

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV091224\
Data File : VV037264.D
Acq On : 12 Sep 2024 13:12
Operator : SY/MD
Sample : PB163300TB 10X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
VLEB300

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 0 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM082624WMA.M

Title : VOC Analysis

Signal : TIC: VV037264.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.085	10	14	21	rVB6	4587	4470	0.09%	0.010%
2	1.172	35	41	49	rBV	5721	9008	0.19%	0.021%
3	1.214	51	54	56	rBV4	2723	1933	0.04%	0.005%
4	1.278	66	74	89	rBV	471622	500250	10.34%	1.169%
5	1.349	92	96	103	rVB	6930	7830	0.16%	0.018%
6	1.413	110	116	123	rBV8	7076	9139	0.19%	0.021%
7	1.532	145	153	167	rVB	369095	439910	9.10%	1.028%
8	2.060	305	317	328	rBV	794167	1145307	23.68%	2.677%
9	2.111	329	333	345	rVB	31346	47587	0.98%	0.111%
10	2.446	428	437	448	rBV2	32584	54893	1.14%	0.128%
11	2.687	507	512	516	rBV7	2533	3023	0.06%	0.007%
12	2.838	556	559	564	rBV7	2545	2263	0.05%	0.005%
13	3.089	629	637	664	rVB3	26233	77806	1.61%	0.182%
14	3.317	703	708	711	rVV7	1821	1907	0.04%	0.004%
15	3.635	795	807	815	rBV7	9711	21917	0.45%	0.051%
16	3.793	846	856	890	rBV	145648	395157	8.17%	0.923%
17	4.243	979	996	1021	rBV	477381	1214705	25.12%	2.839%
18	4.950	1199	1216	1248	rBV2	1219468	3067485	63.43%	7.169%
19	5.198	1280	1293	1335	rBV	237126	536441	11.09%	1.254%
20	5.529	1384	1396	1426	rBV	781647	1604799	33.19%	3.750%
21	5.828	1478	1489	1494	rBV8	9140	16551	0.34%	0.039%
22	5.986	1521	1538	1570	rBV	659737	1376799	28.47%	3.218%
23	6.269	1616	1626	1633	rBV2	71245	132865	2.75%	0.311%
24	6.388	1652	1663	1709	rVB	2709228	4831418	99.91%	11.291%
25	6.571	1717	1720	1736	rVB8	7400	14717	0.30%	0.034%
26	6.908	1814	1825	1858	rBV	387424	733505	15.17%	1.714%
27	7.114	1878	1889	1916	rBV	681888	1165726	24.11%	2.724%
28	7.236	1916	1927	1954	rVV	1416728	2492912	51.55%	5.826%
29	7.545	2011	2023	2066	rBV2	1089107	1933612	39.99%	4.519%
30	7.857	2106	2120	2140	rBV	1015563	1682076	34.78%	3.931%
31	7.969	2140	2155	2160	rVV3	149194	281522	5.82%	0.658%
32	8.018	2160	2170	2192	rVV3	787261	1627638	33.66%	3.804%
33	8.121	2193	2202	2226	rVV	1471824	2338657	48.36%	5.466%
34	8.233	2228	2237	2265	rVB	917657	1455250	30.09%	3.401%
35	8.461	2305	2308	2315	rVB9	1988	1899	0.04%	0.004%
36	8.613	2343	2355	2385	rBV	3202969	4835728	100.00%	11.301%

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

37	8.780	2395	2407	2427	rBV	1222251	2023771	41.85%	4.730%
38	8.950	2458	2460	2465	rVB6	2406	1946	0.04%	0.005%
39	9.053	2481	2492	2510	rBV	33842	69144	1.43%	0.162%
40	9.137	2513	2518	2528	rVB	3752	6287	0.13%	0.015%
41	9.326	2570	2577	2593	rBV	23929	47478	0.98%	0.111%
42	9.484	2615	2626	2649	rBV	227111	391777	8.10%	0.916%
43	9.699	2685	2693	2697	rBV6	5536	7705	0.16%	0.018%
44	9.838	2734	2736	2741	rBV5	2672	2173	0.04%	0.005%
45	10.053	2797	2803	2809	rVB7	3202	4571	0.09%	0.011%
46	10.146	2819	2832	2850	rBV	742798	1188885	24.59%	2.778%
47	10.387	2902	2907	2913	rBV10	1917	2642	0.05%	0.006%
48	10.452	2918	2927	2929	rBV8	7188	9126	0.19%	0.021%
49	10.513	2940	2946	2962	rVB4	19627	33218	0.69%	0.078%
50	10.783	3028	3030	3033	rVV4	2483	1910	0.04%	0.004%
51	10.815	3034	3040	3047	rVB8	5265	7449	0.15%	0.017%
52	10.857	3047	3053	3058	rBV6	4573	5883	0.12%	0.014%
53	10.976	3081	3090	3094	rBV9	4309	7510	0.16%	0.018%
54	11.127	3130	3137	3141	rBV9	2565	3182	0.07%	0.007%
55	11.178	3141	3153	3179	rBV	1259913	1982667	41.00%	4.634%
56	11.310	3191	3194	3200	rVB5	8482	7253	0.15%	0.017%
57	11.358	3206	3209	3212	rVB4	2519	1866	0.04%	0.004%
58	11.394	3212	3220	3229	rBV4	6797	13127	0.27%	0.031%
59	11.503	3243	3254	3257	rBV9	8497	15042	0.31%	0.035%
60	11.555	3259	3270	3299	rVB	1350279	2174991	44.98%	5.083%
61	11.725	3316	3323	3336	rVB3	34252	49602	1.03%	0.116%
62	11.854	3359	3363	3366	rBV6	1779	1812	0.04%	0.004%
63	11.915	3379	3382	3387	rBV6	2479	2795	0.06%	0.007%
64	12.056	3422	3426	3435	rVB10	4582	6951	0.14%	0.016%
65	12.188	3462	3467	3470	rBV7	3445	3931	0.08%	0.009%
66	12.300	3498	3502	3506	rBV7	3687	3696	0.08%	0.009%
67	12.342	3512	3515	3516	rBV3	3310	1828	0.04%	0.004%
68	12.397	3530	3532	3540	rVV7	3215	3049	0.06%	0.007%
69	12.442	3540	3546	3554	rVB10	5681	8589	0.18%	0.020%
70	12.526	3567	3572	3581	rVB10	5745	7260	0.15%	0.017%
71	12.657	3610	3613	3622	rVB10	2854	2376	0.05%	0.006%
72	12.731	3632	3636	3644	rVB10	3585	4626	0.10%	0.011%
73	12.825	3656	3665	3675	rVV2	29976	44542	0.92%	0.104%
74	12.979	3705	3713	3720	rBV4	7845	12358	0.26%	0.029%
75	13.069	3737	3741	3747	rVB8	2939	2731	0.06%	0.006%
76	13.207	3776	3784	3787	rBV9	5173	7792	0.16%	0.018%

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 Title : VOC Analysis

77	13.326	3813	3821	3828	rBV9	11076	19415	0.40%	0.045%
78	13.371	3829	3835	3844	rVB7	9264	15428	0.32%	0.036%
79	13.445	3854	3858	3862	rBV6	3248	3177	0.07%	0.007%
80	13.577	3889	3899	3910	rVB9	10368	22055	0.46%	0.052%
81	13.757	3946	3955	3968	rBV9	12016	23923	0.49%	0.056%
82	13.837	3972	3980	3988	rBV3	30215	43683	0.90%	0.102%
83	13.908	3994	4002	4012	rBV3	5698	12605	0.26%	0.029%
84	14.014	4029	4035	4036	rBV5	8062	7680	0.16%	0.018%
85	14.294	4120	4122	4125	rBV3	2868	1994	0.04%	0.005%
86	14.352	4133	4140	4152	rVB2	15982	24617	0.51%	0.058%
87	14.448	4164	4170	4178	rVB2	7438	10807	0.22%	0.025%
88	14.529	4192	4195	4201	rVV8	5869	5307	0.11%	0.012%
89	14.583	4204	4212	4223	rVB10	16984	38782	0.80%	0.091%
90	14.779	4266	4273	4283	rVB3	32397	51478	1.06%	0.120%
91	14.841	4289	4292	4300	rVB10	6600	6072	0.13%	0.014%
92	15.046	4354	4356	4359	rBV4	3271	1786	0.04%	0.004%
93	15.355	4445	4452	4459	rBV7	14020	18855	0.39%	0.044%
94	15.606	4523	4530	4542	rVV3	21817	39071	0.81%	0.091%
95	15.670	4544	4550	4559	rVB2	29069	39256	0.81%	0.092%
96	15.744	4567	4573	4579	rBV5	20114	31304	0.65%	0.073%
97	16.062	4665	4672	4683	rBV5	11816	23664	0.49%	0.055%
98	16.451	4788	4793	4800	rBV	15398	23003	0.48%	0.054%
99	16.503	4804	4809	4817	rVB3	29934	39942	0.83%	0.093%
100	17.300	5053	5057	5069	rVB8	29420	44926	0.93%	0.105%

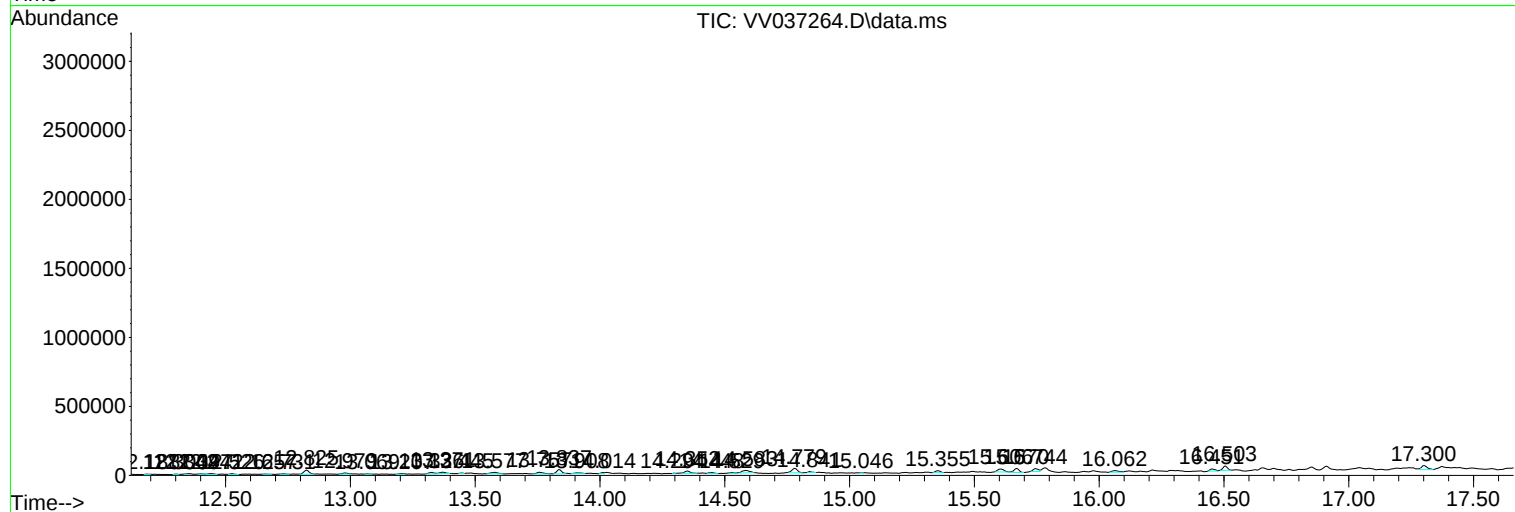
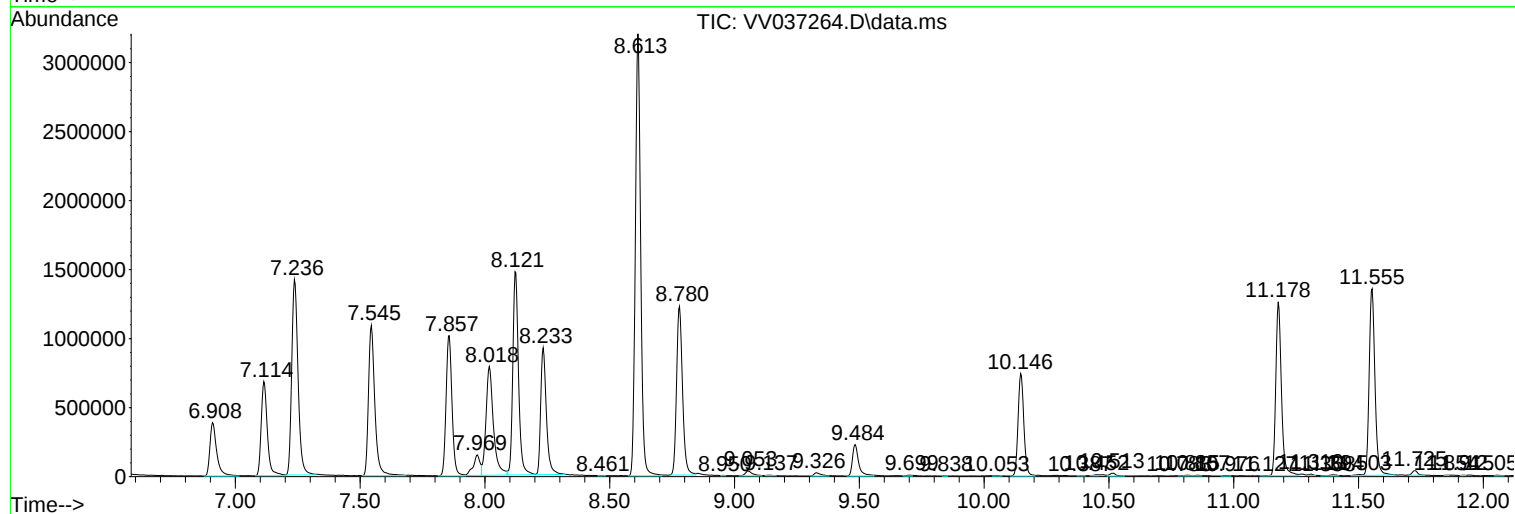
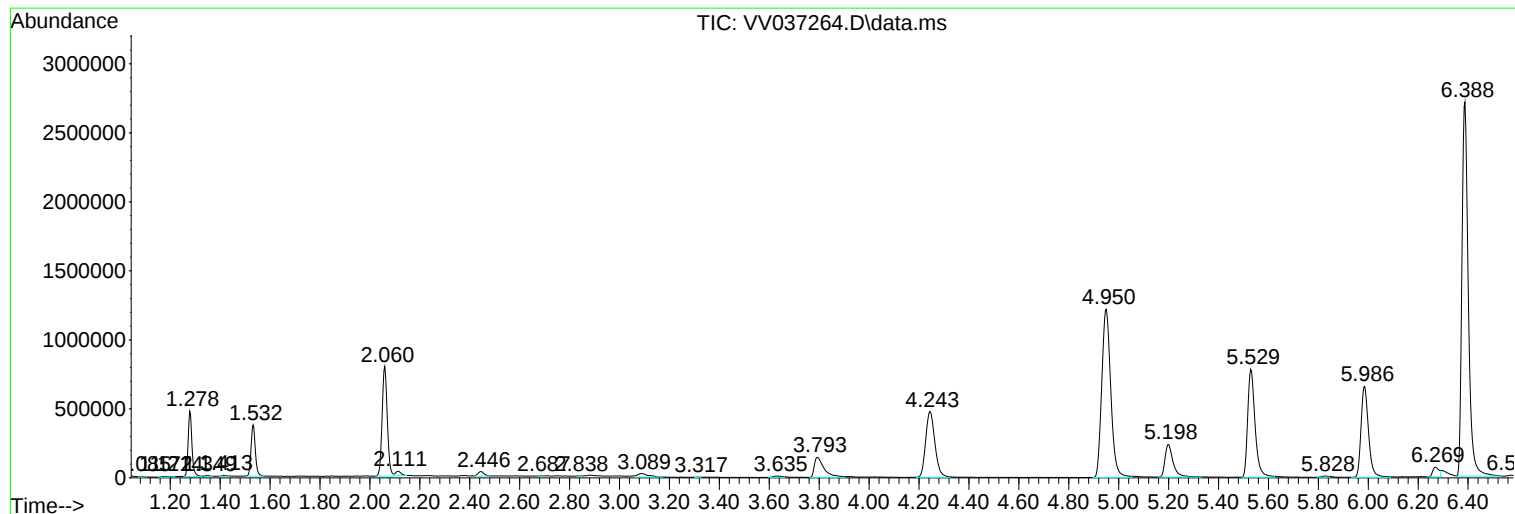
Sum of corrected areas: 42789076

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Instrument :
MSVOA_V
ClientSampleId :
VLEB300

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



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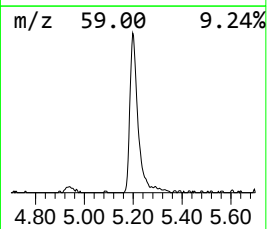
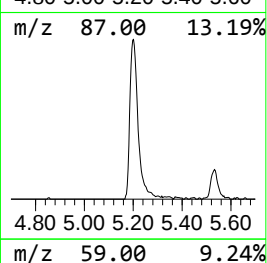
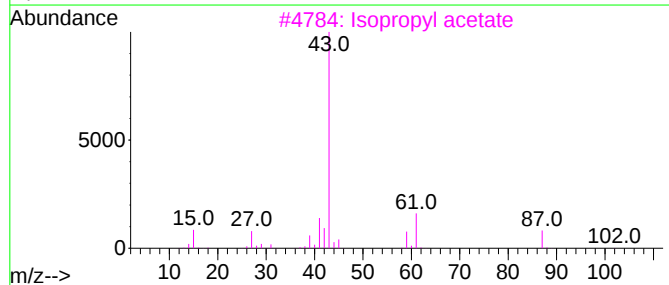
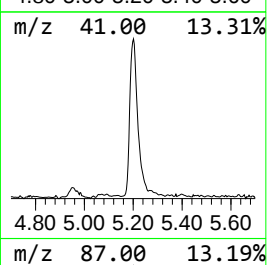
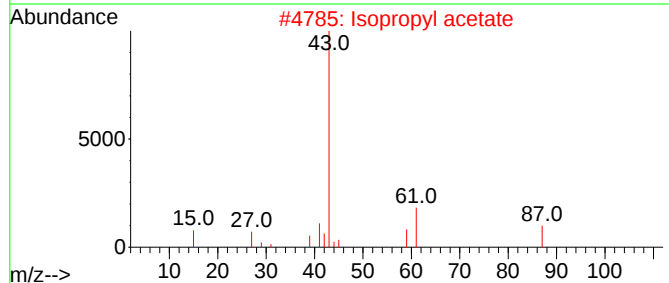
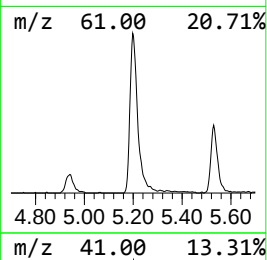
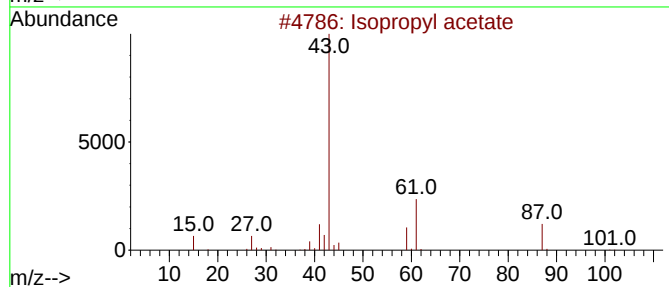
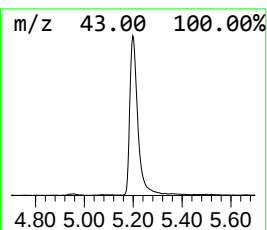
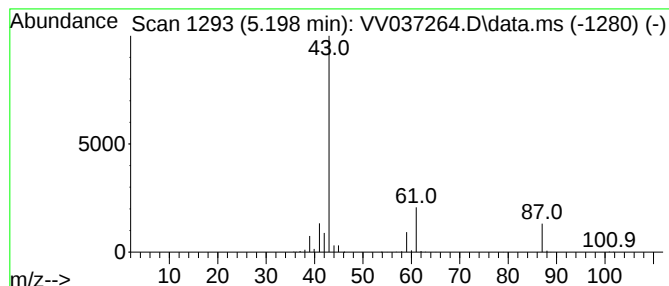
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM082624WMA.M
Quant Title : VOC Analysis

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.
5.198	16.71 ug/L	536441	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	64
4			Isopropyl acetate	102	C5H10O2	000108-21-4	64
5			Propane, 1-(1-methylethoxy)-	102	C6H14O	000627-08-7	23



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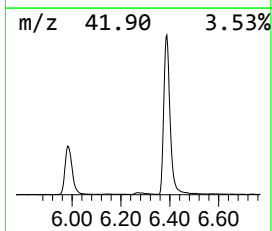
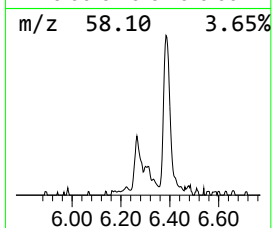
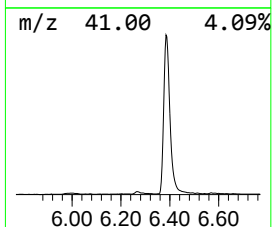
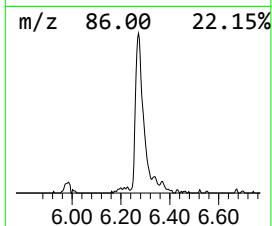
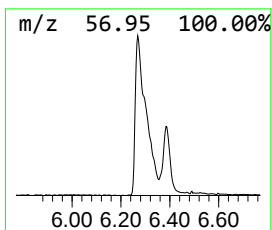
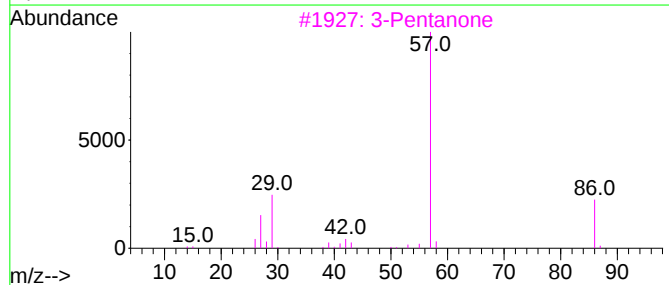
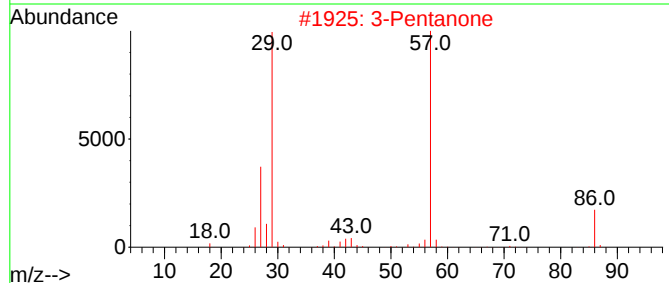
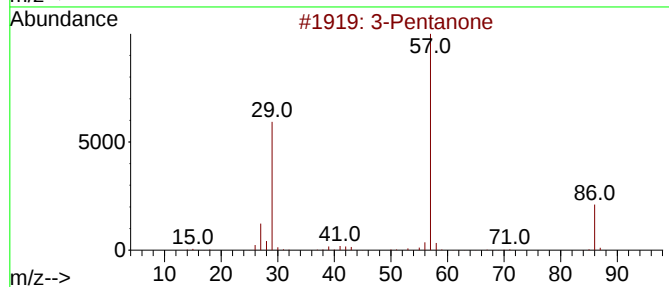
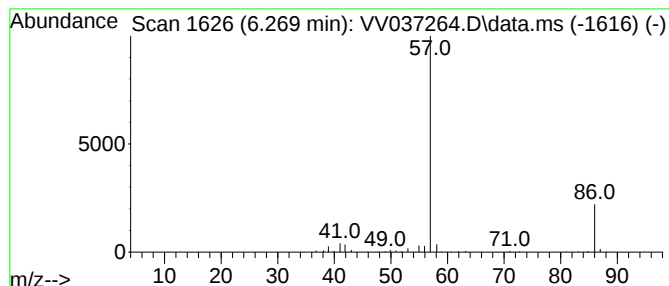
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVLM082624WMA.M
Quant Title : VOC Analysis

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

Peak Number 2 3-Pentanone Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.269	4.14 ug/L	132865	1,4-Difluorobenzene	5.529

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Pentanone			86	C5H10O	000096-22-0	72
2	3-Pentanone			86	C5H10O	000096-22-0	64
3	3-Pentanone			86	C5H10O	000096-22-0	64
4	3-Pentanone			86	C5H10O	000096-22-0	45
5	Propane, 2-bromo-2-methyl-			136	C4H9Br	000507-19-7	23



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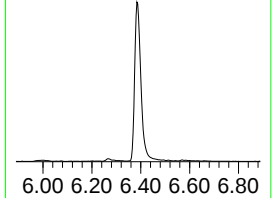
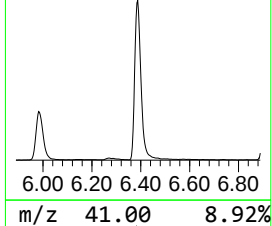
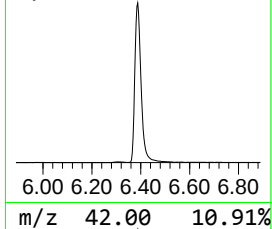
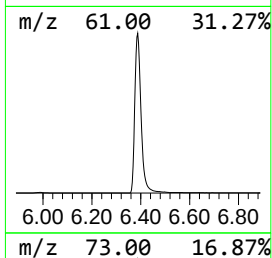
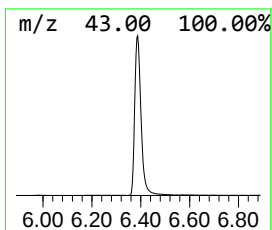
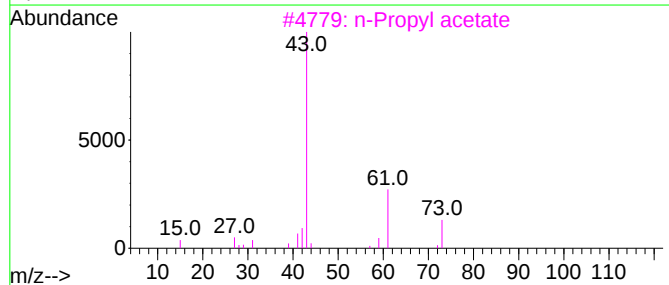
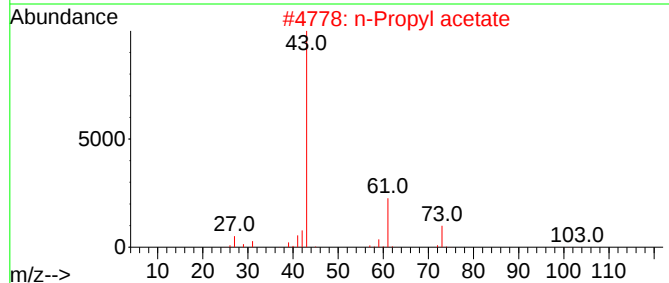
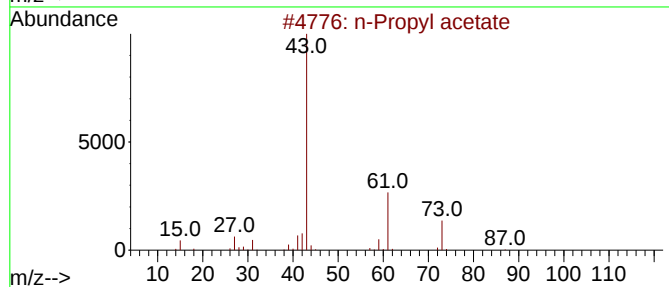
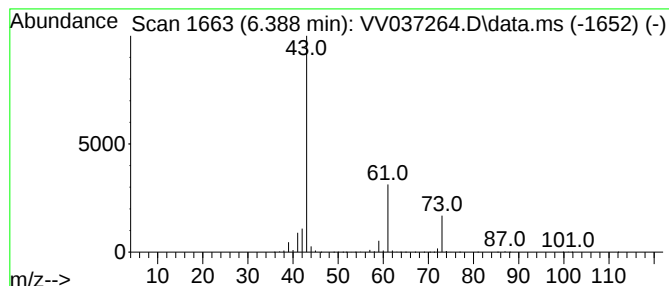
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Quant Title : VOC Analysis

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.
6.388	150.53 ug/L	4831420	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\VV091224\
Data File : VV037264.D
Acq On : 12 Sep 2024 13:12
Operator : SY/MD
Sample : PB163300TB 10X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
VLEB300

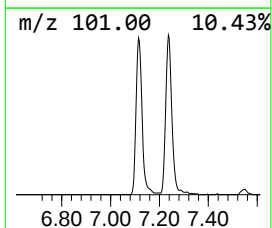
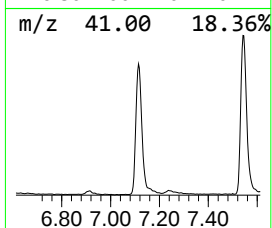
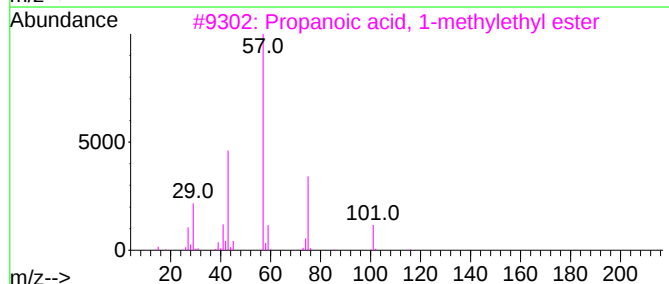
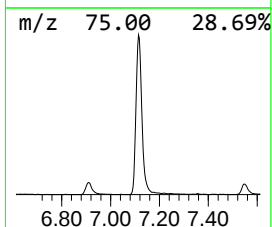
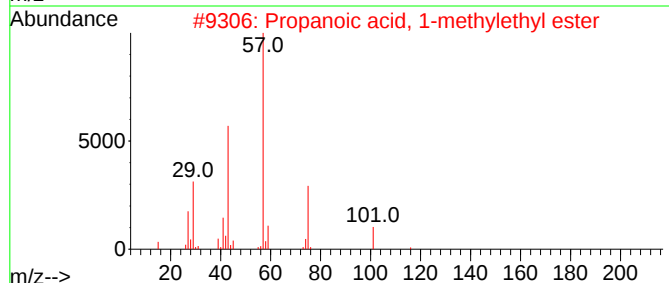
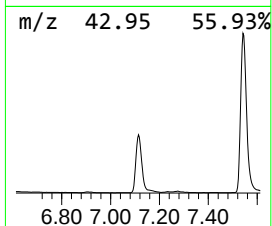
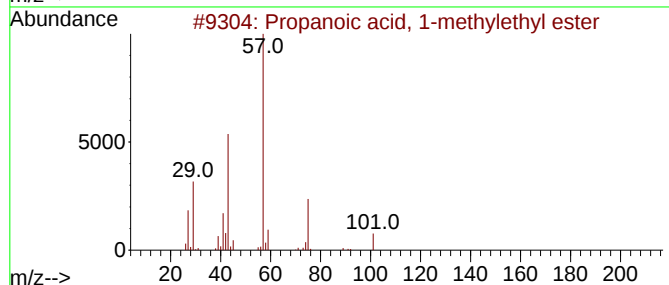
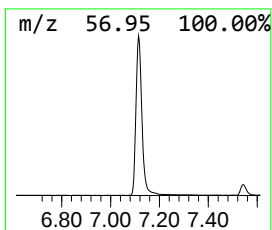
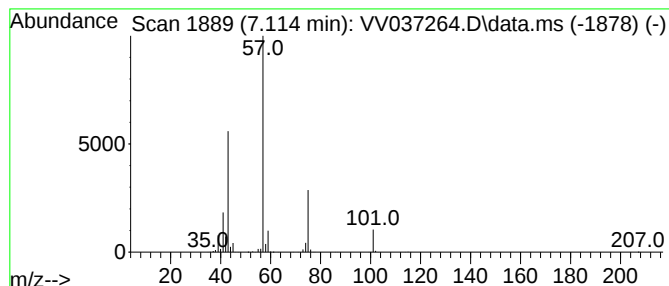
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Quant Title : VOC Analysis

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	36.32 ug/L	1165730	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	78
4			Propanoic acid, 1-methylethyl ester	116	C6H12O2	000637-78-5	74
5			Propanoic acid, propyl ester	116	C6H12O2	000106-36-5	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV091224\
Data File : VV037264.D
Acq On : 12 Sep 2024 13:12
Operator : SY/MD
Sample : PB163300TB 10X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
VLEB300

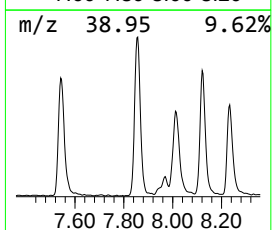
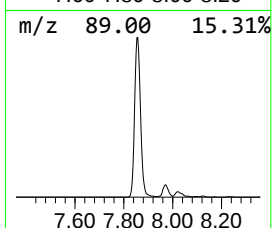
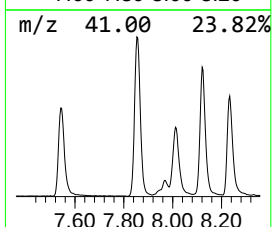
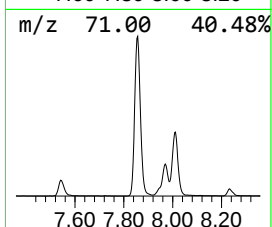
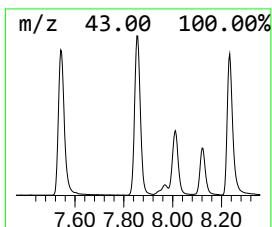
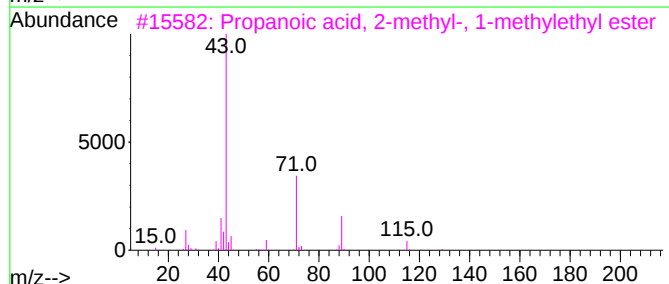
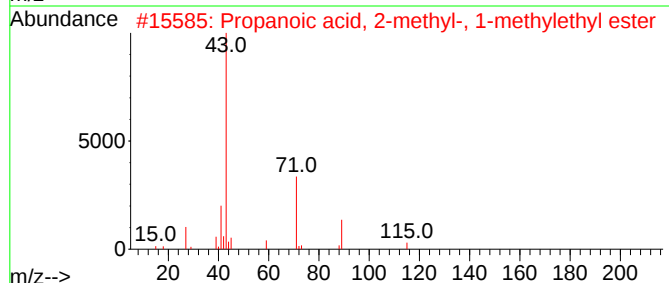
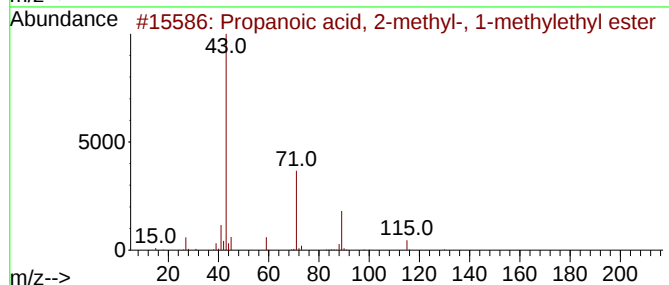
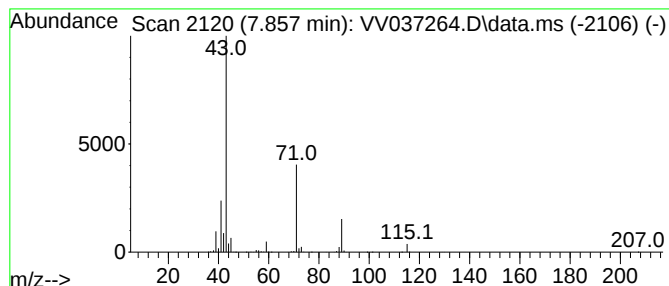
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Quant Title : VOC Analysis

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	41.56 ug/L	1682080	Chlorobenzene-d5	8.780

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_V\Data\VV091224\
Data File : VV037264.D
Acq On : 12 Sep 2024 13:12
Operator : SY/MD
Sample : PB163300TB 10X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
VLEB300

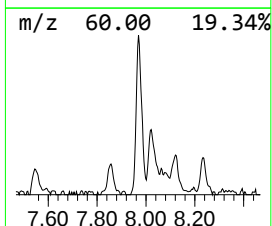
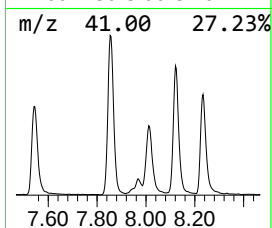
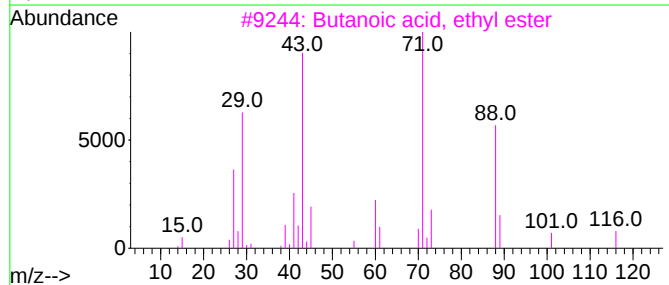
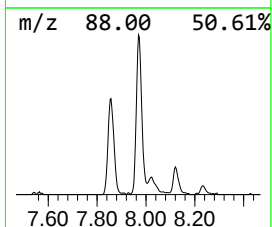
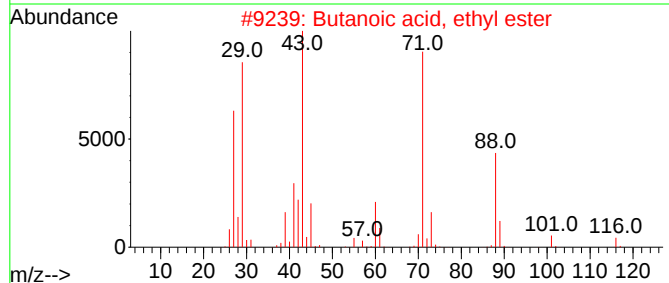
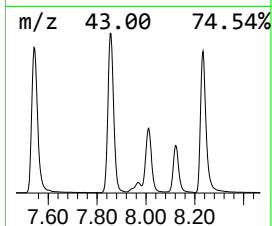
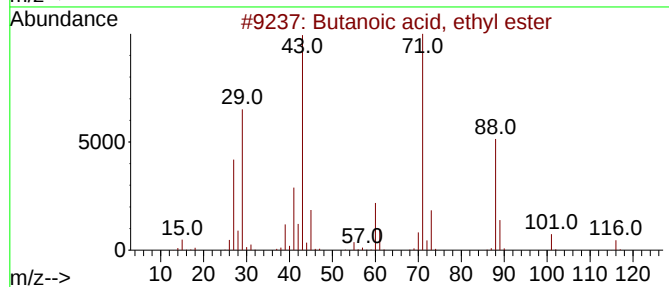
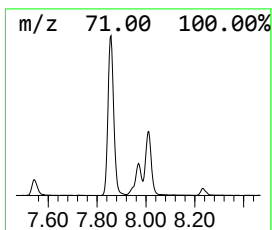
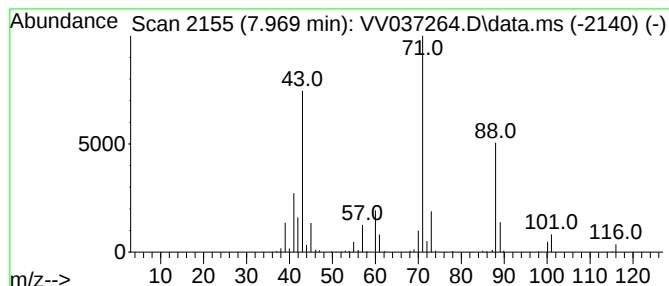
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Quant Title : VOC Analysis

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

Peak Number 7 Butanoic acid, ethyl ester Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.969	6.96 ug/L	281522	Chlorobenzene-d5	8.780

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butanoic acid, ethyl ester	116	C6H12O2	000105-54-4	96
2			Butanoic acid, ethyl ester	116	C6H12O2	000105-54-4	91
3			Butanoic acid, ethyl ester	116	C6H12O2	000105-54-4	90
4			Butanoic acid, ethyl ester	116	C6H12O2	000105-54-4	64
5			Propanoic acid, 2-methyl-, ethyl...	116	C6H12O2	000097-62-1	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV091224\
Data File : VV037264.D
Acq On : 12 Sep 2024 13:12
Operator : SY/MD
Sample : PB163300TB 10X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
VLEB300

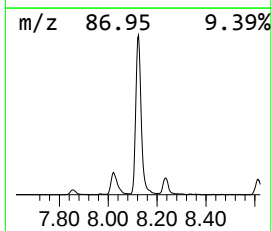
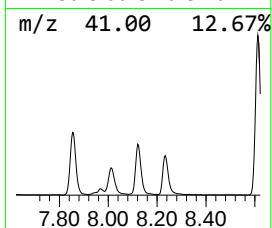
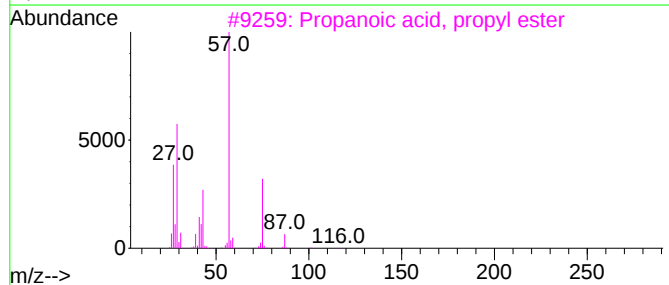
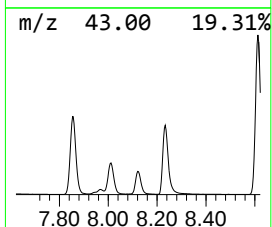
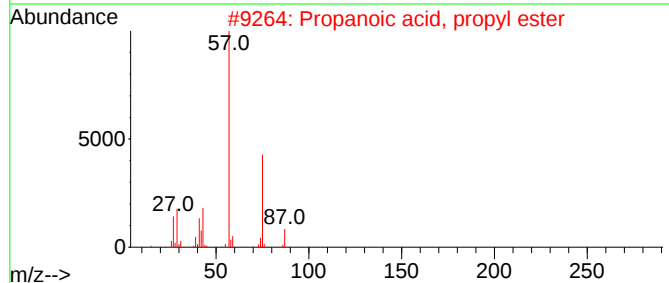
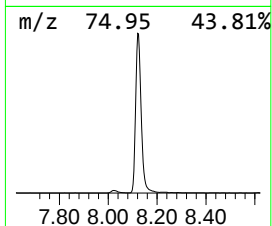
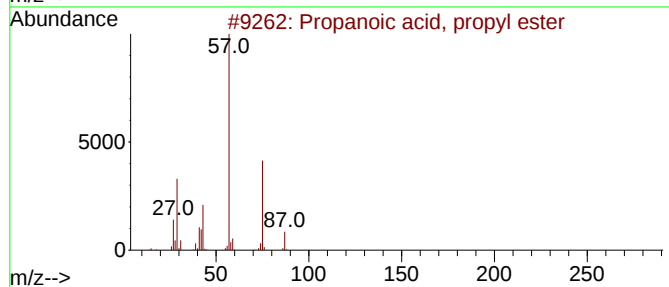
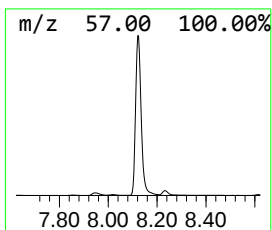
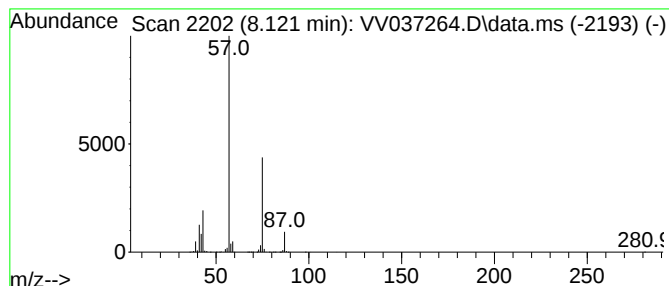
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Quant Title : VOC Analysis

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.
8.121	57.78 ug/L	2338660	Chlorobenzene-d5	8.780

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	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV091224\
Data File : VV037264.D
Acq On : 12 Sep 2024 13:12
Operator : SY/MD
Sample : PB163300TB 10X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
VLEB300

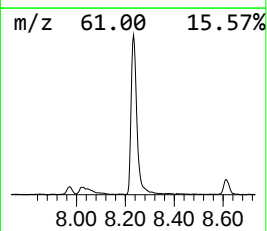
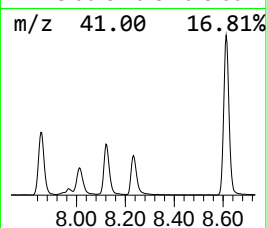
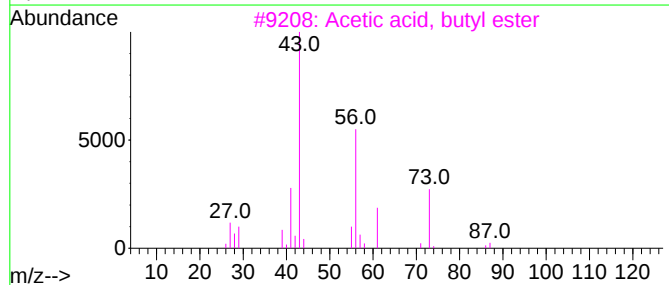
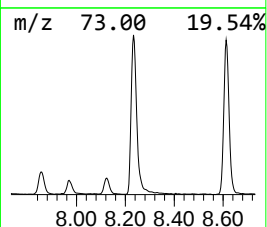
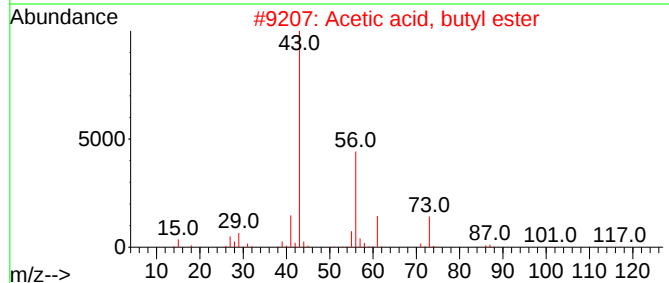
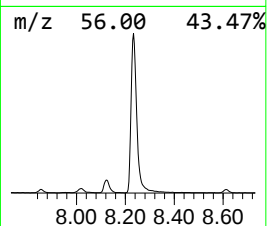
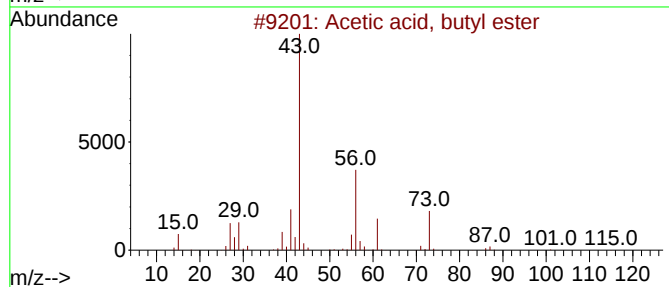
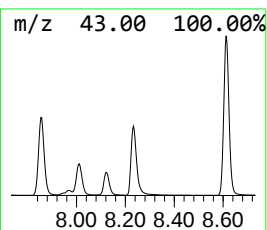
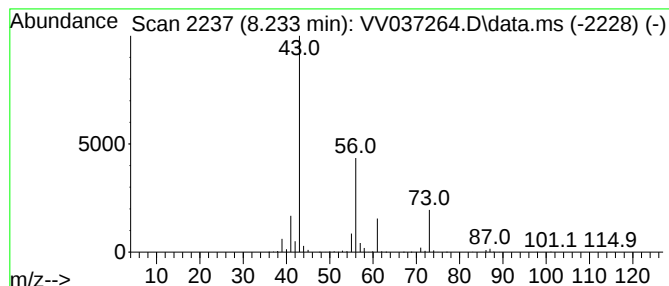
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Quant Title : VOC Analysis

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	35.95 ug/L	1455250	Chlorobenzene-d5	8.780

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	56		



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV091224\
Data File : VV037264.D
Acq On : 12 Sep 2024 13:12
Operator : SY/MD
Sample : PB163300TB 10X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
VLEB300

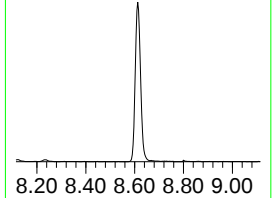
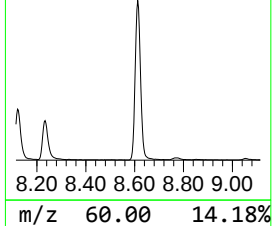
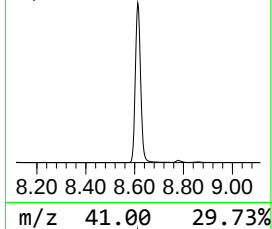
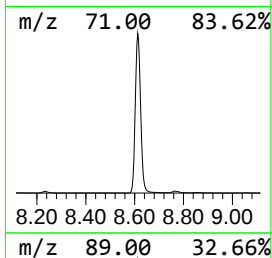
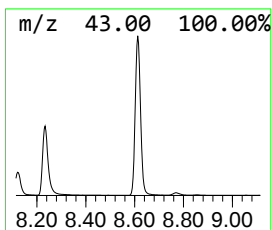
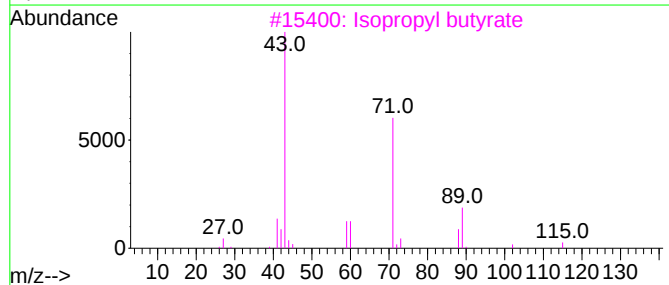
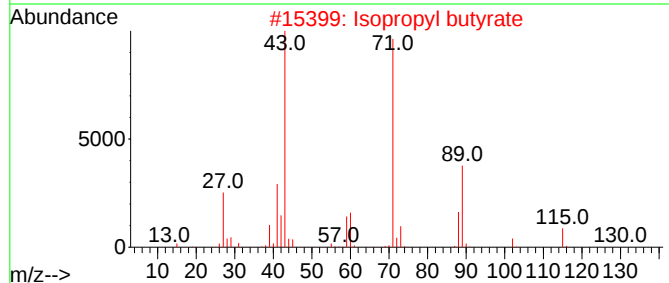
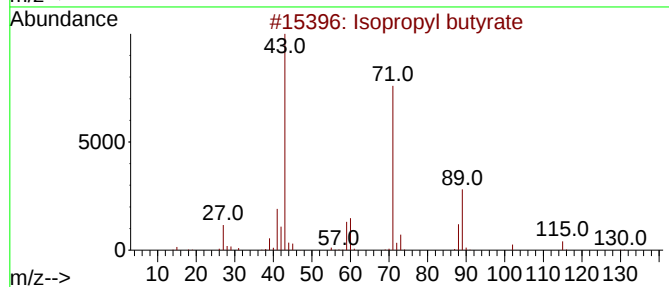
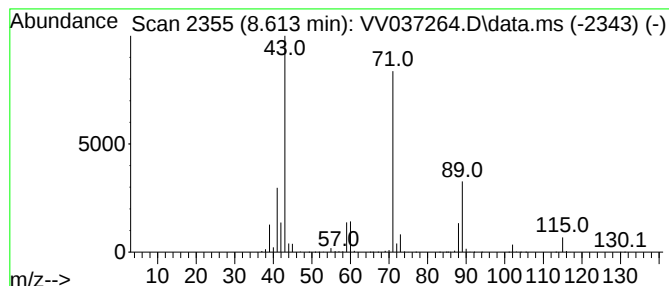
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Quant Title : VOC Analysis

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.
8.613	119.47 ug/L	4835730	Chlorobenzene-d5	8.780

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Isopropyl butyrate	130	C7H14O2	000638-11-9	95
2		Isopropyl butyrate	130	C7H14O2	000638-11-9	91
3		Isopropyl butyrate	130	C7H14O2	000638-11-9	78
4		Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	72
5		Propanoic acid, 2-methyl-, 1-met...	130	C7H14O2	000617-50-5	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV091224\
Data File : VV037264.D
Acq On : 12 Sep 2024 13:12
Operator : SY/MD
Sample : PB163300TB 10X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
VLEB300

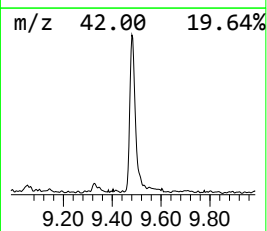
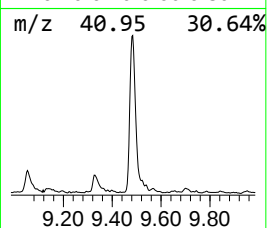
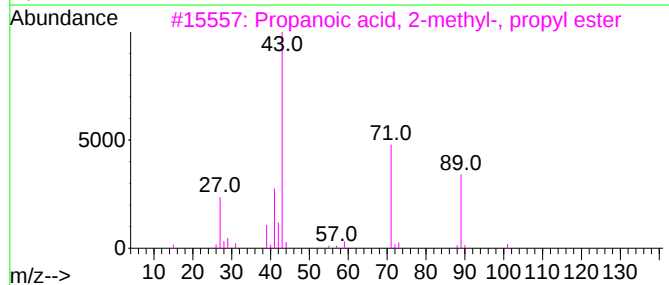
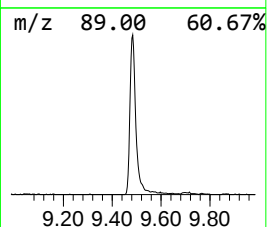
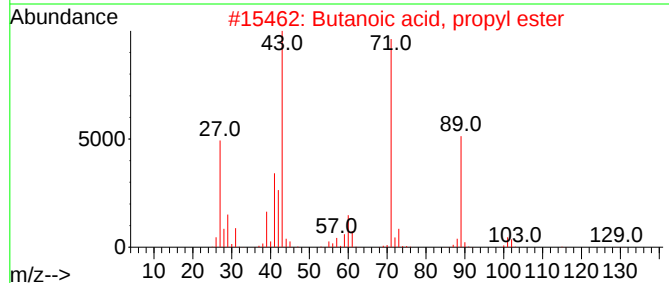
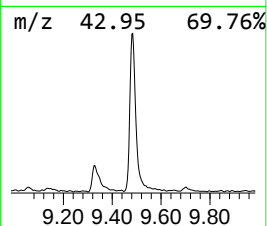
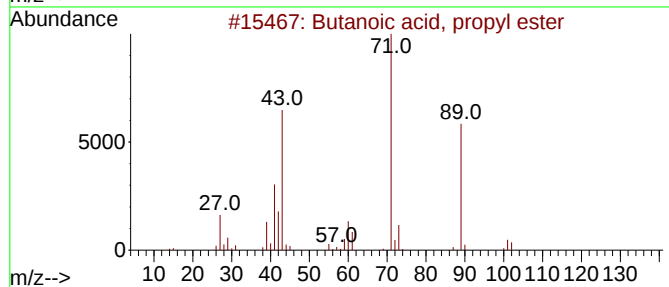
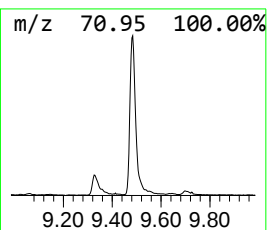
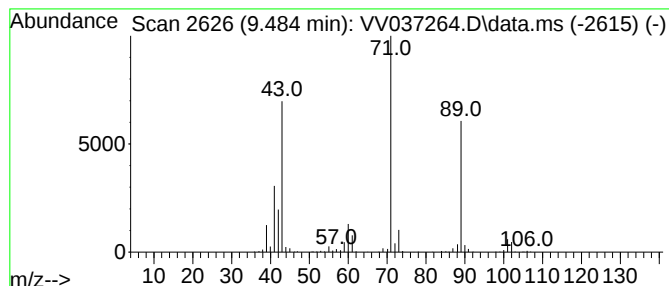
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Quant Title : VOC Analysis

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	9.68 ug/L	391777	Chlorobenzene-d5	8.780

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	86
3			Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	72
4			Propanoic acid, 2-methyl-, propy...	130	C7H14O2	000644-49-5	64
5			Propanoic acid, 2-methyl-, penty...	158	C9H18O2	002445-72-9	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV091224\
Data File : VV037264.D
Acq On : 12 Sep 2024 13:12
Operator : SY/MD
Sample : PB163300TB 10X
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
VLEB300

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	5.198	16.7	ug/L	536441	1	5.529	1604800	50.0
3-Pentanone	6.269	4.1	ug/L	132865	1	5.529	1604800	50.0
n-Propyl acetate	6.388	150.5	ug/L	4831420	1	5.529	1604800	50.0
Propanoic acid,...	7.114	36.3	ug/L	1165730	1	5.529	1604800	50.0
Propanoic acid,...	7.857	41.6	ug/L	1682080	2	8.780	2023770	50.0
Butanoic acid, ...	7.969	7.0	ug/L	281522	2	8.780	2023770	50.0
Propanoic acid,...	8.121	57.8	ug/L	2338660	2	8.780	2023770	50.0
Acetic acid, bu...	8.233	36.0	ug/L	1455250	2	8.780	2023770	50.0
Isopropyl butyrate	8.613	119.5	ug/L	4835730	2	8.780	2023770	50.0
Butanoic acid, ...	9.484	9.7	ug/L	391777	2	8.780	2023770	50.0