

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU061022\
 Data File : VU049042.D
 Acq On : 11 Jun 2022 09:30
 Operator : SY/MD
 Sample : N3249-01
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 56 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 EXYZ2

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\SFAMULM061022WMA.M
 Title : VOC Analysis

Signal : TIC: VU049042.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.600	73	81	101	rBV2	231241	282399	3.08%	1.057%
2	1.925	167	182	199	rBV2	275996	456721	4.98%	1.709%
3	2.076	224	229	235	rVB	4473	5594	0.06%	0.021%
4	2.134	240	247	255	rBV4	3324	5995	0.07%	0.022%
5	2.243	274	281	288	rVB6	2797	3944	0.04%	0.015%
6	2.279	289	292	298	rVB5	1056	891	0.01%	0.003%
7	2.308	298	301	304	rBV3	668	504	0.01%	0.002%
8	2.327	304	307	310	rBV2	717	539	0.01%	0.002%
9	2.472	345	352	357	rBV4	431	621	0.01%	0.002%
10	2.568	364	382	396	rBV4	220075	419442	4.57%	1.569%
11	2.684	415	418	424	rVB2	685	549	0.01%	0.002%
12	2.771	437	445	446	rBV4	494	568	0.01%	0.002%
13	2.780	446	448	454	rVV3	672	614	0.01%	0.002%
14	3.044	520	530	539	rBV5	1681	3325	0.04%	0.012%
15	3.086	539	543	552	rVB2	501	546	0.01%	0.002%
16	3.179	563	572	575	rBV3	2625	3205	0.03%	0.012%
17	3.279	599	603	612	rBV2	421	618	0.01%	0.002%
18	3.353	612	626	644	rVV2	36777	72593	0.79%	0.272%
19	3.694	729	732	741	rVB5	808	1032	0.01%	0.004%
20	3.867	764	786	815	rBV	1082240	2559661	27.91%	9.578%
21	4.526	986	991	992	rBV2	692	540	0.01%	0.002%
22	4.665	1005	1034	1071	rBV	3989787	9171567	100.00%	34.318%
23	5.063	1143	1158	1181	rVB	172052	379205	4.13%	1.419%
24	5.317	1217	1237	1267	rVV	2072058	4658137	50.79%	17.430%
25	5.594	1319	1323	1326	rBV3	611	552	0.01%	0.002%
26	5.722	1343	1363	1378	rBV3	324613	899478	9.81%	3.366%
27	5.886	1410	1414	1415	rBV2	680	516	0.01%	0.002%
28	6.250	1510	1527	1550	rBV	287435	550716	6.00%	2.061%
29	6.330	1550	1552	1558	rVB4	747	576	0.01%	0.002%
30	6.542	1605	1618	1634	rVB	51343	98860	1.08%	0.370%
31	6.690	1651	1664	1679	rBV	216815	424457	4.63%	1.588%
32	6.964	1743	1749	1762	rVB4	1402	2697	0.03%	0.010%
33	7.182	1808	1817	1825	rBV3	1043	2012	0.02%	0.008%
34	7.565	1924	1936	1958	rBV	93891	166224	1.81%	0.622%
35	7.793	1995	2007	2019	rBV2	9297	16392	0.18%	0.061%
36	7.896	2019	2039	2052	rBV	402190	701230	7.65%	2.624%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

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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\SFAMULM061022WMA.M
 Title : VOC Analysis

37	7.970	2052	2062	2074	rVV	72516	124982	1.36%	0.468%
38	8.031	2074	2081	2091	rVB6	3692	6684	0.07%	0.025%
39	8.179	2112	2127	2141	rBV2	70477	124051	1.35%	0.464%
40	8.243	2142	2147	2153	rVB5	1974	2888	0.03%	0.011%
41	8.369	2180	2186	2188	rBV4	974	938	0.01%	0.004%
42	8.401	2190	2196	2205	rVB5	4277	6654	0.07%	0.025%
43	8.475	2208	2219	2228	rVB6	1612	3315	0.04%	0.012%
44	8.552	2232	2243	2256	rBV	125473	206804	2.25%	0.774%
45	8.632	2256	2268	2304	rBV2	308295	603751	6.58%	2.259%
46	8.864	2332	2340	2349	rVB4	4355	6160	0.07%	0.023%
47	8.922	2352	2358	2367	rVB4	2376	3184	0.03%	0.012%
48	9.272	2460	2467	2471	rBV4	1690	1952	0.02%	0.007%
49	9.417	2499	2512	2533	rBV	409643	678552	7.40%	2.539%
50	9.571	2550	2560	2574	rVB	26839	44585	0.49%	0.167%
51	9.693	2586	2598	2616	rBV	54930	90660	0.99%	0.339%
52	10.037	2696	2705	2709	rBV6	1254	1808	0.02%	0.007%
53	10.098	2712	2724	2746	rVB	122975	203037	2.21%	0.760%
54	10.237	2761	2767	2770	rVV3	691	735	0.01%	0.003%
55	10.275	2770	2779	2792	rVB5	5635	8757	0.10%	0.033%
56	10.359	2799	2805	2810	rBV5	1637	2092	0.02%	0.008%
57	10.481	2835	2843	2853	rBV3	4852	7624	0.08%	0.029%
58	10.578	2865	2873	2884	rVB4	1394	2015	0.02%	0.008%
59	10.700	2905	2911	2916	rBV5	1921	2468	0.03%	0.009%
60	10.754	2916	2928	2941	rBV	336890	551622	6.01%	2.064%
61	10.902	2959	2974	2989	rBV2	15373	27902	0.30%	0.104%
62	10.999	2993	3004	3006	rBV2	40473	52023	0.57%	0.195%
63	11.089	3021	3032	3045	rVB	125000	194071	2.12%	0.726%
64	11.230	3066	3076	3083	rBV5	3035	5509	0.06%	0.021%
65	11.291	3083	3095	3112	rVB	153877	239913	2.62%	0.898%
66	11.413	3127	3133	3134	rBV2	1390	1066	0.01%	0.004%
67	11.468	3139	3150	3167	rVB	105776	162756	1.77%	0.609%
68	11.552	3171	3176	3179	rBV2	661	535	0.01%	0.002%
69	11.610	3187	3194	3198	rVV5	3119	4536	0.05%	0.017%
70	11.642	3199	3204	3216	rVB3	7366	10369	0.11%	0.039%
71	11.745	3225	3236	3244	rVV3	18172	27179	0.30%	0.102%
72	11.812	3244	3257	3272	rVV	426180	691914	7.54%	2.589%
73	11.896	3272	3283	3304	rVB	209900	327550	3.57%	1.226%
74	12.005	3310	3317	3328	rVB5	3161	4618	0.05%	0.017%
75	12.095	3334	3345	3363	rBV3	79617	148567	1.62%	0.556%
76	12.192	3363	3375	3397	rVB	480486	791113	8.63%	2.960%

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Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM061022WMA.M
 Title : VOC Analysis

77	12.301	3402	3409	3422	rVB7	6402	10116	0.11%	0.038%
78	12.378	3422	3433	3442	rBV2	17543	25756	0.28%	0.096%
79	12.484	3452	3466	3481	rBV3	65671	135608	1.48%	0.507%
80	12.571	3485	3493	3502	rVV2	27865	44267	0.48%	0.166%
81	12.610	3502	3505	3518	rVB9	4597	6004	0.07%	0.022%
82	12.687	3519	3529	3549	rVB6	15447	40736	0.44%	0.152%
83	12.767	3549	3554	3556	rBV3	1416	1220	0.01%	0.005%
84	12.873	3577	3587	3601	rVB4	13598	21664	0.24%	0.081%
85	12.973	3602	3618	3626	rVV3	16766	28977	0.32%	0.108%
86	13.024	3626	3634	3647	rVB2	25058	39441	0.43%	0.148%
87	13.114	3654	3662	3666	rBV7	2224	2840	0.03%	0.011%
88	13.140	3666	3670	3671	rBV2	1160	837	0.01%	0.003%
89	13.253	3698	3705	3712	rBV5	1388	2314	0.03%	0.009%
90	13.294	3712	3718	3726	rVV6	3649	4664	0.05%	0.017%
91	13.352	3733	3736	3744	rVB6	984	984	0.01%	0.004%
92	13.407	3749	3753	3756	rVV4	1080	1068	0.01%	0.004%
93	13.452	3757	3767	3781	rVB3	26247	45221	0.49%	0.169%
94	13.565	3797	3802	3803	rBV2	685	554	0.01%	0.002%
95	13.626	3811	3821	3833	rVB7	8829	13418	0.15%	0.050%
96	13.738	3853	3856	3863	rVB4	975	966	0.01%	0.004%
97	13.831	3879	3885	3898	rBV8	2205	4600	0.05%	0.017%
98	14.095	3957	3967	3981	rVV2	13926	25562	0.28%	0.096%
99	14.333	4038	4041	4049	rVV3	690	627	0.01%	0.002%
100	15.876	4519	4521	4528	rVB5	741	531	0.01%	0.002%

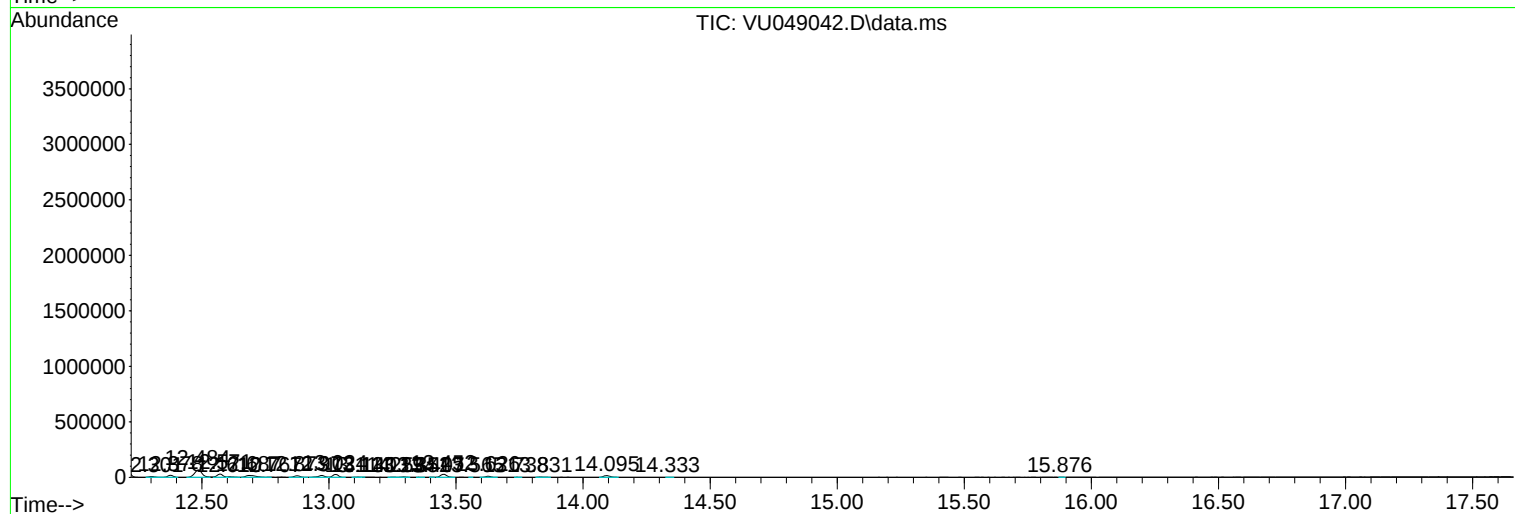
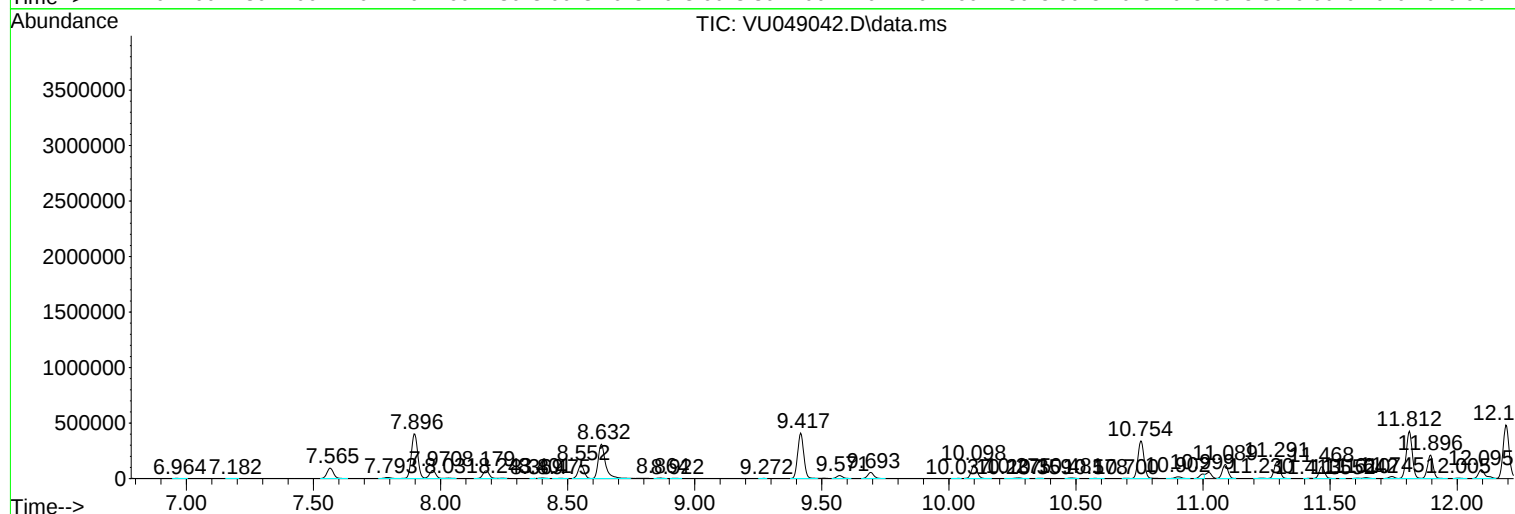
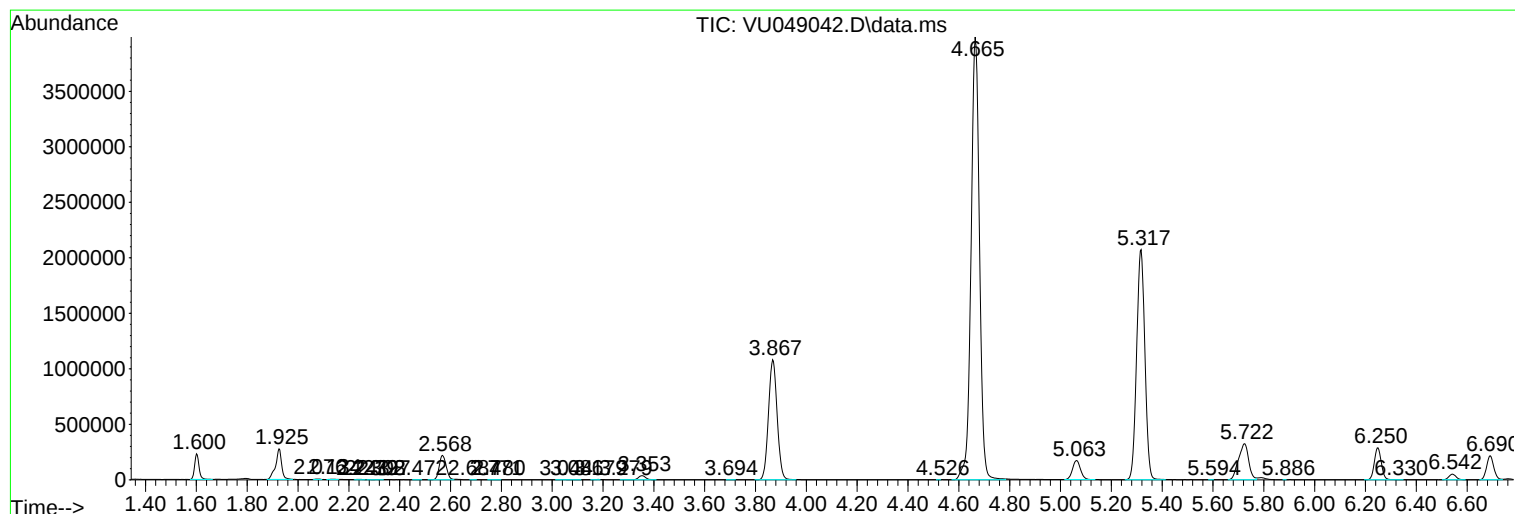
Sum of corrected areas: 26725434

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ALS Vial : 56 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM061022WMA.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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ClientSampleId :
EXYZ2

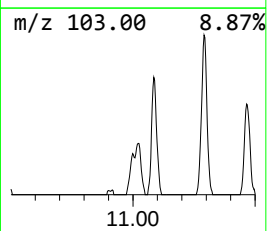
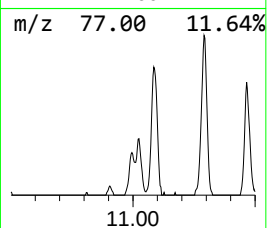
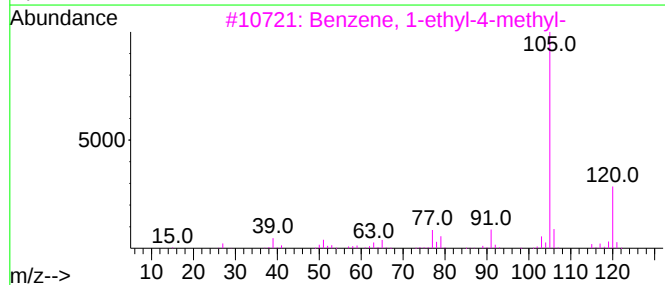
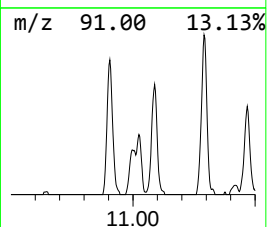
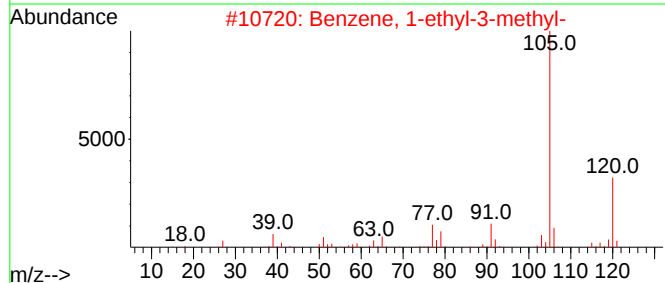
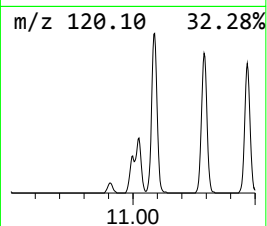
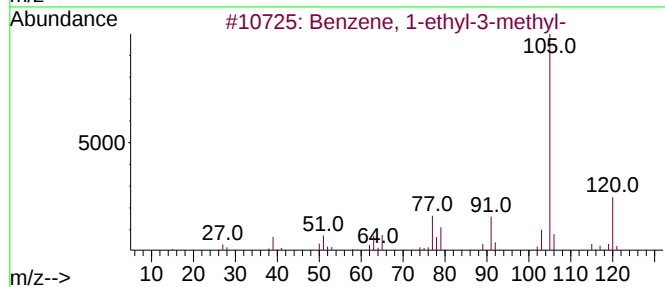
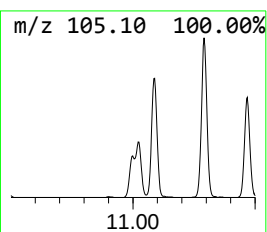
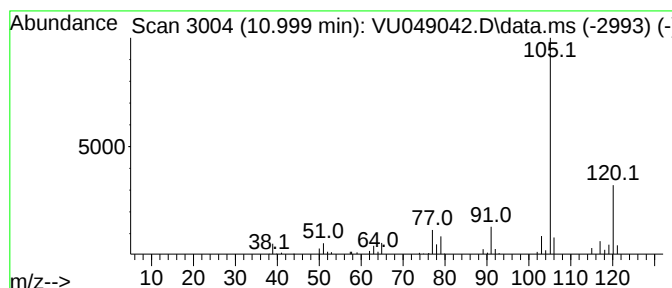
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM061022WMA.M
Quant Title : VOC Analysis

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

Peak Number 2 Benzene, 1-ethyl-3-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.999	3.76 ug/L	52023	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
2			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



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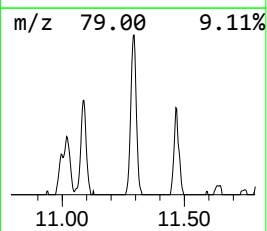
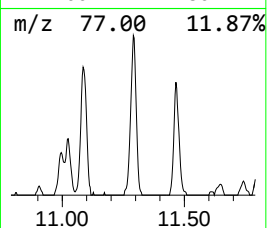
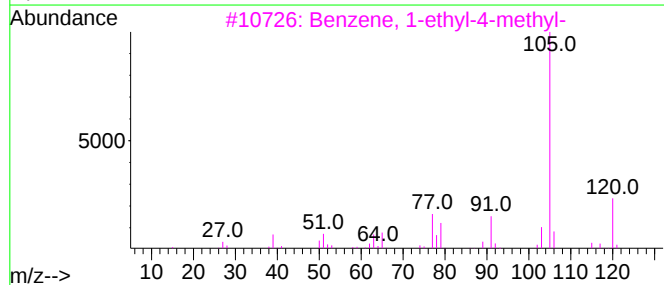
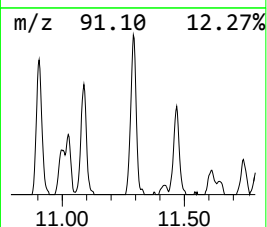
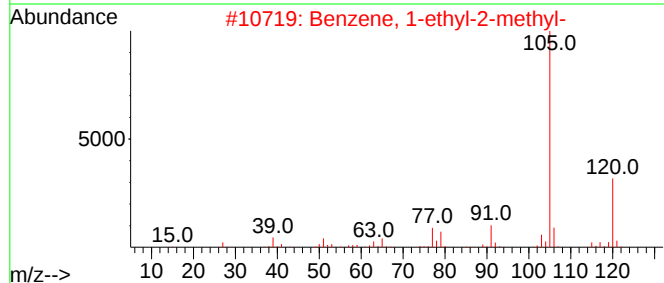
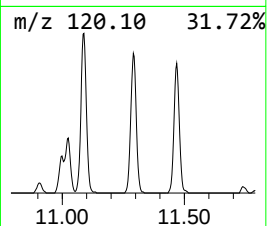
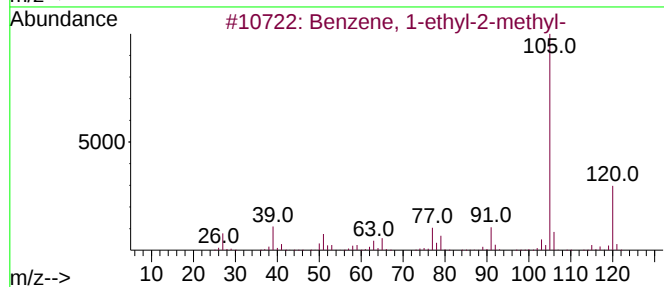
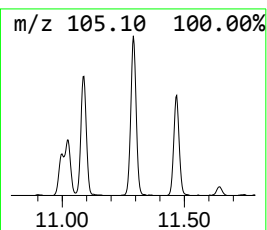
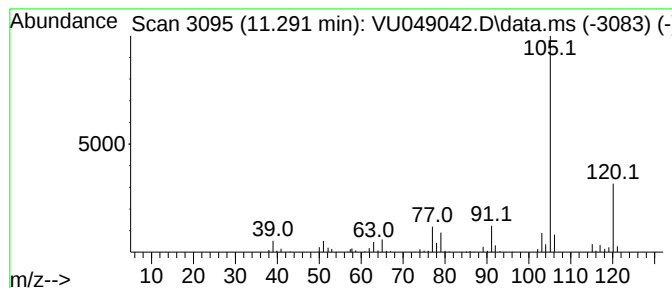
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM061022WMA.M
Quant Title : VOC Analysis

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

Peak Number 3 Benzene, 1-ethyl-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.291	17.34 ug/L	239913	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5			Mesitylene	120	C9H12	000108-67-8	91



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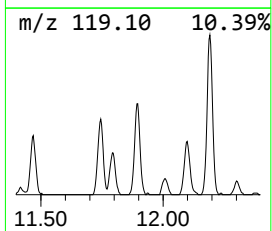
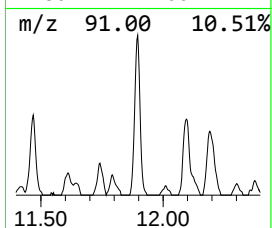
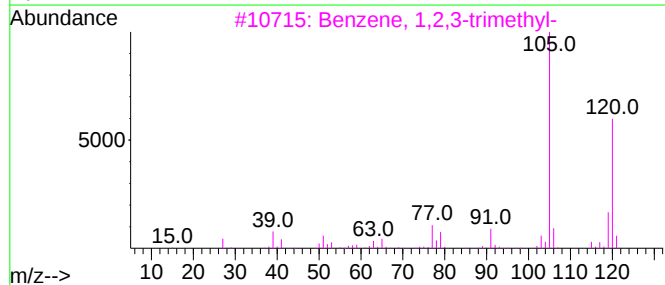
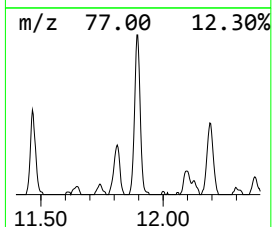
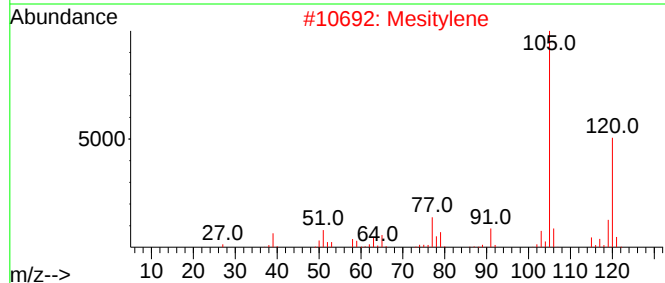
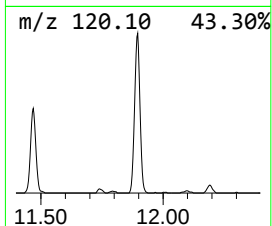
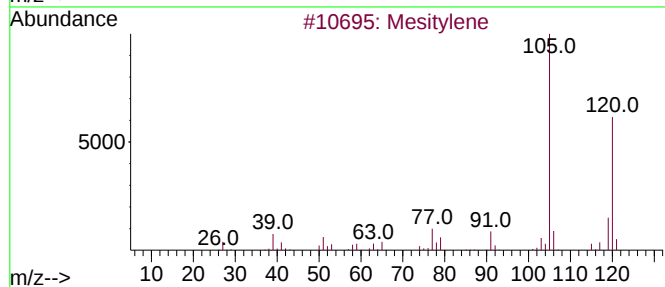
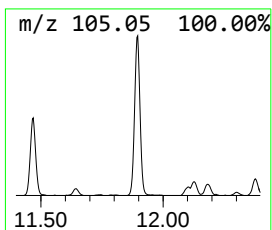
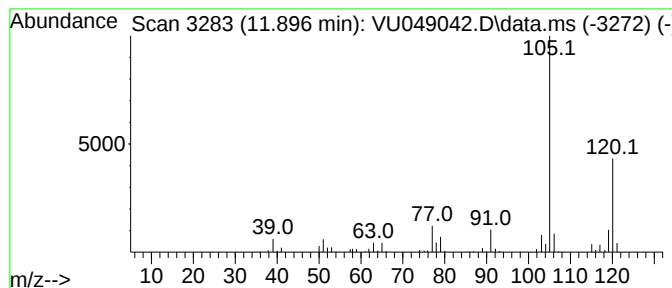
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Quant Title : VOC Analysis

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

Peak Number 4 Benzene, 1,2,3-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.896	23.67 ug/L	327550	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Mesitylene	120	C9H12	000108-67-8	97
2			Mesitylene	120	C9H12	000108-67-8	97
3			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
4			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU061022\
Data File : VU049042.D
Acq On : 11 Jun 2022 09:30
Operator : SY/MD
Sample : N3249-01
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 56 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXYZ2

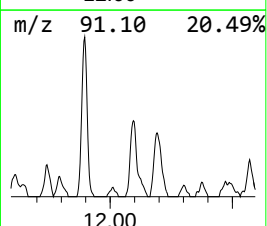
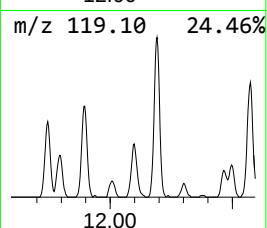
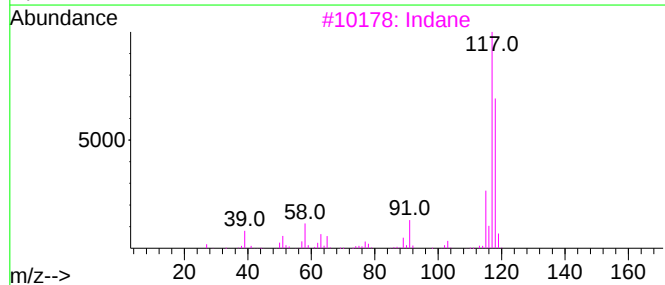
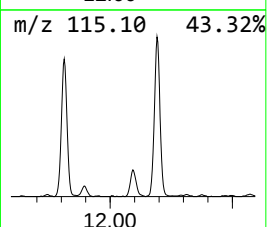
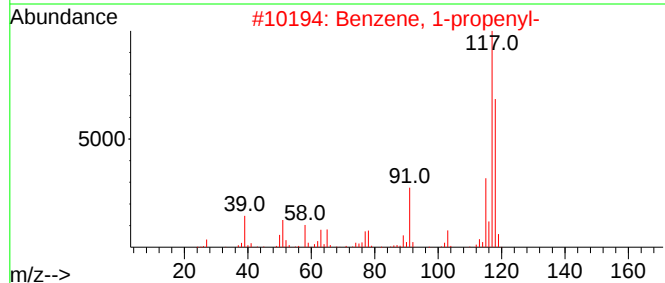
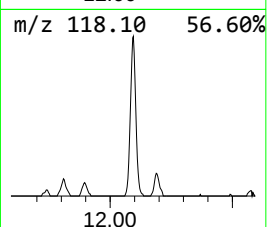
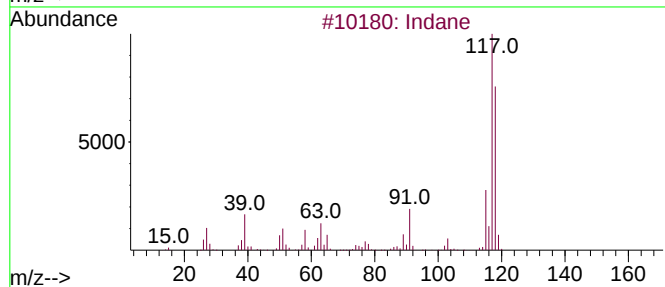
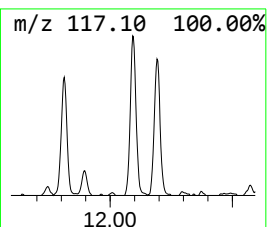
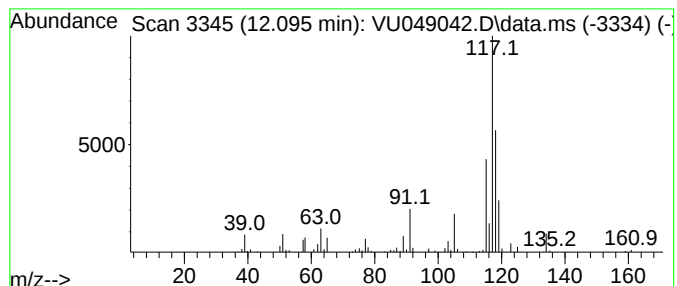
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Quant Title : VOC Analysis

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

Peak Number 5 Indane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.095	10.74 ug/L	148567	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	92
2			Benzene, 1-propenyl-	118	C9H10	000637-50-3	64
3			Indane	118	C9H10	000496-11-7	64
4			Benzene, cyclopropyl-	118	C9H10	000873-49-4	60
5			Bicyclo[4.2.0]octa-1,3,5-triene,...	118	C9H10	055337-80-9	60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU061022\
Data File : VU049042.D
Acq On : 11 Jun 2022 09:30
Operator : SY/MD
Sample : N3249-01
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 56 Sample Multiplier: 1

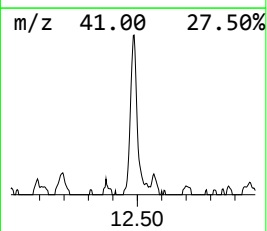
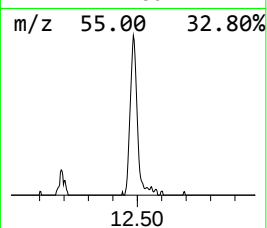
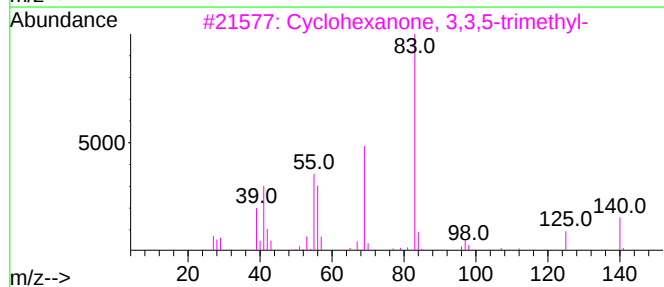
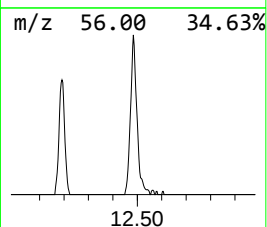
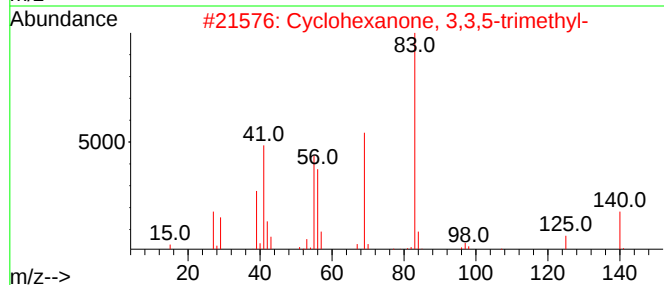
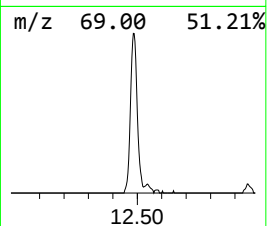
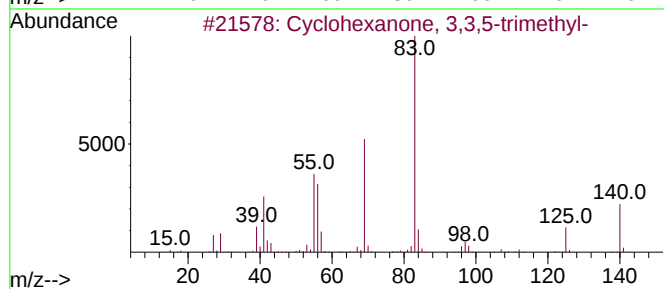
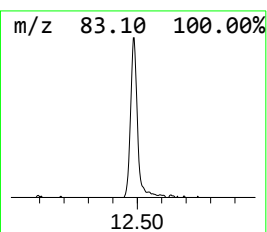
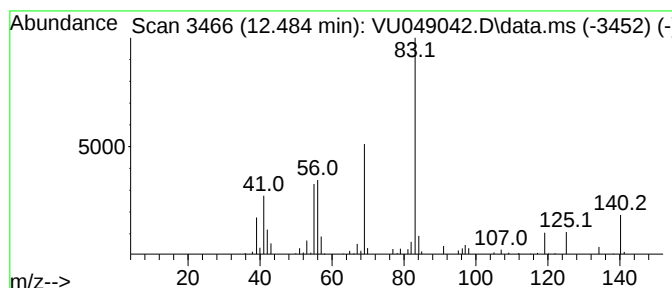
Instrument :
MSVOA_U
ClientSampleId :
EXYZ2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM061022WMA.M
Quant Title : VOC Analysis

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

Peak Number 6 Cyclohexanone, 3,3,5-trimet... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD			R.T.
12.484	9.80 ug/L	135608	1,4-Dichlorobenzene-d4			11.812
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Cyclohexanone, 3,3,5-trimethyl-		140	C9H16O	000873-94-9	95
2	Cyclohexanone, 3,3,5-trimethyl-		140	C9H16O	000873-94-9	93
3	Cyclohexanone, 3,3,5-trimethyl-		140	C9H16O	000873-94-9	93
4	Cyclohexanone, 3,3,5-trimethyl-		140	C9H16O	000873-94-9	87
5	1-Methyl-1H-1,2,4-triazole		83	C3H5N3	006086-21-1	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU061022\
Data File : VU049042.D
Acq On : 11 Jun 2022 09:30
Operator : SY/MD
Sample : N3249-01
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 56 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXYZ2

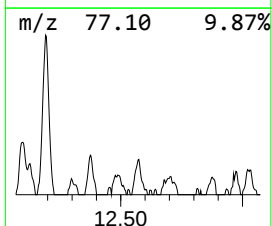
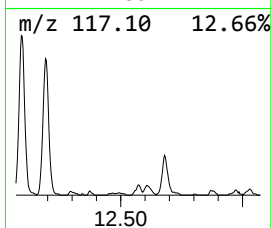
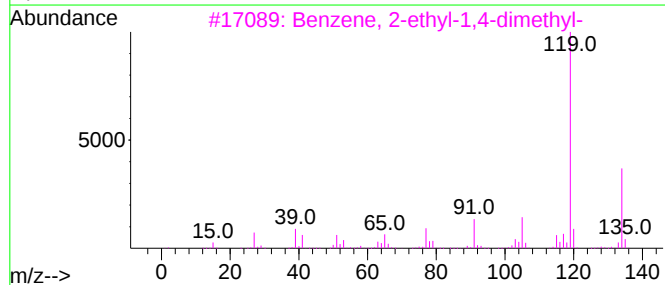
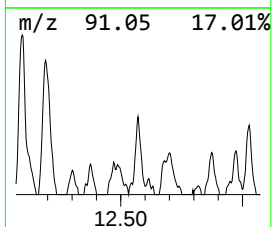
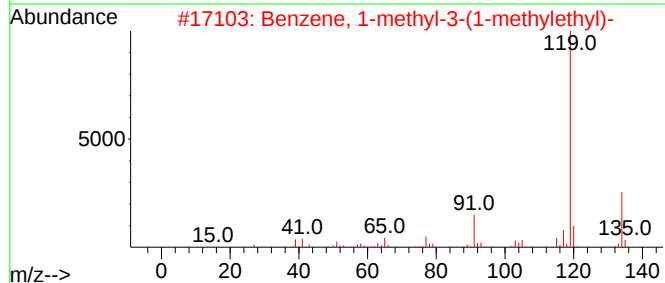
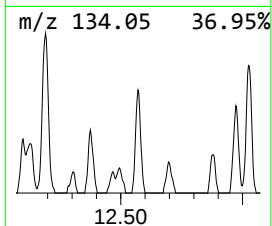
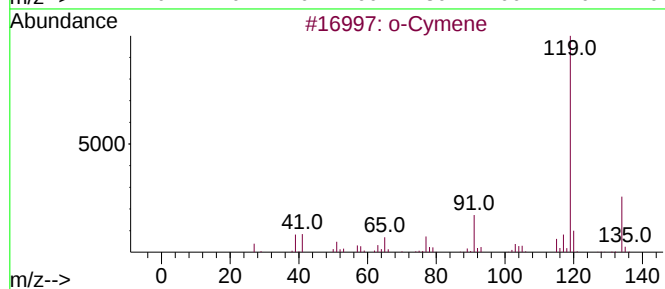
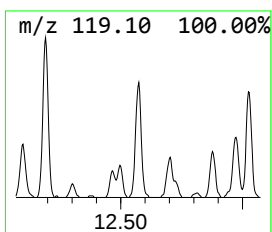
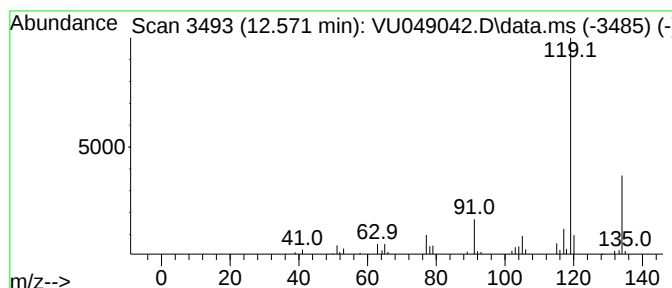
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Quant Title : VOC Analysis

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

Peak Number 7 o-Cymene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.571	3.20 ug/L	44267	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	94
2			Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	93
3			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	93
4			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	93
5			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU061022\
Data File : VU049042.D
Acq On : 11 Jun 2022 09:30
Operator : SY/MD
Sample : N3249-01
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 56 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXYZ2

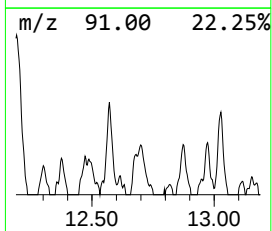
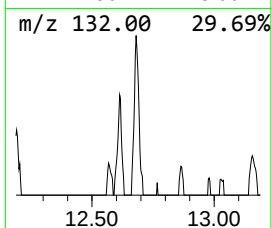
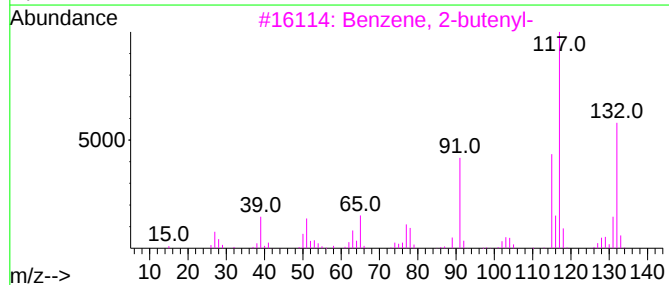
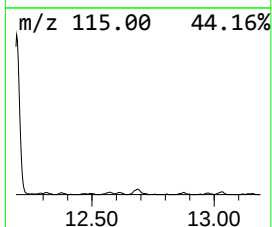
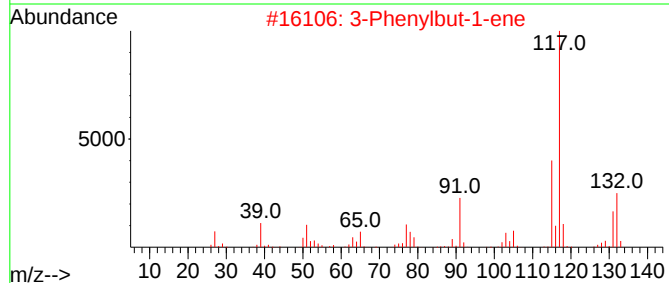
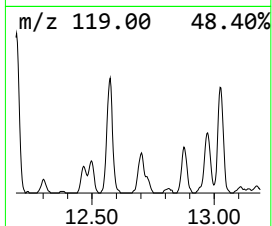
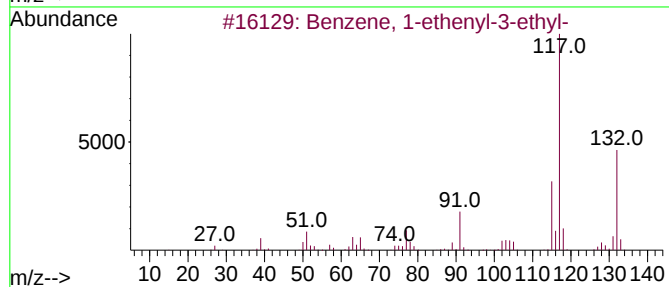
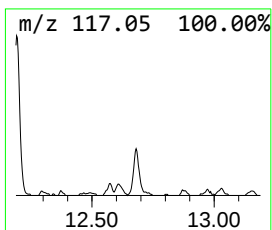
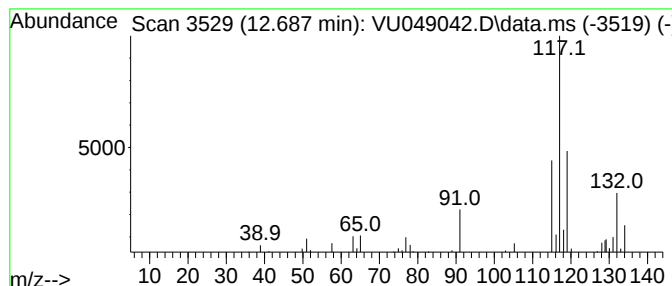
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Quant Title : VOC Analysis

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

Peak Number 8 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.687	2.94 ug/L	40736	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	70
2			3-Phenylbut-1-ene	132	C10H12	000934-10-1	70
3			Benzene, 2-butenyl-	132	C10H12	001560-06-1	68
4			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	68
5			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU061022\
Data File : VU049042.D
Acq On : 11 Jun 2022 09:30
Operator : SY/MD
Sample : N3249-01
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 56 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXYZ2

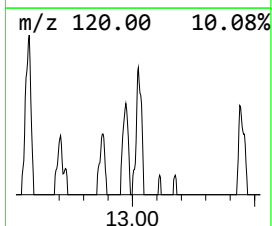
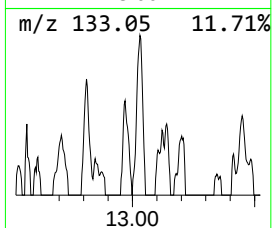
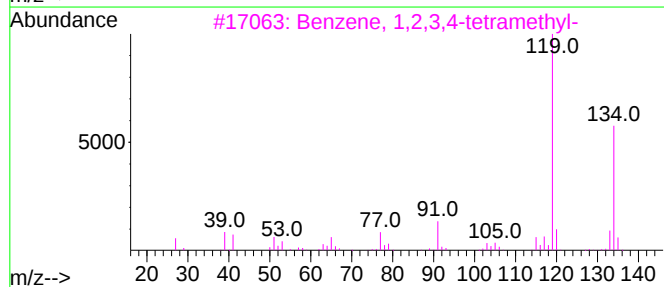
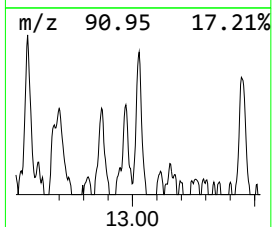
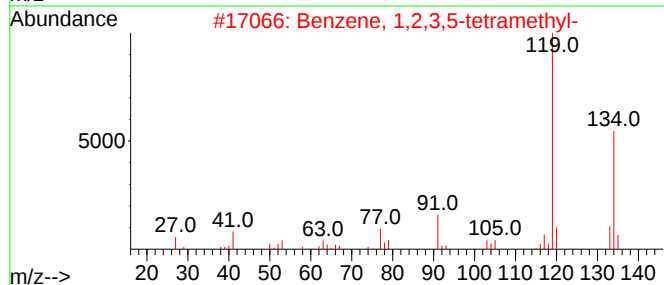
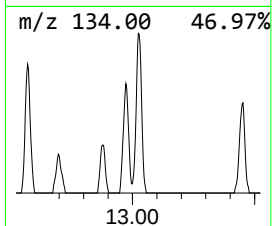
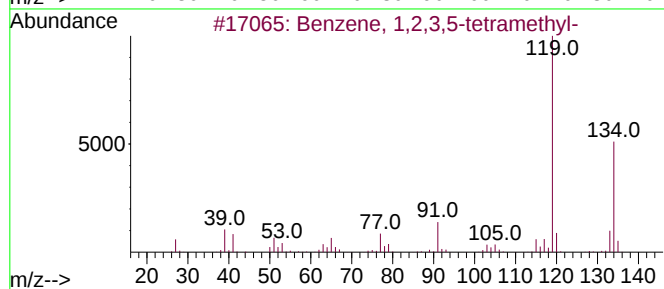
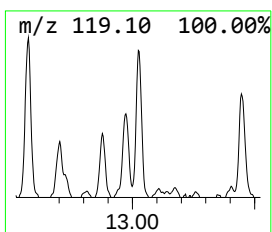
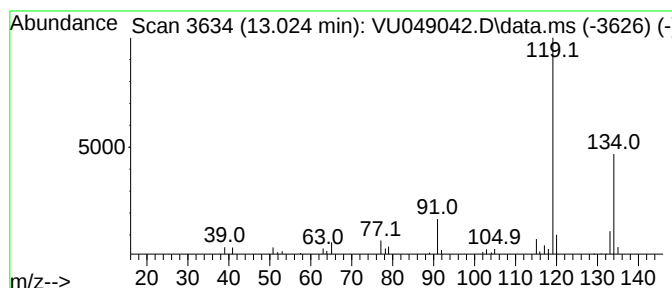
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Quant Title : VOC Analysis

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

Peak Number 9 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.024	2.85 ug/L	39441	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
2			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
3			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
4			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	93
5			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU061022\
Data File : VU049042.D
Acq On : 11 Jun 2022 09:30
Operator : SY/MD
Sample : N3249-01
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 56 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXYZ2

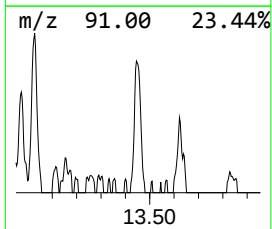
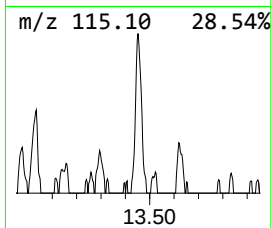
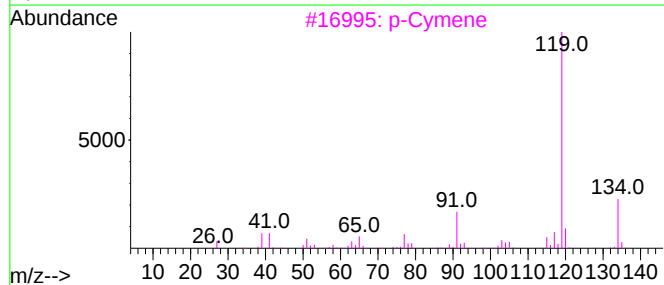
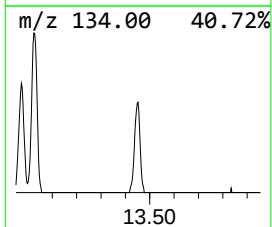
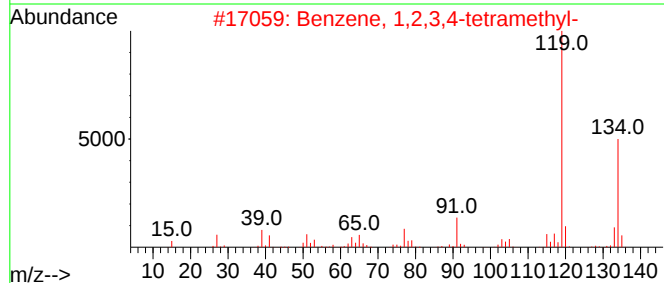
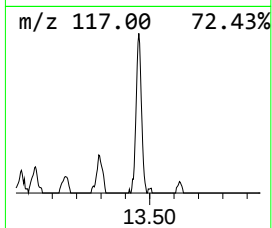
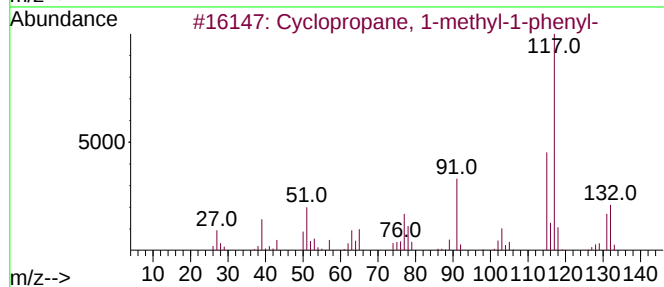
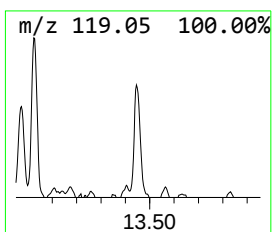
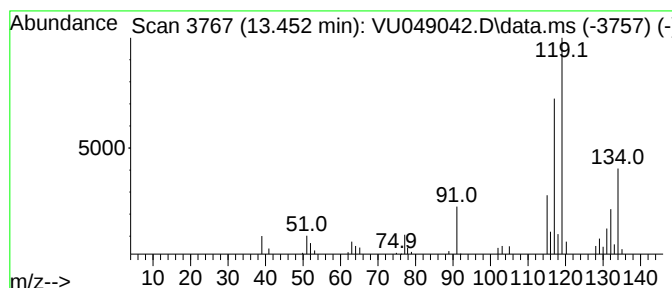
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Quant Title : VOC Analysis

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

Peak Number 10 Cyclopropane, 1-methyl-1-ph... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.452	3.27 ug/L	45221	1,4-Dichlorobenzene-d4	11.812

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopropane, 1-methyl-1-phenyl-	132	C10H12	002214-14-4	84
2		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	49
3		p-Cymene	134	C10H14	000099-87-6	49
4		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	49
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	49



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU061022\
Data File : VU049042.D
Acq On : 11 Jun 2022 09:30
Operator : SY/MD
Sample : N3249-01
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 56 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXYZ2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM061022WMA.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
Benzene, 1-ethy...	10.999	3.8	ug/L	52023	3	11.812	691914	50.0
Benzene, 1-ethy...	11.291	17.3	ug/L	239913	3	11.812	691914	50.0
Benzene, 1,2,3-...	11.896	23.7	ug/L	327550	3	11.812	691914	50.0
Indane	12.095	10.7	ug/L	148567	3	11.812	691914	50.0
Cyclohexanone, ...	12.484	9.8	ug/L	135608	3	11.812	691914	50.0
o-Cymene	12.571	3.2	ug/L	44267	3	11.812	691914	50.0
Benzene, 1-ethe...	12.687	2.9	ug/L	40736	3	11.812	691914	50.0
Benzene, 1,2,3,...	13.024	2.9	ug/L	39441	3	11.812	691914	50.0
Cyclopropane, 1...	13.452	3.3	ug/L	45221	3	11.812	691914	50.0