

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120621\
 Data File : VU046115.D
 Acq On : 06 Dec 2021 16:27
 Operator : SY/MD
 Sample : M4942-05DL 20X
 Misc : 6.58g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BGKQ5DL

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\SFAMULM112921WMA.M

Title : VOC Analysis

Signal : TIC: VU046115.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	93	rVB	98385	107360	8.29%	1.064%
2	1.916	170	179	190	rVB	65214	80905	6.25%	0.802%
3	1.996	201	204	206	rVB	547	312	0.02%	0.003%
4	2.510	360	364	367	rVB	319	208	0.02%	0.002%
5	2.568	370	382	404	rBV	172733	288177	22.25%	2.856%
6	2.800	447	454	457	rVV2	310	357	0.03%	0.004%
7	3.051	522	532	544	rVB3	4562	8034	0.62%	0.080%
8	3.182	565	573	574	rBV2	1779	2045	0.16%	0.020%
9	3.378	630	634	637	rBV	164	201	0.02%	0.002%
10	3.671	723	725	728	rVB	323	187	0.01%	0.002%
11	4.401	950	952	955	rVB2	313	192	0.01%	0.002%
12	4.430	959	961	966	rVV	245	202	0.02%	0.002%
13	4.632	1008	1024	1051	rBV	77203	197758	15.27%	1.960%
14	5.067	1142	1159	1183	rBV	134942	314589	24.29%	3.118%
15	5.186	1194	1196	1202	rBV2	191	233	0.02%	0.002%
16	5.285	1223	1227	1228	rBV	331	222	0.02%	0.002%
17	5.340	1240	1244	1247	rBV	216	236	0.02%	0.002%
18	5.468	1282	1284	1289	rVB	291	218	0.02%	0.002%
19	5.729	1339	1365	1387	rBV2	300085	808003	62.38%	8.008%
20	5.941	1427	1431	1434	rBV2	229	201	0.02%	0.002%
21	5.976	1438	1442	1446	rBV3	240	271	0.02%	0.003%
22	6.034	1456	1460	1462	rBV2	249	216	0.02%	0.002%
23	6.070	1469	1471	1475	rVB	280	245	0.02%	0.002%
24	6.089	1475	1477	1480	rBV	256	183	0.01%	0.002%
25	6.250	1513	1527	1544	rBV	235128	441297	34.07%	4.374%
26	6.404	1572	1575	1579	rBV	296	260	0.02%	0.003%
27	6.533	1606	1615	1625	rBV4	3899	7189	0.56%	0.071%
28	6.693	1651	1665	1682	rBV	184473	354979	27.40%	3.518%
29	6.806	1696	1700	1705	rBV2	505	480	0.04%	0.005%
30	7.214	1824	1827	1832	rVB	196	222	0.02%	0.002%
31	7.247	1832	1837	1841	rBV	253	286	0.02%	0.003%
32	7.568	1924	1937	1953	rBV	106387	186784	14.42%	1.851%
33	7.899	2026	2040	2056	rBV	388859	670738	51.78%	6.648%
34	8.179	2116	2127	2144	rBV	71106	119779	9.25%	1.187%
35	8.240	2144	2146	2149	rVB2	544	276	0.02%	0.003%
36	8.632	2256	2268	2295	rBV	238189	414490	32.00%	4.108%

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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 >

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Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
 Title : VOC Analysis

37	8.790	2315	2317	2320	rVB	427	213	0.02%	0.002%
38	8.870	2334	2342	2348	rBV3	1276	1944	0.15%	0.019%
39	8.918	2348	2357	2359	rBV4	1169	1439	0.11%	0.014%
40	8.954	2365	2368	2370	rVB2	414	250	0.02%	0.002%
41	8.973	2370	2374	2377	rBV2	258	213	0.02%	0.002%
42	9.028	2388	2391	2395	rVB2	233	188	0.01%	0.002%
43	9.057	2395	2400	2403	rBV2	410	384	0.03%	0.004%
44	9.124	2416	2421	2424	rBV5	448	506	0.04%	0.005%
45	9.172	2427	2436	2442	rBV4	1067	1717	0.13%	0.017%
46	9.217	2442	2450	2456	rVB4	1578	2090	0.16%	0.021%
47	9.333	2479	2486	2487	rBV3	1999	1552	0.12%	0.015%
48	9.420	2499	2513	2532	rBV	325839	544979	42.07%	5.401%
49	9.510	2538	2541	2548	rVB5	1224	1194	0.09%	0.012%
50	9.571	2548	2560	2581	rBV	226800	362701	28.00%	3.595%
51	9.693	2584	2598	2615	rBV	793528	1295314	100.00%	12.838%
52	9.886	2650	2658	2659	rBV2	359	395	0.03%	0.004%
53	10.025	2690	2701	2709	rBV6	3433	7515	0.58%	0.074%
54	10.102	2714	2725	2742	rVB	145003	237564	18.34%	2.354%
55	10.250	2760	2771	2786	rVB6	5886	12651	0.98%	0.125%
56	10.314	2786	2791	2795	rBV5	1119	1275	0.10%	0.013%
57	10.356	2797	2804	2813	rBV8	5292	8773	0.68%	0.087%
58	10.484	2830	2844	2854	rBV	9335	14880	1.15%	0.147%
59	10.558	2854	2867	2875	rBV8	3644	7062	0.55%	0.070%
60	10.674	2895	2903	2917	rVB6	3832	7321	0.57%	0.073%
61	10.758	2917	2929	2943	rBV	266277	421278	32.52%	4.175%
62	10.906	2960	2975	2986	rBV	36634	55983	4.32%	0.555%
63	10.999	2994	3004	3021	rBV	82217	186205	14.38%	1.845%
64	11.089	3021	3032	3056	rVB	80644	135452	10.46%	1.342%
65	11.182	3058	3061	3063	rBV3	686	477	0.04%	0.005%
66	11.240	3075	3079	3080	rBV3	649	472	0.04%	0.005%
67	11.295	3086	3096	3113	rBV	40812	69756	5.39%	0.691%
68	11.378	3115	3122	3127	rVV6	4195	5586	0.43%	0.055%
69	11.417	3128	3134	3140	rVV3	13443	17582	1.36%	0.174%
70	11.468	3140	3150	3174	rVB	228526	365002	28.18%	3.618%
71	11.574	3174	3183	3189	rBV2	5303	9568	0.74%	0.095%
72	11.613	3189	3195	3199	rVV2	8595	11923	0.92%	0.118%
73	11.642	3200	3204	3214	rVB	15060	21137	1.63%	0.209%
74	11.706	3215	3224	3229	rBV7	7644	13829	1.07%	0.137%
75	11.745	3229	3236	3244	rVB	22198	33039	2.55%	0.327%
76	11.812	3245	3257	3270	rVV	384062	623483	48.13%	6.179%

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Integration Parameters: LSCINT.P

Integrator: RTE

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Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
 Title : VOC Analysis

77	11.896	3272	3283	3293	rVB	59397	95737	7.39%	0.949%
78	12.098	3334	3346	3350	rBV3	52765	86494	6.68%	0.857%
79	12.195	3363	3376	3390	rVB	469862	794804	61.36%	7.877%
80	12.311	3405	3412	3424	rVB6	12734	24817	1.92%	0.246%
81	12.378	3425	3433	3449	rVB	25389	43367	3.35%	0.430%
82	12.468	3451	3461	3465	rBV	28165	44941	3.47%	0.445%
83	12.571	3486	3493	3501	rVB	35347	49526	3.82%	0.491%
84	12.703	3517	3534	3551	rVV9	13293	44913	3.47%	0.445%
85	12.764	3551	3553	3557	rVB5	1417	812	0.06%	0.008%
86	12.973	3612	3618	3626	rVB	24473	33995	2.62%	0.337%
87	13.028	3627	3635	3650	rVB	34529	57314	4.42%	0.568%
88	13.352	3732	3736	3743	rVB4	2676	3041	0.23%	0.030%
89	13.407	3746	3753	3754	rBV6	2804	3243	0.25%	0.032%
90	13.452	3760	3767	3780	rVB3	21364	39408	3.04%	0.391%
91	13.587	3797	3809	3814	rBV6	5642	9834	0.76%	0.097%
92	13.626	3816	3821	3831	rVB4	9758	14708	1.14%	0.146%
93	13.729	3849	3853	3854	rBV3	1458	960	0.07%	0.010%
94	13.796	3872	3874	3878	rBV4	581	348	0.03%	0.003%
95	14.047	3946	3952	3954	rBV6	2002	2208	0.17%	0.022%
96	14.085	3955	3964	3983	rVV	63532	105297	8.13%	1.044%
97	14.192	3987	3997	4011	rVB10	6547	14632	1.13%	0.145%
98	14.449	4072	4077	4084	rVB4	3741	4441	0.34%	0.044%
99	15.220	4305	4317	4331	rBV	45438	78478	6.06%	0.778%
100	15.410	4368	4376	4394	rVB	26621	46245	3.57%	0.458%

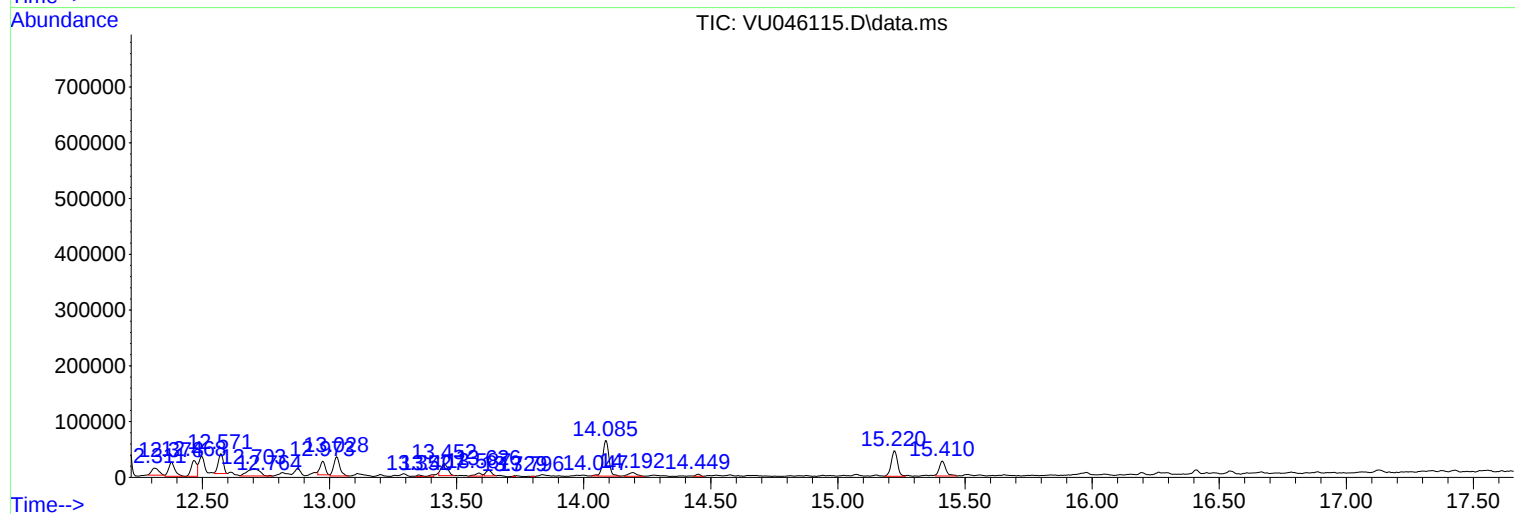
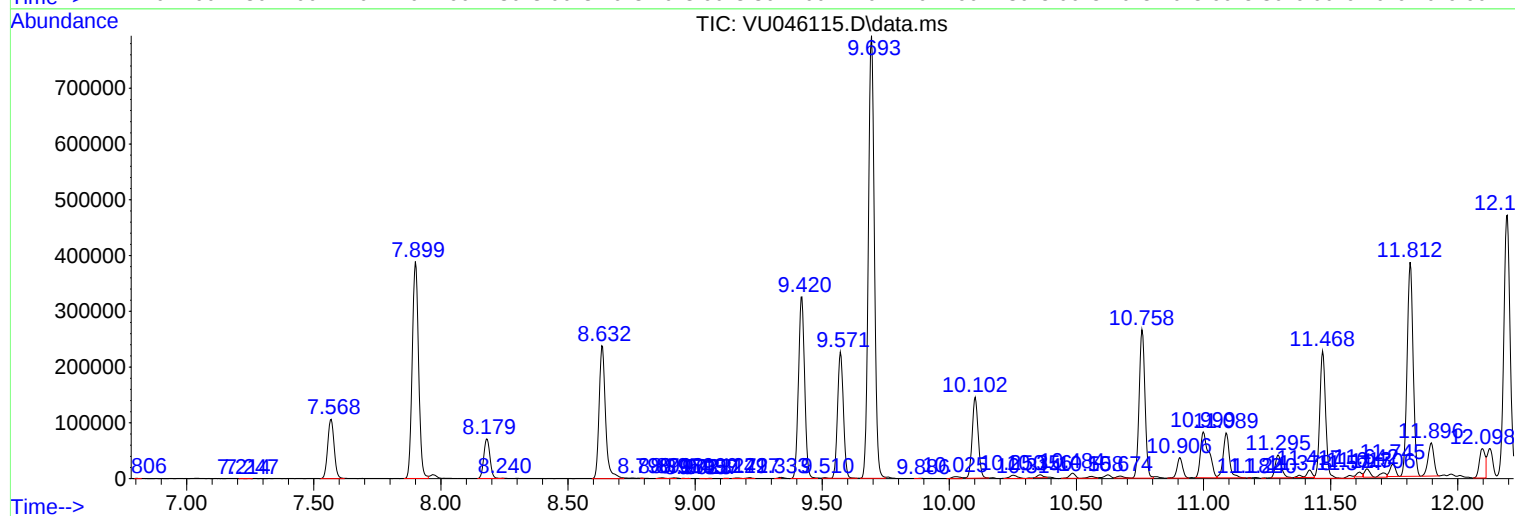
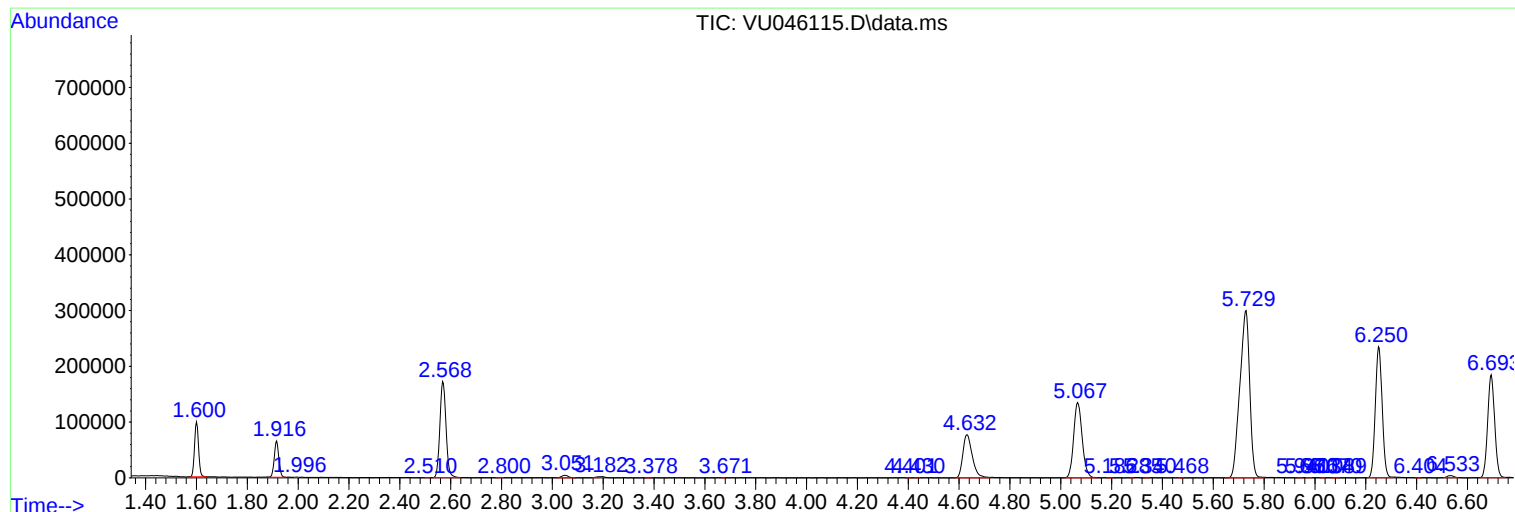
Sum of corrected areas: 10089790

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ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.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 :
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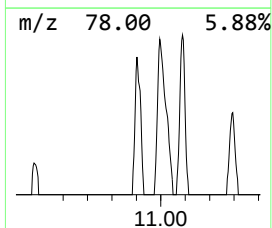
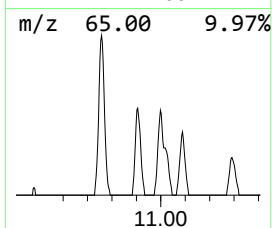
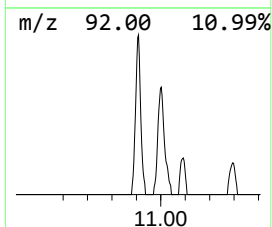
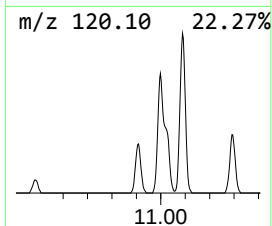
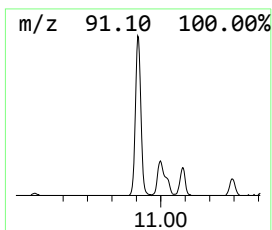
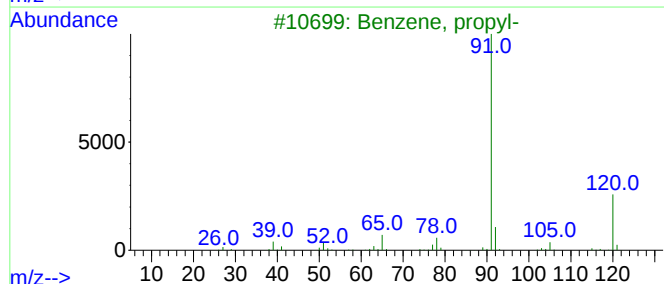
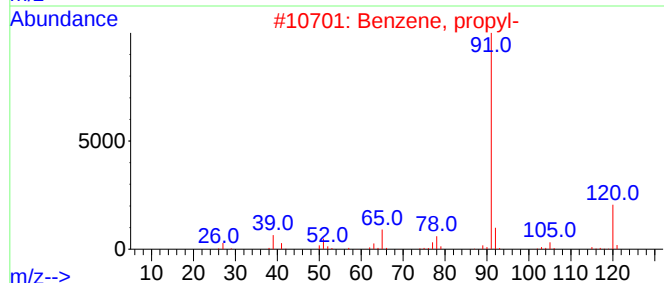
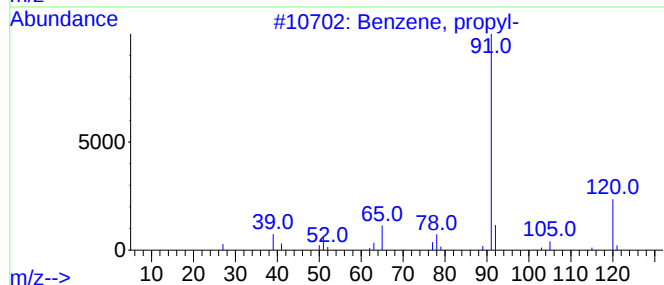
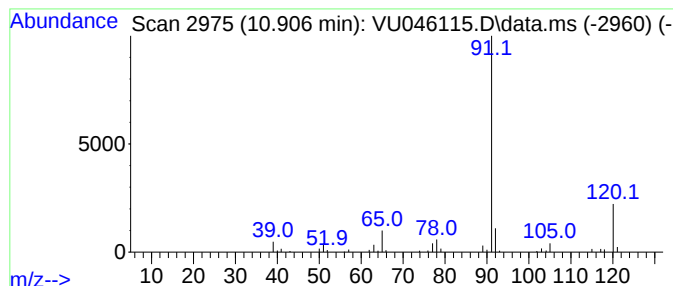
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
Quant Title : VOC Analysis

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

Peak Number 2 Benzene, propyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.906	4.49 ug/L	55983	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, propyl-	120	C9H12	000103-65-1	90
2			Benzene, propyl-	120	C9H12	000103-65-1	87
3			Benzene, propyl-	120	C9H12	000103-65-1	87
4			N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	83
5			N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	78



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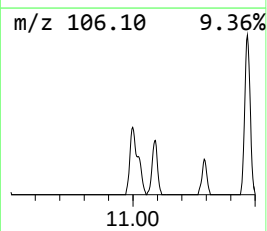
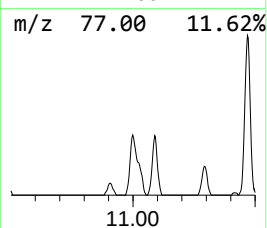
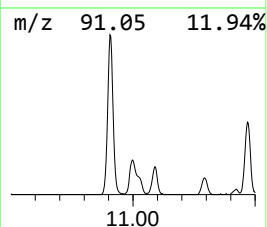
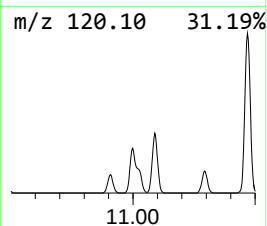
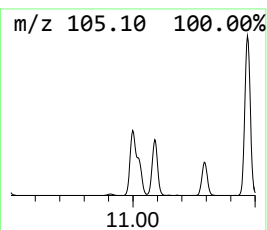
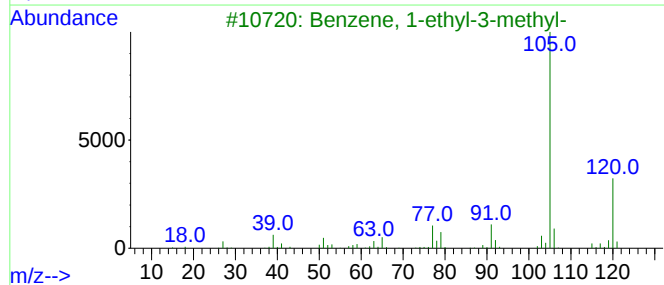
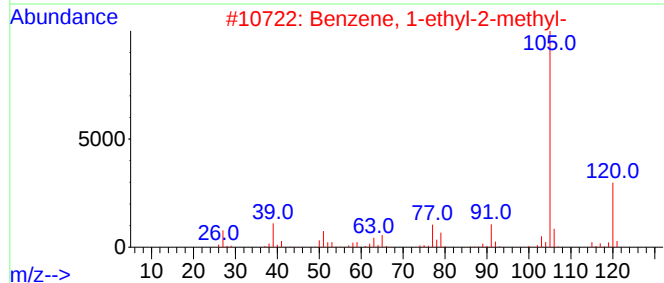
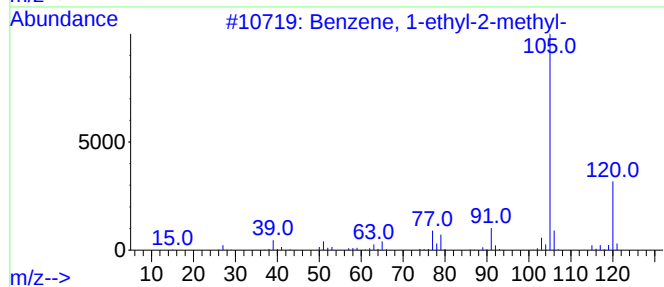
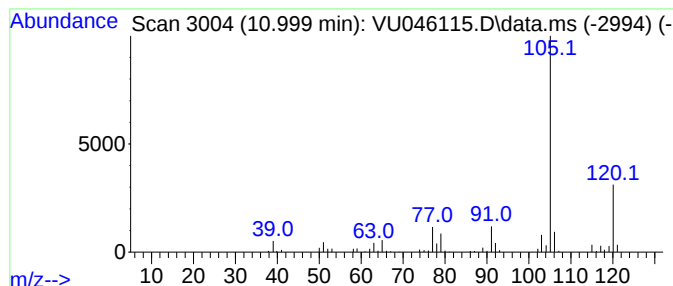
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
Quant Title : VOC Analysis

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.999	14.93 ug/L	186205	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	95
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
5			Mesitylene	120	C9H12	000108-67-8	91



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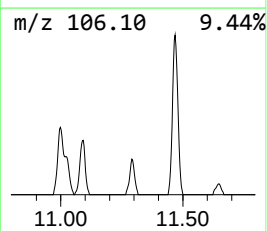
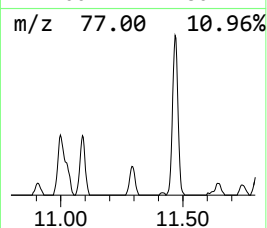
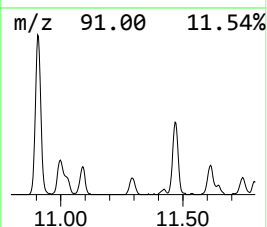
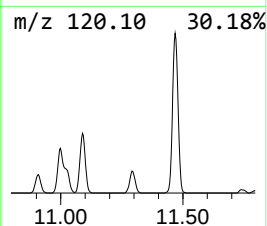
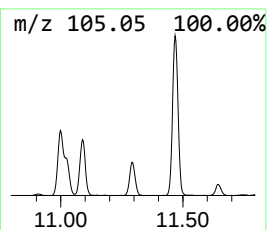
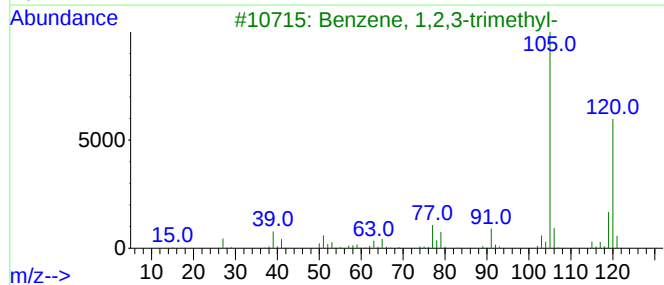
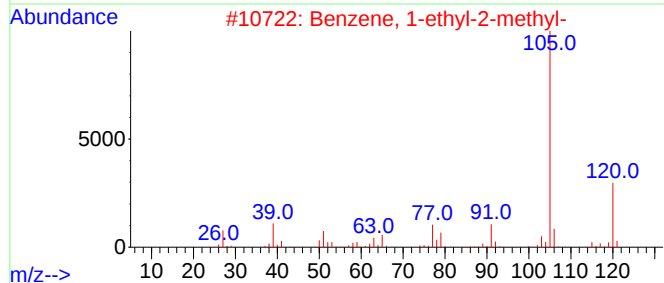
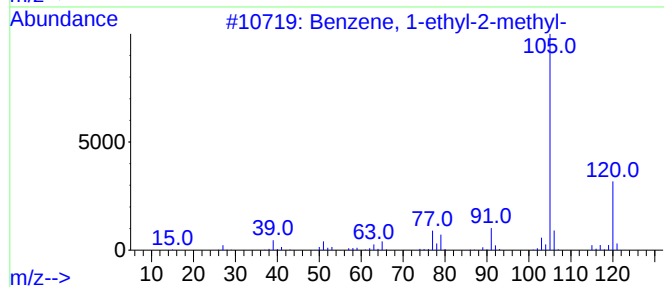
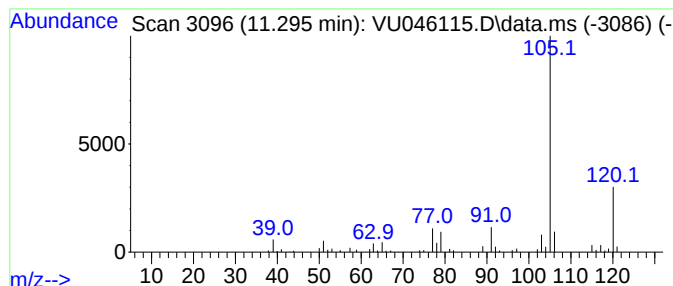
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Quant Title : VOC Analysis

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

Peak Number 4 Benzene, 1-ethyl-2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.295	5.59 ug/L	69756	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	94
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91
3			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
4			Mesitylene	120	C9H12	000108-67-8	91
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120621\
Data File : VU046115.D
Acq On : 06 Dec 2021 16:27
Operator : SY/MD
Sample : M4942-05DL 20X
Misc : 6.58g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BGKQ5DL

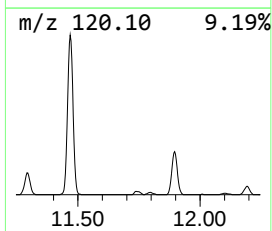
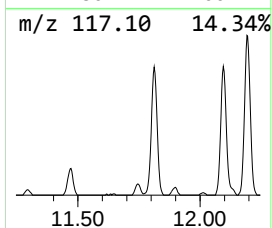
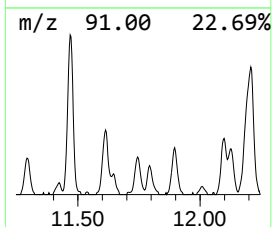
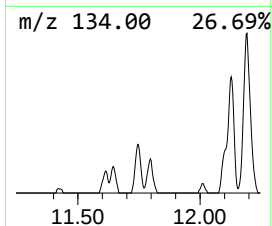
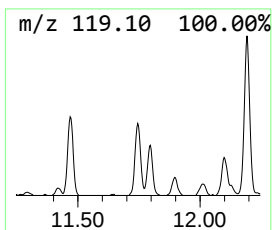
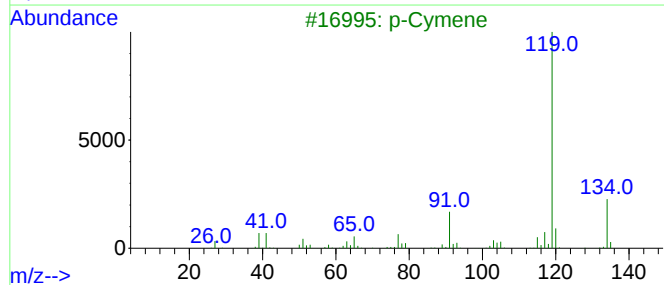
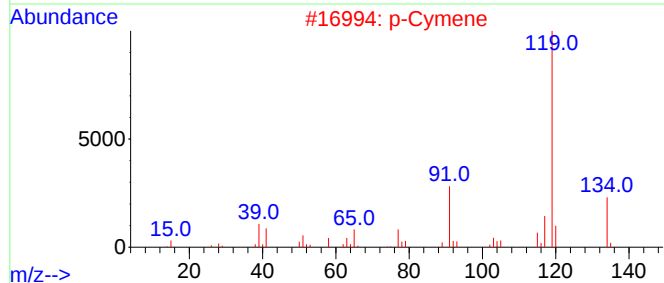
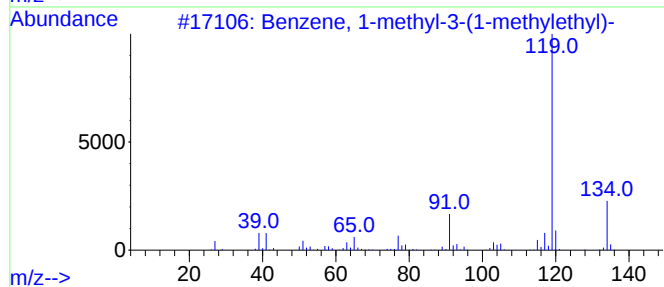
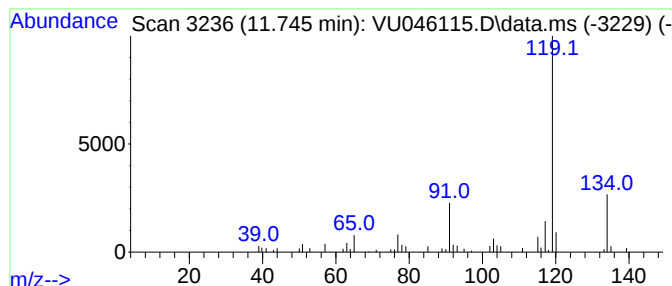
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Quant Title : VOC Analysis

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

Peak Number 5 Benzene, 1-methyl-3-(1-meth... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.745	2.65 ug/L	33039	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
2			p-Cymene	134	C10H14	000099-87-6	95
3			p-Cymene	134	C10H14	000099-87-6	94
4			o-Cymene	134	C10H14	000527-84-4	94
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120621\
Data File : VU046115.D
Acq On : 06 Dec 2021 16:27
Operator : SY/MD
Sample : M4942-05DL 20X
Misc : 6.58g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BGKQ5DL

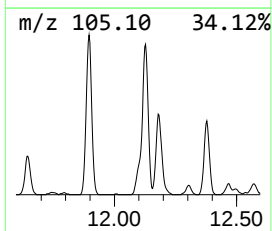
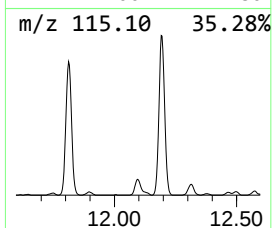
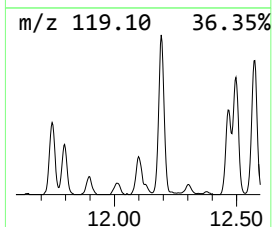
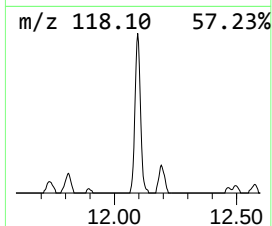
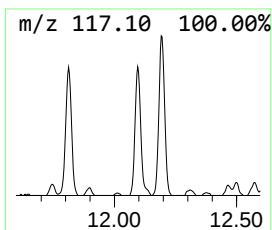
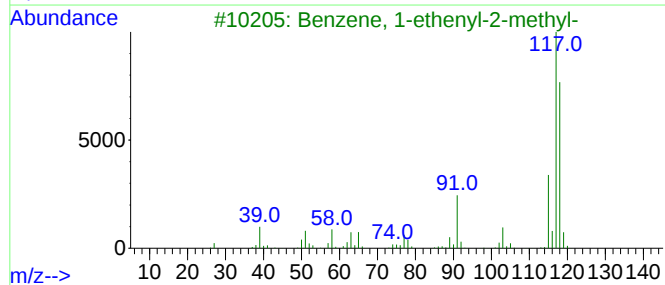
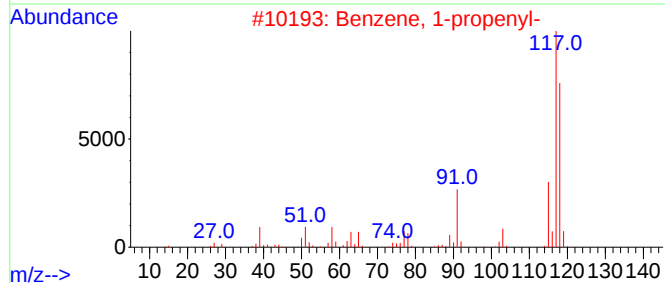
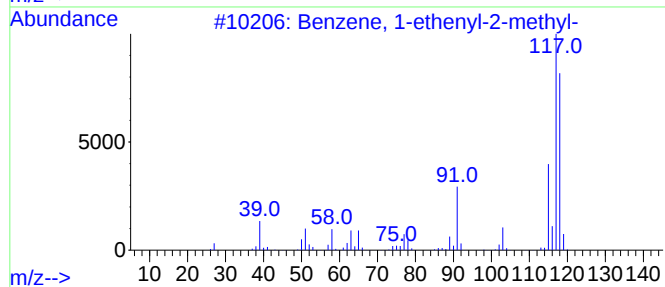
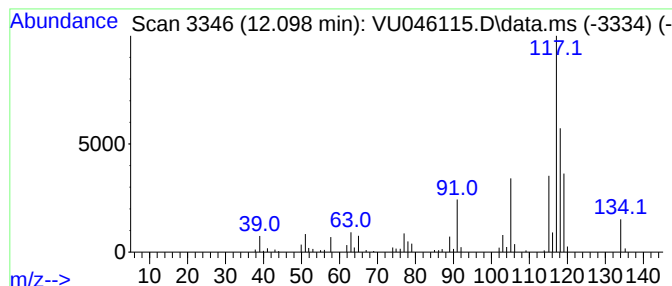
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Quant Title : VOC Analysis

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

Peak Number 7 Benzene, 1-ethenyl-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.098	6.94 ug/L	86494	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	92
2			Benzene, 1-propenyl-	118	C9H10	000637-50-3	92
3			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	91
4			Benzene, 2-propenyl-	118	C9H10	000300-57-2	64
5			Benzene, 2-propenyl-	118	C9H10	000300-57-2	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120621\
Data File : VU046115.D
Acq On : 06 Dec 2021 16:27
Operator : SY/MD
Sample : M4942-05DL 20X
Misc : 6.58g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BGKQ5DL

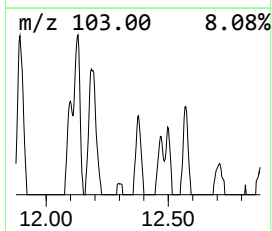
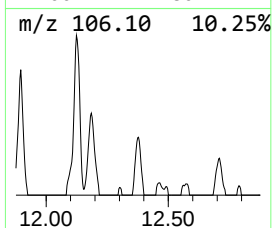
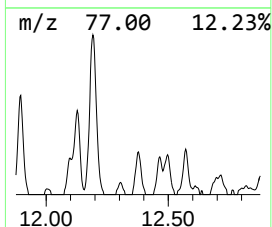
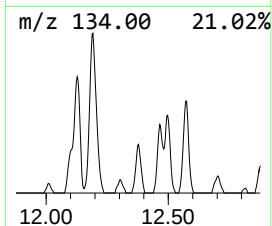
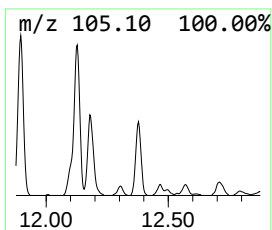
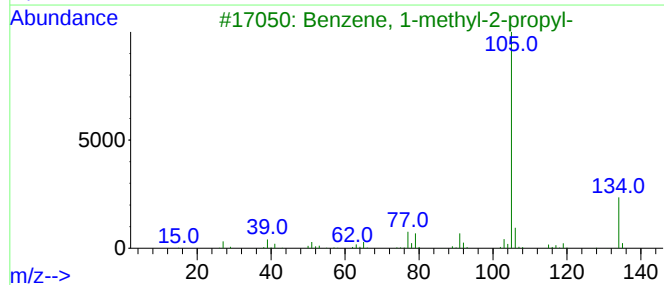
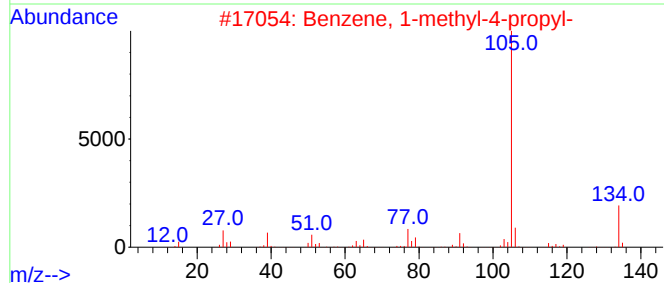
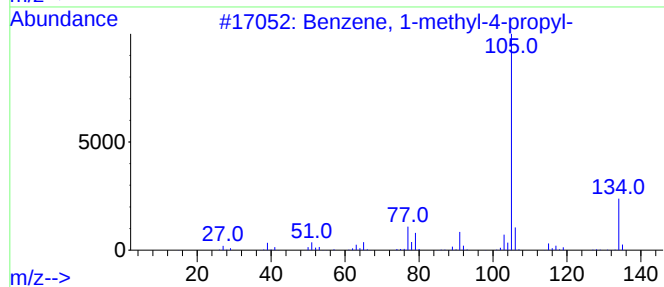
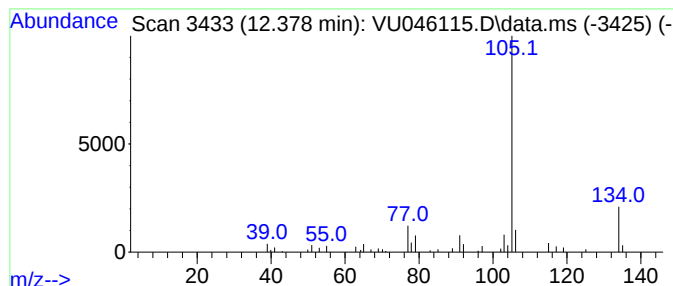
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Quant Title : VOC Analysis

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

Peak Number 8 Benzene, 1-methyl-4-propyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.378	3.48 ug/L	43367	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	95
2			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
3			Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	91
4			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	90
5			Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120621\
Data File : VU046115.D
Acq On : 06 Dec 2021 16:27
Operator : SY/MD
Sample : M4942-05DL 20X
Misc : 6.58g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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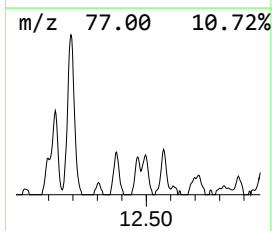
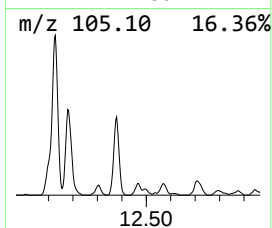
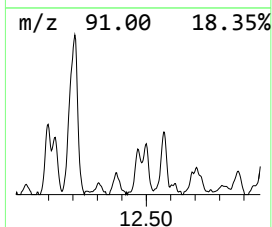
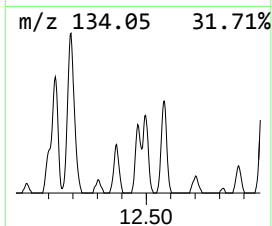
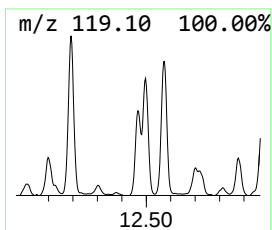
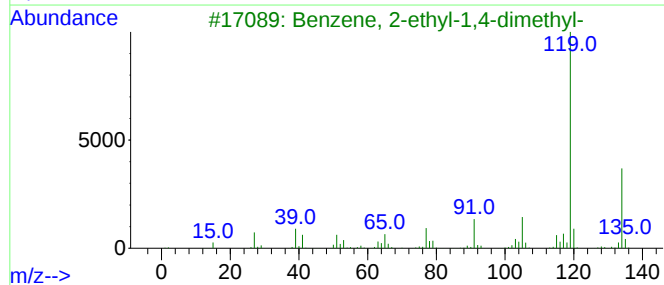
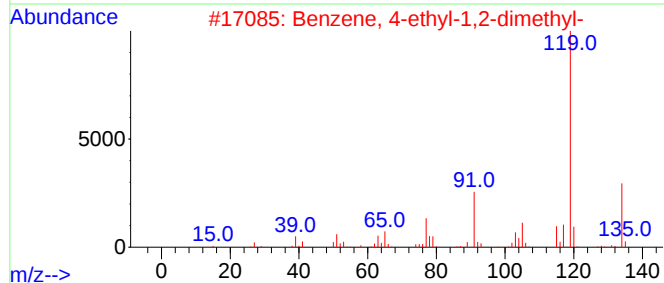
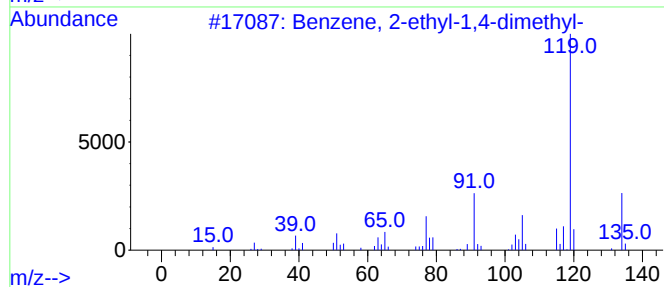
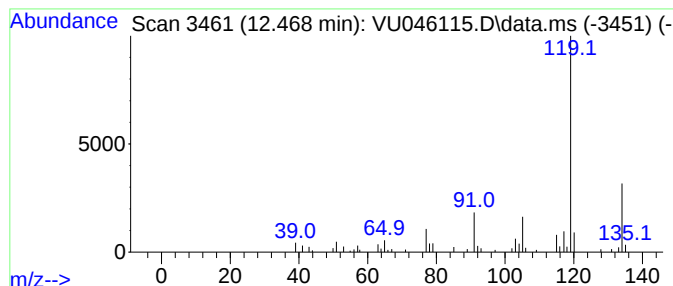
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Quant Title : VOC Analysis

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

Peak Number 9 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.468	3.60 ug/L	44941	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
3			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
4			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120621\
Data File : VU046115.D
Acq On : 06 Dec 2021 16:27
Operator : SY/MD
Sample : M4942-05DL 20X
Misc : 6.58g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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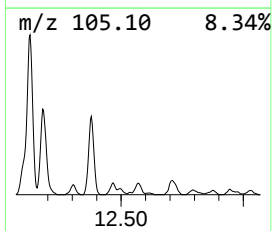
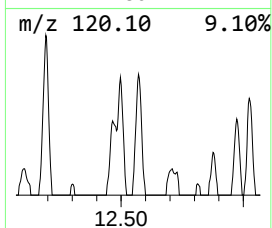
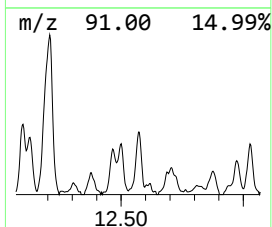
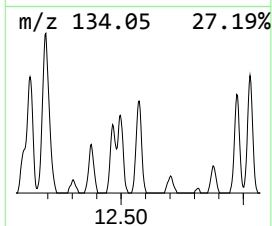
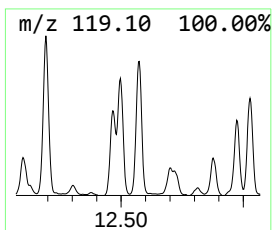
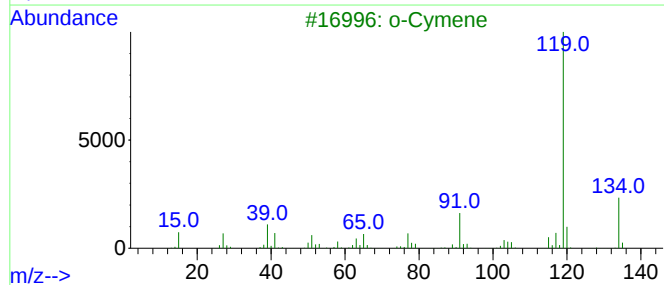
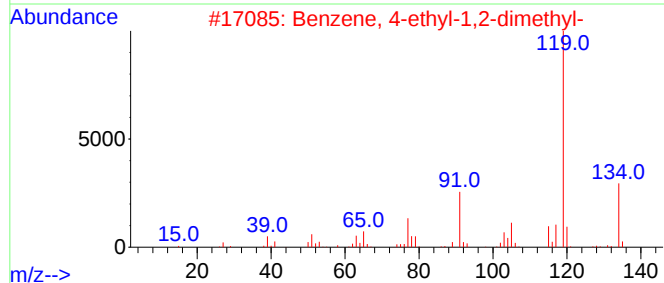
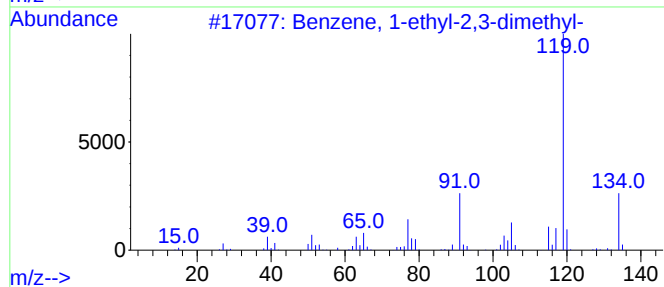
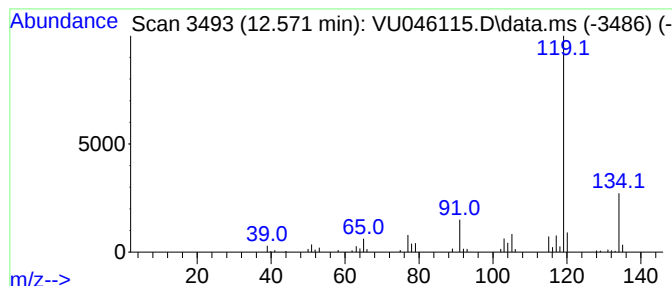
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Quant Title : VOC Analysis

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

Peak Number 10 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.571	3.97 ug/L	49526	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
2			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
3			o-Cymene	134	C10H14	000527-84-4	95
4			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
5			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120621\
Data File : VU046115.D
Acq On : 06 Dec 2021 16:27
Operator : SY/MD
Sample : M4942-05DL 20X
Misc : 6.58g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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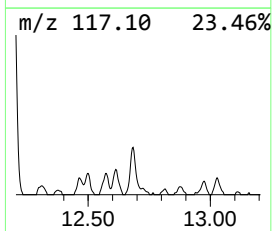
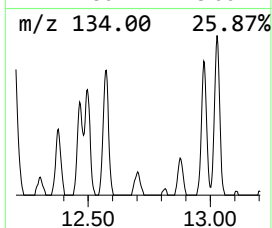
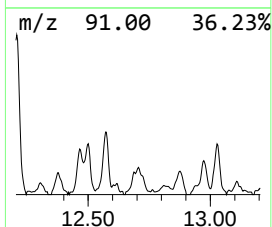
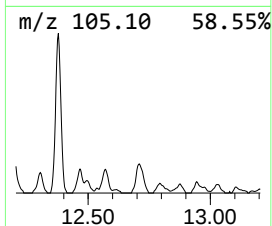
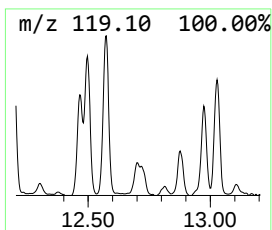
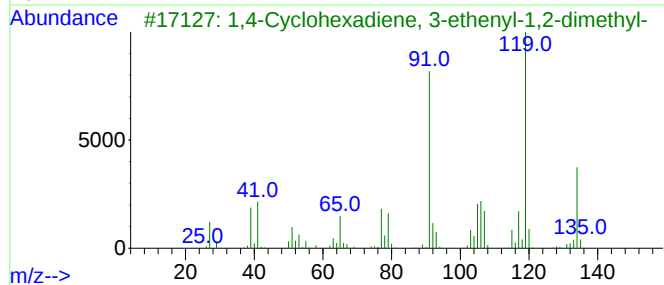
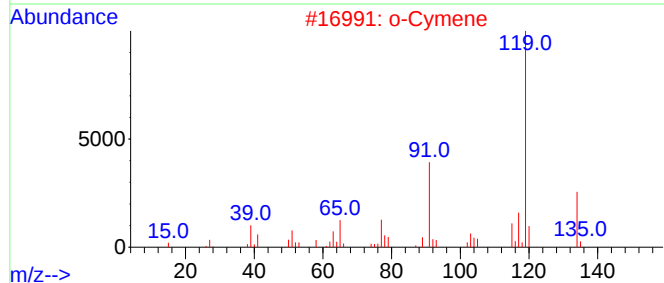
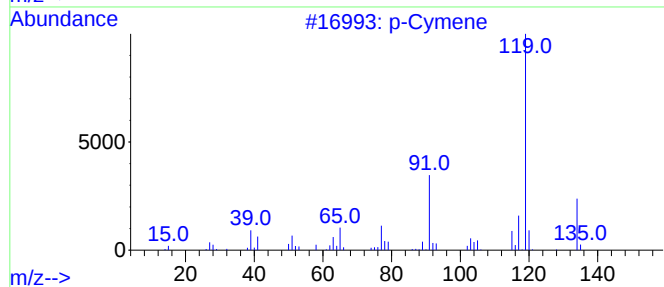
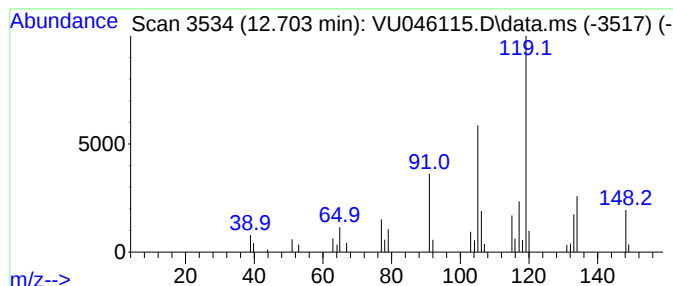
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Quant Title : VOC Analysis

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

Peak Number 11 p-Cymene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.703	3.60 ug/L	44913	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			p-Cymene	134	C10H14	000099-87-6	70
2			o-Cymene	134	C10H14	000527-84-4	64
3			1,4-Cyclohexadiene, 3-ethenyl-1,...	134	C10H14	062338-57-2	64
4			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	58
5			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	58



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120621\
Data File : VU046115.D
Acq On : 06 Dec 2021 16:27
Operator : SY/MD
Sample : M4942-05DL 20X
Misc : 6.58g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BGKQ5DL

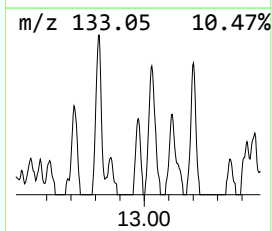
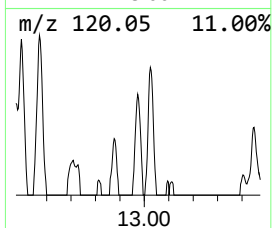
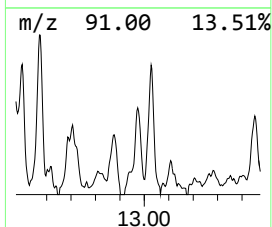
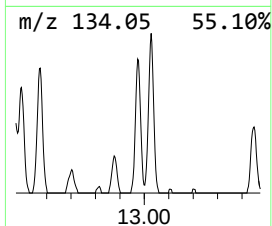
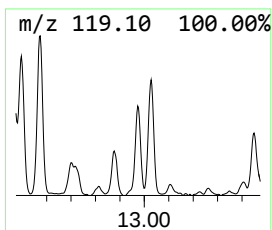
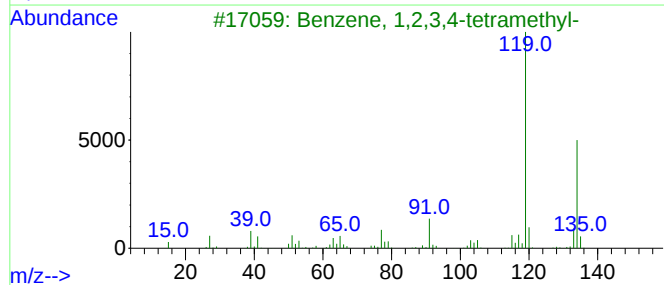
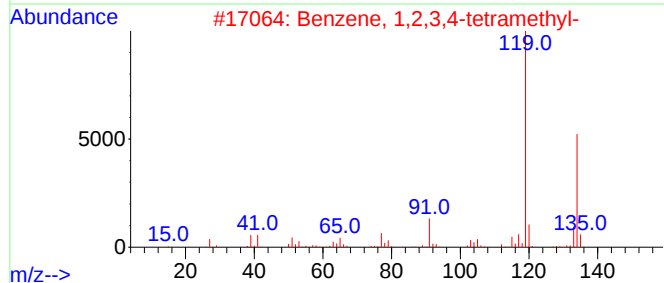
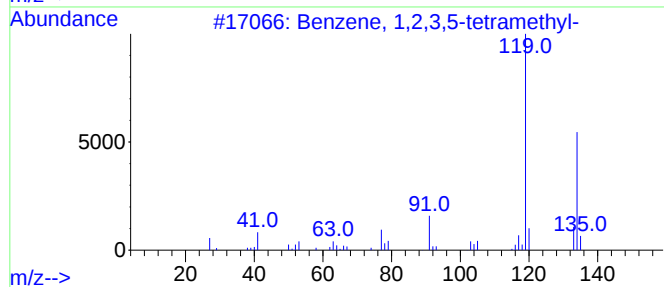
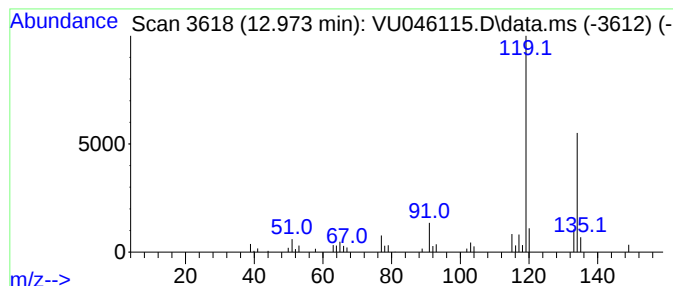
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Quant Title : VOC Analysis

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

Peak Number 12 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.973	2.73 ug/L	33995	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	97
2			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
4			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
5			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120621\
Data File : VU046115.D
Acq On : 06 Dec 2021 16:27
Operator : SY/MD
Sample : M4942-05DL 20X
Misc : 6.58g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BGKQ5DL

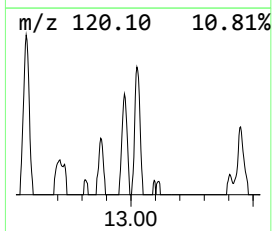
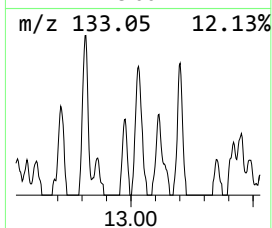
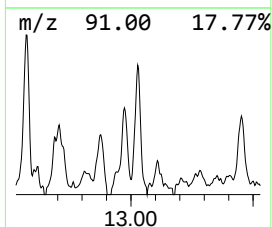
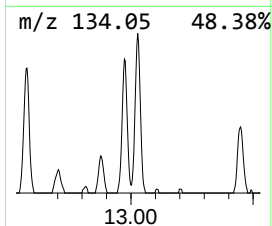
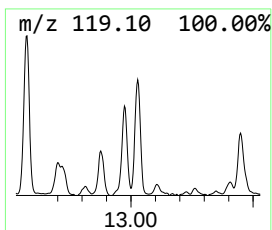
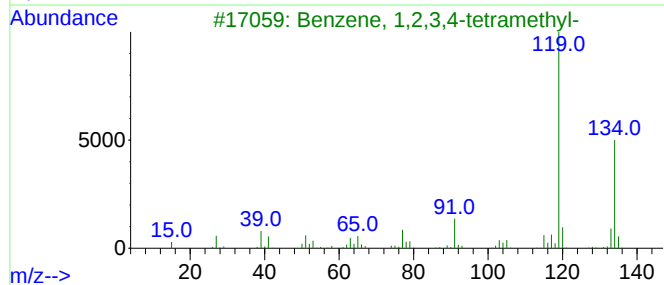
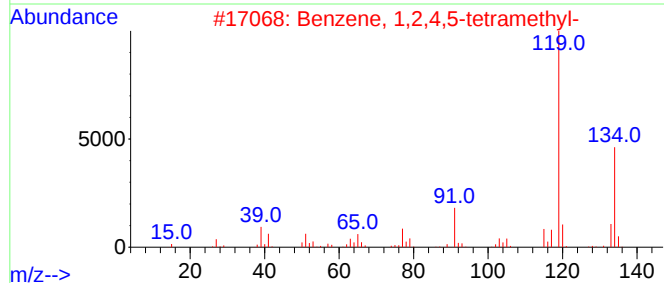
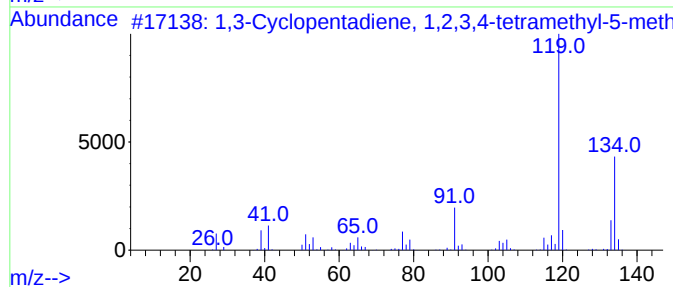
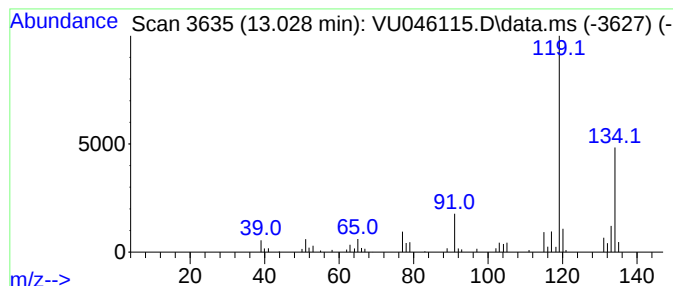
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Quant Title : VOC Analysis

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

Peak Number 13 1,3-Cyclopentadiene, 1,2,3,... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.028	4.60 ug/L	57314	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	97
2			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
3			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
4			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
5			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120621\
Data File : VU046115.D
Acq On : 06 Dec 2021 16:27
Operator : SY/MD
Sample : M4942-05DL 20X
Misc : 6.58g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BGKQ5DL

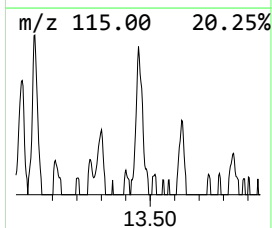
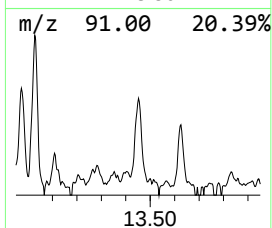
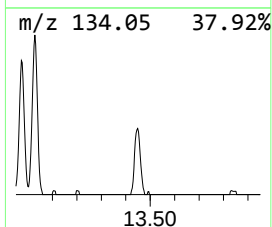
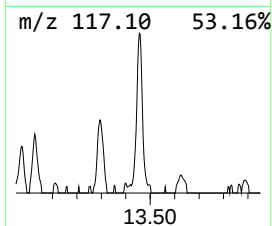
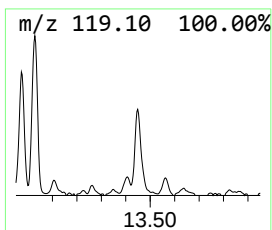
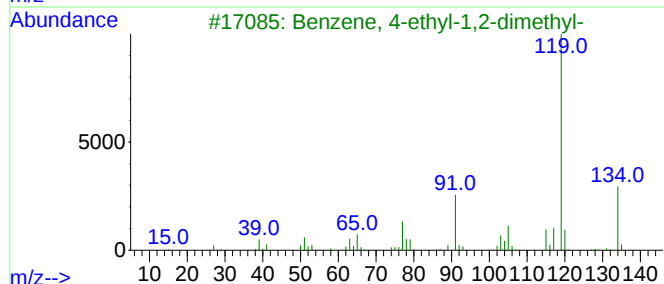
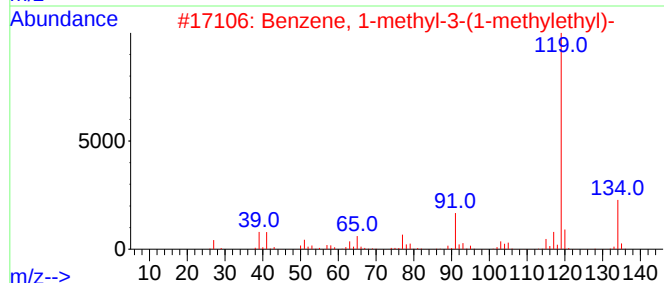
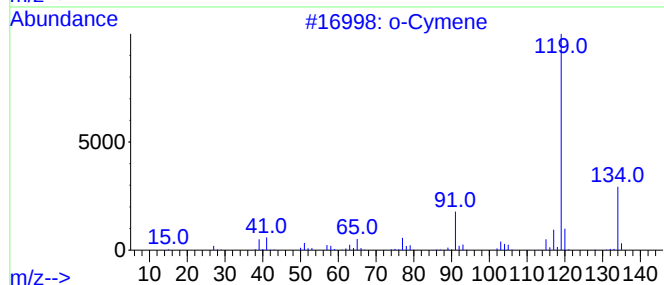
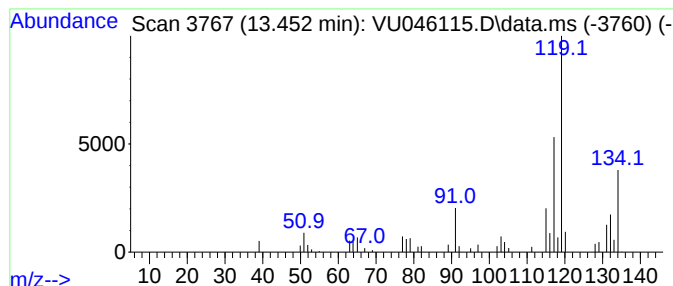
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Quant Title : VOC Analysis

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

Peak Number 14 unknown-01 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.452	3.16 ug/L	39408	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	49
2			Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	49
3			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	49
4			Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	49
5			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	49



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120621\
Data File : VU046115.D
Acq On : 06 Dec 2021 16:27
Operator : SY/MD
Sample : M4942-05DL 20X
Misc : 6.58g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BGKQ5DL

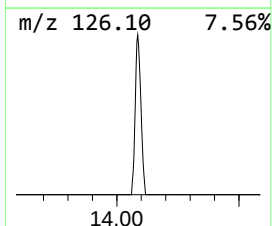
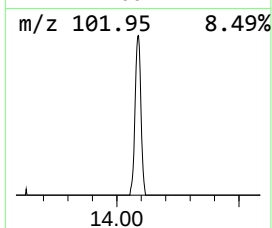
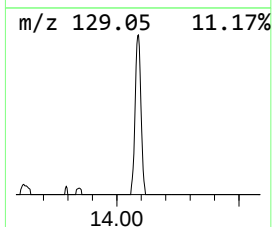
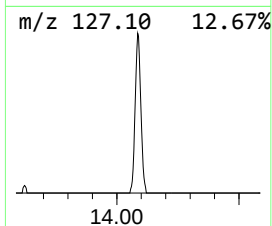
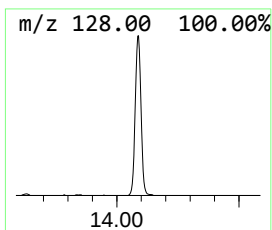
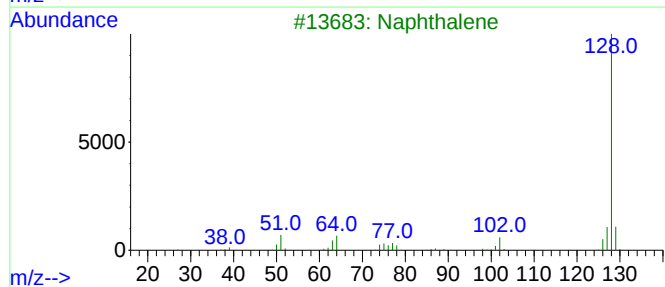
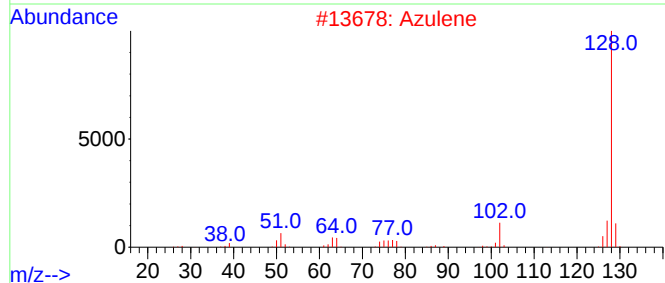
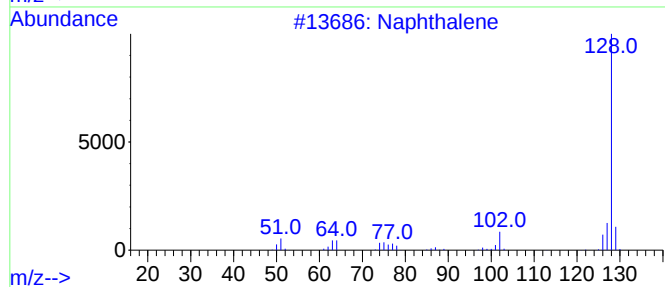
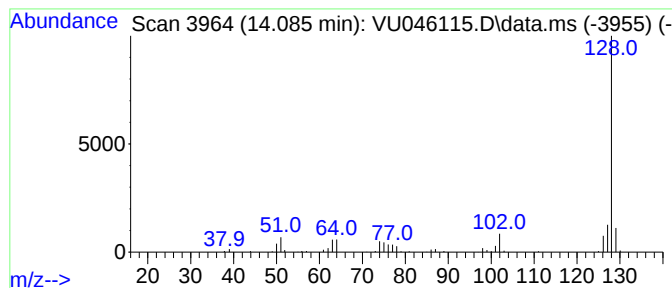
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Quant Title : VOC Analysis

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

Peak Number 15 Azulene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.086	8.44 ug/L	105297	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene	128	C10H8	000091-20-3	95
2			Azulene	128	C10H8	000275-51-4	94
3			Naphthalene	128	C10H8	000091-20-3	94
4			1H-Indene, 1-methylene-	128	C10H8	002471-84-3	94
5			Azulene	128	C10H8	000275-51-4	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120621\
 Data File : VU046115.D
 Acq On : 06 Dec 2021 16:27
 Operator : SY/MD
 Sample : M4942-05DL 20X
 Misc : 6.58g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 BGKQ5DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.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, propyl-	10.906	4.5	ug/L	55983	3	11.812	623483	50.0
Benzene, 1-ethy...	10.999	14.9	ug/L	186205	3	11.812	623483	50.0
Benzene, 1-ethy...	11.295	5.6	ug/L	69756	3	11.812	623483	50.0
Benzene, 1-meth...	11.745	2.6	ug/L	33039	3	11.812	623483	50.0
Benzene, 1-ethe...	12.098	6.9	ug/L	86494	3	11.812	623483	50.0
Benzene, 1-meth...	12.378	3.5	ug/L	43367	3	11.812	623483	50.0
Benzene, 2-ethy...	12.468	3.6	ug/L	44941	3	11.812	623483	50.0
Benzene, 1-ethy...	12.571	4.0	ug/L	49526	3	11.812	623483	50.0
p-Cymene	12.703	3.6	ug/L	44913	3	11.812	623483	50.0
Benzene, 1,2,3,...	12.973	2.7	ug/L	33995	3	11.812	623483	50.0
1,3-Cyclopentad...	13.028	4.6	ug/L	57314	3	11.812	623483	50.0
unknown-01	13.452	3.2	ug/L	39408	3	11.812	623483	50.0
Azulene	14.086	8.4	ug/L	105297	3	11.812	623483	50.0