.002%
16	3.757	434	438	439	rBV	192	287	0.02%	0.002%
17	4.251	512	519	527	rBV4	1820	6164	0.38%	0.049%
18	4.465	546	554	571	rBV	54772	194015	11.86%	1.556%
19	4.739	594	599	600	rBV2	677	987	0.06%	0.008%
20	5.056	639	651	667	rBV	110105	335770	20.52%	2.693%
21	5.391	700	706	715	rVB4	695	1840	0.11%	0.015%
22	5.556	724	733	739	rBV3	1089	2369	0.14%	0.019%
23	5.964	788	800	818	rBV2	288449	842373	51.48%	6.756%
24	6.348	853	863	878	rBV	62633	173456	10.60%	1.391%
25	6.757	921	930	946	rBV	221164	528330	32.29%	4.237%
26	7.031	974	975	977	rBV2	410	424	0.03%	0.003%
27	7.111	985	988	997	rVB5	1127	2432	0.15%	0.020%
28	7.306	1011	1020	1034	rBV	171743	377118	23.05%	3.025%
29	7.482	1043	1049	1056	rBV5	3655	8873	0.54%	0.071%
30	7.647	1064	1076	1082	rBV	35160	78220	4.78%	0.627%
31	7.799	1094	1101	1123	rBV	903795	1636302	100.00%	13.124%
32	7.964	1124	1128	1139	rVB4	4564	11170	0.68%	0.090%
33	8.116	1151	1153	1157	rVB3	438	495	0.03%	0.004%
34	8.171	1161	1162	1166	rVB3	337	313	0.02%	0.003%
35	8.220	1166	1170	1172	rBV	207	324	0.02%	0.003%
36	8.324	1181	1187	1199	rBV	110123	185982	11.37%	1.492%

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ClientSampleId :
VLEB794

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_X\Method\SFAMXLM050924WMA.M
Title : VOC Analysis

37	8.549	1218	1224	1234	rBV	192941	309632	18.92%	2.483%
38	8.647	1234	1240	1250	rBV	451168	693871	42.40%	5.565%
39	8.818	1263	1268	1273	rVB5	1537	2636	0.16%	0.021%
40	8.885	1278	1279	1280	rBV	396	273	0.02%	0.002%
41	8.952	1284	1290	1303	rBV2	328174	516803	31.58%	4.145%
42	9.244	1329	1338	1345	rBV	266915	387426	23.68%	3.107%
43	9.311	1345	1349	1351	rVV	29545	40923	2.50%	0.328%
44	9.342	1351	1354	1356	rVV	43733	59335	3.63%	0.476%
45	9.385	1356	1361	1372	rVV2	385082	643236	39.31%	5.159%
46	9.482	1372	1377	1388	rVV	467697	637090	38.93%	5.110%
47	9.580	1388	1393	1410	rVB	326723	439171	26.84%	3.522%
48	9.817	1427	1432	1434	rBV4	701	1073	0.07%	0.009%
49	9.915	1442	1448	1464	rBV	869276	1135626	69.40%	9.108%
50	10.055	1464	1471	1478	rBV	449004	659022	40.28%	5.286%
51	10.195	1491	1494	1498	rVB3	636	886	0.05%	0.007%
52	10.281	1503	1508	1517	rBV	8892	14856	0.91%	0.119%
53	10.506	1541	1545	1556	rBV	10669	17602	1.08%	0.141%
54	10.640	1559	1567	1571	rVV	65768	85705	5.24%	0.687%
55	10.683	1571	1574	1585	rVV	20715	30079	1.84%	0.241%
56	10.781	1586	1590	1593	rVV2	1918	3166	0.19%	0.025%
57	10.823	1593	1597	1602	rVB3	1737	2965	0.18%	0.024%
58	11.000	1624	1626	1631	rBV3	198	296	0.02%	0.002%
59	11.092	1638	1641	1644	rBV3	747	979	0.06%	0.008%
60	11.189	1652	1657	1669	rBV	341921	447654	27.36%	3.590%
61	11.287	1669	1673	1675	rBV3	1297	2033	0.12%	0.016%
62	11.366	1683	1686	1689	rBV4	192	316	0.02%	0.003%
63	11.439	1696	1698	1699	rVV2	251	265	0.02%	0.002%
64	11.451	1699	1700	1704	rVB2	496	472	0.03%	0.004%
65	11.573	1717	1720	1722	rVB3	214	253	0.02%	0.002%
66	11.628	1725	1729	1736	rVB3	877	1565	0.10%	0.013%
67	11.707	1739	1742	1747	rVB4	685	946	0.06%	0.008%
68	11.762	1747	1751	1757	rBV2	5395	7713	0.47%	0.062%
69	11.847	1761	1765	1771	rVB4	1633	2707	0.17%	0.022%
70	11.933	1776	1779	1781	rBV2	318	446	0.03%	0.004%
71	11.969	1783	1785	1788	rVB2	341	322	0.02%	0.003%
72	12.018	1788	1793	1803	rBV	472600	613287	37.48%	4.919%
73	12.225	1823	1827	1828	rBV2	973	1139	0.07%	0.009%
74	12.250	1829	1831	1837	rVB5	1355	1851	0.11%	0.015%
75	12.317	1837	1842	1860	rBV	506150	653872	39.96%	5.244%
76	12.573	1880	1884	1891	rVB2	1627	2493	0.15%	0.020%

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Instrument :
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ClientSampleId :
VLEB794

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
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Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM050924WMA.M
Title : VOC Analysis

77	12.695	1903	1904	1907	rVV	347	280	0.02%	0.002%
78	12.774	1914	1917	1919	rBV	302	340	0.02%	0.003%
79	12.799	1919	1921	1925	rVB2	323	377	0.02%	0.003%
80	12.866	1929	1932	1937	rVV4	581	935	0.06%	0.007%
81	13.116	1969	1973	1976	rBV3	437	624	0.04%	0.005%
82	13.597	2049	2052	2054	rBV3	516	618	0.04%	0.005%
83	13.695	2064	2068	2069	rBV3	432	496	0.03%	0.004%
84	13.786	2078	2083	2090	rVB3	416	857	0.05%	0.007%
85	13.920	2101	2105	2108	rBV3	201	307	0.02%	0.002%
86	13.969	2110	2113	2117	rVB2	514	614	0.04%	0.005%
87	14.115	2134	2137	2138	rBV2	303	289	0.02%	0.002%
88	14.176	2143	2147	2148	rBV2	204	255	0.02%	0.002%
89	14.603	2214	2217	2220	rBV3	244	283	0.02%	0.002%
90	14.993	2277	2281	2289	rVV4	679	979	0.06%	0.008%
91	15.073	2292	2294	2298	rVB3	219	252	0.02%	0.002%
92	15.298	2328	2331	2333	rBV3	415	388	0.02%	0.003%
93	15.530	2366	2369	2373	rVB3	342	551	0.03%	0.004%
94	15.566	2373	2375	2376	rBV	345	354	0.02%	0.003%
95	15.737	2402	2403	2405	rBV	416	315	0.02%	0.003%
96	15.877	2423	2426	2427	rBV2	342	355	0.02%	0.003%
97	15.926	2432	2434	2435	rBV	529	377	0.02%	0.003%
98	16.200	2477	2479	2480	rBV	334	297	0.02%	0.002%
99	16.310	2495	2497	2499	rBV2	420	407	0.02%	0.003%
100	16.365	2504	2506	2509	rBV2	487	501	0.03%	0.004%

Sum of corrected areas: 12468468

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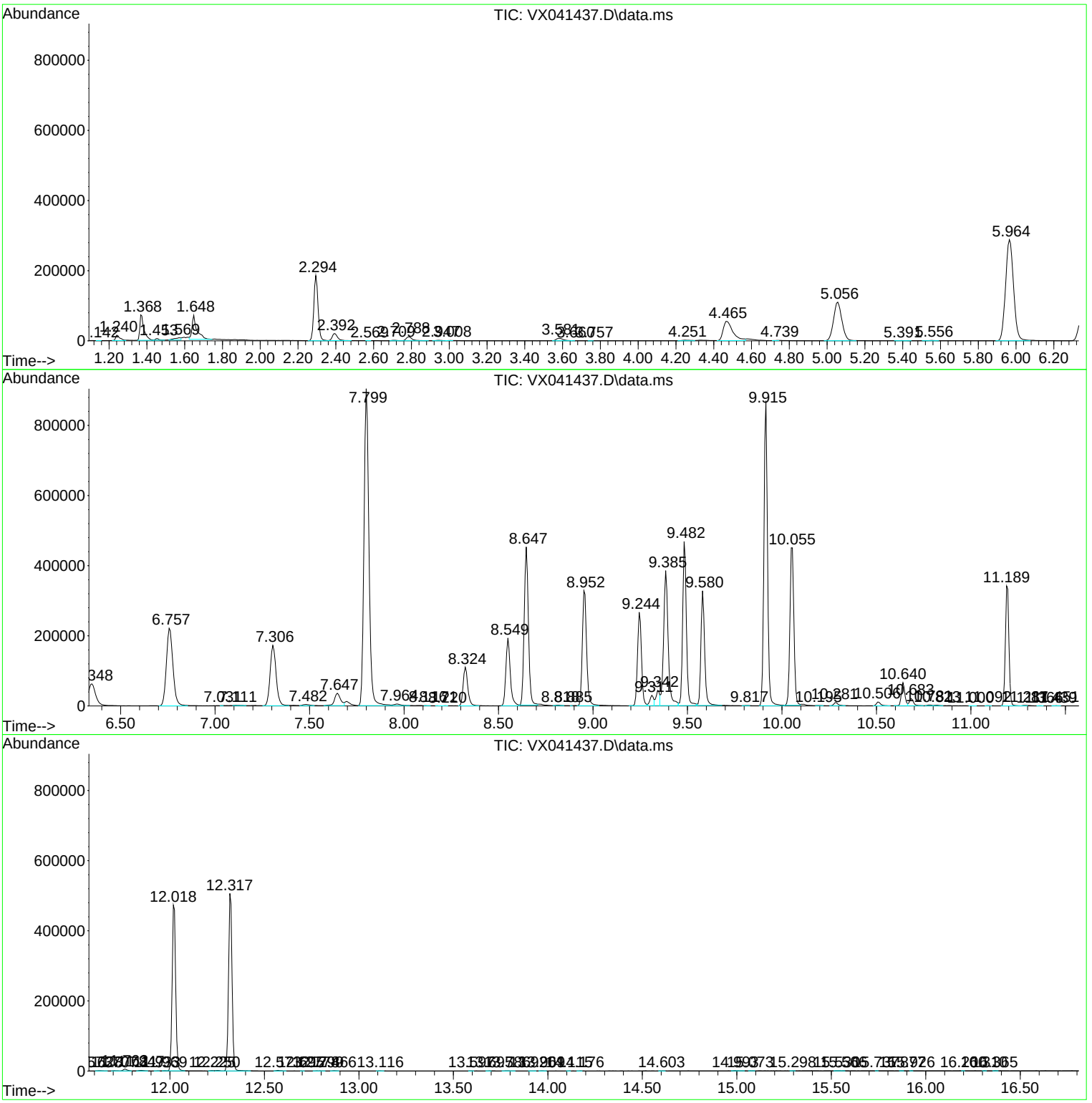
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
VLEB794

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM050924WMA.M
Quant Title : VOC Analysis

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



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ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
VLEB794

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM050924WMA.M
Quant Title : VOC Analysis

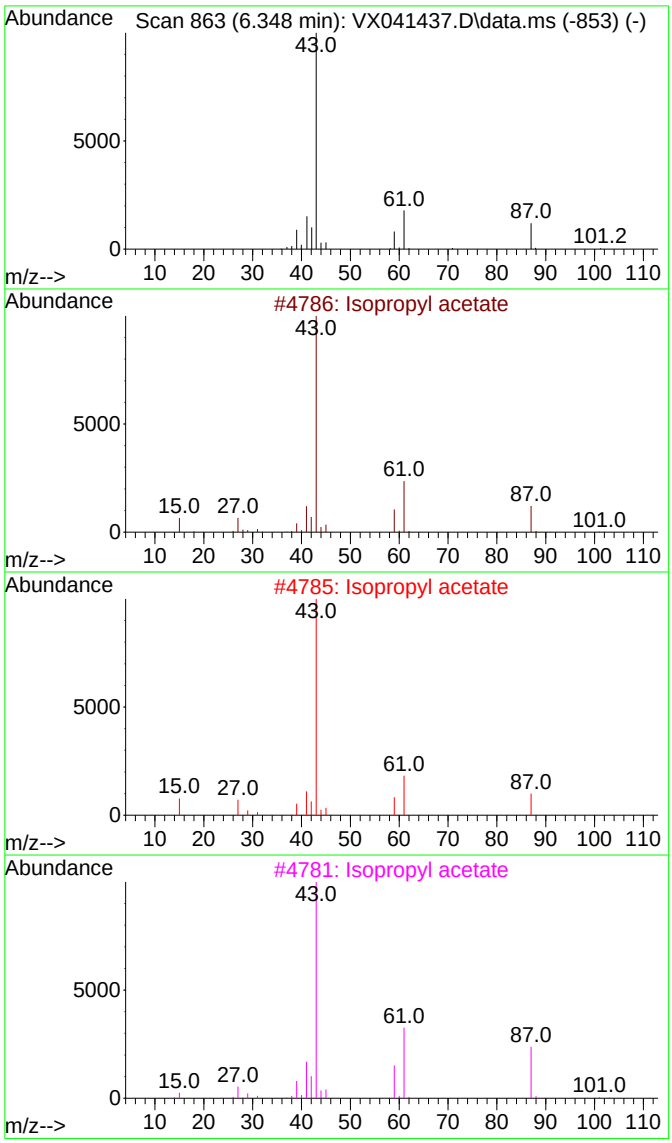
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

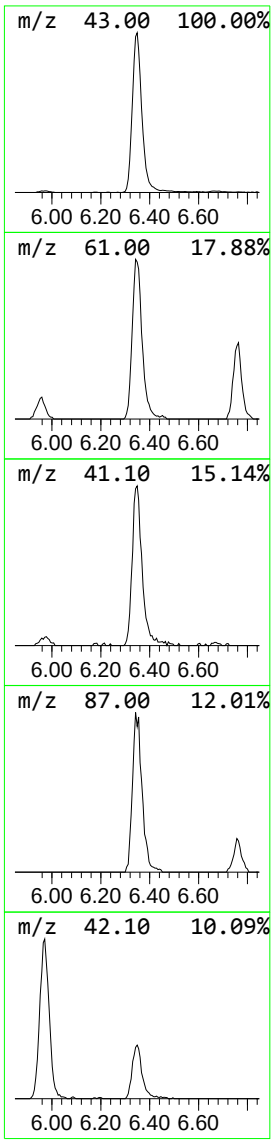
****

Peak Number 2 Isopropyl acetate Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.348	16.42 ug/L	173456	1,4-Difluorobenzene	6.757

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Isopropyl acetate	102	C5H10O2	000108-21-4	83
2	Isopropyl acetate	102	C5H10O2	000108-21-4	72
3	Isopropyl acetate	102	C5H10O2	000108-21-4	59
4	Isopropyl acetate	102	C5H10O2	000108-21-4	56
5	Acetaldehyde, O-ethyloxime-	87	C4H9NO	1000288-34-6	16





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX050924\
Data File : VX041437.D
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Sample : PB160794TB 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
VLEB794

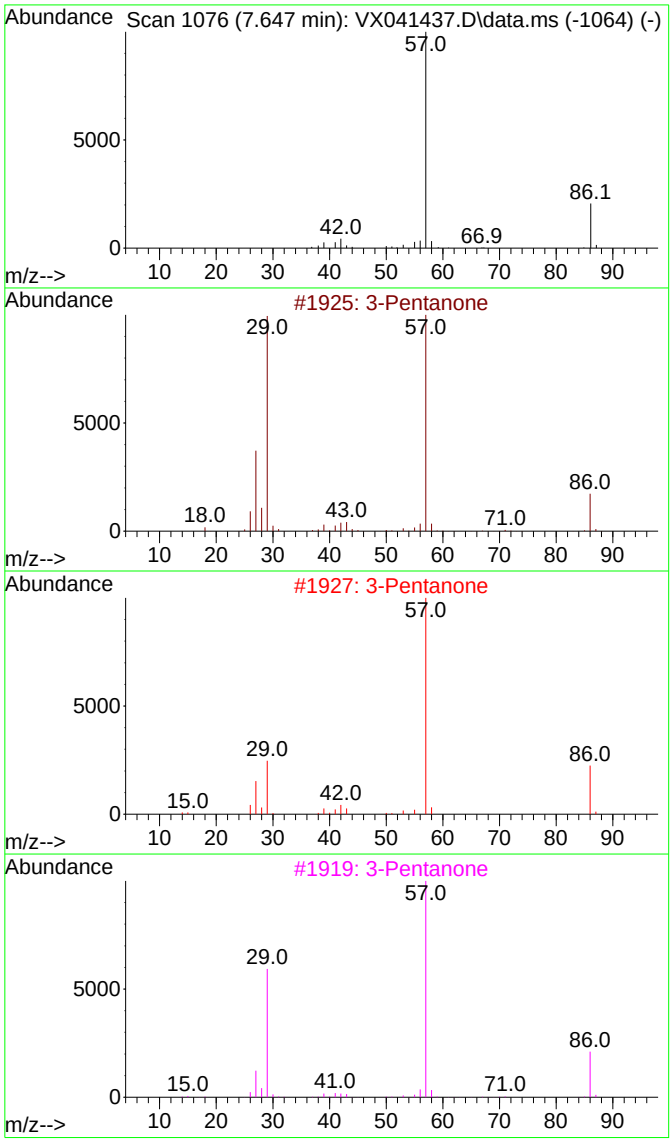
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM050924WMA.M
Quant Title : VOC Analysis

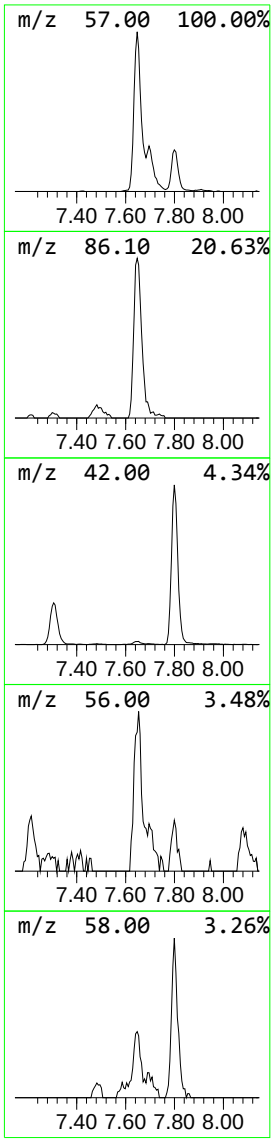
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 3-Pentanone Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.647	7.40 ug/L	78220	1,4-Difluorobenzene	6.757

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Pentanone	86	C5H10O	000096-22-0	86
2	3-Pentanone	86	C5H10O	000096-22-0	80
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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ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM050924WMA.M
Quant Title : VOC Analysis

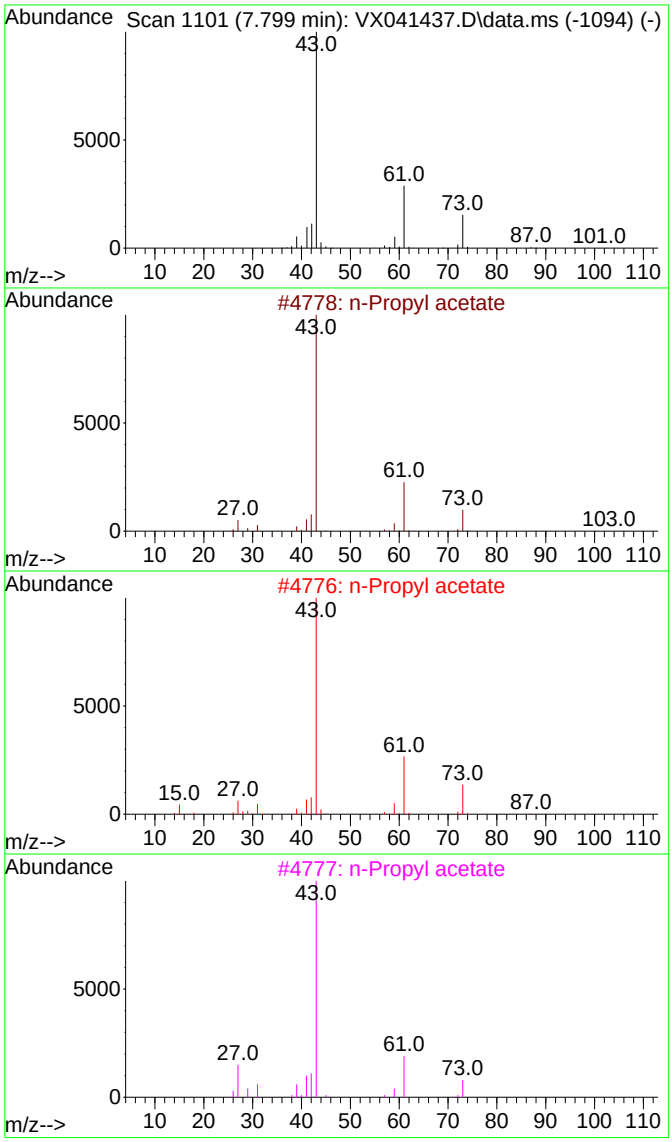
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

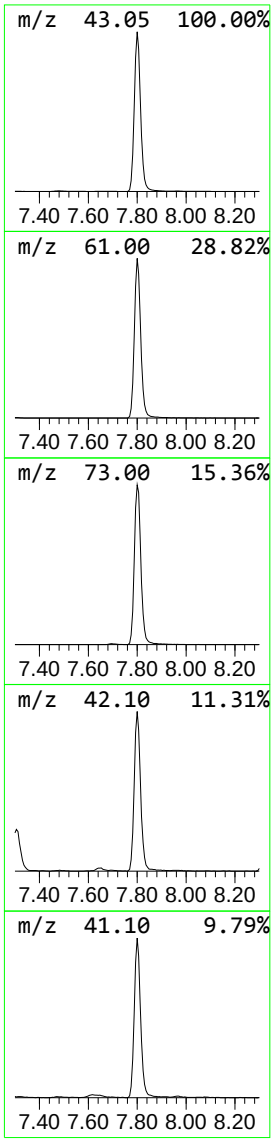
****

Peak Number 4 n-Propyl acetate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.799	154.86 ug/L	1636300	1,4-Difluorobenzene	6.757

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Propyl acetate	102	C5H10O2	000109-60-4	83
2	n-Propyl acetate	102	C5H10O2	000109-60-4	78
3	n-Propyl acetate	102	C5H10O2	000109-60-4	78
4	n-Propyl acetate	102	C5H10O2	000109-60-4	78
5	n-Propyl acetate	102	C5H10O2	000109-60-4	64





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Instrument :
MSVOA_X
ClientSampleId :
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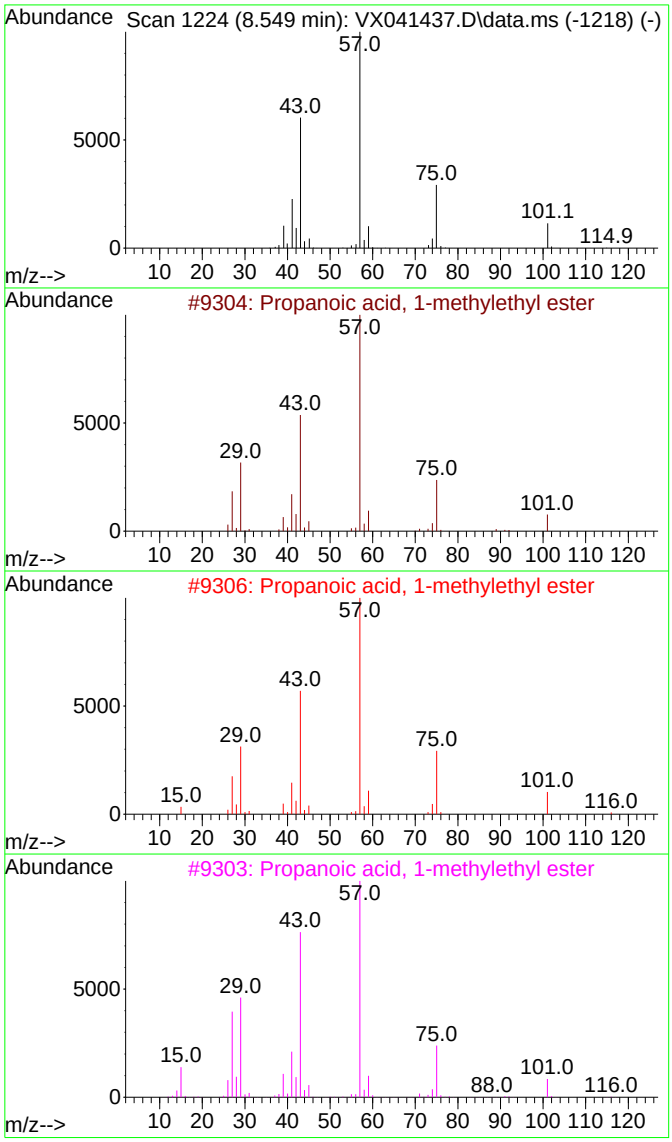
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM050924WMA.M
Quant Title : VOC Analysis

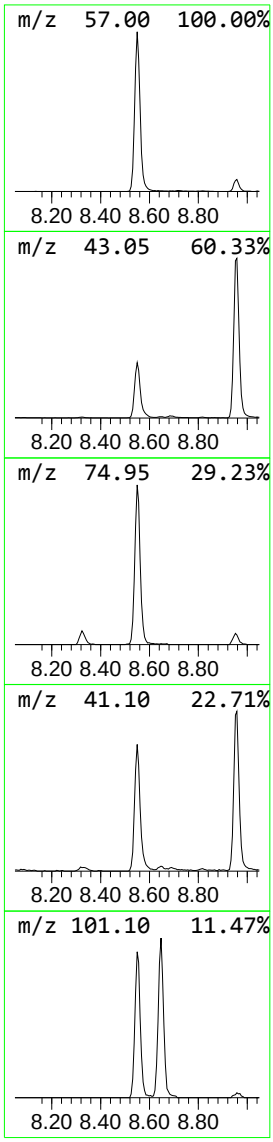
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Propanoic acid, 1-methyleth... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.549	23.49 ug/L	309632	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propanoic acid, 1-methylethyl ester	116	C6H12O2	000637-78-5	90
2	Propanoic acid, 1-methylethyl ester	116	C6H12O2	000637-78-5	83
3	Propanoic acid, 1-methylethyl ester	116	C6H12O2	000637-78-5	74
4	Propanoic acid, 1-methylethyl ester	116	C6H12O2	000637-78-5	59
5	Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	53





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Operator : JC/MD
Sample : PB160794TB 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
VLEB794

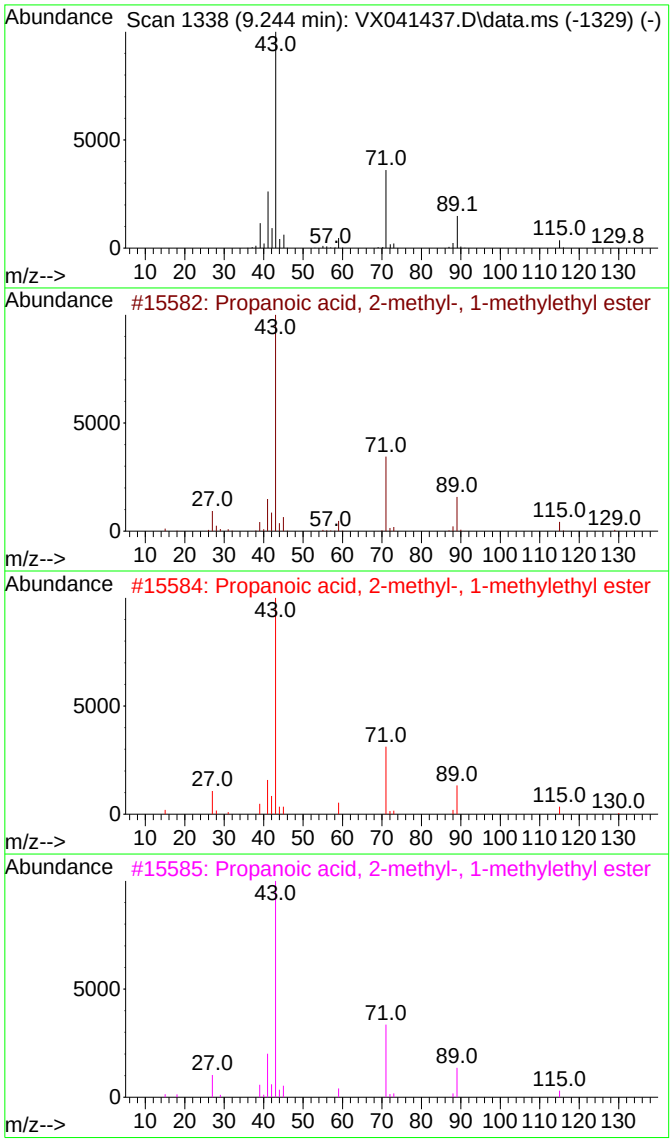
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM050924WMA.M
Quant Title : VOC Analysis

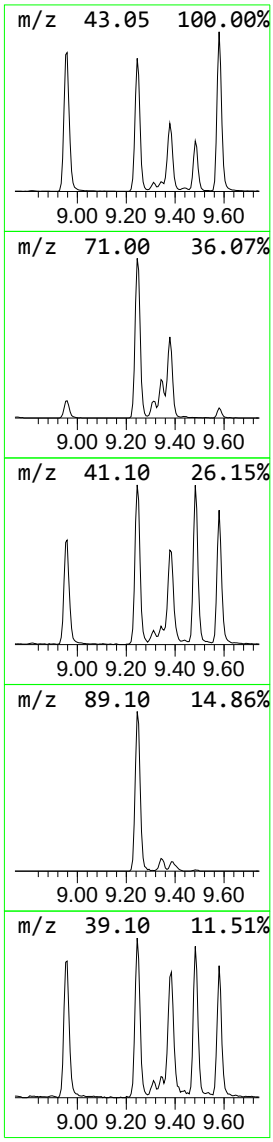
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Propanoic acid, 2-methyl-, ... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.244	29.39 ug/L	387426	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	83
2	Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	83
3	Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	83
4	Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	78
5	Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	72





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX050924\
Data File : VX041437.D
Acq On : 09 May 2024 16:20
Operator : JC/MD
Sample : PB160794TB 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
VLEB794

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM050924WMA.M
Quant Title : VOC Analysis

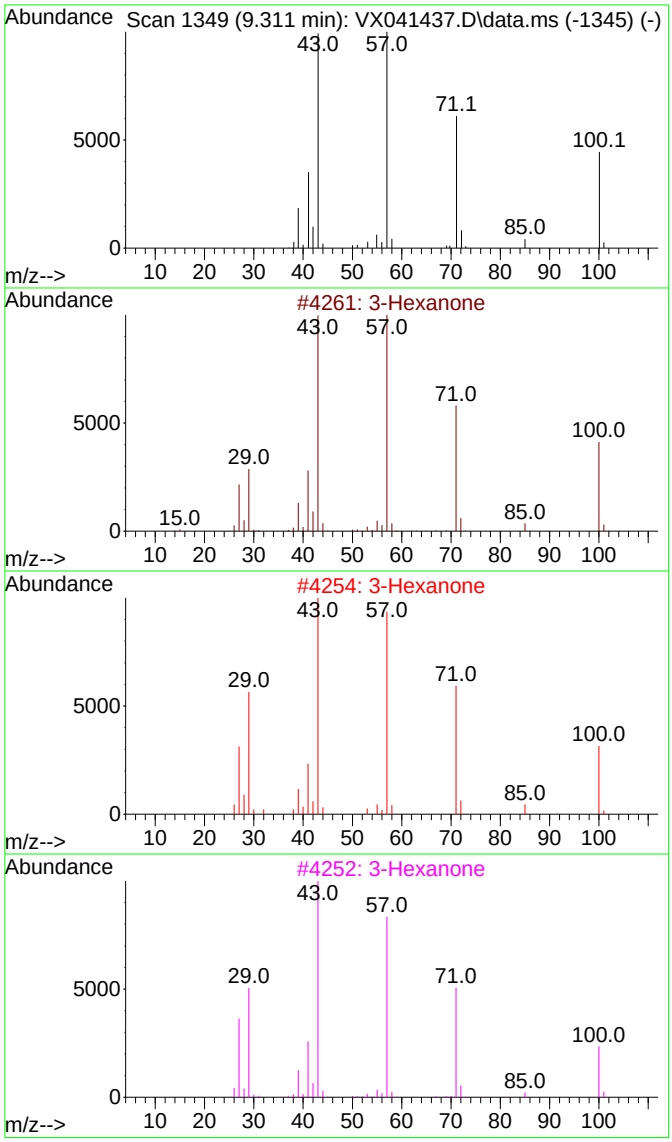
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

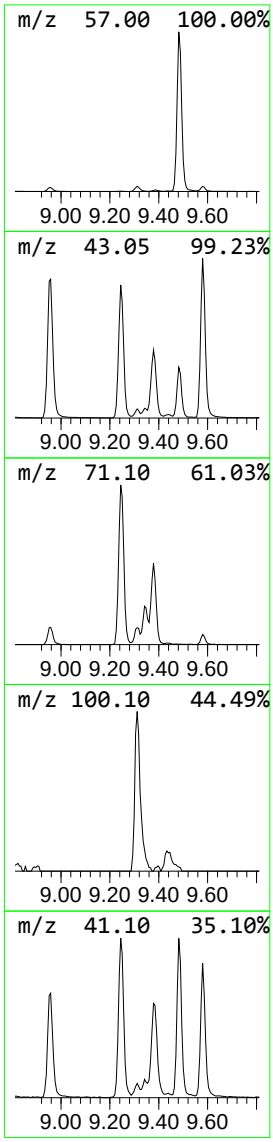
****

Peak Number 8 3-Hexanone Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.311	3.10 ug/L	40923	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Hexanone	100	C6H12O	000589-38-8	91
2	3-Hexanone	100	C6H12O	000589-38-8	59
3	3-Hexanone	100	C6H12O	000589-38-8	59
4	3-Hexanone	100	C6H12O	000589-38-8	59
5	3-Hexanone	100	C6H12O	000589-38-8	53





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Data File : VX041437.D
Acq On : 09 May 2024 16:20
Operator : JC/MD
Sample : PB160794TB 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
VLEB794

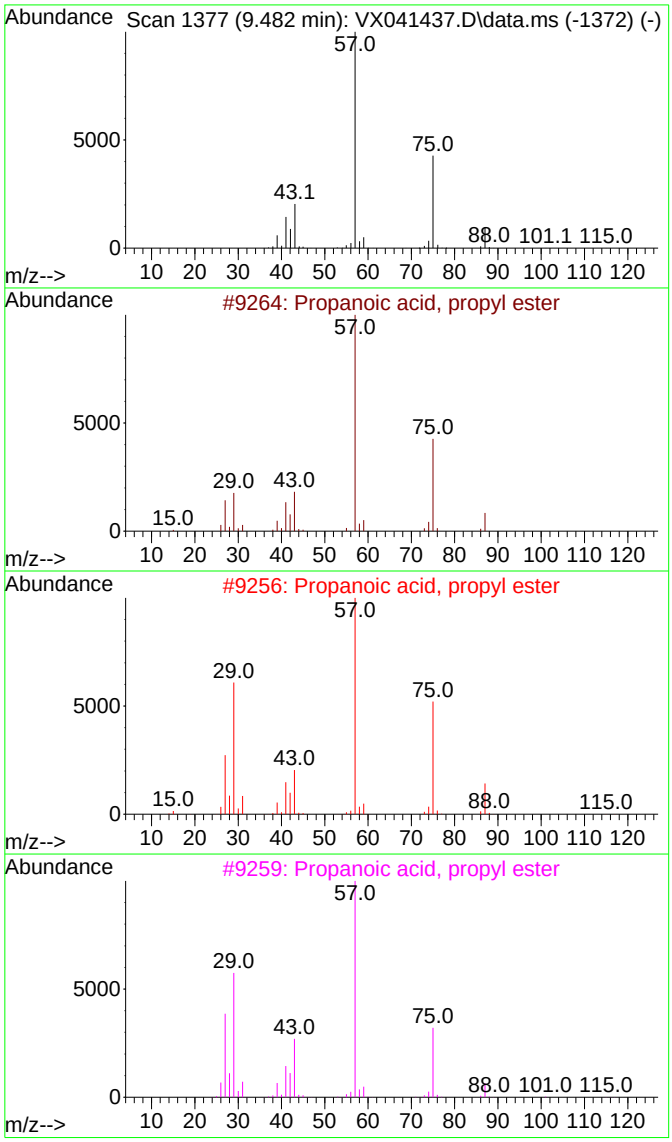
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM050924WMA.M
Quant Title : VOC Analysis

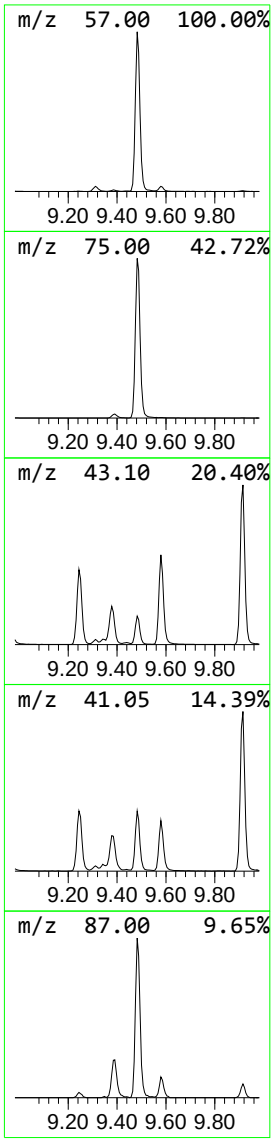
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Propanoic acid, propyl ester Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.482	48.34 ug/L	637090	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	83
2	Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	83
3	Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	83
4	Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	83
5	Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	83





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Data File : VX041437.D
Acq On : 09 May 2024 16:20
Operator : JC/MD
Sample : PB160794TB 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
VLEB794

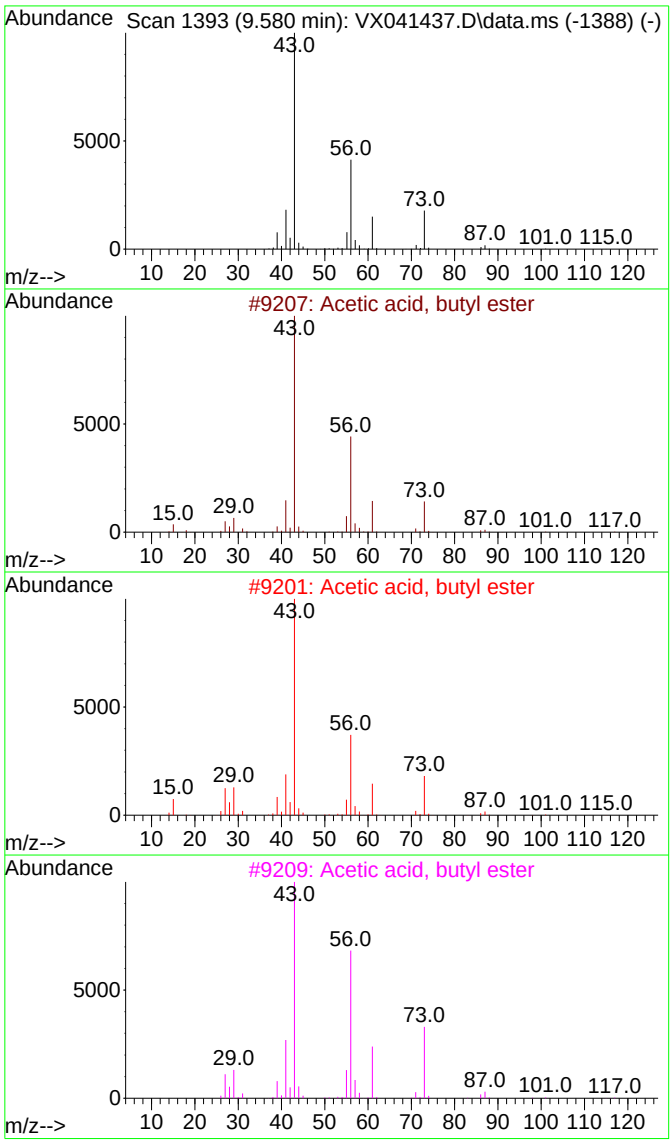
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM050924WMA.M
Quant Title : VOC Analysis

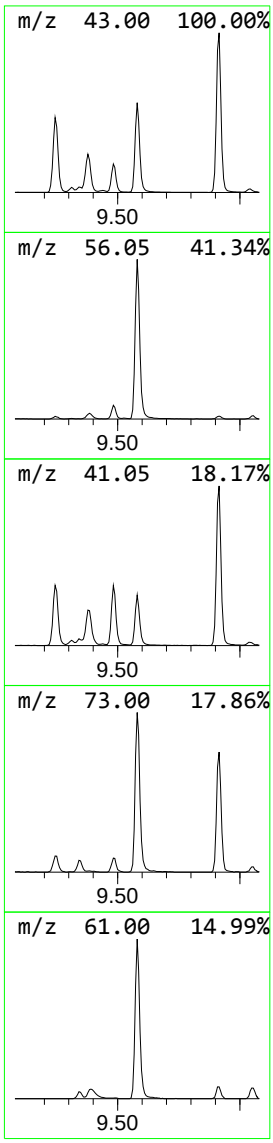
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Acetic acid, butyl ester Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.580	33.32 ug/L	439171	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetic acid, butyl ester	116	C6H12O2	000123-86-4	83
2	Acetic acid, butyl ester	116	C6H12O2	000123-86-4	83
3	Acetic acid, butyl ester	116	C6H12O2	000123-86-4	72
4	Acetic acid, butyl ester	116	C6H12O2	000123-86-4	72
5	Isobutyl acetate	116	C6H12O2	000110-19-0	40





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX050924\
Data File : VX041437.D
Acq On : 09 May 2024 16:20
Operator : JC/MD
Sample : PB160794TB 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
VLEB794

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM050924WMA.M
Quant Title : VOC Analysis

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

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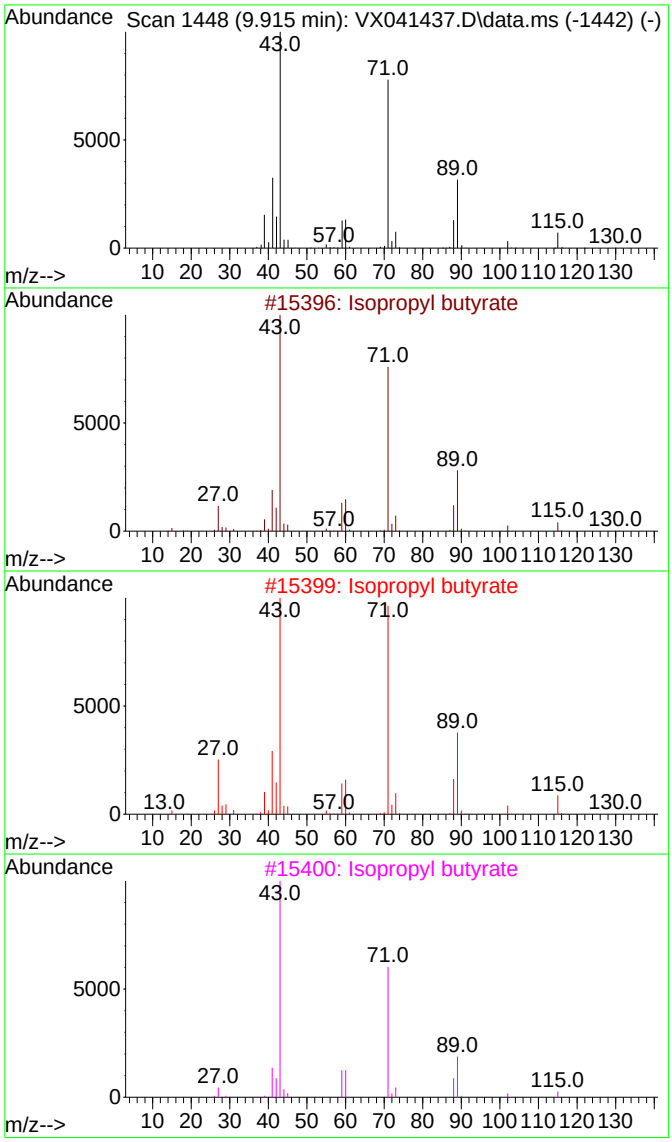
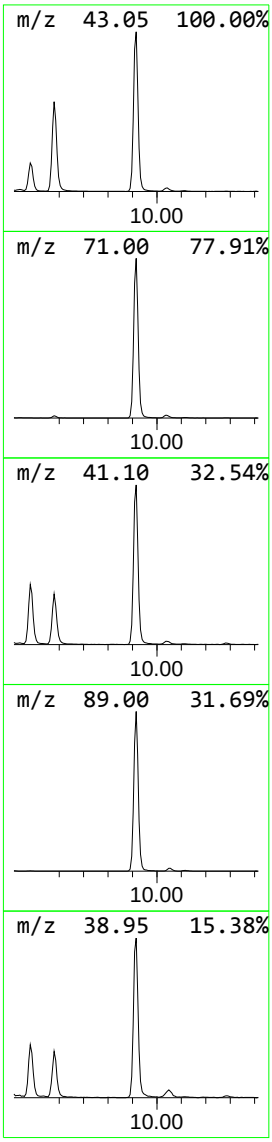
Peak Number 11 Isopropyl butyrate Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.915	86.16 ug/L	1135630	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Isopropyl butyrate	130	C7H14O2	000638-11-9	97
2	Isopropyl butyrate	130	C7H14O2	000638-11-9	91
3	Isopropyl butyrate	130	C7H14O2	000638-11-9	78
4	Butanoic acid, propyl ester	130	C7H14O2	000105-66-8	64
5	Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX050924\
Data File : VX041437.D
Acq On : 09 May 2024 16:20
Operator : JC/MD
Sample : PB160794TB 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
VLEB794

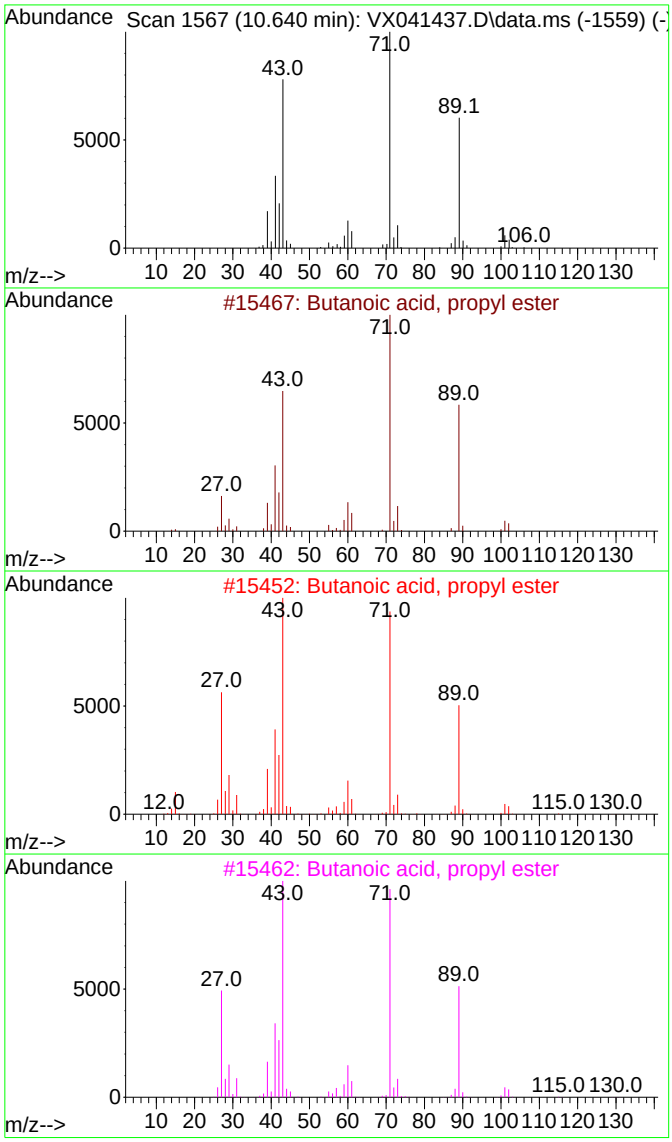
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM050924WMA.M
Quant Title : VOC Analysis

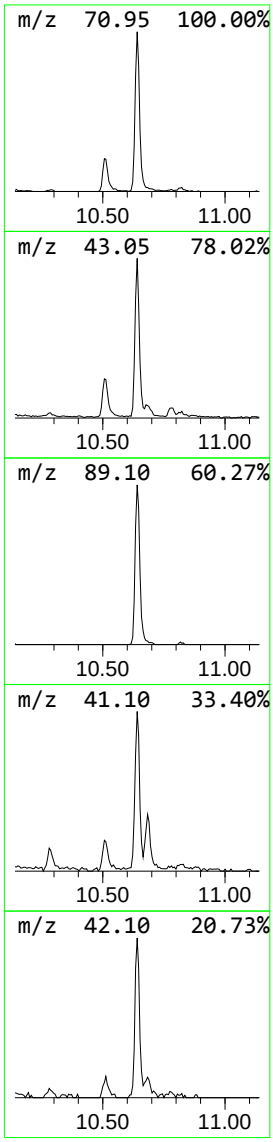
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Butanoic acid, propyl ester Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.640	6.50 ug/L	85705	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butanoic acid, propyl ester	130	C7H14O2	000105-66-8	90
2	Butanoic acid, propyl ester	130	C7H14O2	000105-66-8	90
3	Butanoic acid, propyl ester	130	C7H14O2	000105-66-8	90
4	Butanoic acid, propyl ester	130	C7H14O2	000105-66-8	90
5	Butanoic acid, hexyl ester	172	C10H20O2	002639-63-6	83





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX050924\
Data File : VX041437.D
Acq On : 09 May 2024 16:20
Operator : JC/MD
Sample : PB160794TB 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
VLEB794

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM050924WMA.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
Isopropyl acetate	6.348	16.4	ug/L	173456	1	6.757	528330	50.0
3-Pentanone	7.647	7.4	ug/L	78220	1	6.757	528330	50.0
n-Propyl acetate	7.799	154.9	ug/L	1636300	1	6.757	528330	50.0
Propanoic acid,...	8.549	23.5	ug/L	309632	2	10.055	659022	50.0
Propanoic acid,...	9.244	29.4	ug/L	387426	2	10.055	659022	50.0
3-Hexanone	9.311	3.1	ug/L	40923	2	10.055	659022	50.0
Propanoic acid,...	9.482	48.3	ug/L	637090	2	10.055	659022	50.0
Acetic acid, bu...	9.580	33.3	ug/L	439171	2	10.055	659022	50.0
Isopropyl butyrate	9.915	86.2	ug/L	1135630	2	10.055	659022	50.0
Butanoic acid, ...	10.640	6.5	ug/L	85705	2	10.055	659022	50.0