

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV082624\
 Data File : VV037024.D
 Acq On : 26 Aug 2024 14:06
 Operator : SY/MD
 Sample : PB162923TB 10X
 Misc : 5.00mL/MSVOA_V/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 VLEB923

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_V\Method\SFAMVLM082624WMA.M

Title : VOC Analysis

Signal : TIC: VV037024.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.278	67	74	91	rBV	380870	408599	8.43%	1.054%
2	1.532	145	153	171	rVB	313034	379823	7.83%	0.980%
3	2.059	306	317	328	rBV	660097	958801	19.77%	2.474%
4	2.442	430	436	446	rVB	30606	47686	0.98%	0.123%
5	3.336	712	714	717	rBV2	1339	726	0.01%	0.002%
6	3.407	734	736	741	rVB5	1548	1216	0.03%	0.003%
7	3.468	751	755	757	rBV5	1178	719	0.01%	0.002%
8	3.490	759	762	765	rVV5	972	687	0.01%	0.002%
9	3.526	768	773	776	rBV5	828	894	0.02%	0.002%
10	3.674	815	819	820	rBV4	1422	672	0.01%	0.002%
11	3.709	825	830	834	rVB5	1156	936	0.02%	0.002%
12	3.747	838	842	844	rBV5	1448	1268	0.03%	0.003%
13	3.796	848	857	889	rBV	149712	429856	8.86%	1.109%
14	4.143	960	965	968	rBV5	2081	2145	0.04%	0.006%
15	4.243	979	996	1026	rBV2	415928	1081597	22.30%	2.791%
16	4.374	1034	1037	1039	rBV4	1589	968	0.02%	0.002%
17	4.474	1065	1068	1073	rVB6	1191	1107	0.02%	0.003%
18	4.503	1073	1077	1081	rBV6	965	835	0.02%	0.002%
19	4.555	1090	1093	1096	rVB4	1146	633	0.01%	0.002%
20	4.590	1100	1104	1105	rBV4	1207	750	0.02%	0.002%
21	4.715	1139	1143	1148	rBV5	1783	1766	0.04%	0.005%
22	4.809	1168	1172	1177	rBV6	1705	1943	0.04%	0.005%
23	4.950	1197	1216	1250	rBV2	1083590	2729118	56.28%	7.042%
24	5.201	1282	1294	1323	rBV2	220703	505568	10.43%	1.305%
25	5.529	1385	1396	1432	rBV	750321	1515421	31.25%	3.910%
26	5.931	1516	1521	1525	rBV5	2211	2772	0.06%	0.007%
27	5.982	1525	1537	1570	rVV	598196	1242395	25.62%	3.206%
28	6.268	1618	1626	1652	rBV	74927	225238	4.64%	0.581%
29	6.387	1653	1663	1706	rVV	2726253	4849186	100.00%	12.512%
30	6.908	1815	1825	1848	rBV	335259	619629	12.78%	1.599%
31	7.034	1862	1864	1869	rVB4	1383	1100	0.02%	0.003%
32	7.114	1878	1889	1916	rBV	612376	1078409	22.24%	2.783%
33	7.236	1916	1927	1955	rVV	1254369	2198438	45.34%	5.673%
34	7.545	2012	2023	2048	rBV2	1013905	1768793	36.48%	4.564%
35	7.857	2109	2120	2140	rBV	915454	1491189	30.75%	3.848%
36	7.969	2141	2155	2160	rVV3	136505	256443	5.29%	0.662%

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Integrator: RTE

Smoothing : OFF

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Stop Thrs : 0

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Max Peaks: 100

Peak Location: TOP

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Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM082624WMA.M
Title : VOC Analysis

37	8.017	2161	2170	2193	rVV3	735072	1520985	31.37%	3.925%
38	8.123	2194	2203	2228	rVV	1384619	2198468	45.34%	5.673%
39	8.233	2228	2237	2265	rVB	881510	1400926	28.89%	3.615%
40	8.612	2344	2355	2379	rBV	2788843	4269948	88.05%	11.018%
41	8.779	2395	2407	2427	rBV	1196412	1975404	40.74%	5.097%
42	9.053	2481	2492	2512	rBV	31582	65520	1.35%	0.169%
43	9.329	2568	2578	2599	rBV2	23165	50429	1.04%	0.130%
44	9.484	2616	2626	2647	rBV	200182	344606	7.11%	0.889%
45	9.660	2678	2681	2684	rBV4	1738	1407	0.03%	0.004%
46	9.702	2689	2694	2705	rVB8	4700	7764	0.16%	0.020%
47	9.857	2737	2742	2743	rBV4	1249	1157	0.02%	0.003%
48	10.056	2802	2804	2810	rVB7	1296	1217	0.03%	0.003%
49	10.149	2821	2833	2851	rBV	751936	1193979	24.62%	3.081%
50	10.281	2871	2874	2880	rVB7	2117	2115	0.04%	0.005%
51	10.471	2930	2933	2942	rBV8	2545	2967	0.06%	0.008%
52	10.535	2950	2953	2954	rBV3	1330	809	0.02%	0.002%
53	10.570	2961	2964	2968	rBV6	1186	1342	0.03%	0.003%
54	10.606	2971	2975	2980	rBV7	1166	1224	0.03%	0.003%
55	10.760	3020	3023	3025	rBV4	1721	1237	0.03%	0.003%
56	10.840	3045	3048	3050	rVV3	1819	1306	0.03%	0.003%
57	10.860	3050	3054	3061	rVV7	3665	4660	0.10%	0.012%
58	10.895	3063	3065	3069	rVB4	1533	989	0.02%	0.003%
59	11.030	3105	3107	3110	rVB4	1107	714	0.01%	0.002%
60	11.130	3125	3138	3143	rBV10	6290	12574	0.26%	0.032%
61	11.181	3143	3154	3183	rVV	1193390	1861956	38.40%	4.804%
62	11.377	3213	3215	3218	rVB4	1996	947	0.02%	0.002%
63	11.554	3257	3270	3297	rBV	1217177	1932739	39.86%	4.987%
64	11.873	3367	3369	3374	rVB5	1912	1077	0.02%	0.003%
65	11.966	3393	3398	3402	rBV6	1145	999	0.02%	0.003%
66	12.004	3405	3410	3411	rBV4	745	637	0.01%	0.002%
67	12.310	3503	3505	3509	rVV4	1230	1019	0.02%	0.003%
68	12.329	3509	3511	3514	rVV4	1155	700	0.01%	0.002%
69	12.358	3518	3520	3523	rVV4	1430	1063	0.02%	0.003%
70	12.448	3545	3548	3551	rBV2	891	640	0.01%	0.002%
71	12.464	3551	3553	3556	rBV3	1312	874	0.02%	0.002%
72	12.583	3582	3590	3592	rBV9	5709	6484	0.13%	0.017%
73	12.705	3624	3628	3630	rBV4	1344	732	0.02%	0.002%
74	12.725	3633	3634	3640	rVB4	2040	1615	0.03%	0.004%
75	12.773	3640	3649	3650	rBV6	1200	1338	0.03%	0.003%
76	12.828	3661	3666	3672	rBV8	1738	2223	0.05%	0.006%

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Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM082624WMA.M
 Title : VOC Analysis

77	12.873	3678	3680	3685	rVB4	1618	1089	0.02%	0.003%
78	12.898	3685	3688	3692	rVB6	1071	783	0.02%	0.002%
79	13.088	3745	3747	3749	rBV	1457	751	0.02%	0.002%
80	13.127	3756	3759	3763	rVB6	1153	897	0.02%	0.002%
81	13.207	3774	3784	3795	rBV8	9801	19445	0.40%	0.050%
82	13.310	3813	3816	3820	rVB4	2078	1380	0.03%	0.004%
83	13.358	3828	3831	3836	rVB5	951	961	0.02%	0.002%
84	13.445	3851	3858	3873	rBV2	14052	27165	0.56%	0.070%
85	13.606	3903	3908	3910	rBV4	1715	1377	0.03%	0.004%
86	13.647	3918	3921	3923	rBV4	1397	1018	0.02%	0.003%
87	13.676	3926	3930	3931	rBV3	5970	3942	0.08%	0.010%
88	13.760	3954	3956	3959	rVB4	1648	685	0.01%	0.002%
89	13.782	3959	3963	3964	rBV3	1357	951	0.02%	0.002%
90	13.863	3985	3988	3991	rBV5	1515	1102	0.02%	0.003%
91	14.024	4035	4038	4042	rBV5	1272	1326	0.03%	0.003%
92	14.098	4059	4061	4064	rBV3	1188	774	0.02%	0.002%
93	14.342	4134	4137	4139	rBV3	1410	884	0.02%	0.002%
94	14.397	4152	4154	4159	rBV4	1150	978	0.02%	0.003%
95	14.419	4159	4161	4166	rBV5	1798	1150	0.02%	0.003%
96	14.715	4249	4253	4256	rBV4	1225	907	0.02%	0.002%
97	14.856	4295	4297	4302	rVB6	1327	961	0.02%	0.002%
98	14.911	4311	4314	4316	rBV4	1328	815	0.02%	0.002%
99	15.290	4430	4432	4436	rVB4	1445	756	0.02%	0.002%
100	15.660	4544	4547	4550	rBV4	2199	1316	0.03%	0.003%

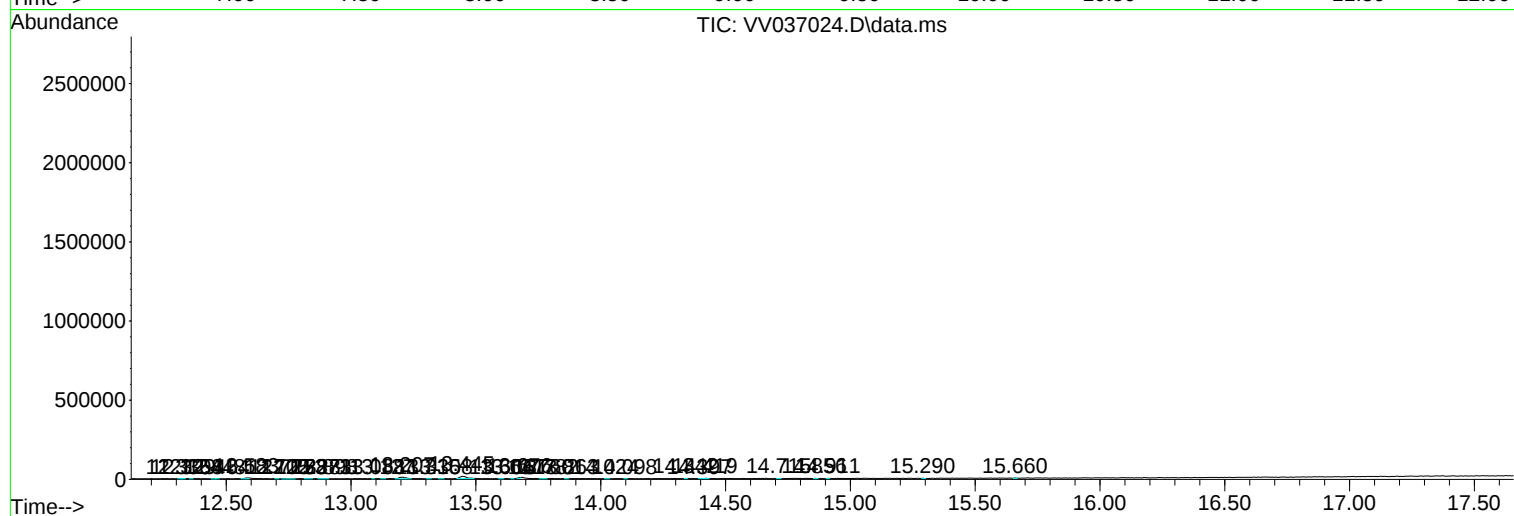
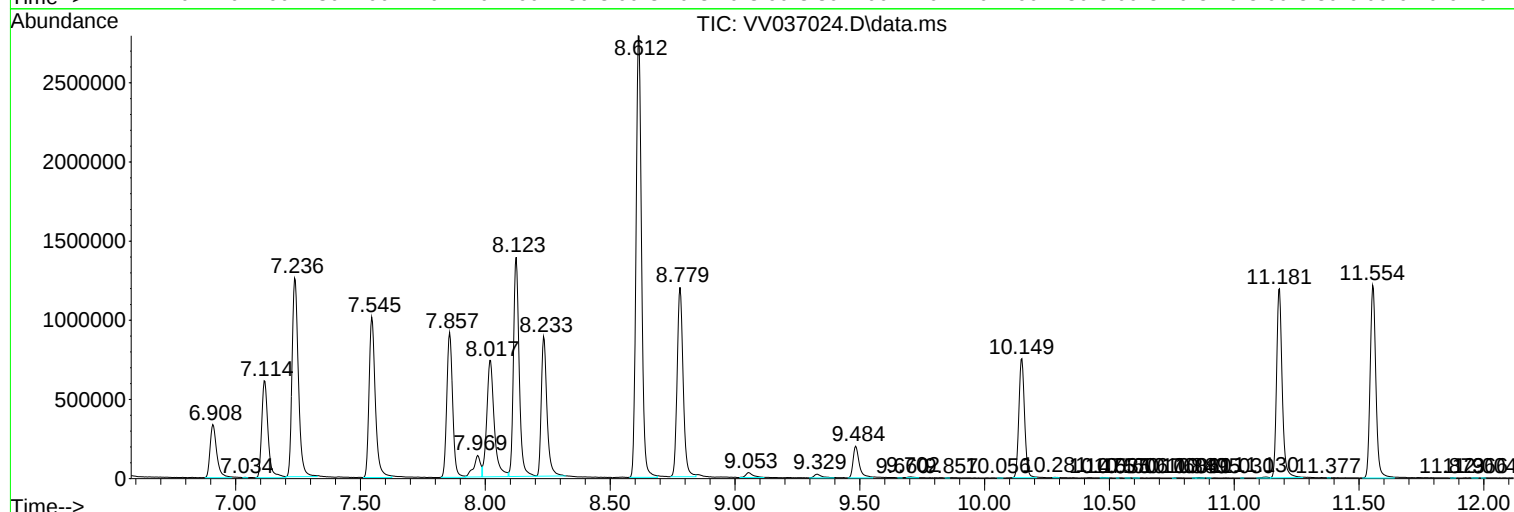
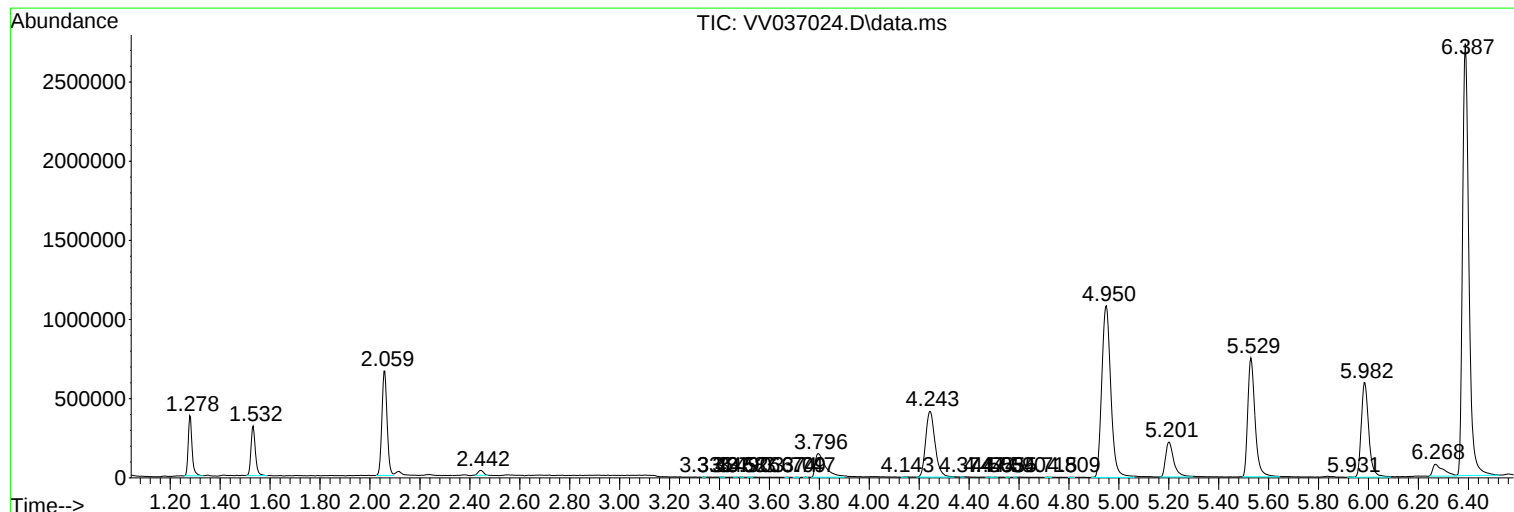
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ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
VLEB923

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM082624WMA.M
Quant Title : VOC Analysis

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV082624\
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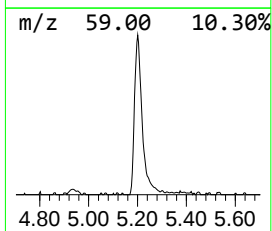
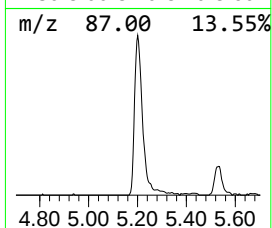
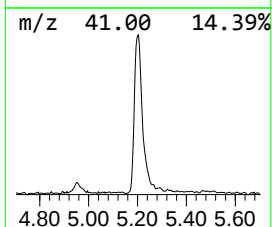
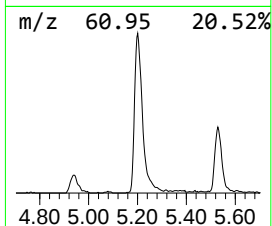
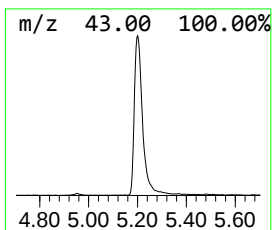
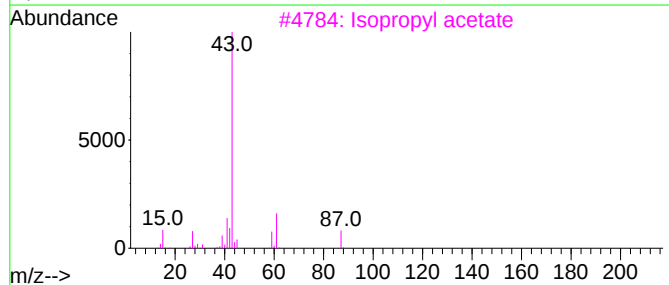
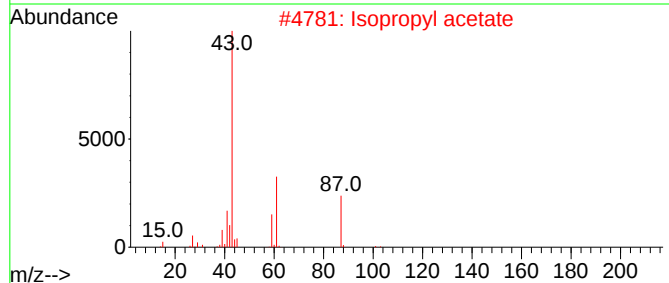
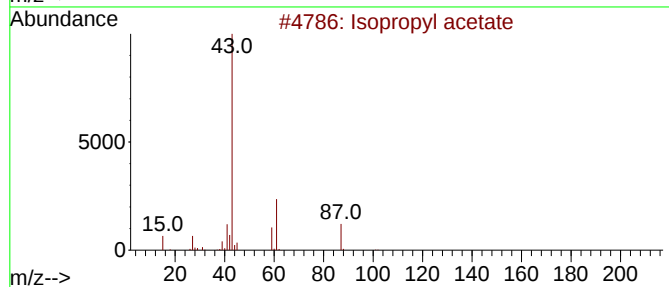
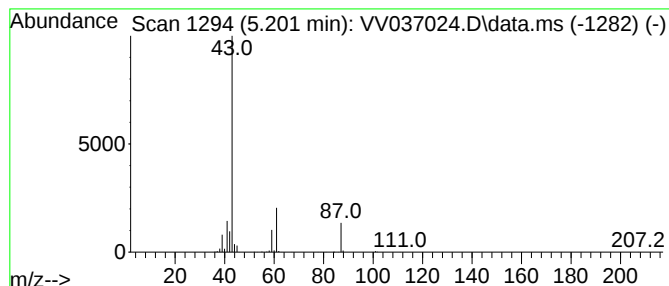
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM082624WMA.M
Quant Title : VOC Analysis

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

Peak Number 1 Isopropyl acetate Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.201	16.68 ug/L	505568	1,4-Difluorobenzene	5.529

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	78
3			Isopropyl acetate	102	C5H10O2	000108-21-4	74
4			Isopropyl acetate	102	C5H10O2	000108-21-4	72
5			Propane, 1-(1-methylethoxy)-	102	C6H14O	000627-08-7	40



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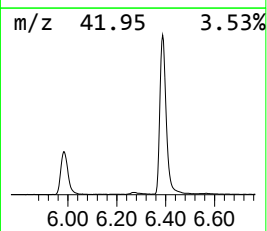
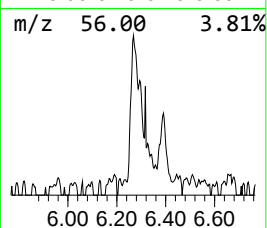
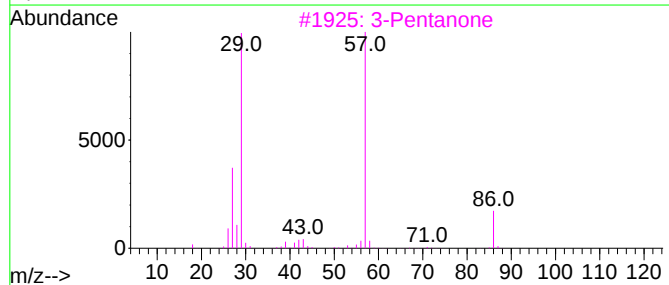
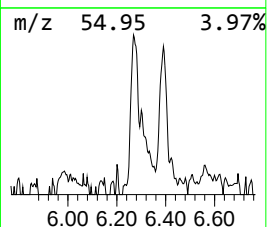
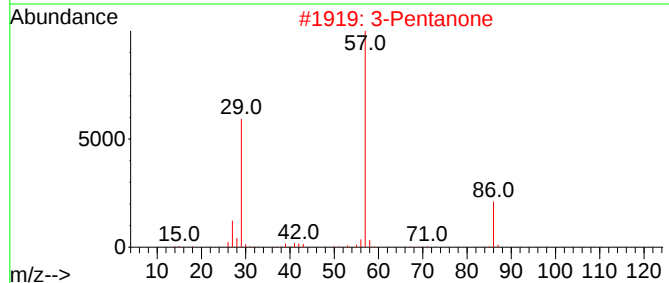
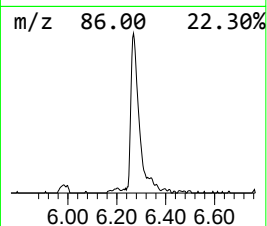
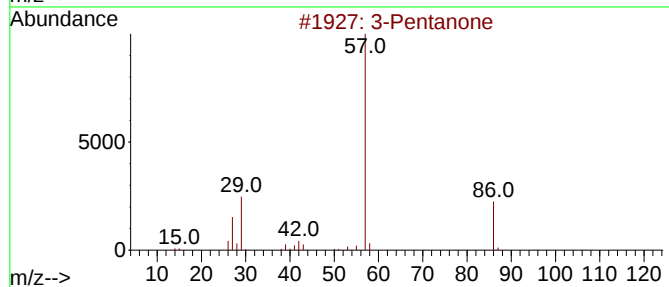
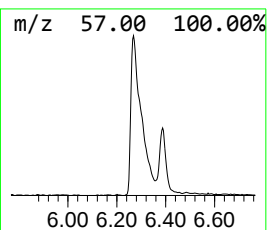
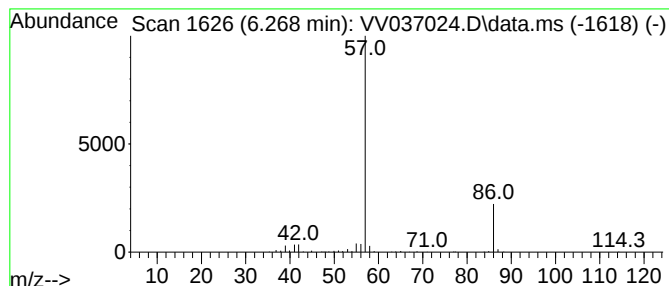
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM082624WMA.M
Quant Title : VOC Analysis

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

Peak Number 2 3-Pentanone Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.268	7.43 ug/L	225238	1,4-Difluorobenzene	5.529

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Pentanone	86	C5H10O	000096-22-0	59
2			3-Pentanone	86	C5H10O	000096-22-0	59
3			3-Pentanone	86	C5H10O	000096-22-0	59
4			3-Pentanone	86	C5H10O	000096-22-0	45
5			Propanoic acid, 2-propenyl ester	114	C6H10O2	002408-20-0	9



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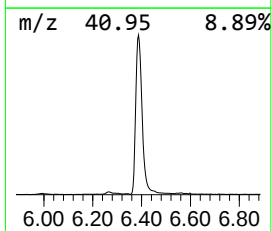
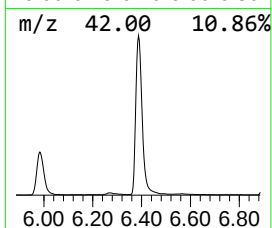
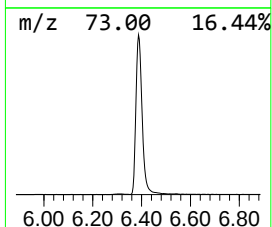
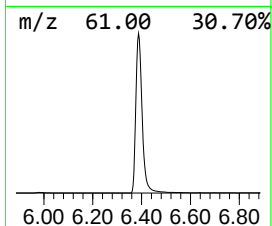
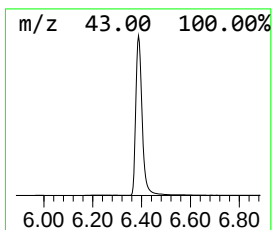
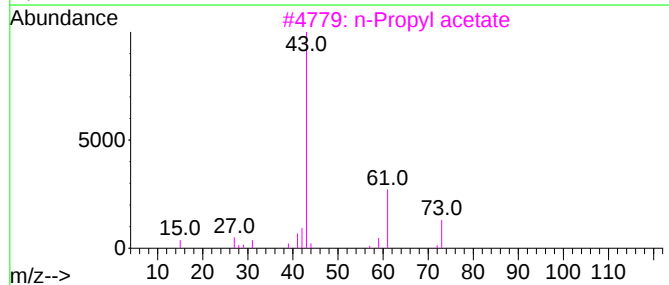
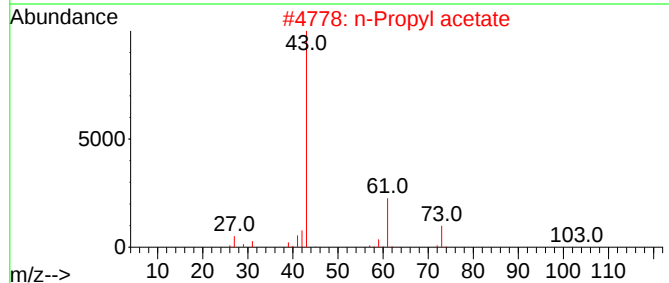
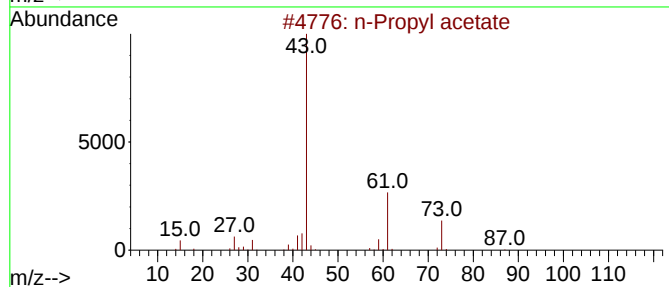
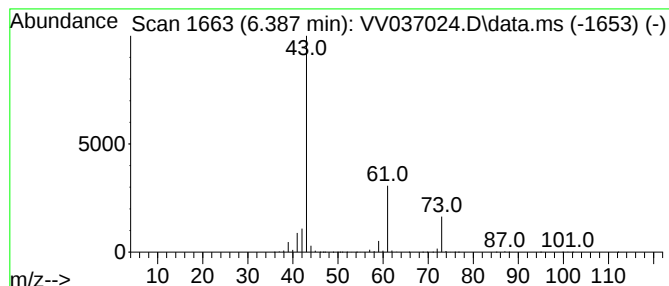
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Quant Title : VOC Analysis

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

Peak Number 3 n-Propyl acetate Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.387	160.00 ug/L	4849190	1,4-Difluorobenzene	5.529

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	83
3	n-Propyl acetate			102	C5H10O2	000109-60-4	78
4	n-Propyl acetate			102	C5H10O2	000109-60-4	64
5	Isopropyl acetate			102	C5H10O2	000108-21-4	36



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV082624\
Data File : VV037024.D
Acq On : 26 Aug 2024 14:06
Operator : SY/MD
Sample : PB162923TB 10X
Misc : 5.00mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
VLEB923

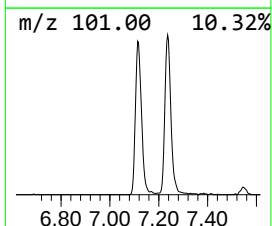
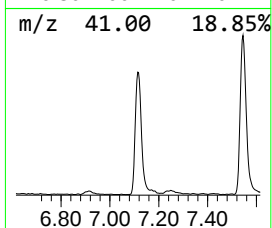
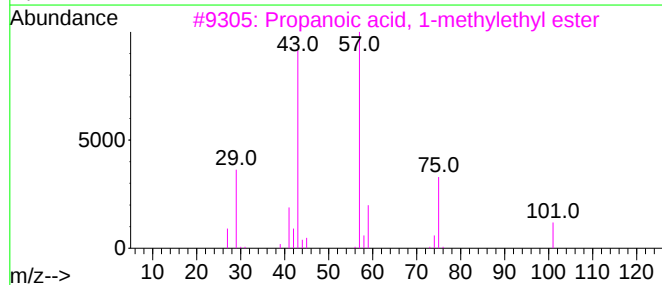
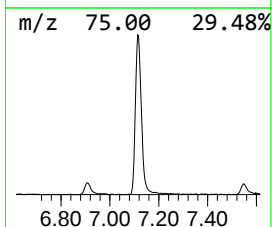
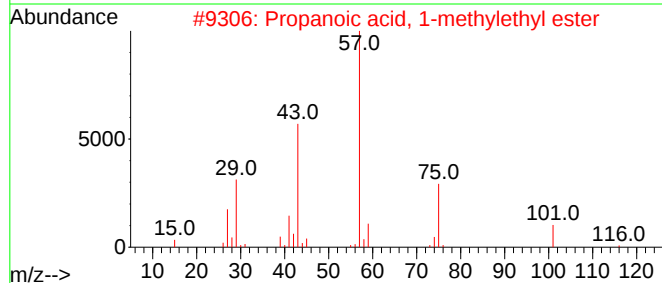
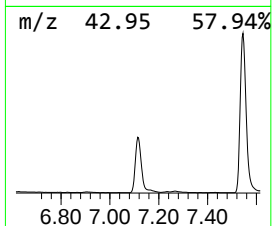
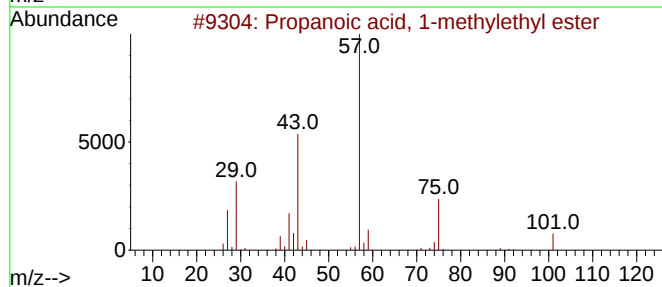
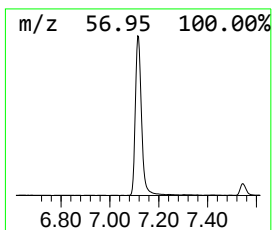
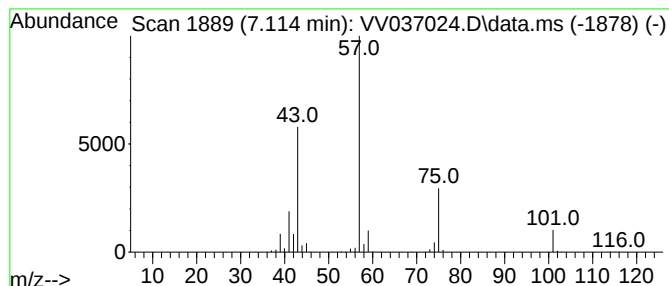
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Quant Title : VOC Analysis

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

Peak Number 5 Propanoic acid, 1-methyleth... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.114	35.58 ug/L	1078410	1,4-Difluorobenzene	5.529

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	53
4			Propanoic acid, 1-methylethyl ester	116	C6H12O2	000637-78-5	50
5			Propanoic acid, 1-methylethyl ester	116	C6H12O2	000637-78-5	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV082624\
Data File : VV037024.D
Acq On : 26 Aug 2024 14:06
Operator : SY/MD
Sample : PB162923TB 10X
Misc : 5.00mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
VLEB923

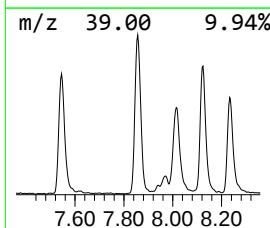
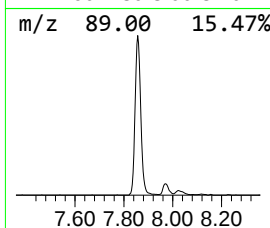
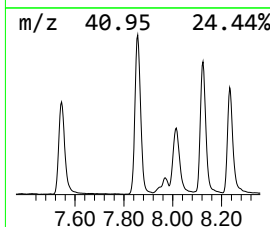
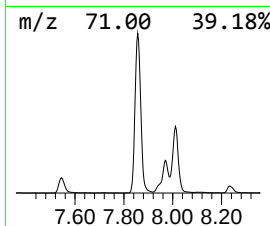
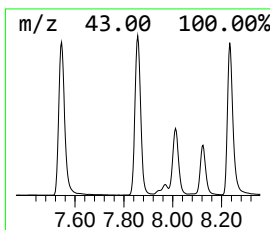
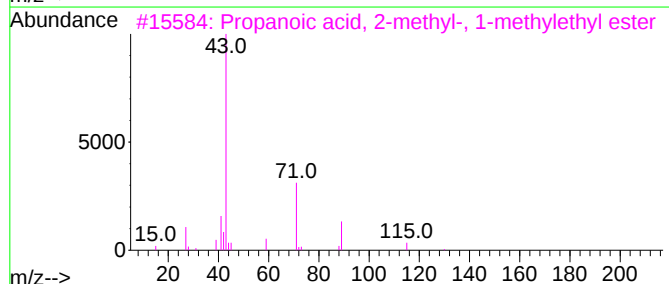
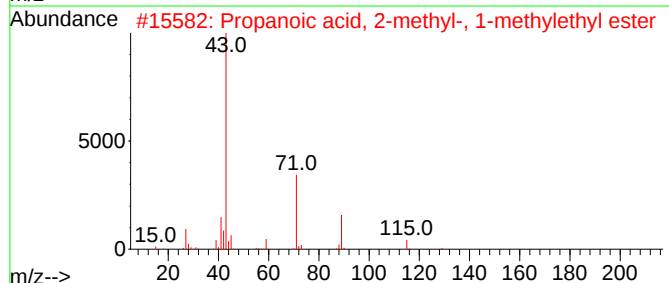
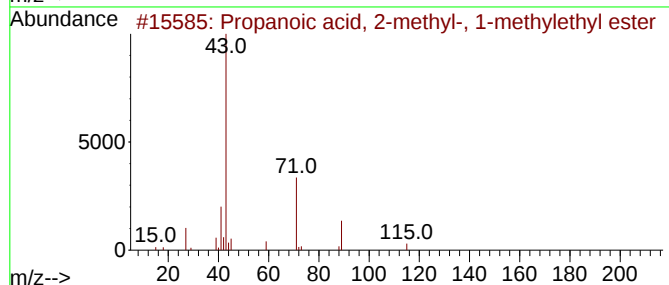
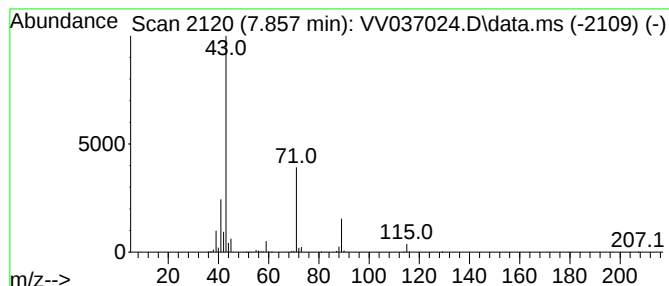
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Quant Title : VOC Analysis

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

Peak Number 6 Propanoic acid, 2-methyl-, ... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.857	37.74 ug/L	1491190	Chlorobenzene-d5	8.779

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	78
4			Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	78
5			Isopropyl butyrate	130	C7H14O2	000638-11-9	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV082624\
Data File : VV037024.D
Acq On : 26 Aug 2024 14:06
Operator : SY/MD
Sample : PB162923TB 10X
Misc : 5.00mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
VLEB923

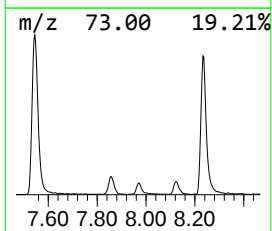
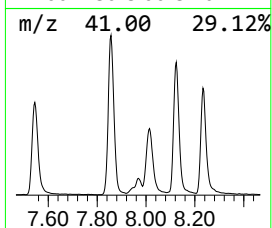
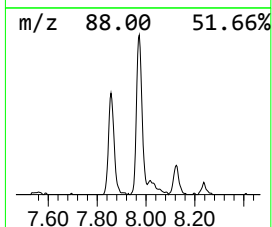
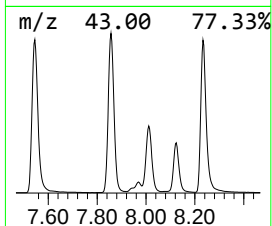
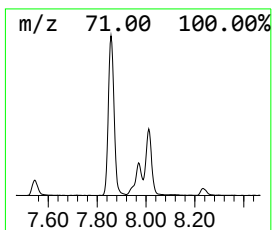
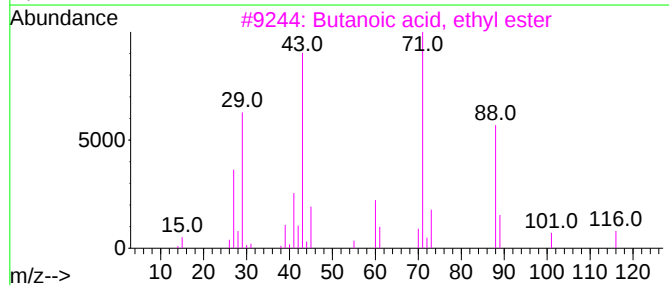
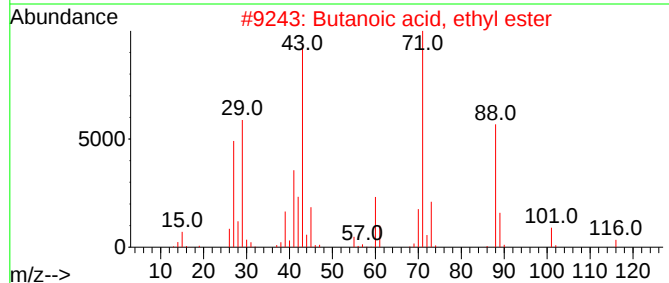
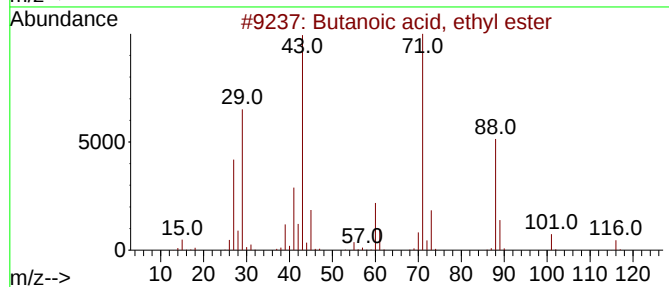
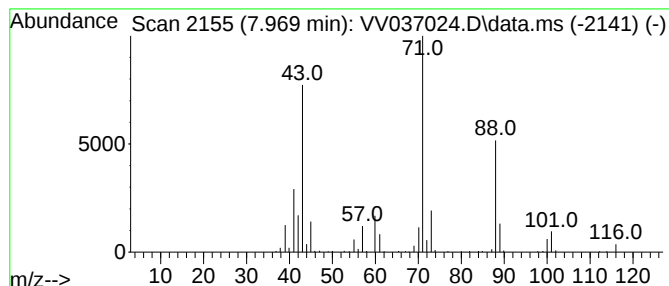
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Quant Title : VOC Analysis

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

Peak Number 7 Butanoic acid, ethyl ester Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.969	6.49 ug/L	256443	Chlorobenzene-d5	8.779

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butanoic acid, ethyl ester	116	C6H12O2	000105-54-4	97
2			Butanoic acid, ethyl ester	116	C6H12O2	000105-54-4	93
3			Butanoic acid, ethyl ester	116	C6H12O2	000105-54-4	90
4			Butanoic acid, ethyl ester	116	C6H12O2	000105-54-4	80
5			Butanoic acid, ethyl ester	116	C6H12O2	000105-54-4	46



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV082624\
Data File : VV037024.D
Acq On : 26 Aug 2024 14:06
Operator : SY/MD
Sample : PB162923TB 10X
Misc : 5.00mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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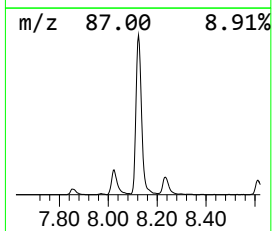
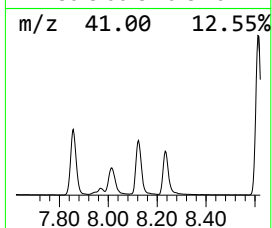
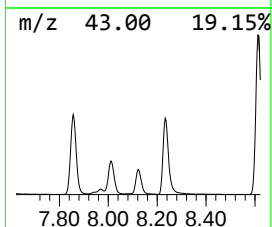
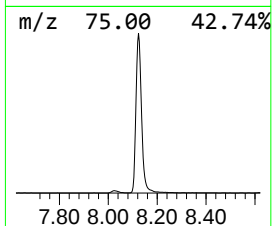
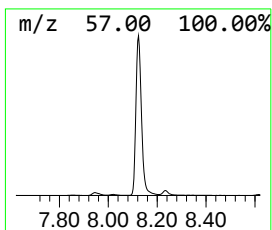
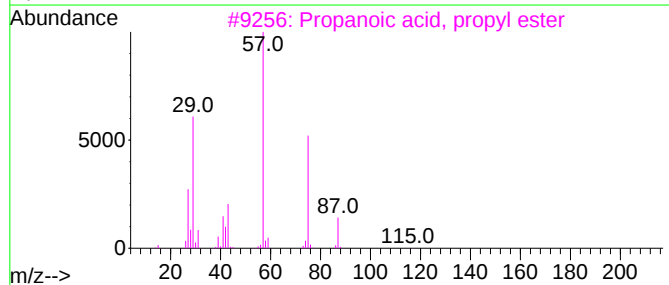
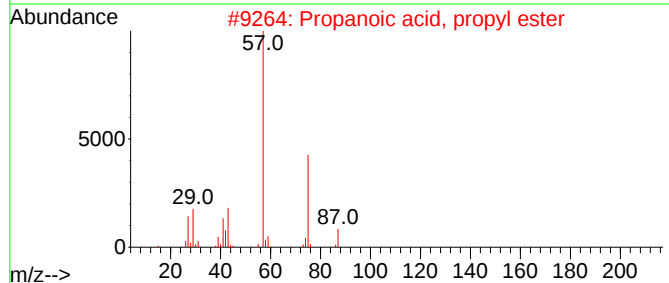
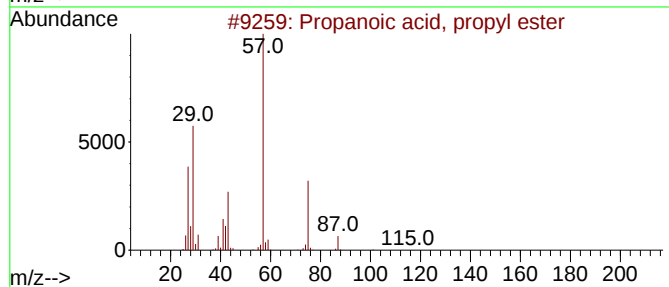
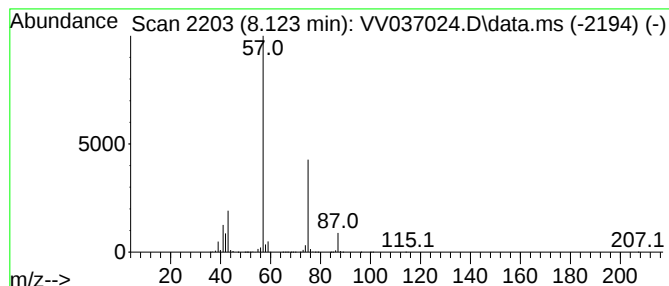
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Quant Title : VOC Analysis

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

Peak Number 8 Propanoic acid, propyl ester Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.123	55.65 ug/L	2198470	Chlorobenzene-d5	8.779

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV082624\
Data File : VV037024.D
Acq On : 26 Aug 2024 14:06
Operator : SY/MD
Sample : PB162923TB 10X
Misc : 5.00mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
VLEB923

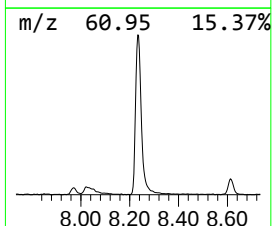
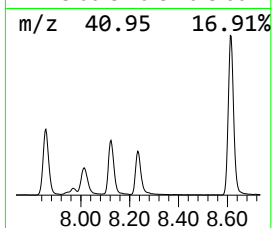
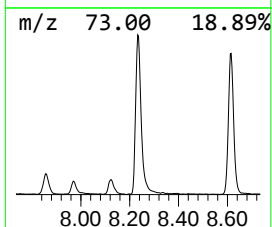
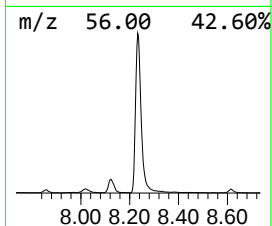
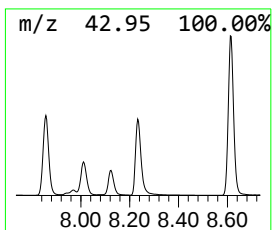
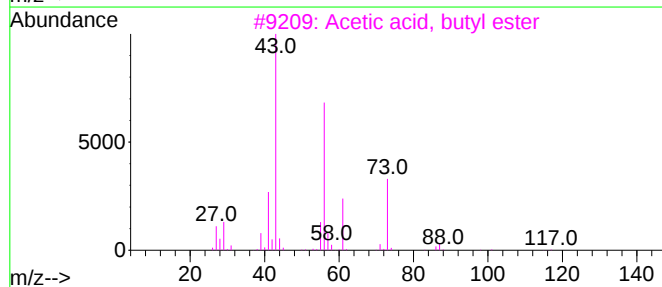
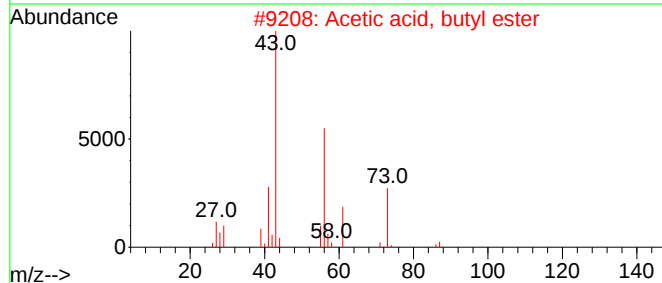
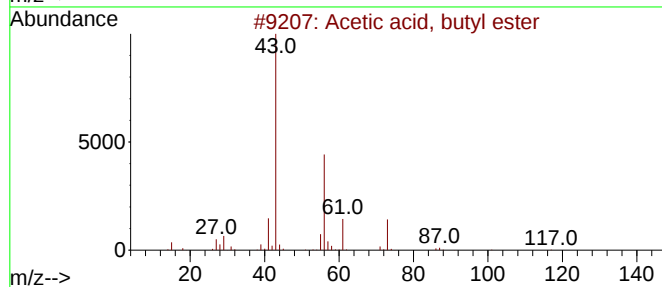
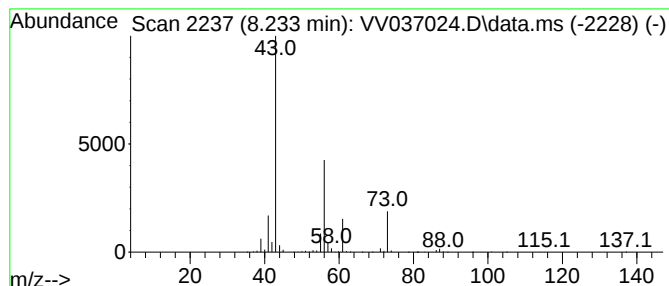
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Quant Title : VOC Analysis

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

Peak Number 9 Acetic acid, butyl ester Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.233	35.46 ug/L	1400930	Chlorobenzene-d5	8.779

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetic acid, butyl ester			116	C6H12O2	000123-86-4	74
2	Acetic acid, butyl ester			116	C6H12O2	000123-86-4	72
3	Acetic acid, butyl ester			116	C6H12O2	000123-86-4	72
4	Acetic acid, butyl ester			116	C6H12O2	000123-86-4	59
5	Isobutyl acetate			116	C6H12O2	000110-19-0	40



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV082624\
Data File : VV037024.D
Acq On : 26 Aug 2024 14:06
Operator : SY/MD
Sample : PB162923TB 10X
Misc : 5.00mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
VLEB923

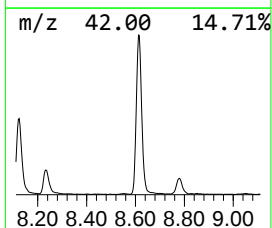
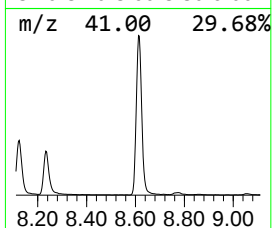
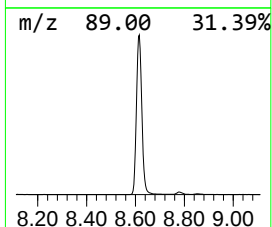
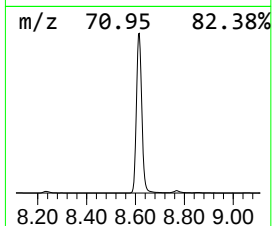
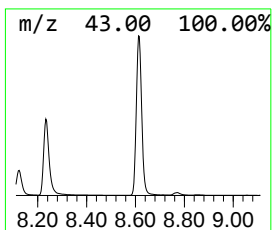
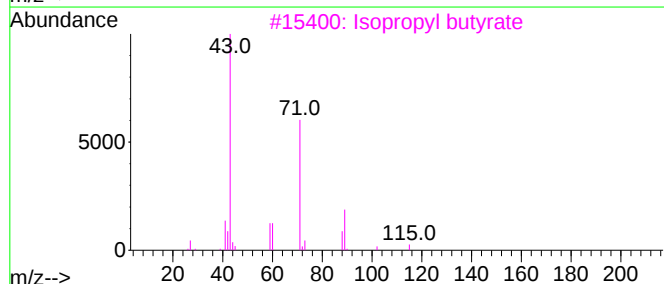
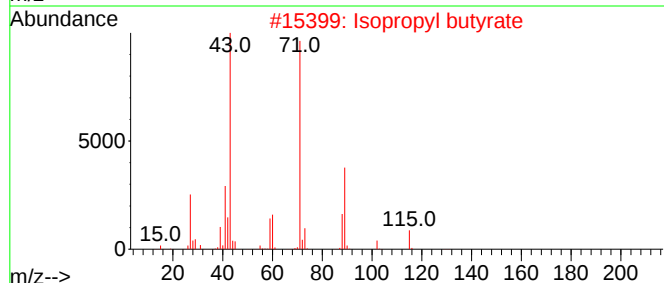
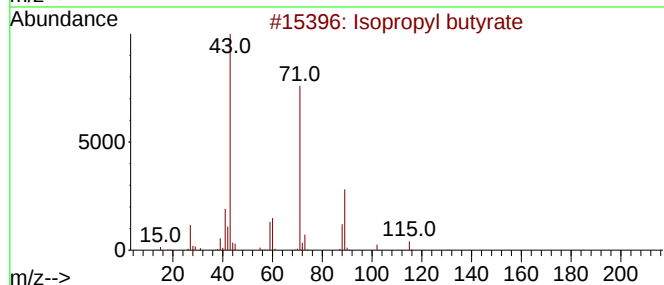
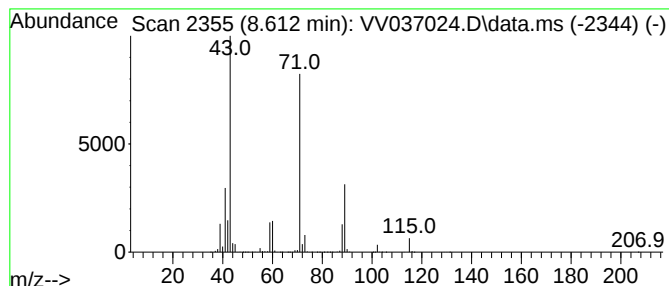
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Quant Title : VOC Analysis

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

Peak Number 10 Isopropyl butyrate Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.612	108.08 ug/L	4269950	Chlorobenzene-d5	8.779

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Isopropyl butyrate	130	C7H14O2	000638-11-9	94
2			Isopropyl butyrate	130	C7H14O2	000638-11-9	91
3			Isopropyl butyrate	130	C7H14O2	000638-11-9	83
4			Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	72
5			Butanoic acid, pentyl ester	158	C9H18O2	000540-18-1	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV082624\
Data File : VV037024.D
Acq On : 26 Aug 2024 14:06
Operator : SY/MD
Sample : PB162923TB 10X
Misc : 5.00mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
VLEB923

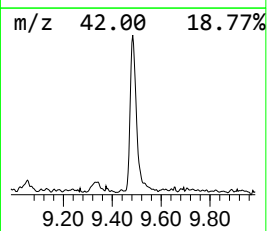
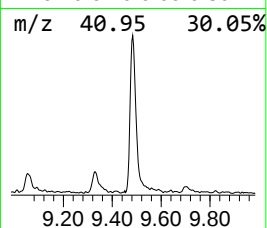
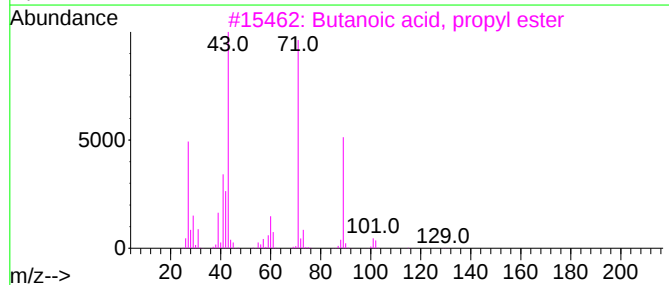
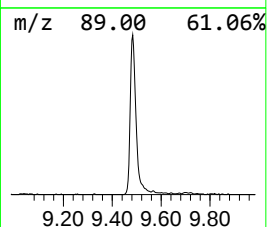
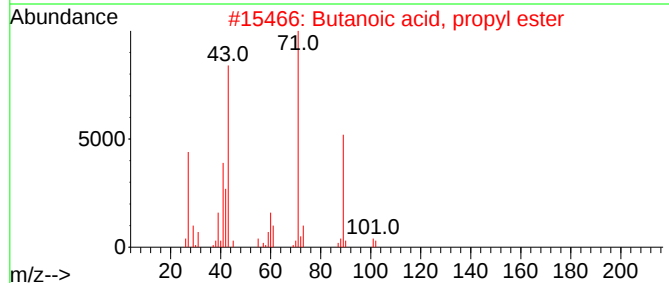
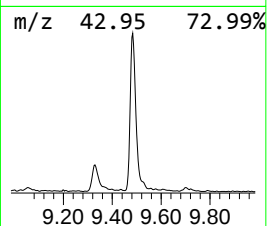
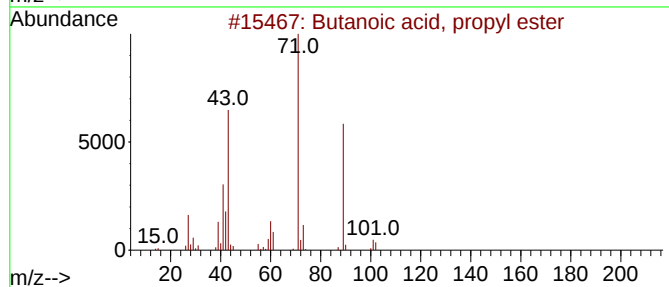
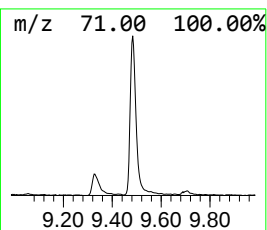
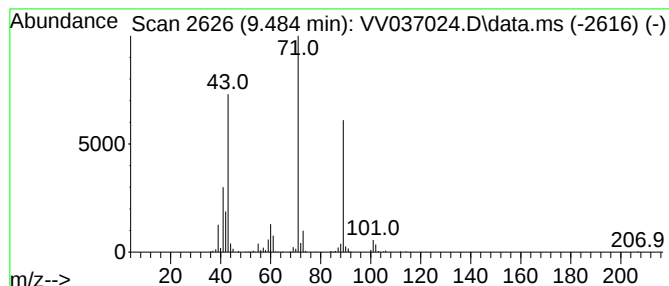
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Quant Title : VOC Analysis

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

Peak Number 11 Butanoic acid, propyl ester Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.484	8.72 ug/L	344606	Chlorobenzene-d5	8.779

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	80
3			Butanoic acid, propyl ester	130	C7H14O2	000105-66-8	80
4			Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	72
5			Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV082624\
Data File : VV037024.D
Acq On : 26 Aug 2024 14:06
Operator : SY/MD
Sample : PB162923TB 10X
Misc : 5.00mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
VLEB923

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM082624WMA.M
Quant Title : VOC Analysis

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Isopropyl acetate	5.201	16.7	ug/L	505568	1	5.529	1515420	50.0
3-Pentanone	6.268	7.4	ug/L	225238	1	5.529	1515420	50.0
n-Propyl acetate	6.387	160.0	ug/L	4849190	1	5.529	1515420	50.0
Propanoic acid,...	7.114	35.6	ug/L	1078410	1	5.529	1515420	50.0
Propanoic acid,...	7.857	37.7	ug/L	1491190	2	8.779	1975400	50.0
Butanoic acid, ...	7.969	6.5	ug/L	256443	2	8.779	1975400	50.0
Propanoic acid,...	8.123	55.6	ug/L	2198470	2	8.779	1975400	50.0
Acetic acid, bu...	8.233	35.5	ug/L	1400930	2	8.779	1975400	50.0
Isopropyl butyrate	8.612	108.1	ug/L	4269950	2	8.779	1975400	50.0
Butanoic acid, ...	9.484	8.7	ug/L	344606	2	8.779	1975400	50.0