

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120821\
 Data File : VU046193.D
 Acq On : 08 Dec 2021 22:13
 Operator : SY/MD
 Sample : M4983-04
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 EW5N3

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: VU046193.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.607	74	83	101	rVB2	185372	223008	2.04%	0.755%
2	1.938	171	186	204	rBV	789055	1110184	10.15%	3.758%
3	2.478	349	354	364	rVB6	937	1429	0.01%	0.005%
4	2.578	371	385	404	rBV3	205233	411266	3.76%	1.392%
5	3.051	524	532	535	rBV5	856	1091	0.01%	0.004%
6	3.186	564	574	586	rVB2	2065	4540	0.04%	0.015%
7	3.308	606	612	615	rBV2	410	519	0.00%	0.002%
8	3.359	615	628	643	rBV2	28152	54897	0.50%	0.186%
9	3.485	663	667	672	rVB	466	458	0.00%	0.002%
10	3.774	749	757	759	rBV2	389	460	0.00%	0.002%
11	3.877	767	789	815	rVV	1082622	2578154	23.57%	8.726%
12	4.456	963	969	973	rVB	590	697	0.01%	0.002%
13	4.504	980	984	987	rVB3	534	463	0.00%	0.002%
14	4.671	1011	1036	1066	rBV	2355716	5423859	49.58%	18.358%
15	5.067	1145	1159	1179	rVB	130663	290528	2.66%	0.983%
16	5.321	1217	1238	1258	rBV	4845253	10940511	100.00%	37.030%
17	5.729	1345	1365	1380	rBV2	262777	713189	6.52%	2.414%
18	5.906	1409	1420	1428	rVB3	2938	5918	0.05%	0.020%
19	5.980	1439	1443	1445	rBV3	589	480	0.00%	0.002%
20	6.253	1511	1528	1545	rBV	240640	461310	4.22%	1.561%
21	6.459	1589	1592	1597	rVB3	607	598	0.01%	0.002%
22	6.546	1605	1619	1636	rBV	44296	81941	0.75%	0.277%
23	6.620	1636	1642	1648	rBV2	456	587	0.01%	0.002%
24	6.694	1651	1665	1680	rBV	164636	319729	2.92%	1.082%
25	7.038	1769	1772	1778	rVB3	469	472	0.00%	0.002%
26	7.137	1801	1803	1809	rBV2	453	445	0.00%	0.002%
27	7.568	1923	1937	1953	rBV	98441	173395	1.58%	0.587%
28	7.726	1981	1986	1991	rBV2	517	615	0.01%	0.002%
29	7.899	2026	2040	2052	rVV	335275	580297	5.30%	1.964%
30	7.970	2052	2062	2087	rVB	293564	508900	4.65%	1.722%
31	8.179	2114	2127	2143	rBV	70507	118667	1.08%	0.402%
32	8.404	2190	2197	2205	rVB3	962	1425	0.01%	0.005%
33	8.475	2210	2219	2229	rVV2	7430	13464	0.12%	0.046%
34	8.555	2232	2244	2257	rVV	115877	196910	1.80%	0.666%
35	8.636	2257	2269	2302	rVB	245401	446038	4.08%	1.510%
36	8.870	2331	2342	2353	rBV3	1589	3099	0.03%	0.010%

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

37	8.925	2353	2359	2366	rVB3	868	1140	0.01%	0.004%
38	9.041	2390	2395	2398	rBV2	712	636	0.01%	0.002%
39	9.186	2437	2440	2444	rVB	641	488	0.00%	0.002%
40	9.417	2499	2512	2536	rVV	350106	586844	5.36%	1.986%
41	9.571	2549	2560	2576	rVB	72395	118083	1.08%	0.400%
42	9.694	2583	2598	2616	rBV	192819	330313	3.02%	1.118%
43	10.102	2713	2725	2743	rVB	113915	188683	1.72%	0.639%
44	10.275	2771	2779	2788	rVB4	3120	5180	0.05%	0.018%
45	10.343	2793	2800	2801	rBV2	784	606	0.01%	0.002%
46	10.484	2835	2844	2860	rVB	13595	20930	0.19%	0.071%
47	10.571	2865	2871	2872	rBV2	1360	1229	0.01%	0.004%
48	10.639	2887	2892	2899	rBV4	1604	2361	0.02%	0.008%
49	10.700	2907	2911	2917	rVB4	1574	1896	0.02%	0.006%
50	10.758	2917	2929	2944	rBV	272235	428391	3.92%	1.450%
51	10.902	2958	2974	2987	rBV2	37452	68570	0.63%	0.232%
52	10.999	2992	3004	3022	rBV	106347	240690	2.20%	0.815%
53	11.089	3022	3032	3045	rVB	64472	98244	0.90%	0.333%
54	11.234	3069	3077	3085	rVV3	3456	5160	0.05%	0.017%
55	11.292	3085	3095	3111	rVV	88266	138764	1.27%	0.470%
56	11.369	3114	3119	3126	rVB2	613	824	0.01%	0.003%
57	11.423	3126	3136	3137	rBV	1758	2134	0.02%	0.007%
58	11.468	3139	3150	3167	rVB	294765	452686	4.14%	1.532%
59	11.610	3187	3194	3198	rBV2	3746	5203	0.05%	0.018%
60	11.645	3199	3205	3215	rVB	10438	15469	0.14%	0.052%
61	11.697	3215	3221	3226	rBV2	781	1029	0.01%	0.003%
62	11.745	3226	3236	3244	rBV2	16369	23349	0.21%	0.079%
63	11.812	3244	3257	3271	rVB	419549	666062	6.09%	2.254%
64	11.893	3271	3282	3301	rVB	133562	207585	1.90%	0.703%
65	11.973	3301	3307	3313	rBV5	1073	1559	0.01%	0.005%
66	12.012	3314	3319	3329	rVB4	2553	3432	0.03%	0.012%
67	12.095	3329	3345	3352	rBV3	57881	100118	0.92%	0.339%
68	12.192	3363	3375	3392	rVB	415288	697186	6.37%	2.360%
69	12.304	3401	3410	3423	rBV8	3832	6245	0.06%	0.021%
70	12.375	3423	3432	3443	rVB2	14279	20879	0.19%	0.071%
71	12.481	3450	3465	3484	rBV3	51525	130361	1.19%	0.441%
72	12.571	3485	3493	3501	rVV	24993	37403	0.34%	0.127%
73	12.613	3501	3506	3513	rVB4	4396	5341	0.05%	0.018%
74	12.690	3519	3530	3546	rBV4	13215	34227	0.31%	0.116%
75	12.771	3548	3555	3563	rBV2	6054	8911	0.08%	0.030%
76	12.877	3577	3588	3597	rBV2	9002	14413	0.13%	0.049%

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 Title : VOC Analysis

77	12.973	3603	3618	3626	rBV2	11397	17857	0.16%	0.060%
78	13.025	3626	3634	3646	rVB	20471	29744	0.27%	0.101%
79	13.108	3656	3660	3667	rBV5	1599	1989	0.02%	0.007%
80	13.253	3700	3705	3711	rVV4	1922	2386	0.02%	0.008%
81	13.295	3711	3718	3727	rVV4	3843	5624	0.05%	0.019%
82	13.333	3727	3730	3733	rVV2	601	433	0.00%	0.001%
83	13.407	3746	3753	3756	rBV4	2000	2582	0.02%	0.009%
84	13.449	3757	3766	3778	rVV3	25907	43123	0.39%	0.146%
85	13.523	3782	3789	3796	rVV7	1955	3012	0.03%	0.010%
86	13.558	3796	3800	3809	rVV2	1134	1689	0.02%	0.006%
87	13.623	3813	3820	3833	rVB3	7608	11168	0.10%	0.038%
88	13.783	3864	3870	3876	rBV2	991	981	0.01%	0.003%
89	13.838	3879	3887	3892	rBV6	2037	2462	0.02%	0.008%
90	14.086	3952	3964	3982	rVB	35423	56023	0.51%	0.190%
91	14.253	4011	4016	4018	rBV2	513	429	0.00%	0.001%
92	14.497	4086	4092	4102	rVB4	450	633	0.01%	0.002%
93	14.578	4107	4117	4128	rVB3	2962	5384	0.05%	0.018%
94	14.668	4139	4145	4147	rBV3	525	527	0.00%	0.002%
95	15.018	4252	4254	4258	rVB2	612	422	0.00%	0.001%
96	15.224	4309	4318	4328	rVB3	5874	8781	0.08%	0.030%
97	15.307	4340	4344	4348	rBV2	472	499	0.00%	0.002%
98	15.410	4368	4376	4387	rBV2	4301	6069	0.06%	0.021%
99	15.812	4498	4501	4509	rBV3	440	564	0.01%	0.002%
100	16.442	4695	4697	4702	rVB5	924	549	0.01%	0.002%

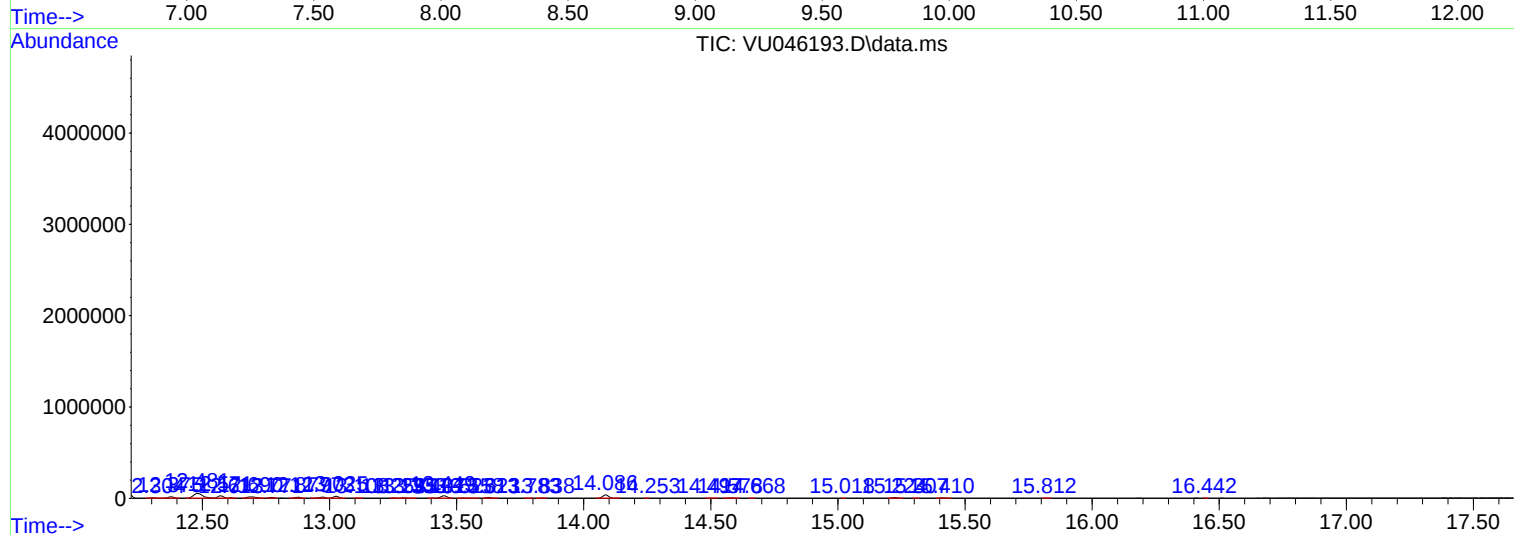
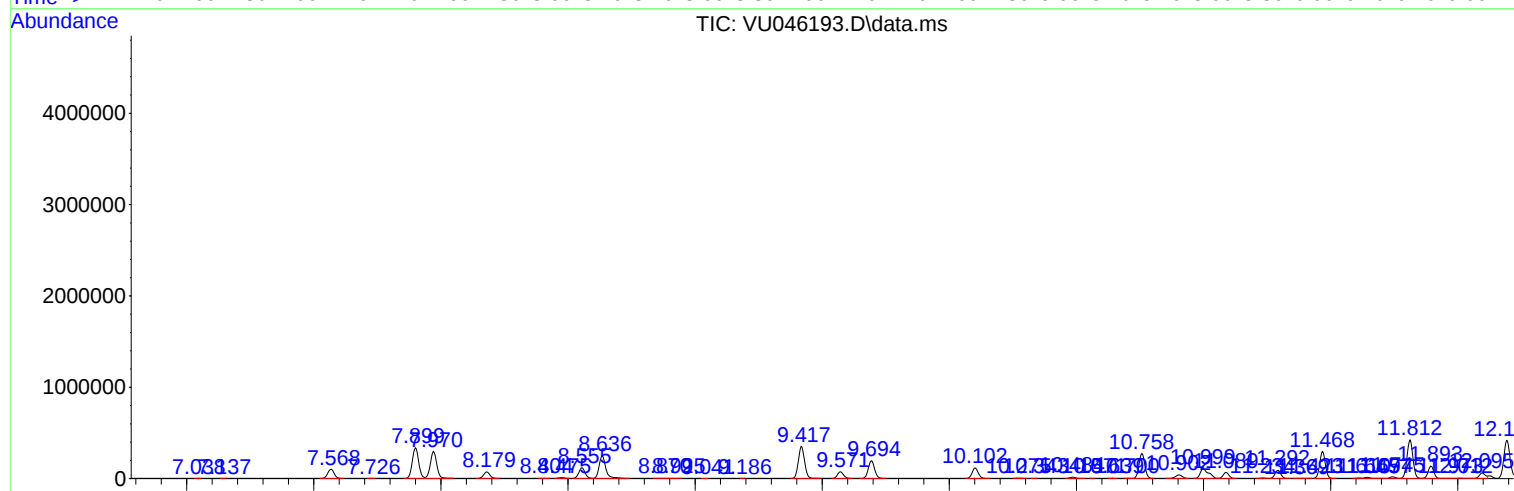
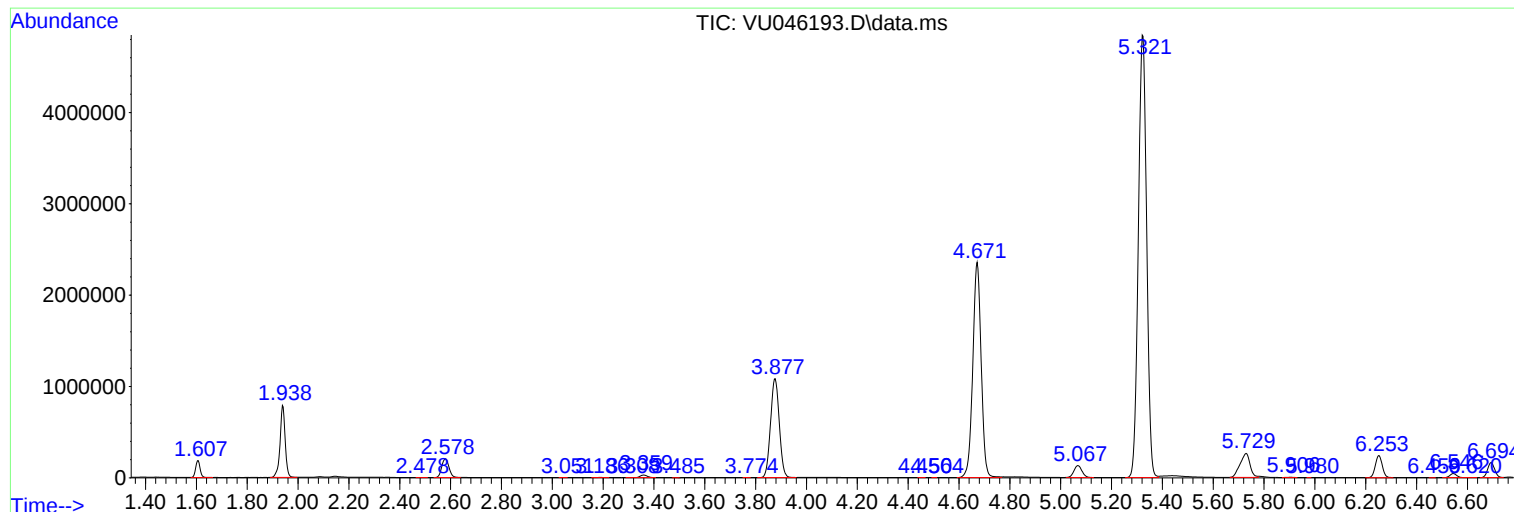
Sum of corrected areas: 29545097

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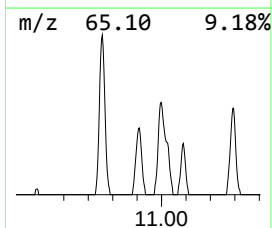
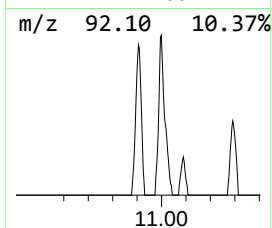
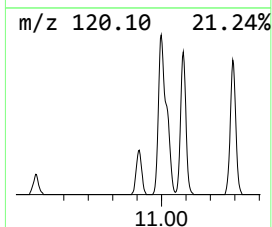
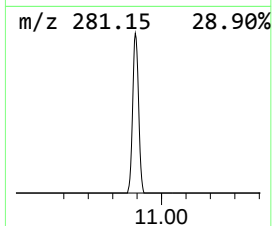
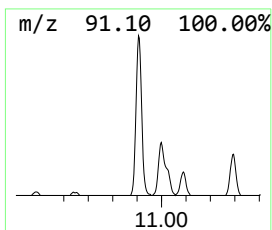
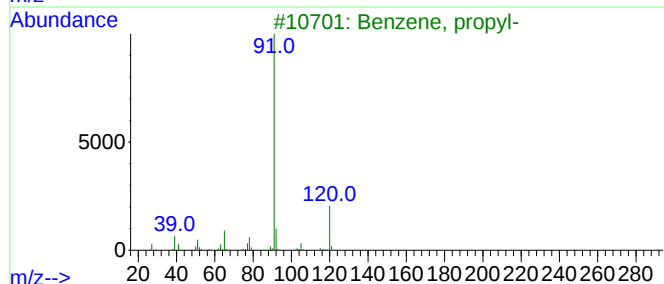
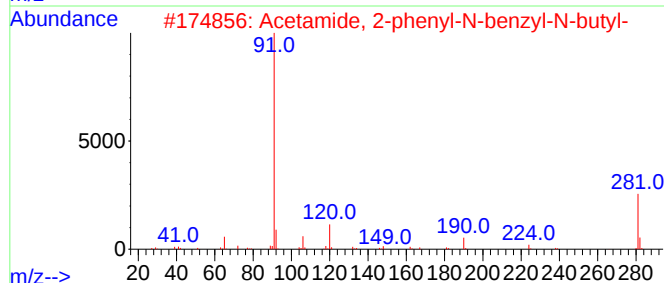
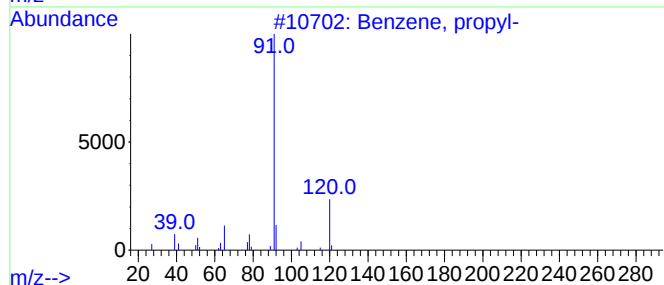
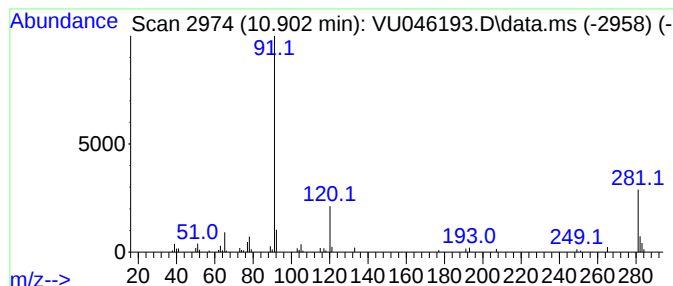
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
Quant Title : VOC Analysis

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

Peak Number 2 Benzene, propyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.903	5.15 ug/L	68570	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			Acetamide, 2-phenyl-N-benzyl-N-b...	281	C19H23NO	1010446-79-6	64
3			Benzene, propyl-	120	C9H12	000103-65-1	60
4			Benzene, propyl-	120	C9H12	000103-65-1	60
5			Benzene, propyl-	120	C9H12	000103-65-1	53



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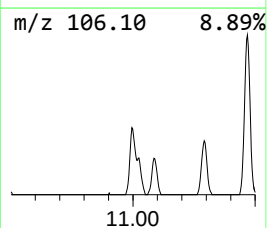
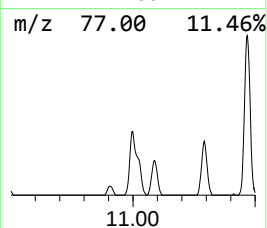
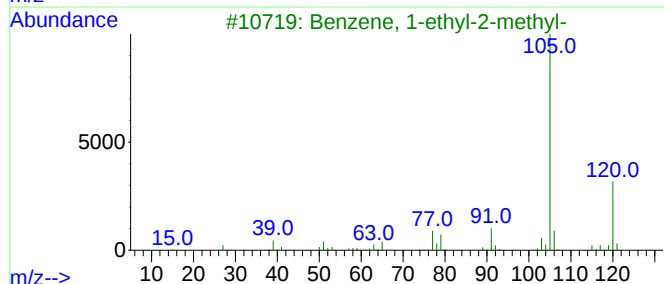
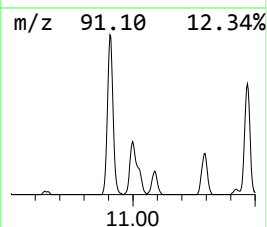
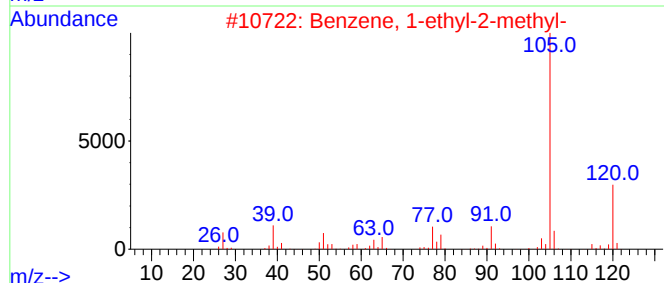
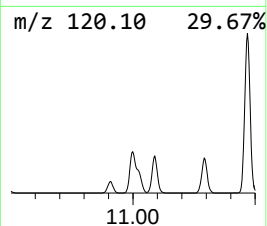
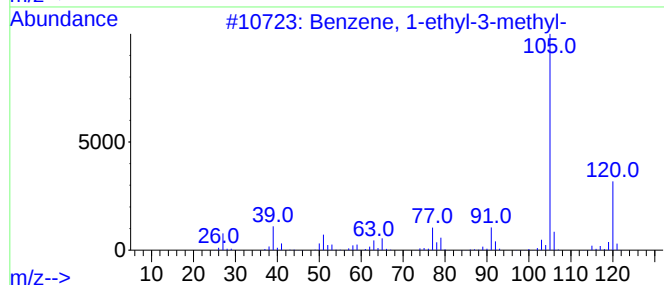
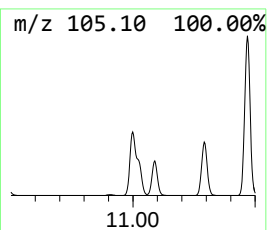
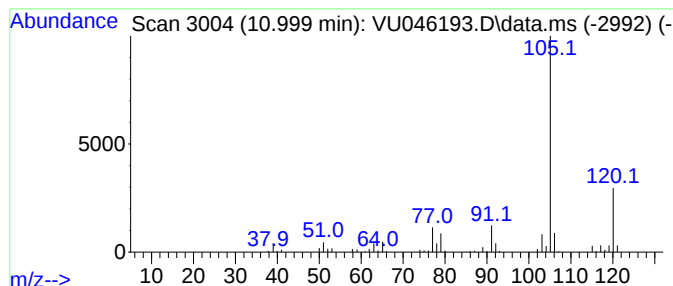
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM112921WMA.M
Quant Title : VOC Analysis

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	18.07 ug/L	240690	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	95
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



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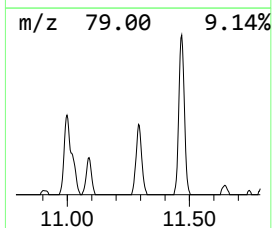
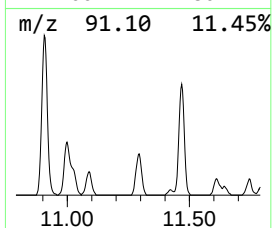
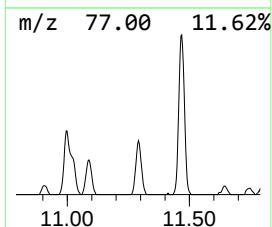
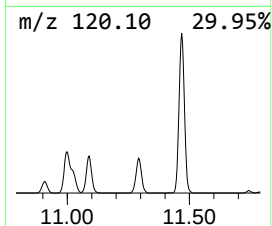
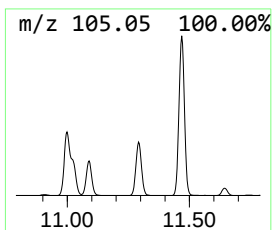
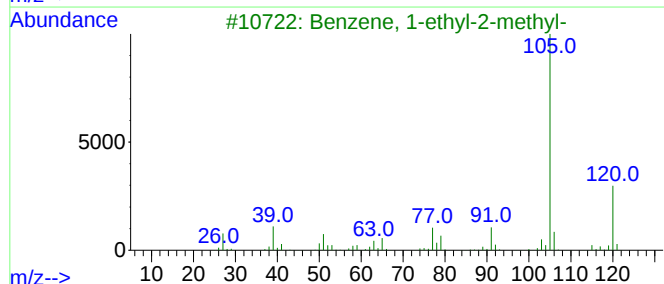
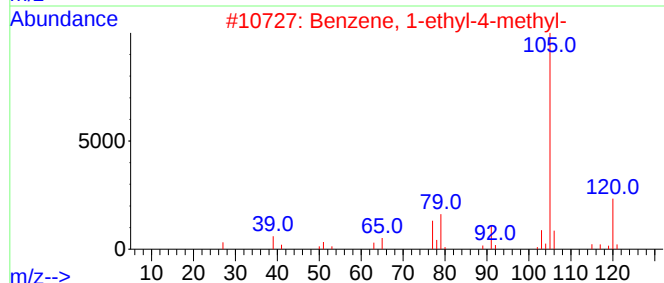
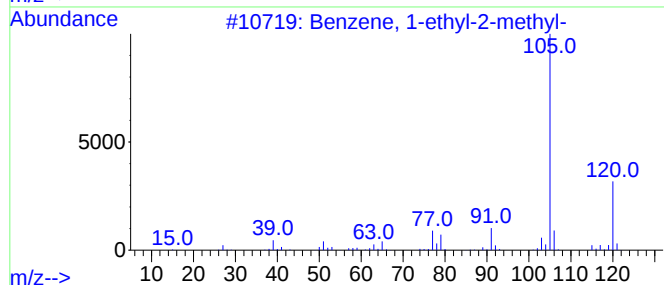
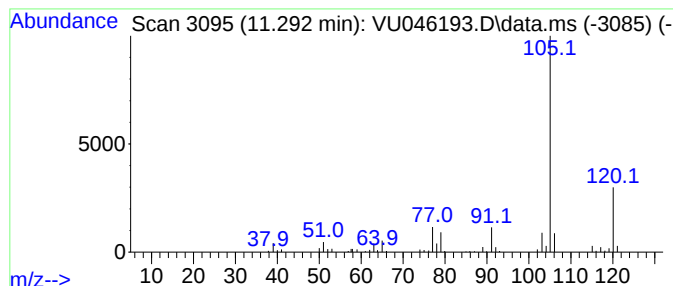
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Quant Title : VOC Analysis

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.291	10.42 ug/L	138764	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-4-methyl-	120	C9H12	000622-96-8	91
3			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120821\
Data File : VU046193.D
Acq On : 08 Dec 2021 22:13
Operator : SY/MD
Sample : M4983-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5N3

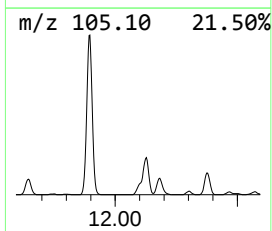
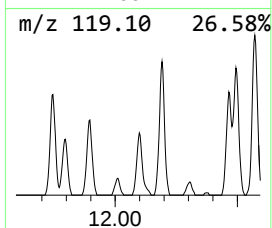
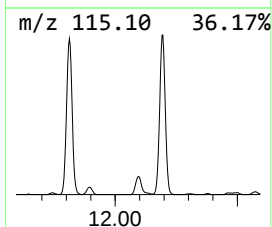
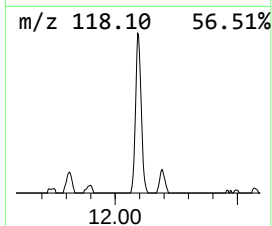
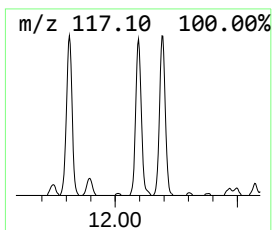
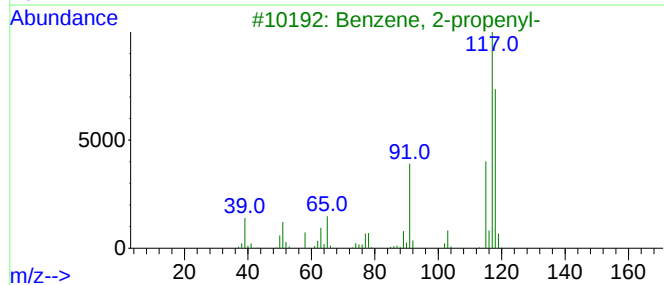
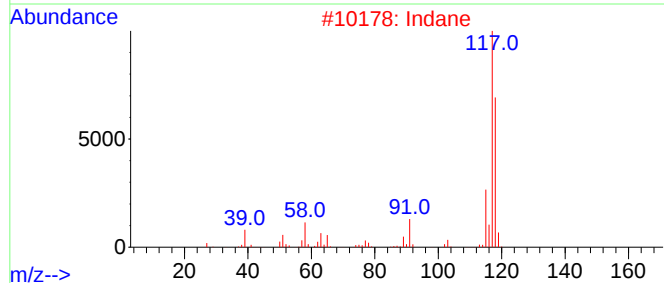
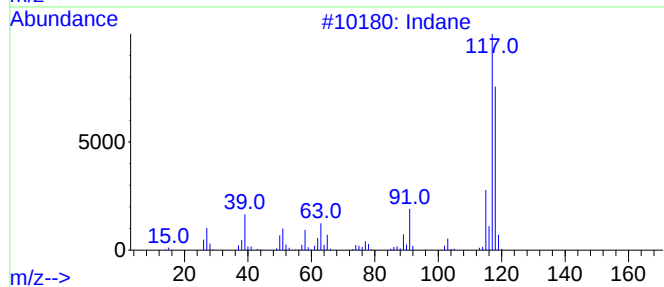
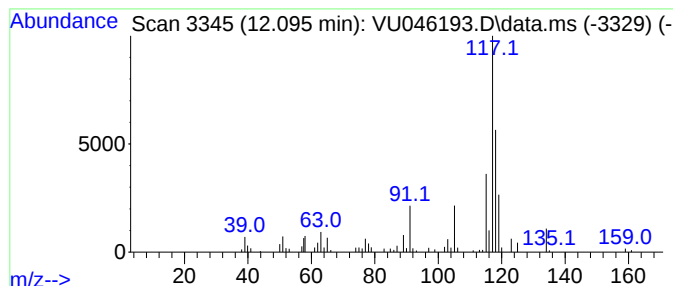
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Quant Title : VOC Analysis

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

Peak Number 6 Indane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.095	7.52 ug/L	100118	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	92
2			Indane	118	C9H10	000496-11-7	64
3			Benzene, 2-propenyl-	118	C9H10	000300-57-2	60
4			Benzene, 2-propenyl-	118	C9H10	000300-57-2	60
5			Benzene, cyclopropyl-	118	C9H10	000873-49-4	60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120821\
Data File : VU046193.D
Acq On : 08 Dec 2021 22:13
Operator : SY/MD
Sample : M4983-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5N3

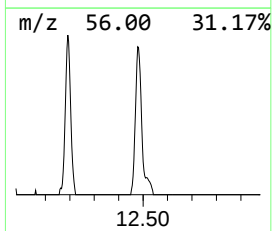
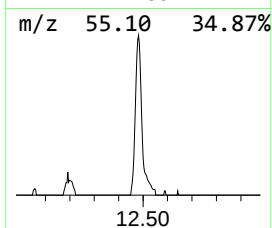
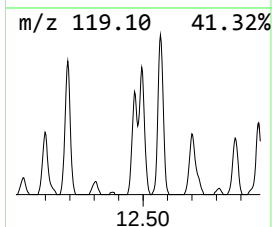
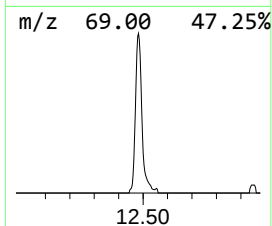
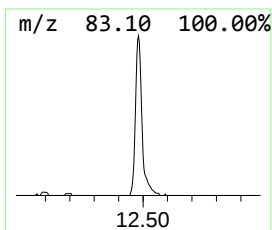
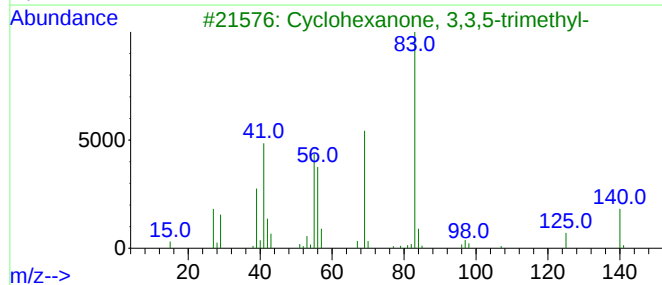
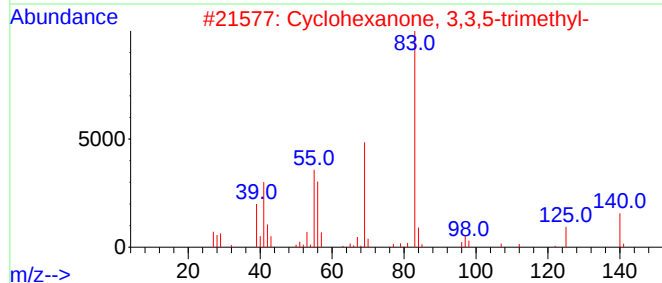
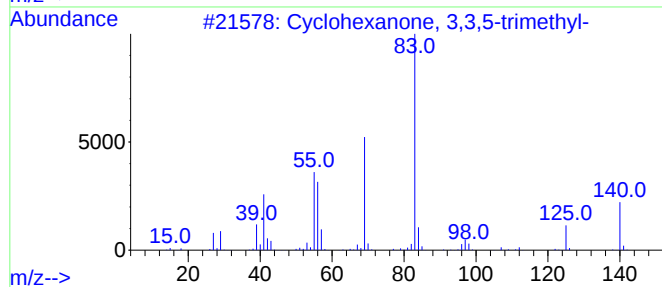
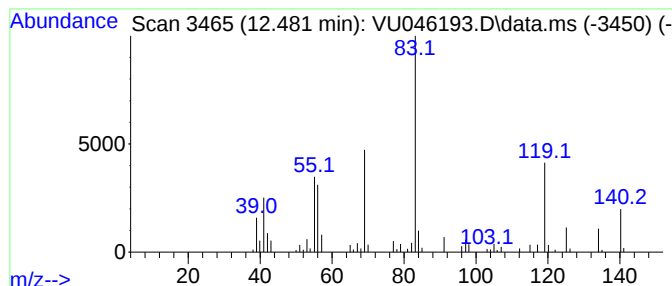
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Quant Title : VOC Analysis

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

Peak Number 7 Cyclohexanone, 3,3,5-trimet... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.481	9.79 ug/L	130361	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	94
3			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	70
4			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	70
5			3-Acetamidofuran	125	C6H7NO2	059445-85-1	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120821\
Data File : VU046193.D
Acq On : 08 Dec 2021 22:13
Operator : SY/MD
Sample : M4983-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5N3

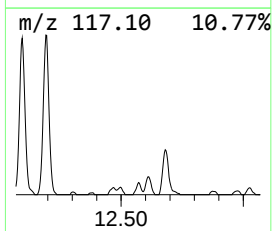
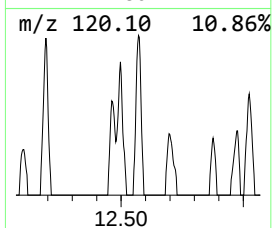
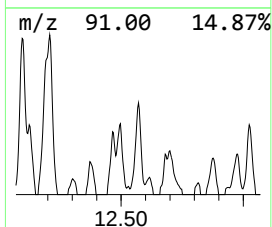
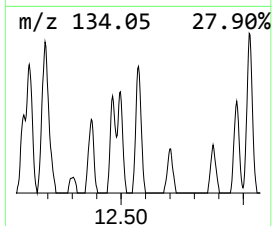
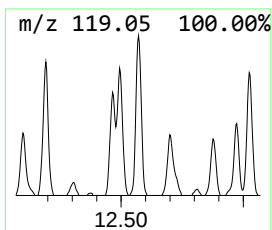
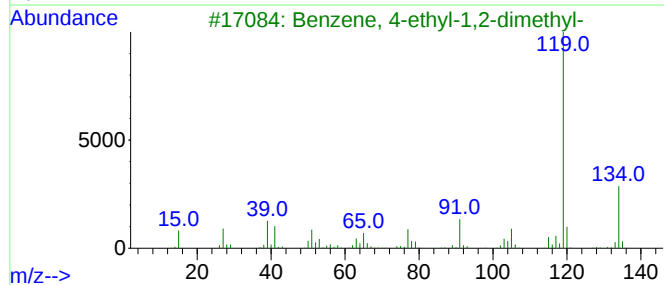
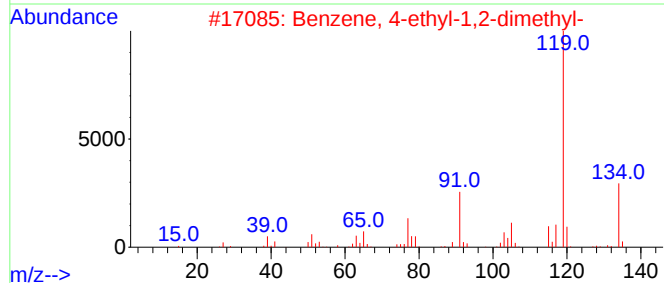
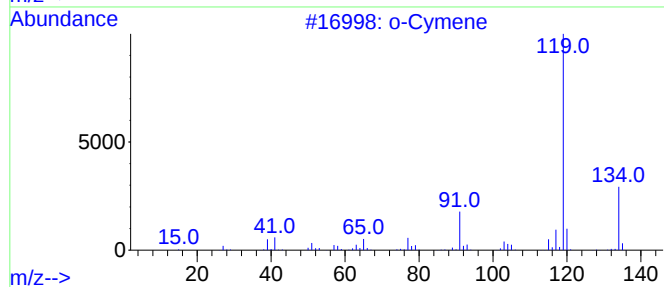
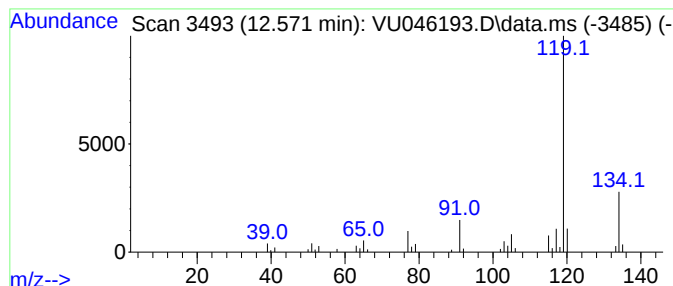
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Quant Title : VOC Analysis

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

Peak Number 8 o-Cymene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.571	2.81 ug/L	37403	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	97
2			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
3			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	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\VU120821\
Data File : VU046193.D
Acq On : 08 Dec 2021 22:13
Operator : SY/MD
Sample : M4983-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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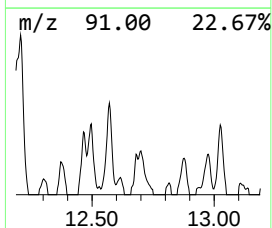
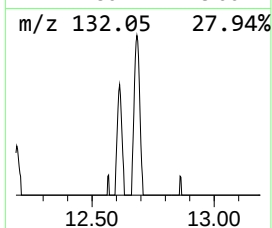
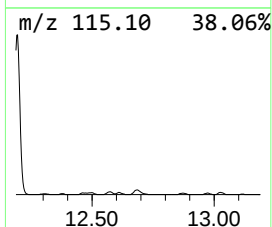
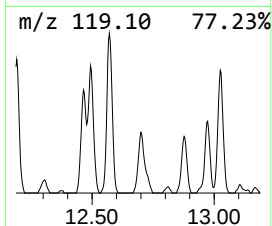
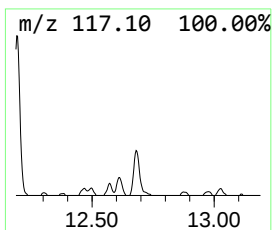
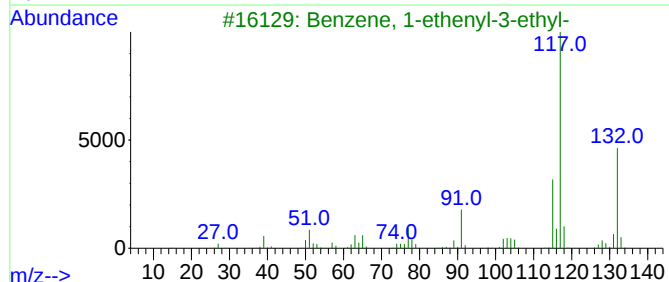
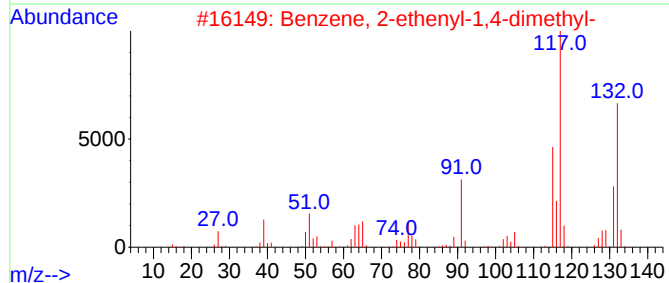
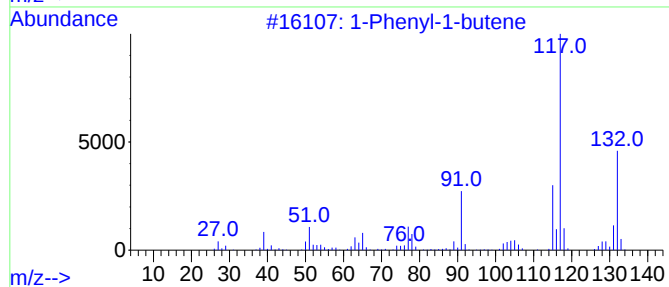
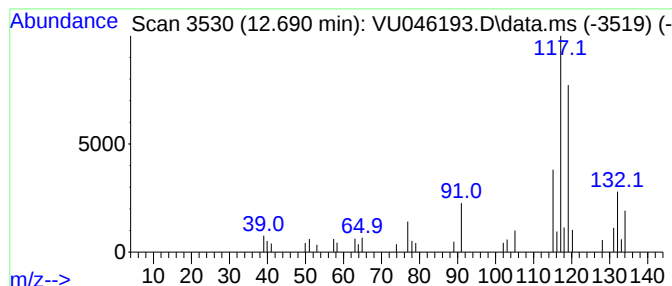
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Quant Title : VOC Analysis

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

Peak Number 9 1-Phenyl-1-butene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.690	2.57 ug/L	34227	1,4-Dichlorobenzene-d4	11.812

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Phenyl-1-butene	132	C10H12	000824-90-8	60
2		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	55
3		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	55
4		1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	53
5		3-Phenylbut-1-ene	132	C10H12	000934-10-1	49



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120821\
Data File : VU046193.D
Acq On : 08 Dec 2021 22:13
Operator : SY/MD
Sample : M4983-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5N3

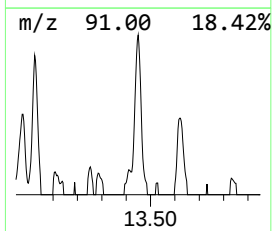
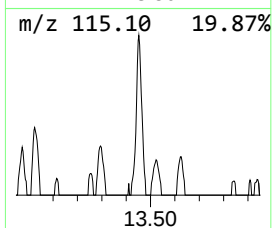
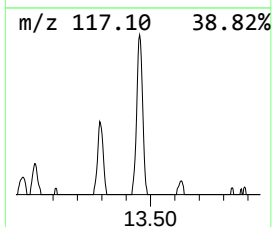
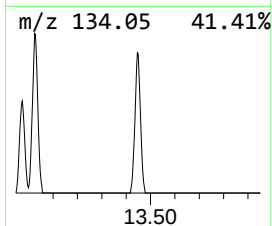
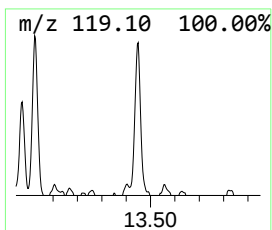
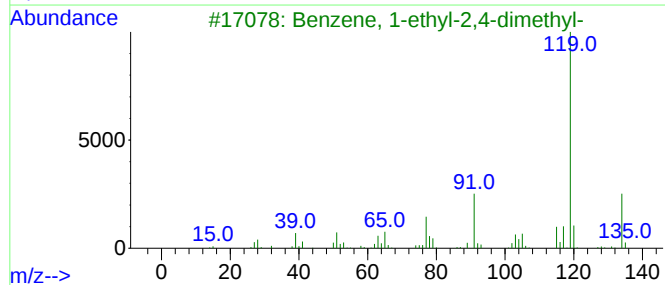
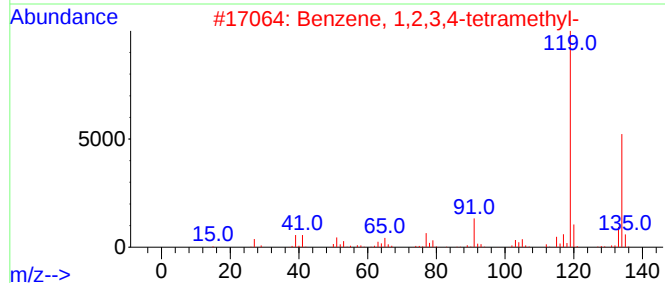
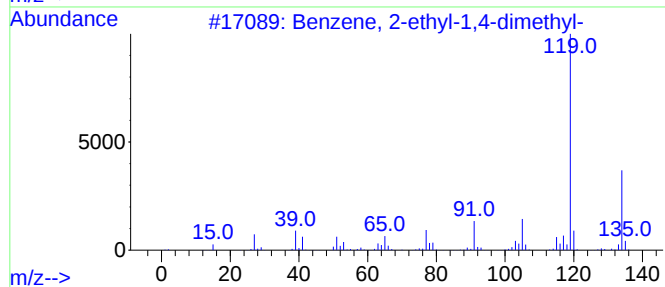
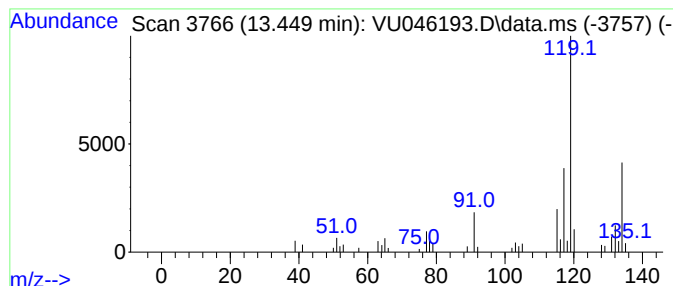
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Quant Title : VOC Analysis

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

Peak Number 10 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.449	3.24 ug/L	43123	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	70
2			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	70
3			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	64
4			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	64
5			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	62



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120821\
Data File : VU046193.D
Acq On : 08 Dec 2021 22:13
Operator : SY/MD
Sample : M4983-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5N3

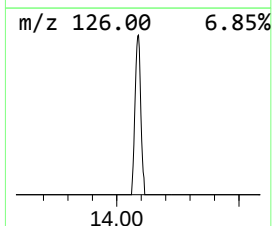
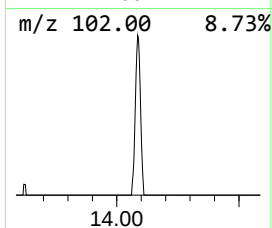
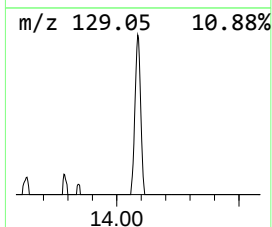
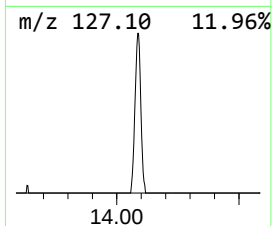
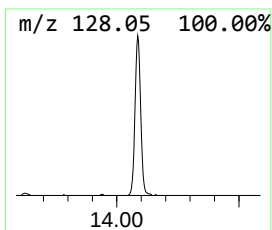
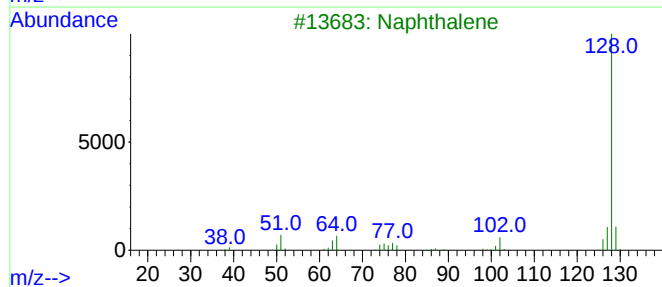
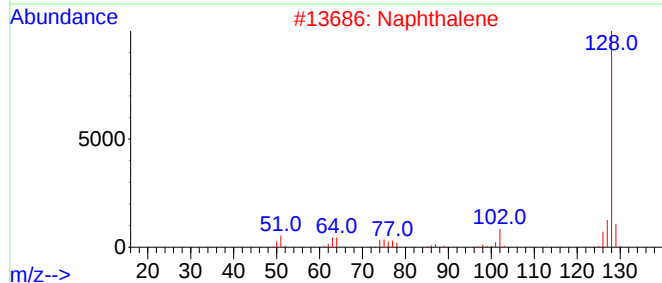
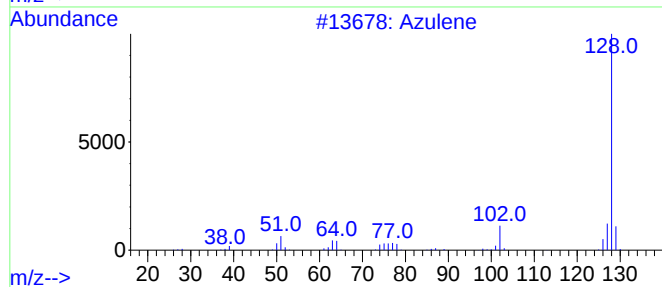
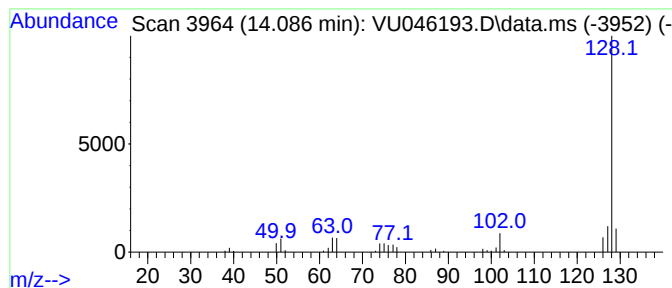
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Quant Title : VOC Analysis

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

Peak Number 11 Azulene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.086	4.21 ug/L	56023	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Azulene	128	C10H8	000275-51-4	95
2			Naphthalene	128	C10H8	000091-20-3	95
3			Naphthalene	128	C10H8	000091-20-3	95
4			Azulene	128	C10H8	000275-51-4	91
5			Naphthalene	128	C10H8	000091-20-3	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU120821\
Data File : VU046193.D
Acq On : 08 Dec 2021 22:13
Operator : SY/MD
Sample : M4983-04
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 32 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EW5N3

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
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Benzene, 1-ethy...	10.999	18.1	ug/L	240690	3	11.812	666062	50.0
Benzene, 1-ethy...	11.291	10.4	ug/L	138764	3	11.812	666062	50.0
Indane	12.095	7.5	ug/L	100118	3	11.812	666062	50.0
Cyclohexanone, ...	12.481	9.8	ug/L	130361	3	11.812	666062	50.0
o-Cymene	12.571	2.8	ug/L	37403	3	11.812	666062	50.0
1-Phenyl-1-butene	12.690	2.6	ug/L	34227	3	11.812	666062	50.0
Benzene, 2-ethy...	13.449	3.2	ug/L	43123	3	11.812	666062	50.0
Azulene	14.086	4.2	ug/L	56023	3	11.812	666062	50.0