

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
 Data File : VV020147.D
 Acq On : 22 Jan 2021 19:47
 Operator : SY/MD
 Sample : M1201-06
 Misc : 5.0mL/MSVOA_V/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 BG501

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM012121WMA.M

Title : VOC Analysis

Signal : TIC: VV020147.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.110	3	6	19	rVB	165831	140129	1.88%	0.197%
2	1.309	58	68	79	rVB2	289351	337515	4.52%	0.475%
3	1.386	79	92	99	rBV3	72230	156880	2.10%	0.221%
4	1.566	141	148	162	rVB	231049	268713	3.60%	0.378%
5	2.110	305	317	330	rBV	385925	572328	7.67%	0.805%
6	2.534	426	449	466	rVB	102149	211576	2.84%	0.298%
7	2.766	505	521	542	rBV	147599	295674	3.96%	0.416%
8	3.045	594	608	630	rVV	418620	830603	11.13%	1.168%
9	3.293	670	685	703	rVV2	68512	162020	2.17%	0.228%
10	3.672	782	803	817	rBV2	71299	193473	2.59%	0.272%
11	3.769	818	833	854	rVB	200441	515548	6.91%	0.725%
12	3.881	857	868	895	rBV	119132	320799	4.30%	0.451%
13	4.354	996	1015	1054	rBV	264150	713949	9.57%	1.004%
14	4.672	1096	1114	1129	rBV	265296	626692	8.40%	0.882%
15	4.769	1129	1144	1170	rVB	247096	640666	8.58%	0.901%
16	4.913	1170	1189	1213	rBV	368686	885471	11.87%	1.246%
17	5.052	1214	1232	1259	rBV2	651351	1691641	22.67%	2.380%
18	5.183	1261	1273	1279	rVV3	114394	251612	3.37%	0.354%
19	5.238	1280	1290	1307	rVV2	325635	858347	11.50%	1.207%
20	5.331	1307	1319	1332	rVB3	62711	137881	1.85%	0.194%
21	5.495	1356	1370	1386	rBV	286943	581635	7.79%	0.818%
22	5.621	1397	1409	1437	rBV	482433	1004441	13.46%	1.413%
23	6.074	1536	1550	1560	rBV2	391679	867620	11.63%	1.220%
24	6.203	1581	1590	1599	rBV	103258	184609	2.47%	0.260%
25	6.261	1599	1608	1618	rVV2	135778	245741	3.29%	0.346%
26	6.370	1633	1642	1656	rVB3	72581	136781	1.83%	0.192%
27	6.466	1657	1672	1686	rVB4	112738	259150	3.47%	0.365%
28	6.659	1713	1732	1749	rVB	344249	791388	10.60%	1.113%
29	6.781	1750	1770	1781	rBV2	456903	1011712	13.56%	1.423%
30	6.843	1781	1789	1799	rVB	126294	209774	2.81%	0.295%
31	6.923	1804	1814	1820	rBV2	207609	347605	4.66%	0.489%
32	6.961	1821	1826	1829	rVV	86791	112309	1.50%	0.158%
33	6.990	1830	1835	1845	rVB	215080	334070	4.48%	0.470%
34	7.084	1856	1864	1873	rBV	273636	449872	6.03%	0.633%
35	7.187	1889	1896	1908	rVB4	72569	137286	1.84%	0.193%
36	7.270	1908	1922	1928	rBV3	316357	615731	8.25%	0.866%

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Integrator: RTE

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Sampling : 1

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Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM012121WMA.M
 Title : VOC Analysis

37	7.318	1929	1937	1953	rVV	868724	1571436	21.06%	2.211%
38	7.444	1968	1976	1983	rVV4	128314	225164	3.02%	0.317%
39	7.489	1984	1990	2006	rVV3	108250	283227	3.80%	0.398%
40	7.572	2006	2016	2027	rVV3	214820	506890	6.79%	0.713%
41	7.624	2028	2032	2038	rVV	148388	240077	3.22%	0.338%
42	7.659	2039	2043	2053	rVB3	152727	237857	3.19%	0.335%
43	7.794	2075	2085	2094	rVB3	177973	311274	4.17%	0.438%
44	7.849	2094	2102	2110	rBV	103913	161541	2.16%	0.227%
45	8.090	2167	2177	2190	rBV	531115	926340	12.41%	1.303%
46	8.238	2209	2223	2232	rVB5	126262	268970	3.60%	0.378%
47	8.296	2233	2241	2248	rBV	122735	181567	2.43%	0.255%
48	8.518	2298	2310	2321	rBV4	79259	161452	2.16%	0.227%
49	8.672	2351	2358	2381	rVB4	70888	154912	2.08%	0.218%
50	8.794	2388	2396	2405	rVB	78618	128591	1.72%	0.181%
51	8.855	2405	2415	2434	rBV	840736	1366715	18.31%	1.923%
52	8.939	2434	2441	2457	rVB5	75447	152781	2.05%	0.215%
53	9.209	2517	2525	2553	rVB8	74689	227449	3.05%	0.320%
54	9.476	2598	2608	2628	rVB4	88029	199951	2.68%	0.281%
55	9.711	2664	2681	2688	rBV9	70359	181358	2.43%	0.255%
56	9.807	2703	2711	2725	rVB8	53242	110342	1.48%	0.155%
57	10.222	2829	2840	2858	rVB	579206	956098	12.81%	1.345%
58	11.064	3089	3102	3106	rBV	169097	260503	3.49%	0.366%
59	11.093	3107	3111	3125	rVB	176114	253833	3.40%	0.357%
60	11.199	3135	3144	3152	rBV	332076	484858	6.50%	0.682%
61	11.254	3152	3161	3185	rVB2	1045099	1636397	21.93%	2.302%
62	11.556	3240	3255	3257	rBV2	1255832	1605356	21.51%	2.258%
63	11.582	3257	3263	3271	rVV	2425695	3583166	48.01%	5.040%
64	11.643	3271	3282	3302	rVB5	3358614	7462643	100.00%	10.498%
65	11.752	3306	3316	3327	rVB	184333	275106	3.69%	0.387%
66	11.826	3327	3339	3356	rVB2	1041239	1602386	21.47%	2.254%
67	11.919	3356	3368	3372	rBV	1260191	1863903	24.98%	2.622%
68	11.948	3372	3377	3388	rVB	1289675	1822057	24.42%	2.563%
69	12.022	3388	3400	3421	rVB	2407895	3727582	49.95%	5.244%
70	12.125	3421	3432	3437	rBV	286159	475030	6.37%	0.668%
71	12.167	3438	3445	3462	rVB5	294124	683698	9.16%	0.962%
72	12.267	3462	3476	3483	rBV5	106974	227379	3.05%	0.320%
73	12.321	3483	3493	3505	rVB2	280859	520829	6.98%	0.733%
74	12.421	3505	3524	3532	rBV	1258290	1976904	26.49%	2.781%
75	12.473	3532	3540	3556	rVB	2026413	3022519	40.50%	4.252%
76	12.559	3556	3567	3572	rBV	568233	836043	11.20%	1.176%

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Title : VOC Analysis

77	12.595	3572	3578	3582	rVV	395310	518230	6.94%	0.729%
78	12.624	3583	3587	3595	rVB	307802	356937	4.78%	0.502%
79	12.691	3598	3608	3614	rBV3	196255	379668	5.09%	0.534%
80	12.733	3614	3621	3630	rVB	537599	712276	9.54%	1.002%
81	12.804	3636	3643	3650	rVB2	282007	366795	4.92%	0.516%
82	12.858	3650	3660	3664	rBV3	464925	791714	10.61%	1.114%
83	12.891	3665	3670	3680	rVB2	463279	719529	9.64%	1.012%
84	13.009	3699	3707	3718	rBV	382807	545747	7.31%	0.768%
85	13.177	3747	3759	3765	rBV	440581	681255	9.13%	0.958%
86	13.225	3765	3774	3782	rVB	750869	1117916	14.98%	1.573%
87	13.276	3782	3790	3797	rBV2	1133871	1766287	23.67%	2.485%
88	13.344	3807	3811	3815	rBV2	120363	120275	1.61%	0.169%
89	13.376	3815	3821	3827	rVV	477937	579861	7.77%	0.816%
90	13.408	3828	3831	3843	rVB	371608	455299	6.10%	0.640%
91	13.472	3843	3851	3868	rBV4	50731	135542	1.82%	0.191%
92	13.556	3868	3877	3889	rBV2	95599	163991	2.20%	0.231%
93	13.620	3889	3897	3905	rBV2	69412	107702	1.44%	0.152%
94	13.720	3918	3928	3939	rBV3	311593	618919	8.29%	0.871%
95	13.778	3939	3946	3956	rVB2	282754	436773	5.85%	0.614%
96	13.919	3979	3990	4003	rVB	542974	816215	10.94%	1.148%
97	14.099	4036	4046	4059	rBV	401372	829068	11.11%	1.166%
98	14.241	4081	4090	4100	rVV2	198322	304428	4.08%	0.428%
99	14.305	4100	4110	4123	rVB3	230280	443311	5.94%	0.624%
100	14.450	4142	4155	4166	rBV6	62223	144922	1.94%	0.204%

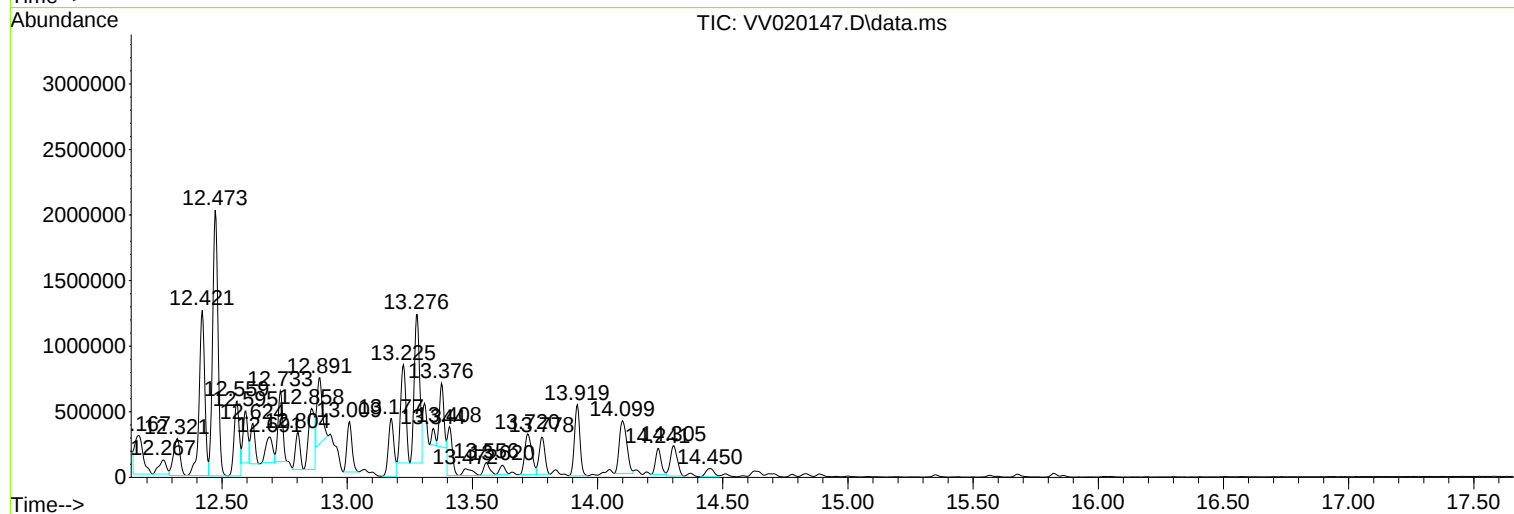
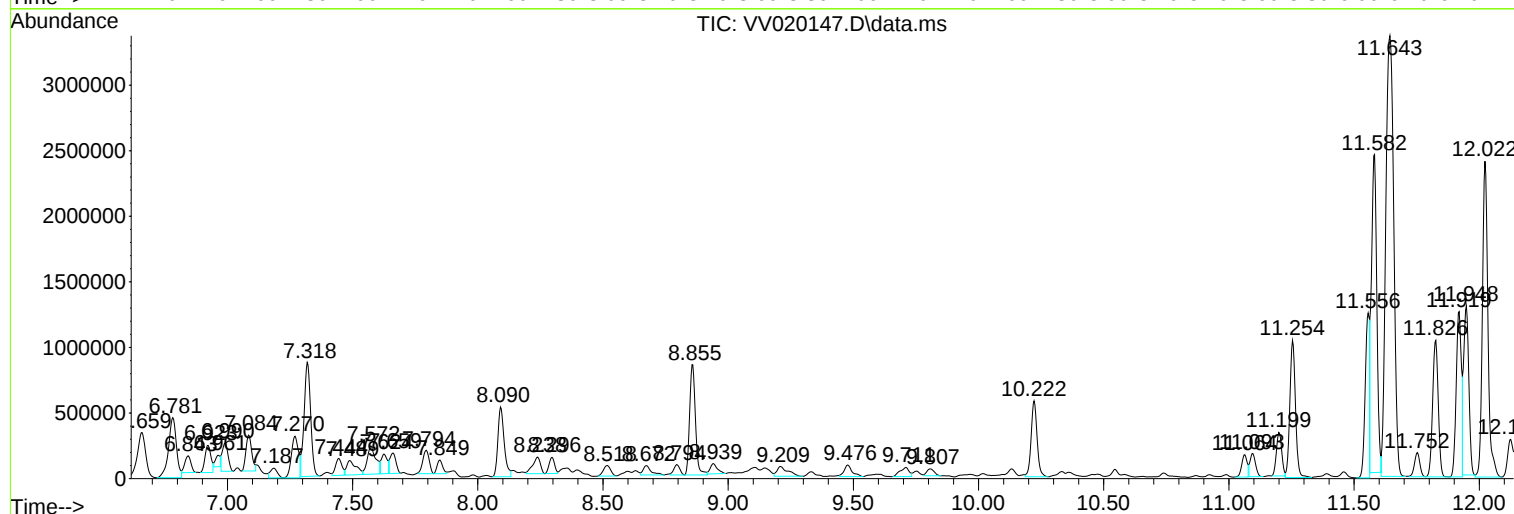
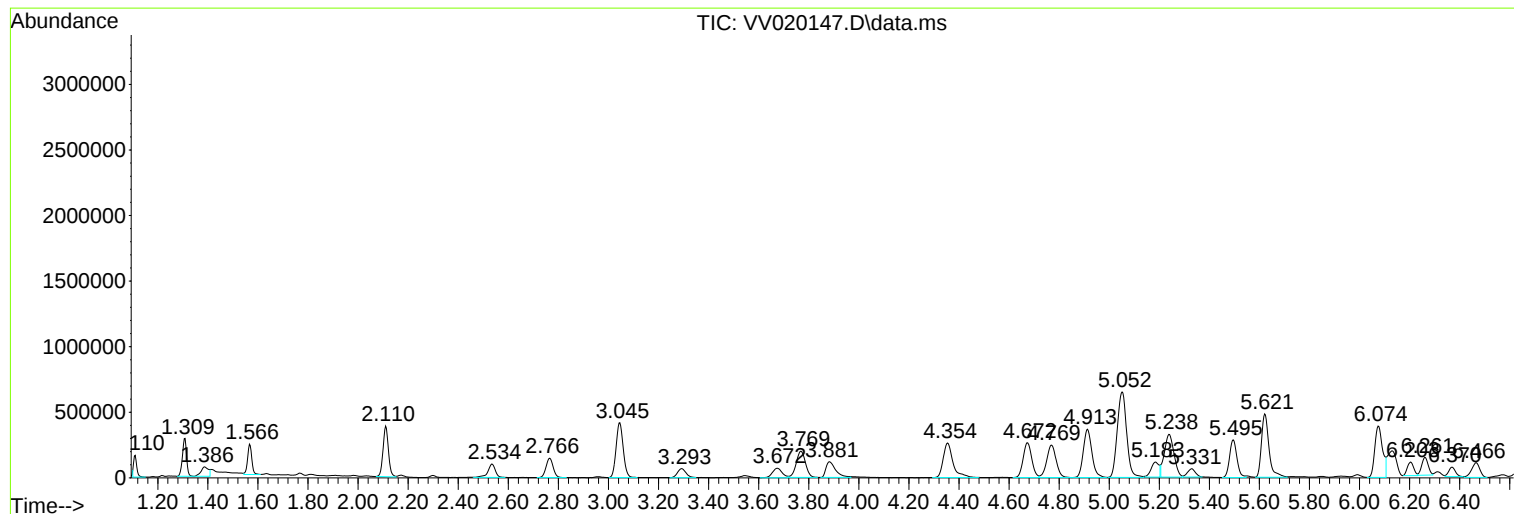
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ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM012121WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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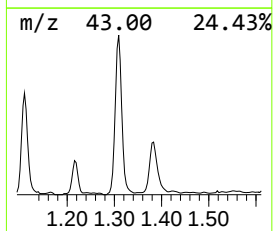
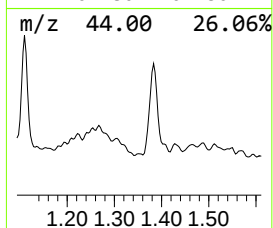
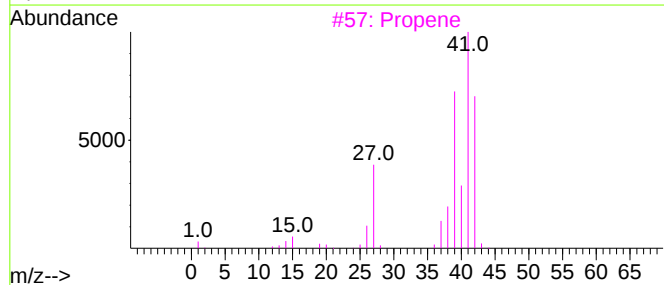
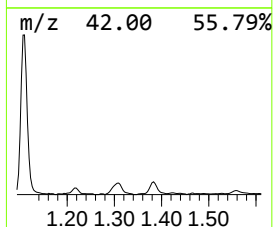
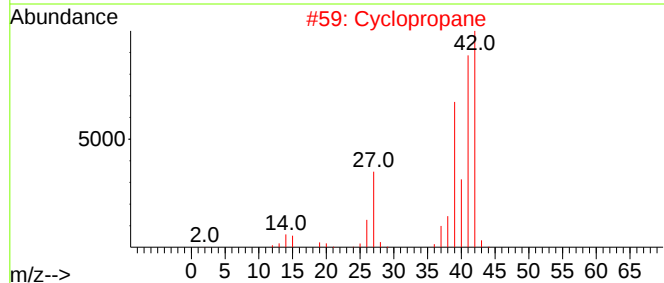
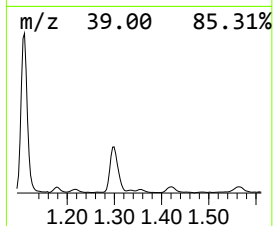
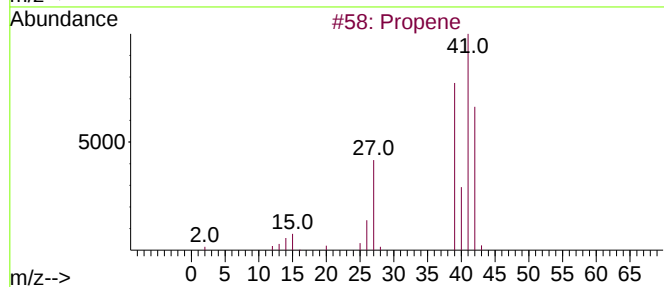
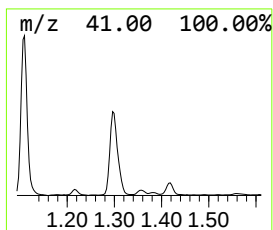
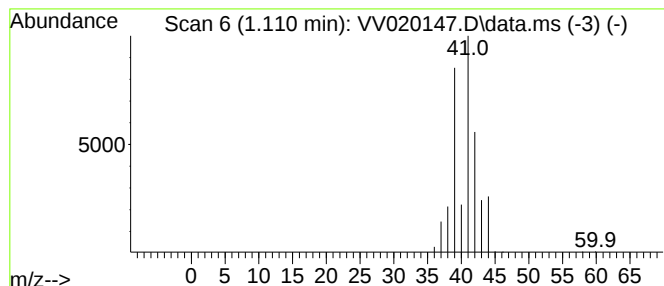
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM012121WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 63

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.110	6.98 ug/L	140129	1,4-Difluorobenzene	5.621

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propene	42	C3H6	000115-07-1	42
2			Cyclopropane	42	C3H6	000075-19-4	40
3			Propene	42	C3H6	000115-07-1	30
4			Cyclopropene	40	C3H4	002781-85-3	9
5			Cyclopropane	42	C3H6	000075-19-4	7



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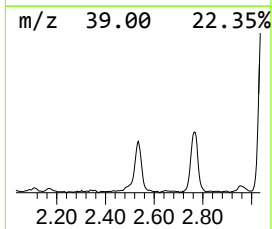
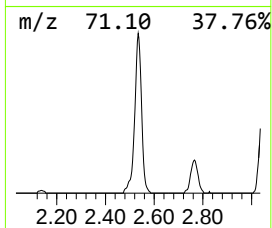
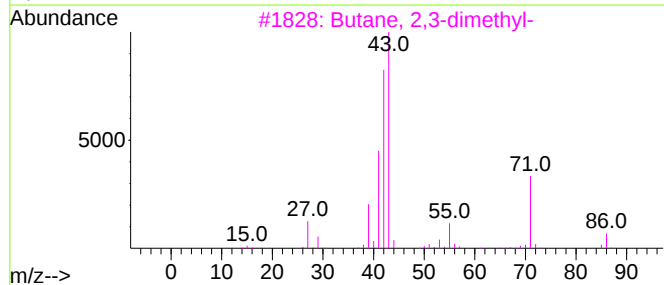
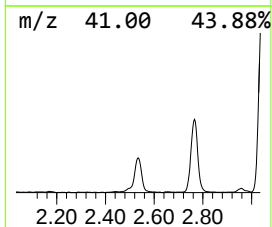
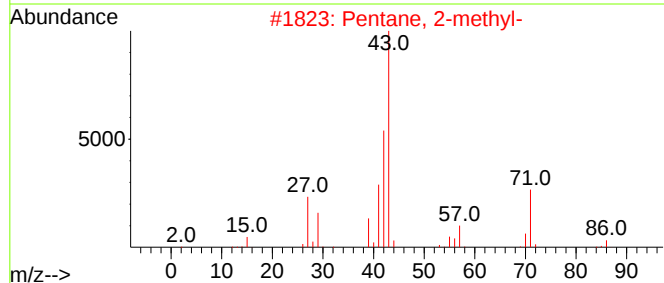
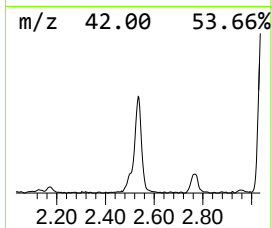
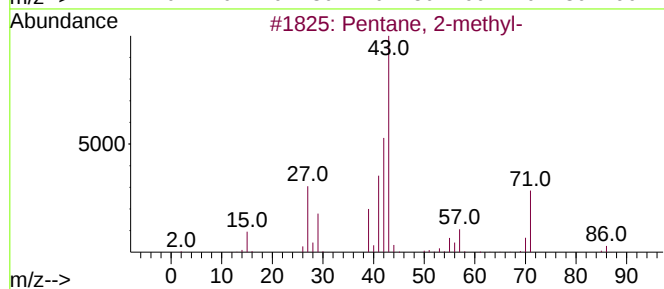
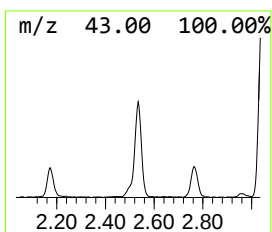
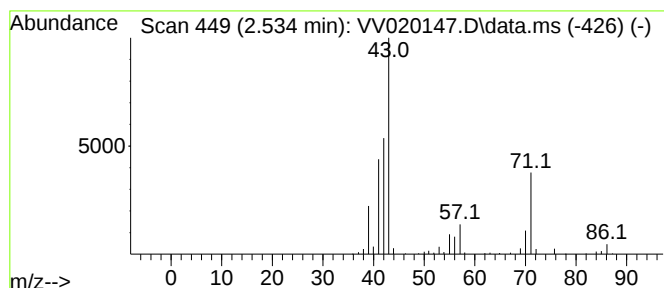
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM012121WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 48

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.534	10.53 ug/L	211576	1,4-Difluorobenzene	5.621

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2-methyl-	86	C6H14	000107-83-5	87
2			Pentane, 2-methyl-	86	C6H14	000107-83-5	86
3			Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	58
4			Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	53
5			Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	42



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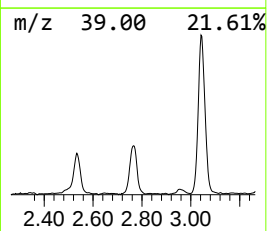
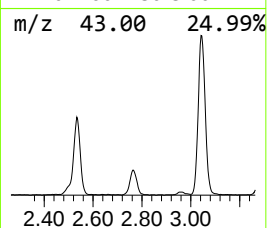
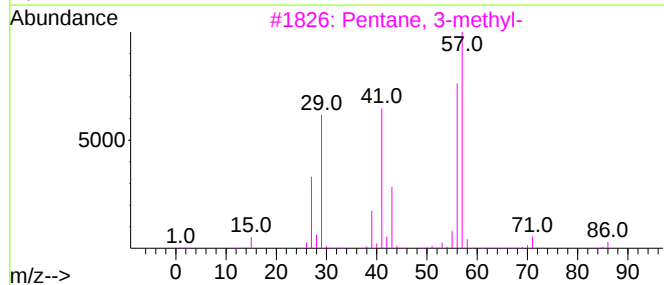
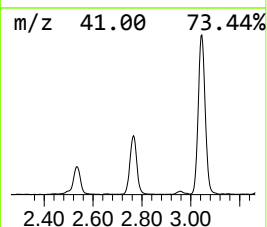
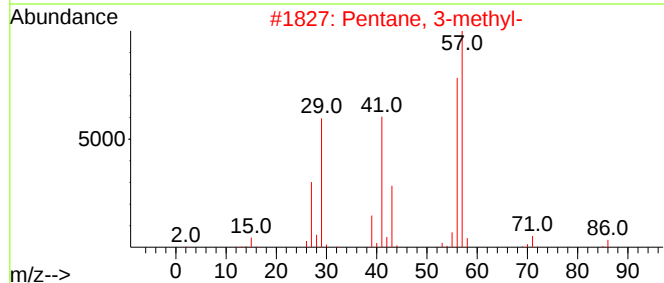
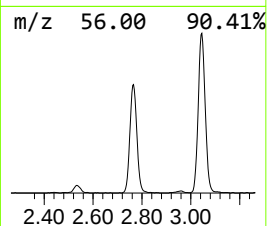
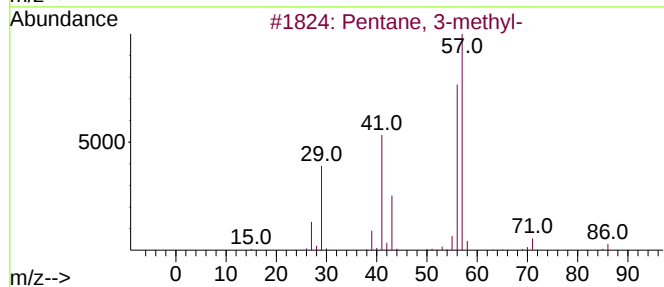
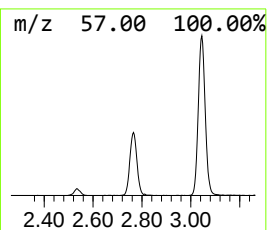
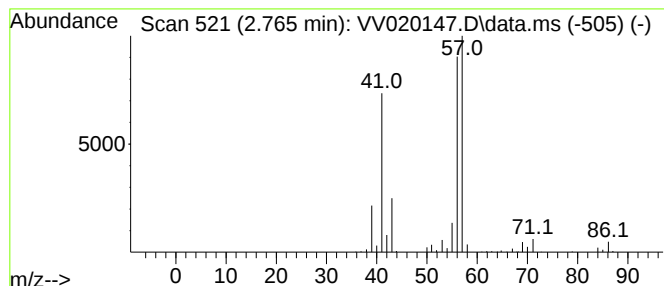
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.765	14.72 ug/L	295674	1,4-Difluorobenzene	5.621

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 3-methyl-	86	C6H14	000096-14-0	86
2	Pentane, 3-methyl-	86	C6H14	000096-14-0	86
3	Pentane, 3-methyl-	86	C6H14	000096-14-0	86
4	Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	64
5	Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

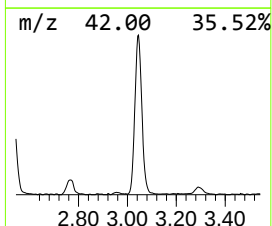
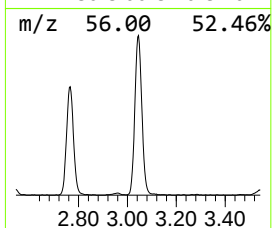
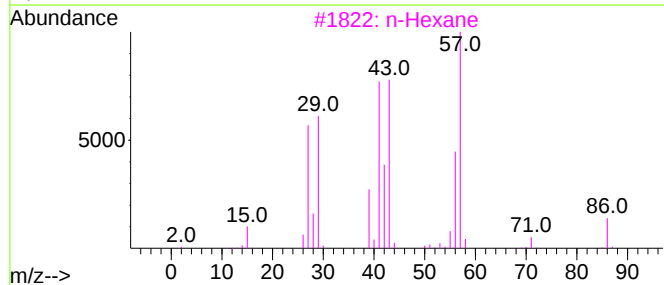
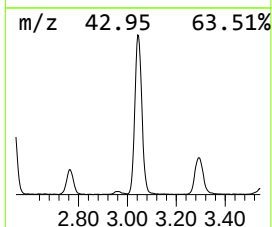
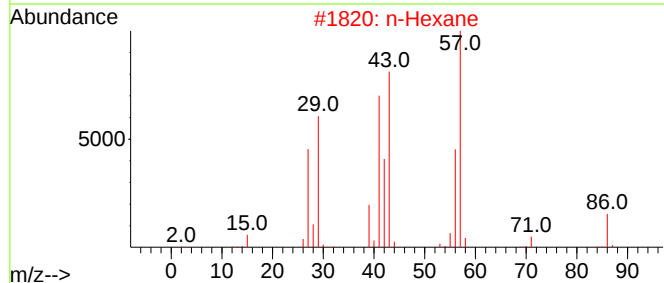
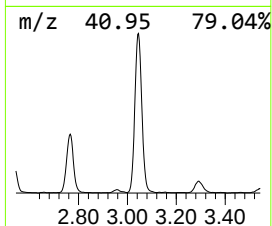
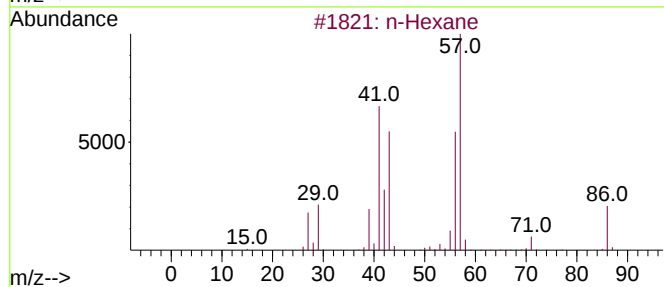
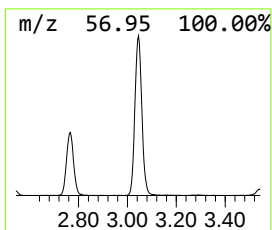
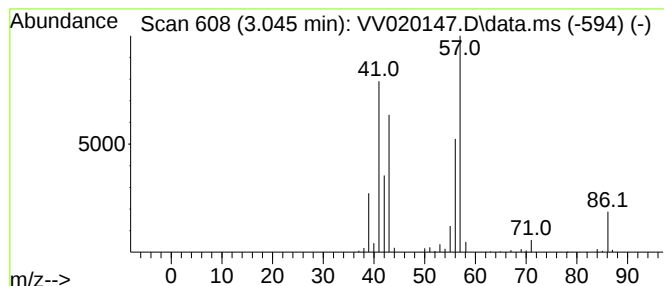
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.045	41.35 ug/L	830603	1,4-Difluorobenzene	5.621

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexane	86	C6H14	000110-54-3	91
2	n-Hexane	86	C6H14	000110-54-3	83
3	n-Hexane	86	C6H14	000110-54-3	72
4	Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	53
5	1-Pentene, 4-methyl-	84	C6H12	000691-37-2	40



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

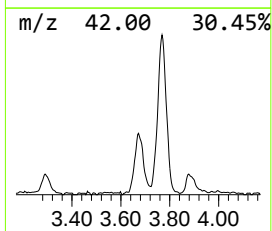
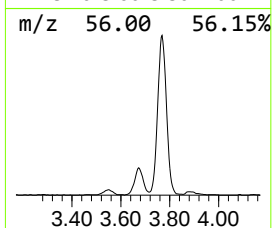
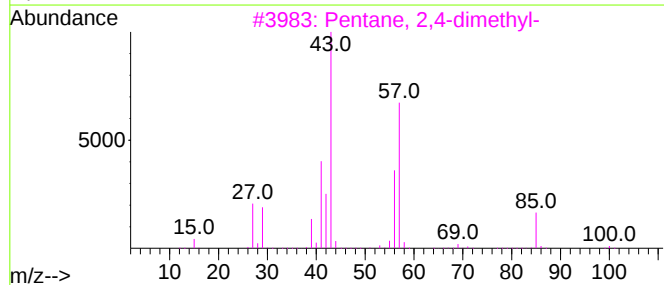
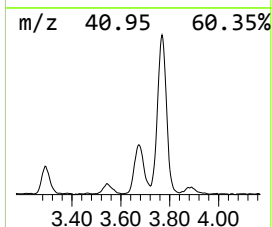
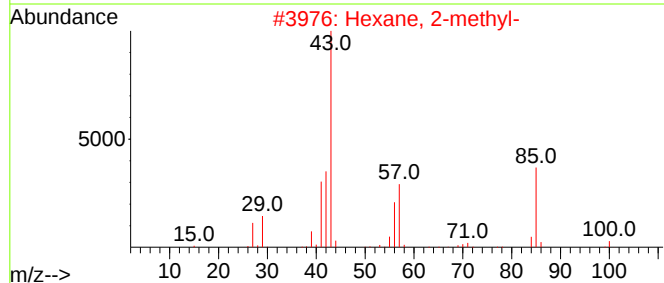
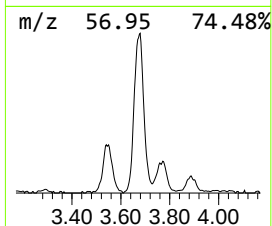
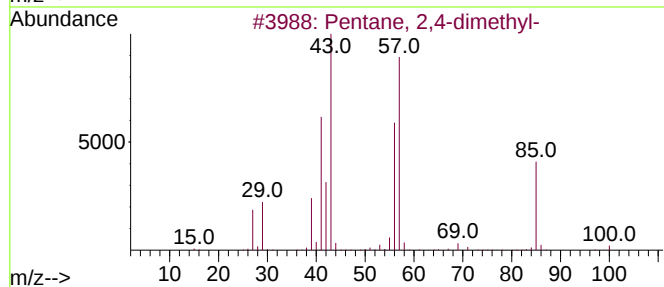
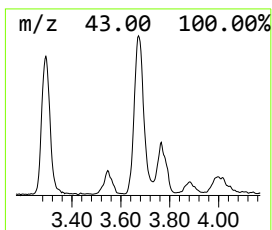
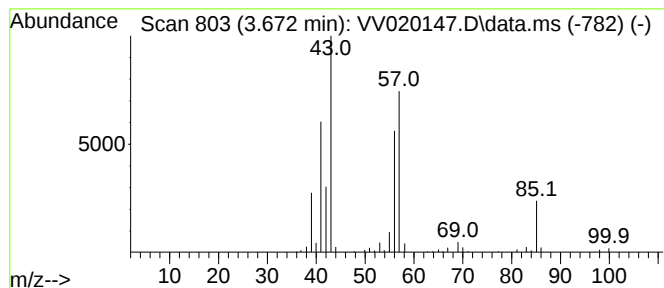
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 52

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.672	9.63 ug/L	193473	1,4-Difluorobenzene	5.621

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	81
2		Hexane, 2-methyl-	100	C7H16	000591-76-4	78
3		Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	78
4		Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	64
5		Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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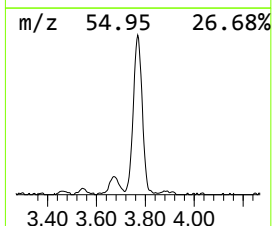
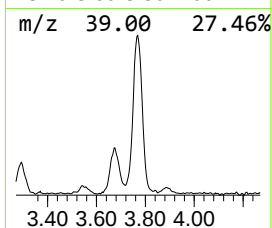
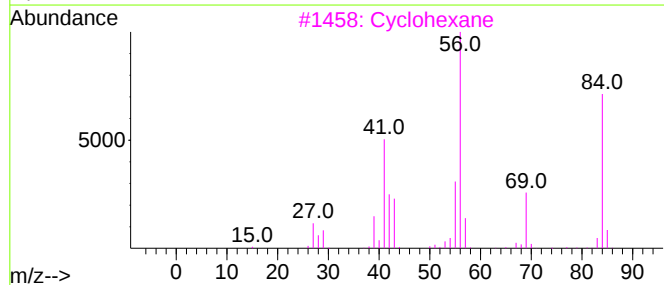
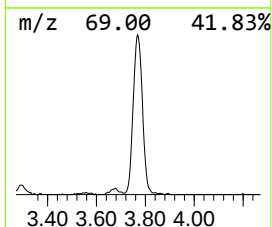
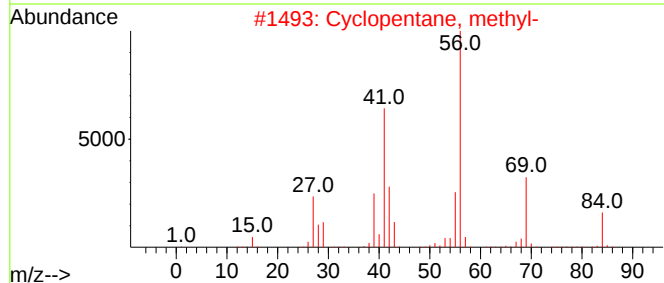
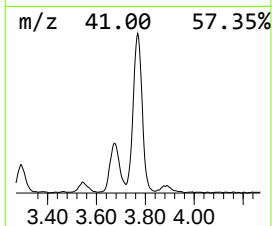
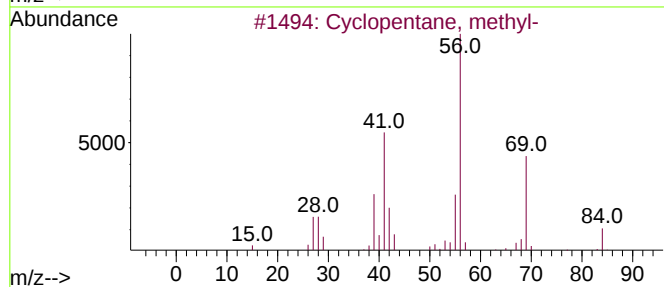
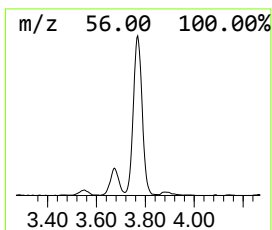
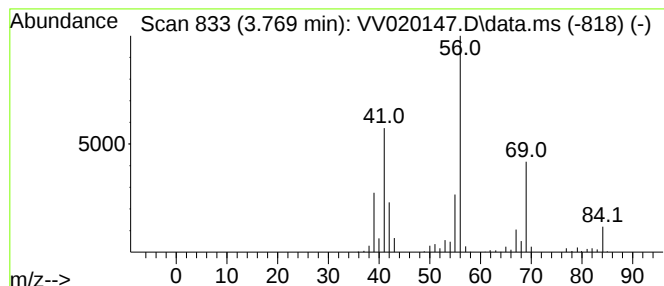
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 (DEL) Alkane: Cyclic3.769 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.769	25.66 ug/L	515548	1,4-Difluorobenzene	5.621

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2			Cyclopentane, methyl-	84	C6H12	000096-37-7	87
3			Cyclohexane	84	C6H12	000110-82-7	78
4			1-Pentene, 2-methyl-	84	C6H12	000763-29-1	64
5			1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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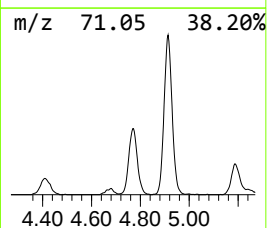
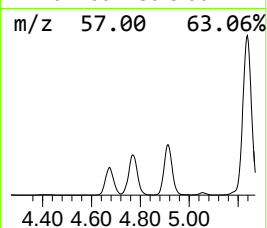
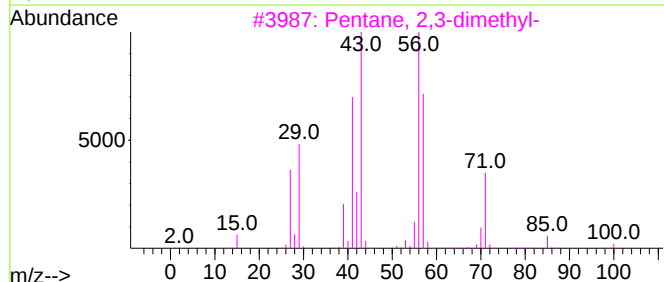
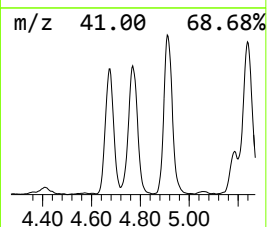
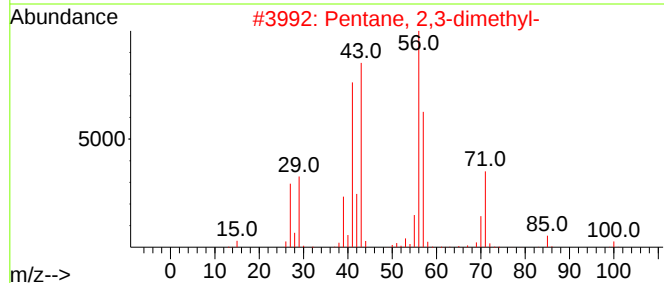
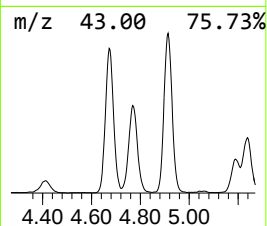
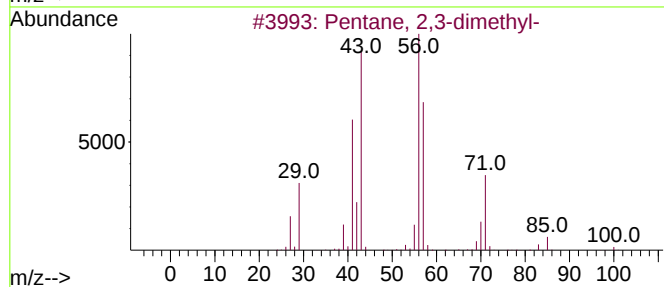
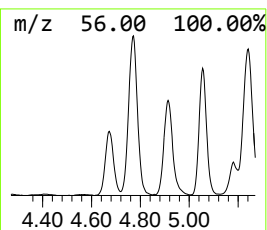
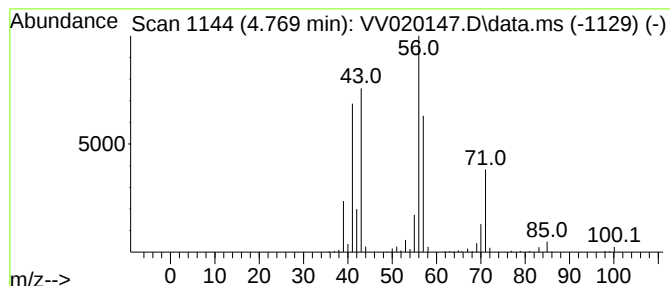
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.769	31.89 ug/L	640666	1,4-Difluorobenzene	5.621

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	94
2		Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	94
3		Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	91
4		Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	91
5		Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	45



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

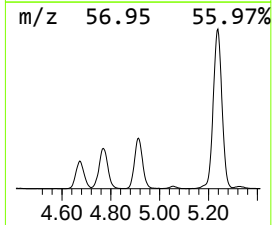
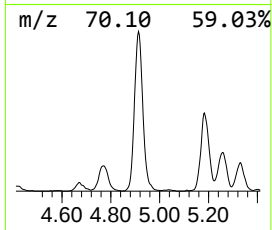
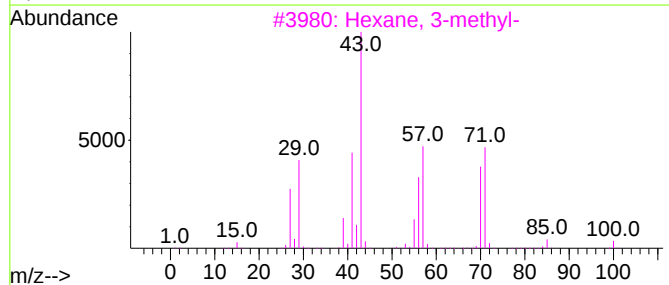
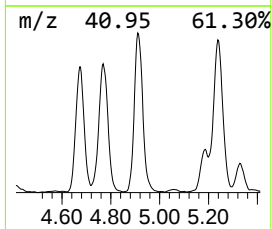
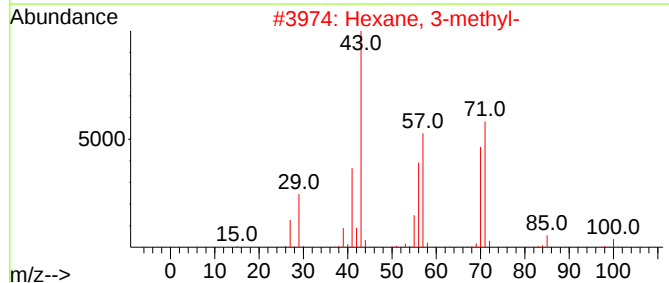
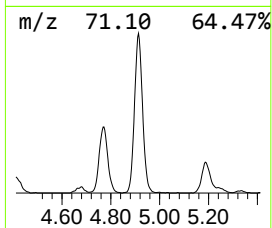
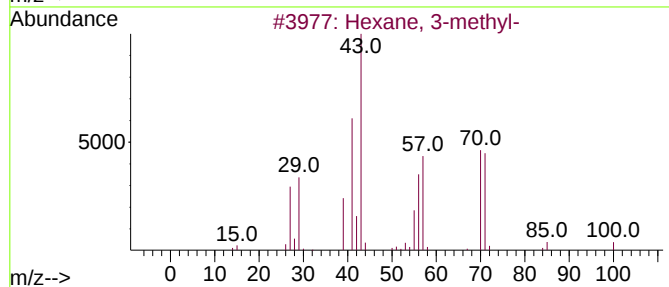
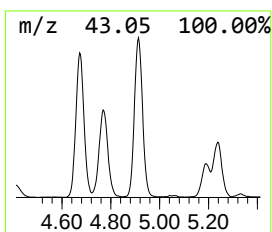
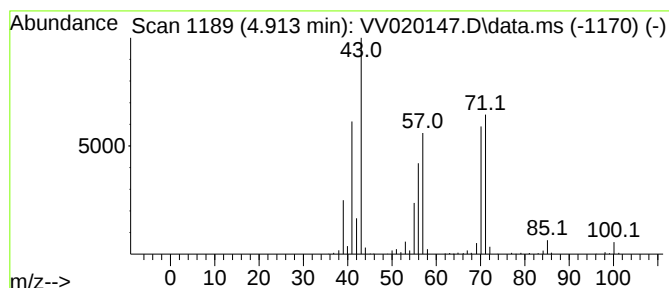
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.913	44.08 ug/L	885471	1,4-Difluorobenzene	5.621

Hit#	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 3-methyl-		100	C7H16	000589-34-4	94
2	Hexane, 3-methyl-		100	C7H16	000589-34-4	91
3	Hexane, 3-methyl-		100	C7H16	000589-34-4	87
4	Heptane		100	C7H16	000142-82-5	80
5	Heptane		100	C7H16	000142-82-5	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

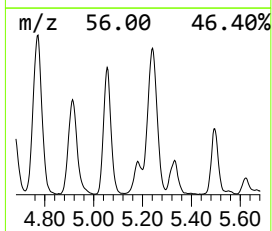
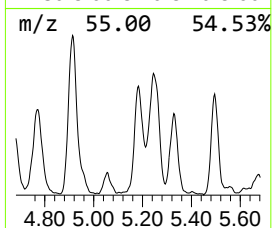
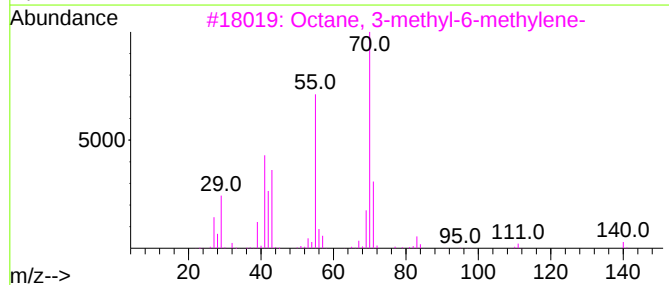
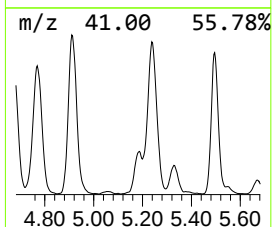
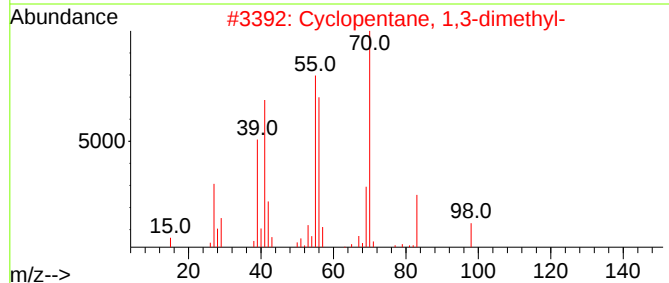
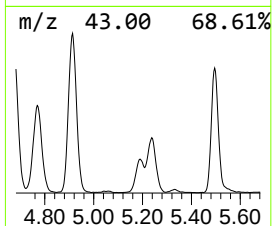
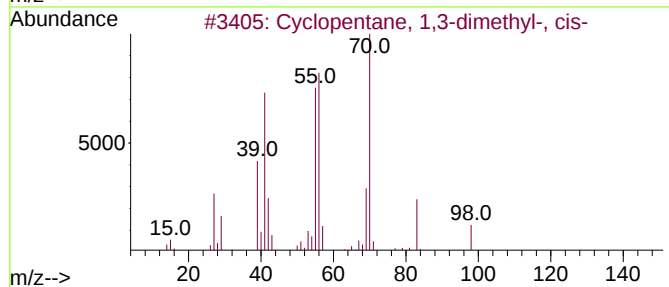
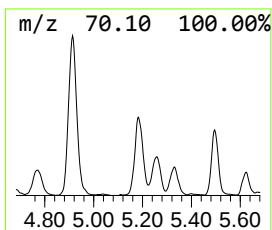
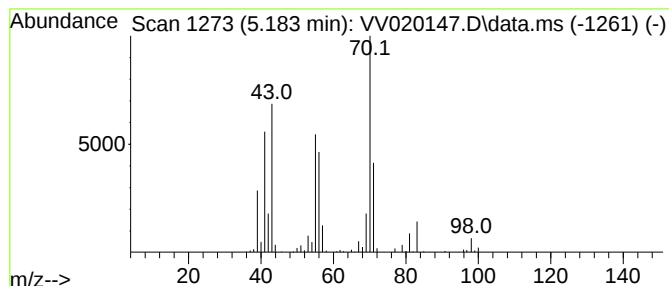
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 (DEL) Alkane: Cyclic5.183 Concentration Rank 42

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.183	12.53 ug/L	251612	1,4-Difluorobenzene	5.621

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	60
2		Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	60
3		Octane, 3-methyl-6-methylene-	140	C10H20	074630-07-2	59
4		Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14	001192-18-3	53
5		Heptane, 3-ethyl-5-methylene-	140	C10H20	1000160-41-7	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

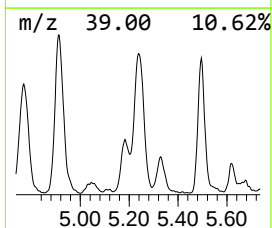
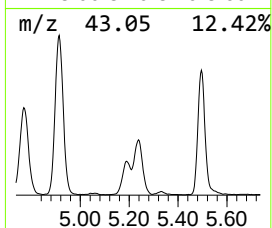
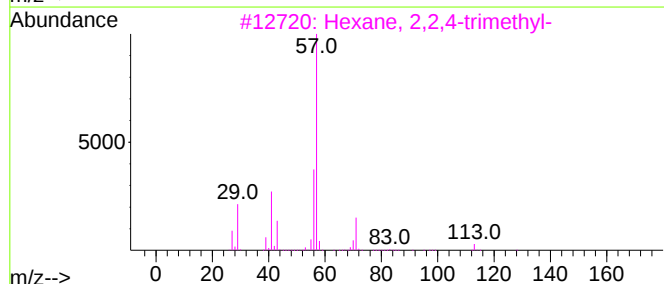
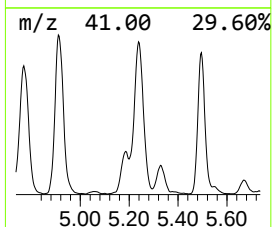
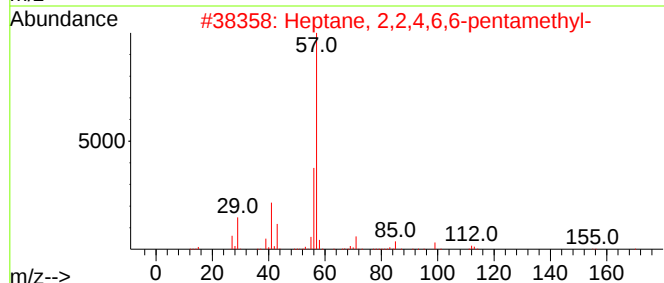
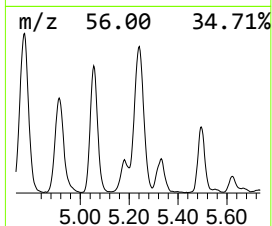
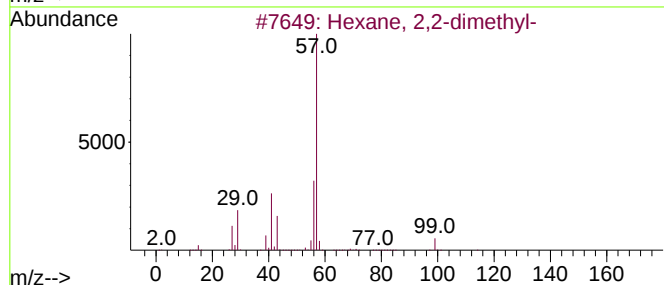
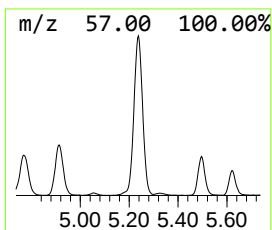
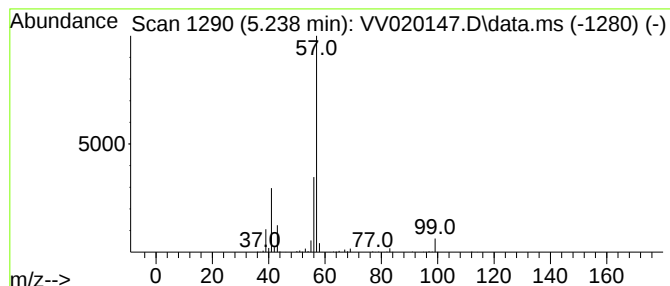
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.238	42.73 ug/L	858347	1,4-Difluorobenzene	5.621

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	72
2		Heptane, 2,2,4,6,6-pentamethyl-	170	C12H26	013475-82-6	72
3		Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	64
4		Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	59
5		Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

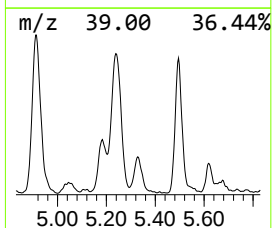
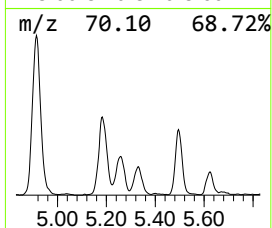
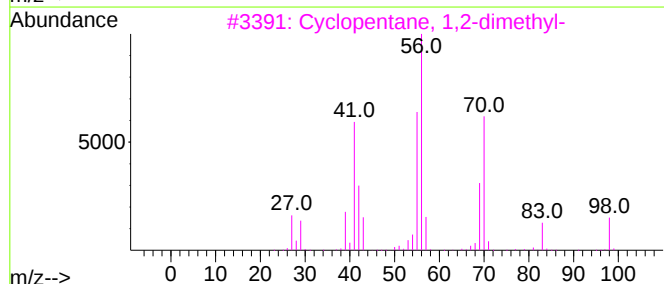
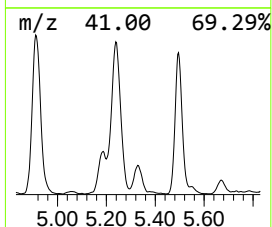
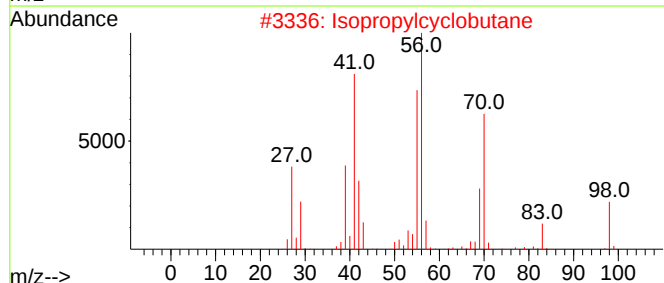
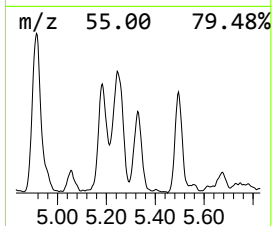
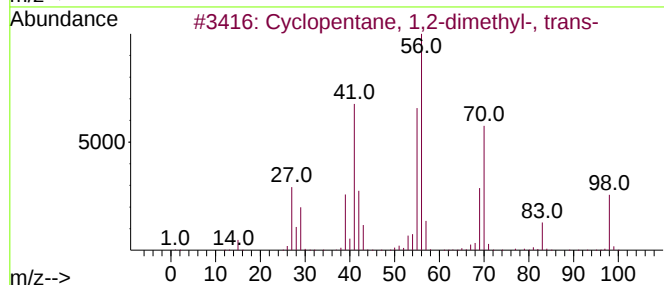
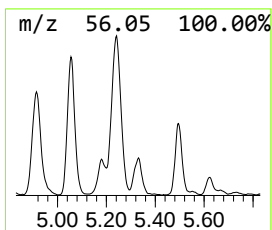
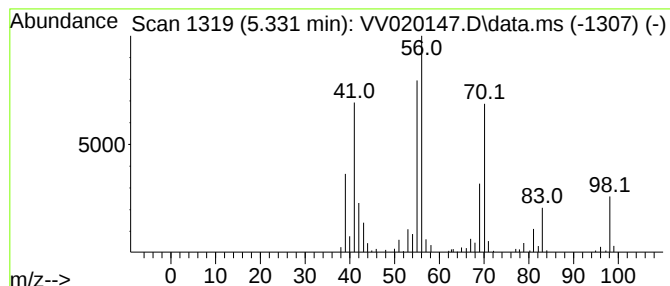
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 (DEL) Alkane: Cyclic5.331 Concentration Rank 65

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.331	6.86 ug/L	137881	1,4-Difluorobenzene	5.621

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopentane, 1,2-dimethyl-, trans-	98	C7H14	000822-50-4	90
2		Isopropylcyclobutane	98	C7H14	000872-56-0	83
3		Cyclopentane, 1,2-dimethyl-	98	C7H14	002452-99-5	72
4		Cyclopentane, 1,2-dimethyl-, trans-	98	C7H14	000822-50-4	72
5		1-Heptene	98	C7H14	000592-76-7	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

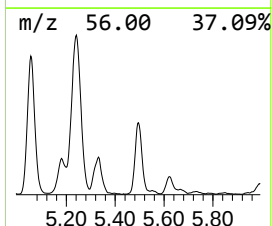
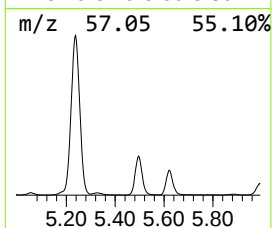
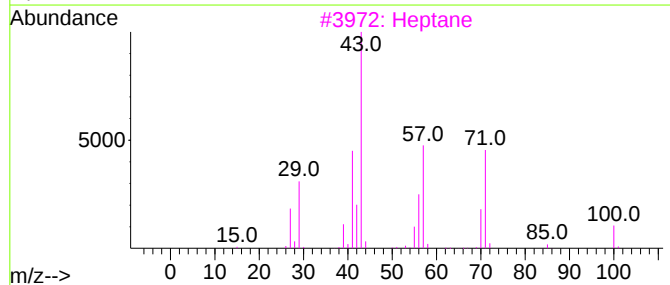
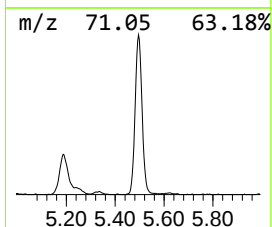
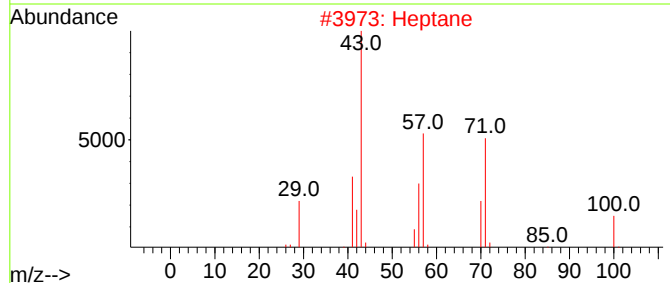
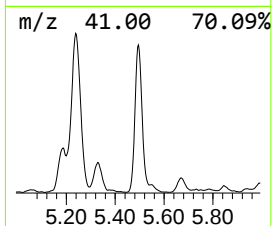
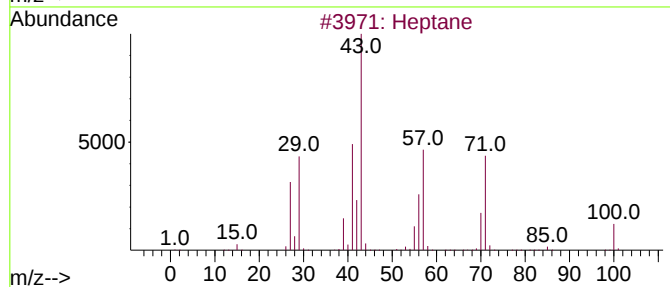
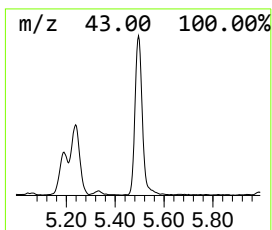
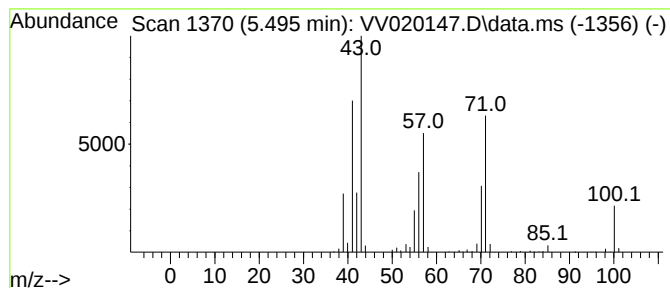
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.495	28.95 ug/L	581635	1,4-Difluorobenzene	5.621

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane			100	C7H16	000142-82-5	87
2	Heptane			100	C7H16	000142-82-5	86
3	Heptane			100	C7H16	000142-82-5	68
4	Heptane			100	C7H16	000142-82-5	58
5	Hexane, 3-methyl-			100	C7H16	000589-34-4	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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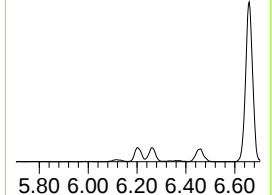
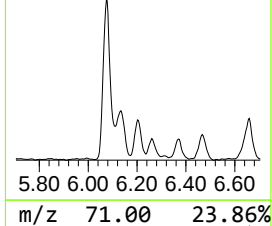
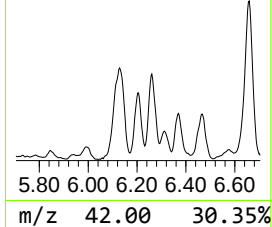
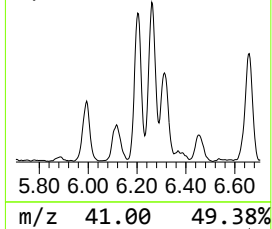
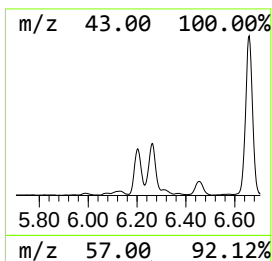
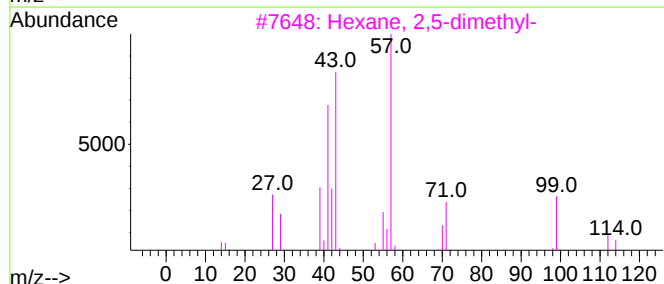
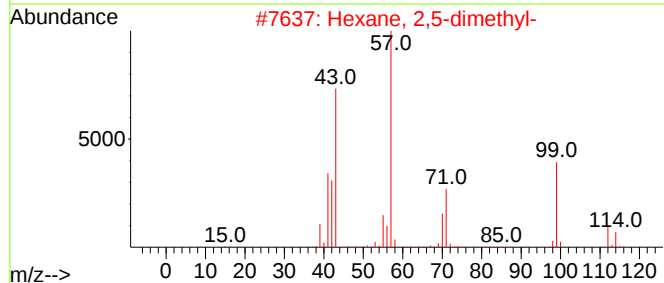
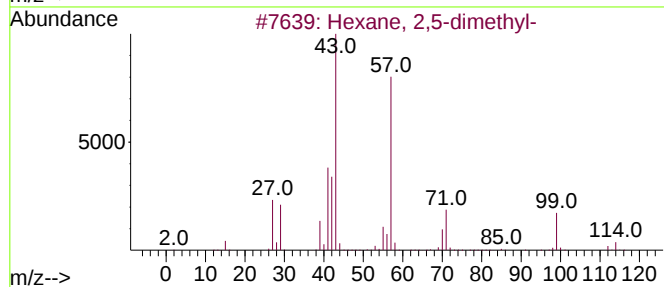
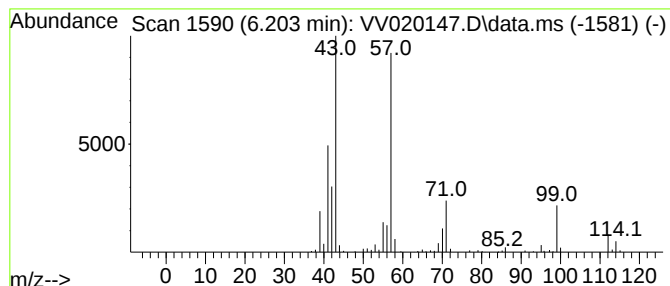
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 54

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.203	9.19 ug/L	184609	1,4-Difluorobenzene	5.621

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 2,5-dimethyl-	114	C8H18	000592-13-2	91
2			Hexane, 2,5-dimethyl-	114	C8H18	000592-13-2	87
3			Hexane, 2,5-dimethyl-	114	C8H18	000592-13-2	87
4			Octane, 1,1'-oxybis-	242	C16H34O	000629-82-3	53
5			Heptane, 2-methyl-	114	C8H18	000592-27-8	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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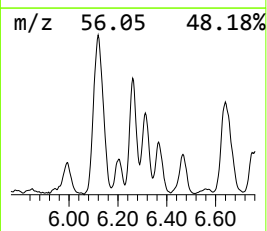
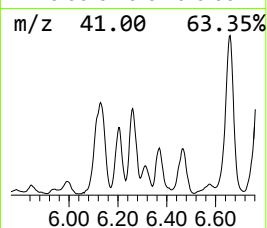
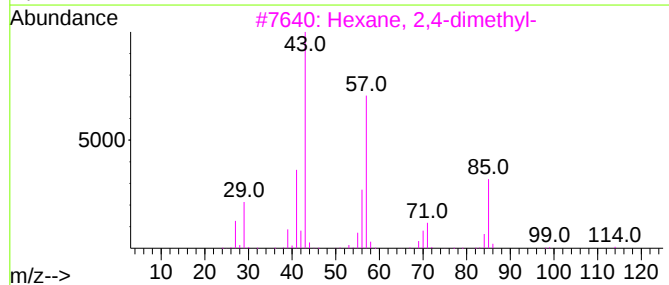
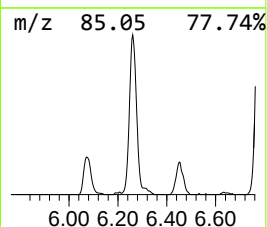
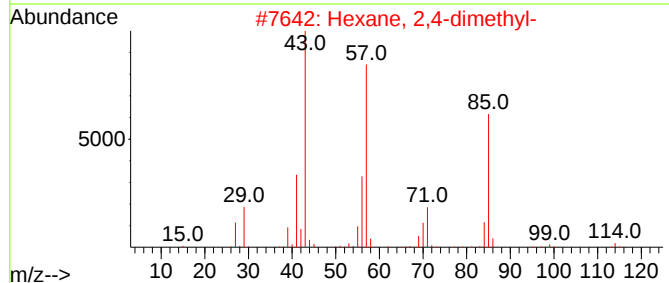
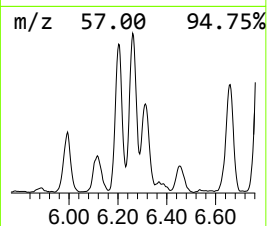
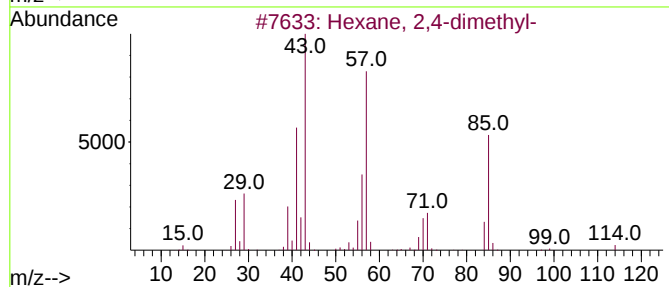
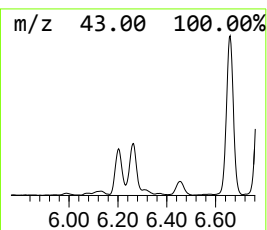
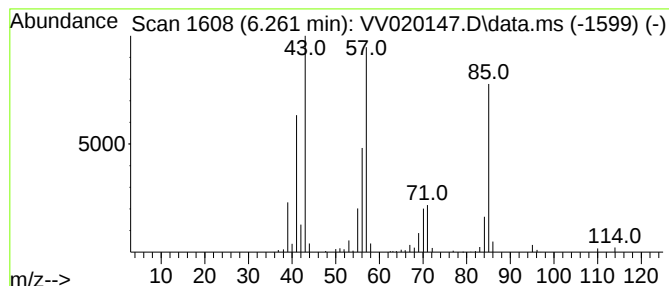
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 43

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.261	12.23 ug/L	245741	1,4-Difluorobenzene	5.621

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	96
2			Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	91
3			Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	91
4			Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	91
5			Pentane, 2,2,3,4-tetramethyl-	128	C9H20	001186-53-4	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

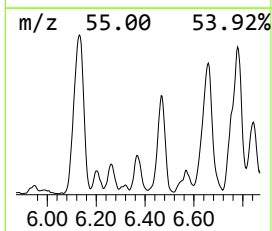
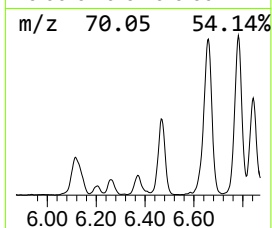
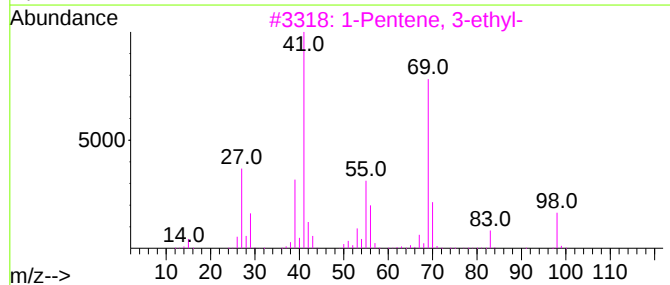
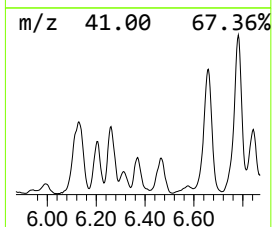
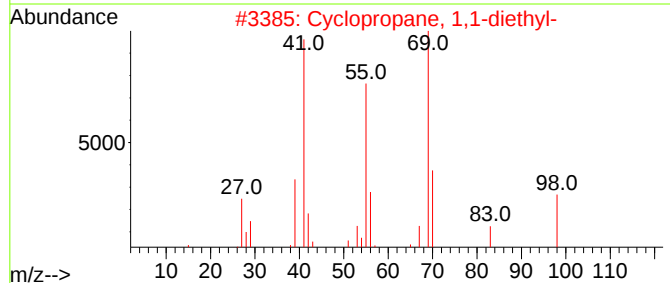
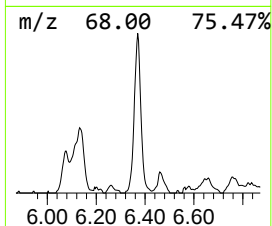
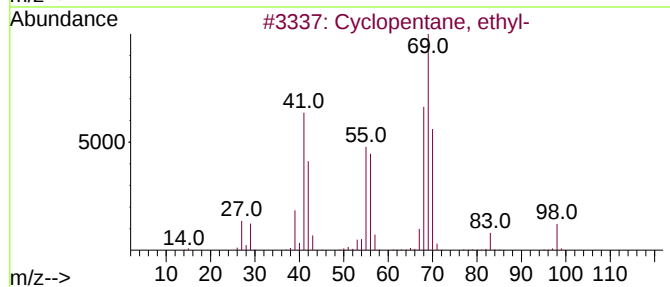
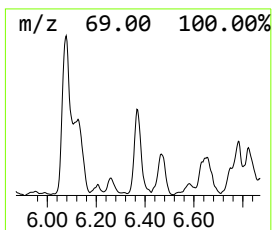
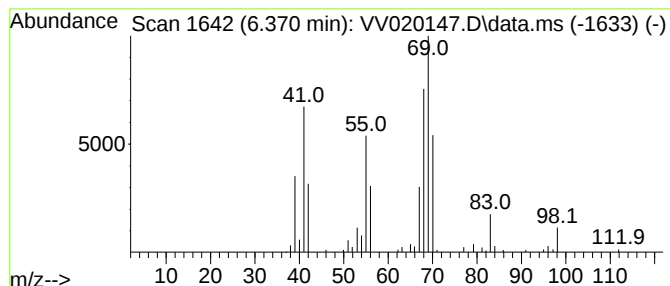
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 (DEL) Alkane: Cyclic6.370 Concentration Rank 67

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.370	6.81 ug/L	136781	1,4-Difluorobenzene	5.621

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopentane, ethyl-	98	C7H14	001640-89-7	80
2		Cyclopropane, 1,1-diethyl-	98	C7H14	001003-19-6	46
3		1-Pentene, 3-ethyl-	98	C7H14	004038-04-4	46
4		2-Hexene, 3-methyl-, (Z)-	98	C7H14	010574-36-4	43
5		3-Methyl-2-hexene	98	C7H14	017618-77-8	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

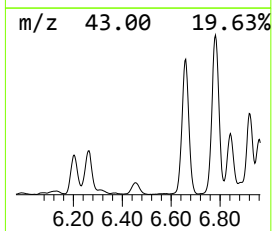
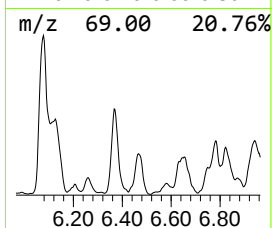
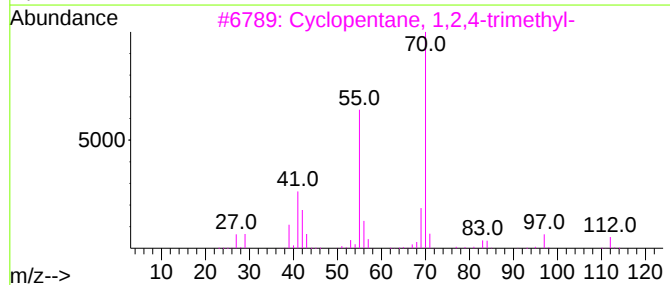
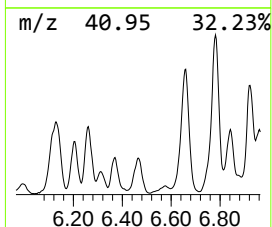
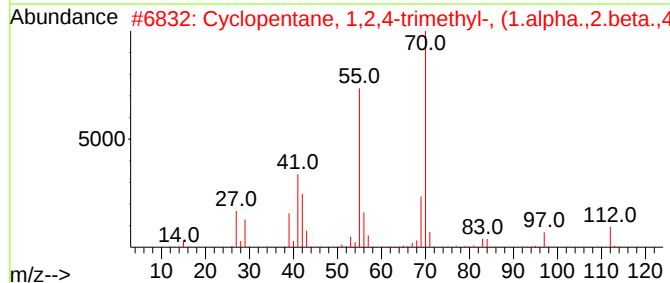
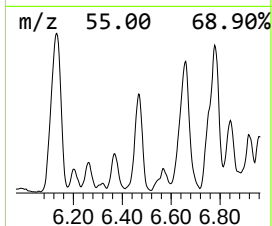
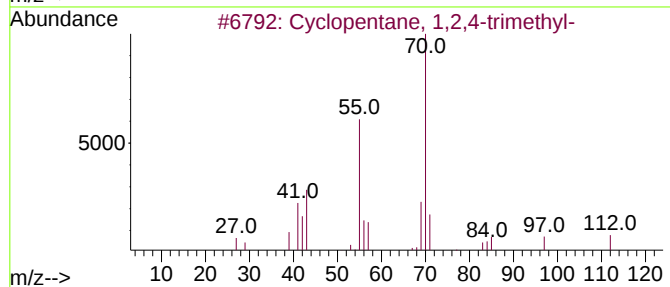
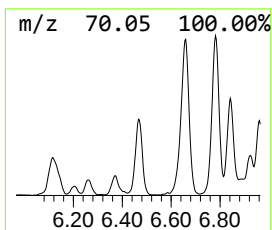
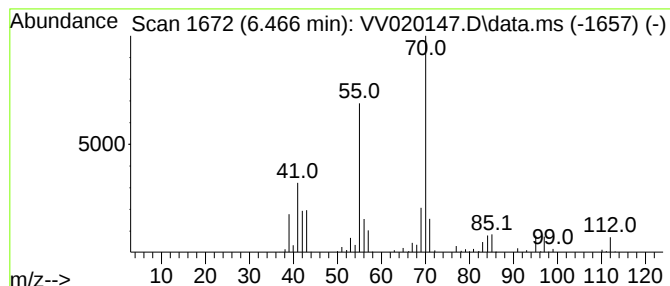
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 18 (DEL) Alkane: Cyclic6.466 Concentration Rank 41

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.466	12.90 ug/L	259150	1,4-Difluorobenzene	5.621

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, 1,2,4-trimethyl-	112	C8H16	002815-58-9	90
2			Cyclopentane, 1,2,4-trimethyl-, ...	112	C8H16	016883-48-0	87
3			Cyclopentane, 1,2,4-trimethyl-	112	C8H16	002815-58-9	83
4			Hexane, 2-methyl-4-methylene-	112	C8H16	003404-80-6	72
5			Cyclopentane, 1,2,3-trimethyl-, ...	112	C8H16	002613-69-6	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

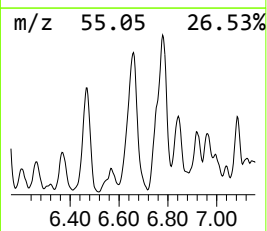
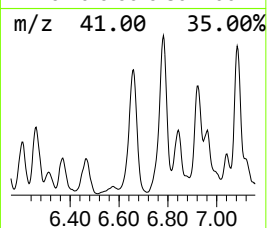
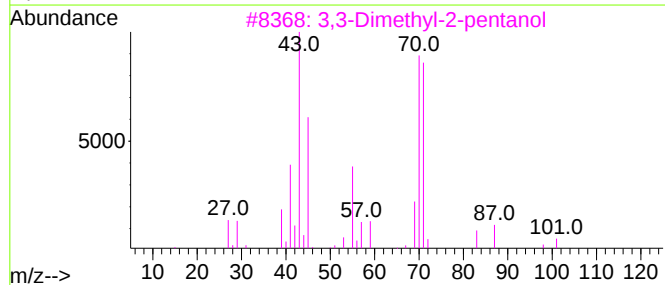
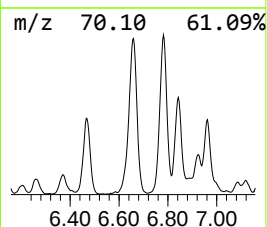
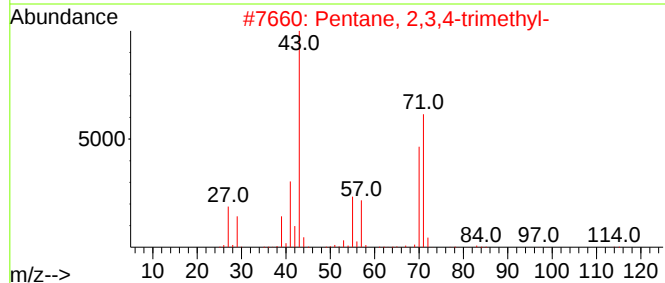
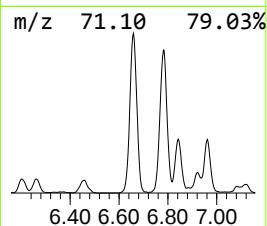
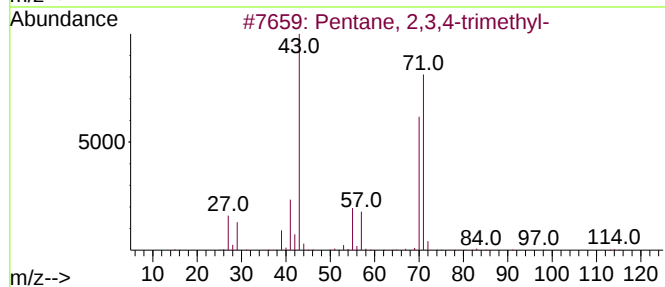
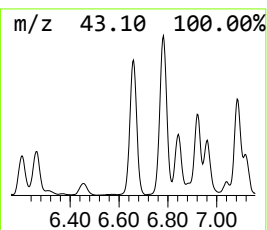
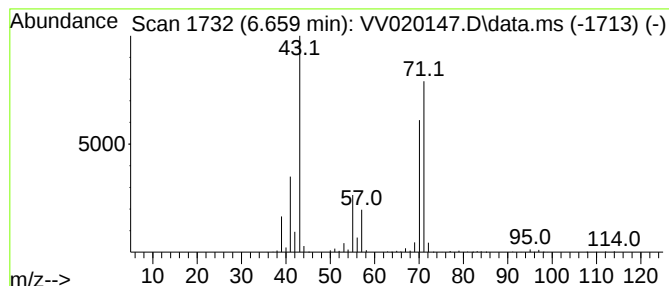
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.659	39.39 ug/L	791388	1,4-Difluorobenzene	5.621

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	91
2		Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	87
3		3,3-Dimethyl-2-pentanol	116	C7H16O	019781-24-9	78
4		Pentane, 3-ethyl-	100	C7H16	000617-78-7	78
5		Heptane, 4-methyl-	114	C8H18	000589-53-7	74



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

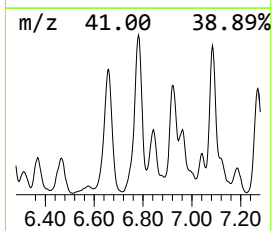
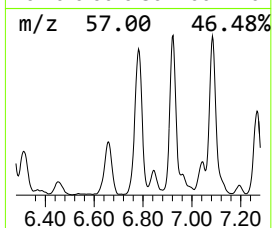
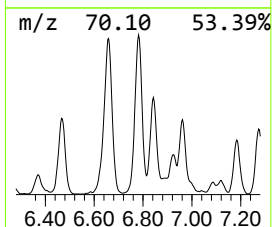
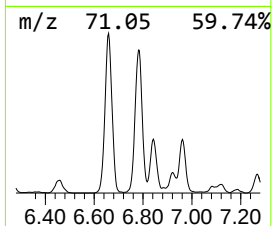
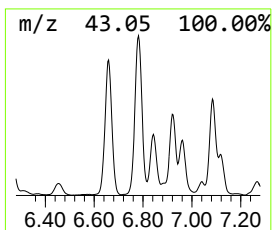
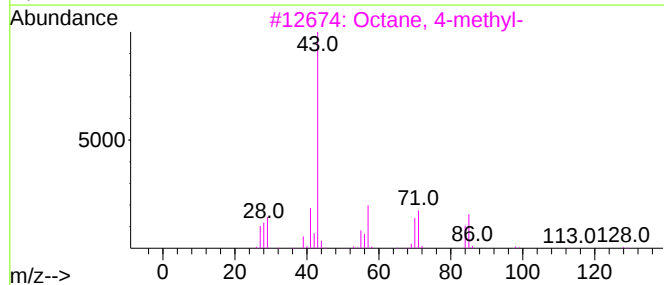
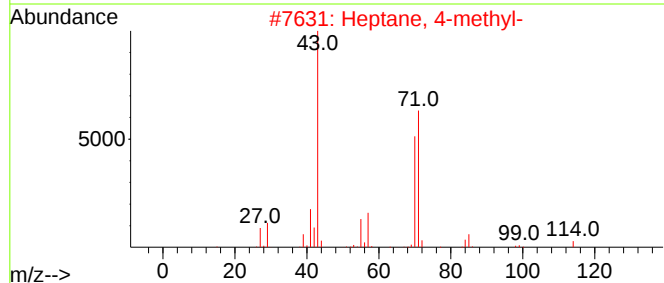
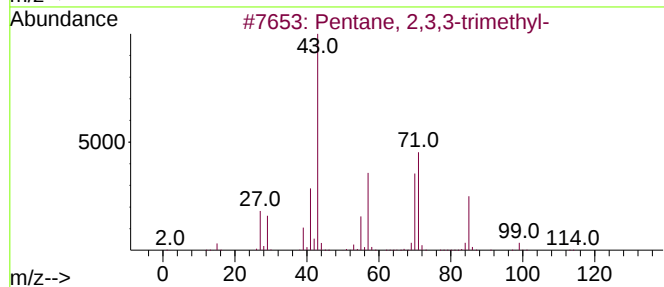
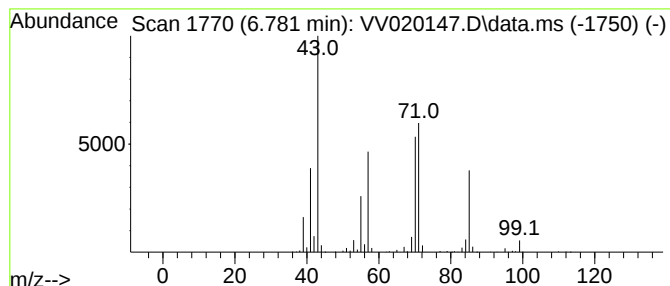
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.781	50.36 ug/L	1011710	1,4-Difluorobenzene	5.621

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	90
2		Heptane, 4-methyl-	114	C8H18	000589-53-7	64
3		Octane, 4-methyl-	128	C9H20	002216-34-4	64
4		Hexane, 3-methyl-	100	C7H16	000589-34-4	59
5		Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

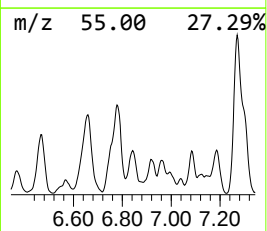
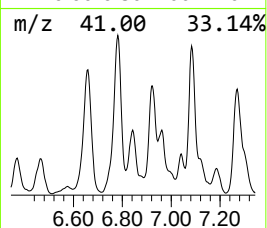
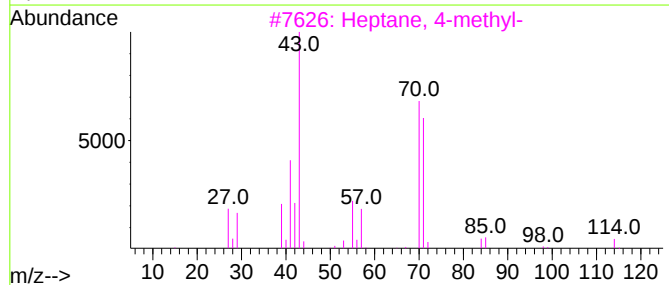
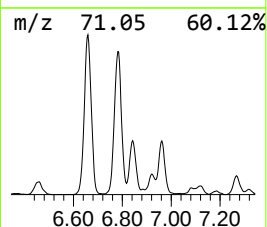
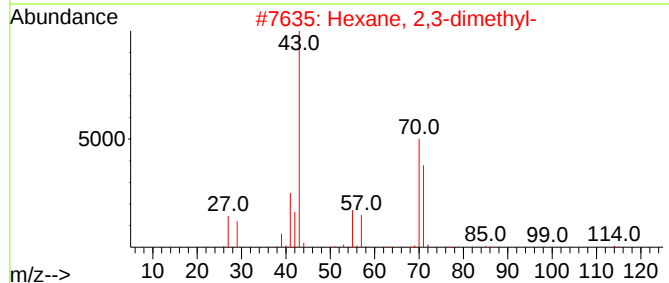
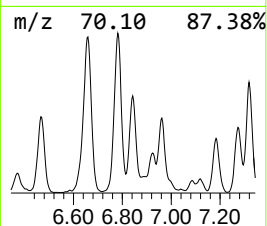
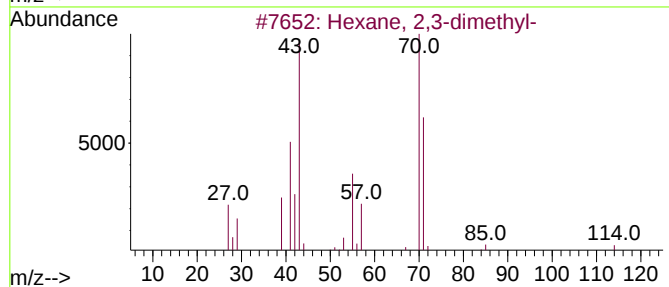
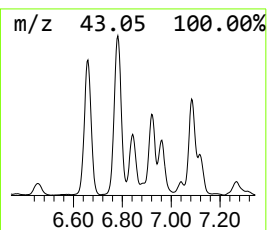
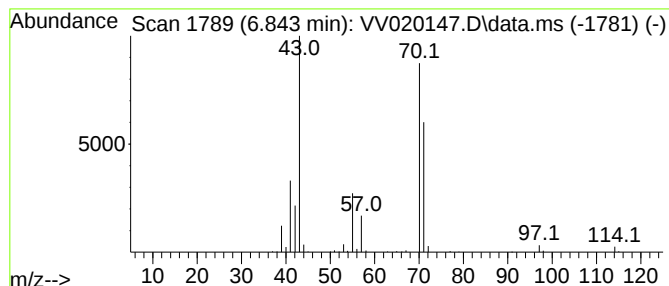
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 21 (DEL) Alkane: Straight-Chai... Concentration Rank 49

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.843	10.44 ug/L	209774	1,4-Difluorobenzene	5.621

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	87
2			Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	78
3			Heptane, 4-methyl-	114	C8H18	000589-53-7	74
4			Pyrrolidine	71	C4H9N	000123-75-1	72
5			Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	68



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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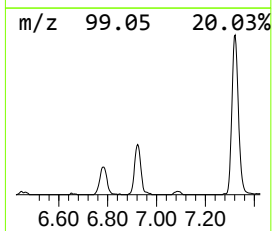
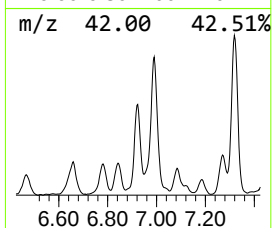
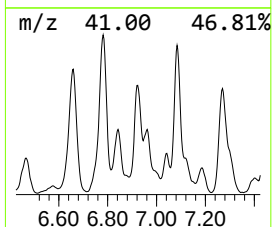
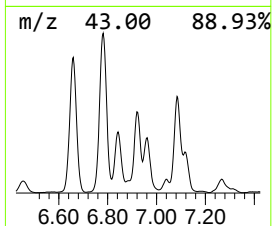
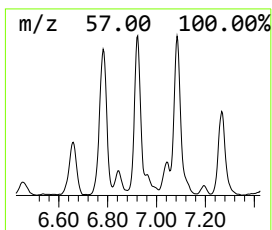
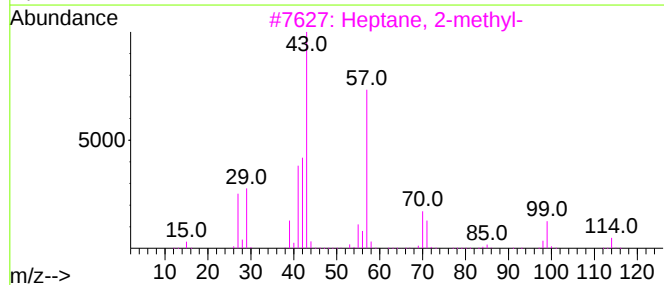
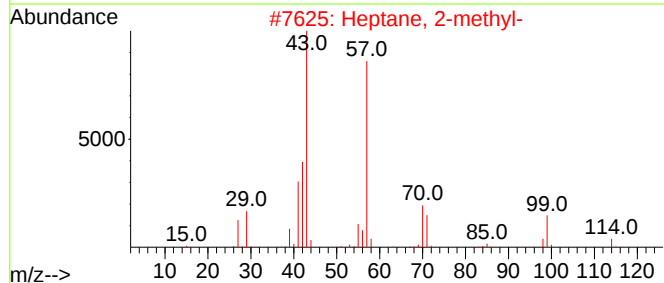
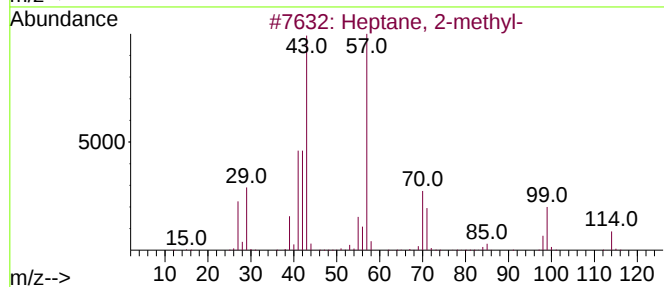
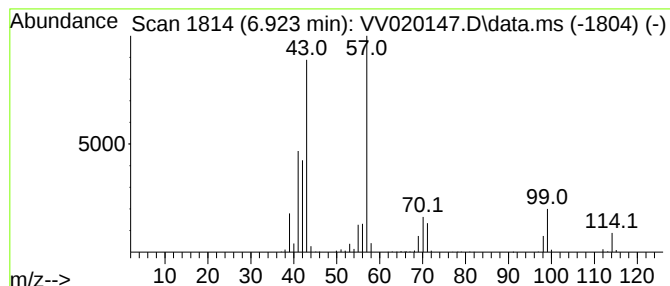
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.923	17.30 ug/L	347605	1,4-Difluorobenzene	5.621

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptane, 2-methyl-	114	C8H18	000592-27-8	80
2		Heptane, 2-methyl-	114	C8H18	000592-27-8	72
3		Heptane, 2-methyl-	114	C8H18	000592-27-8	72
4		Hexane, 2,5-dimethyl-	114	C8H18	000592-13-2	64
5		Hexane, 4-ethyl-2-methyl-	128	C9H20	003074-75-7	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
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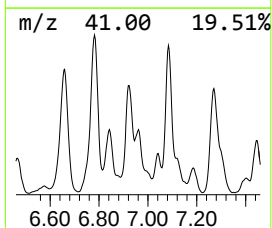
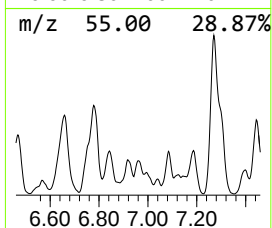
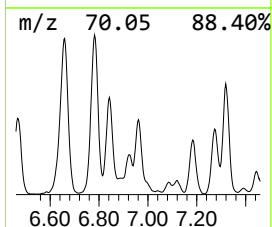
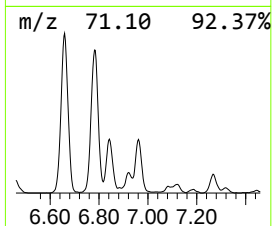
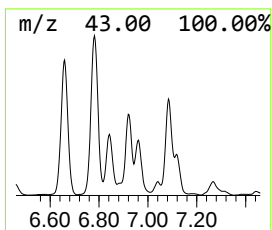
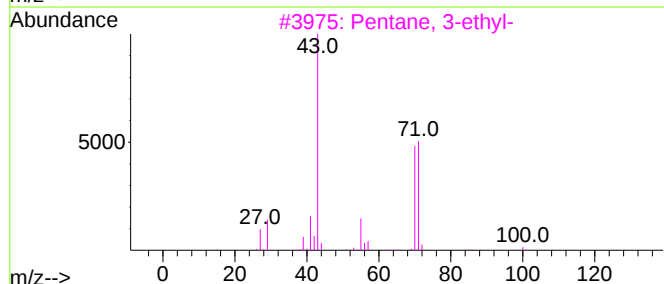
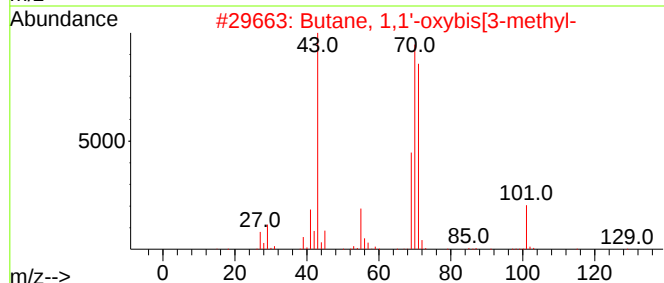
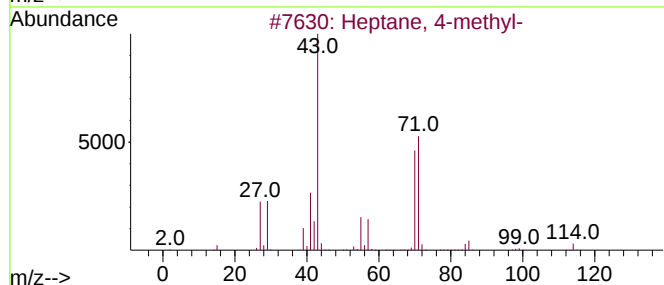
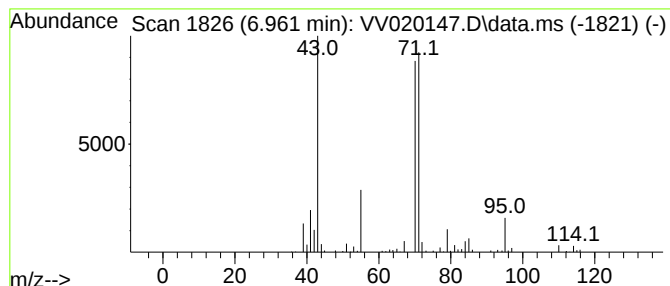
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 23 (DEL) Alkane: Straight-Chai... Concentration Rank 73

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.961	5.59 ug/L	112309	1,4-Difluorobenzene	5.621

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptane, 4-methyl-	114	C8H18	000589-53-7	86
2		Butane, 1,1'-oxybis[3-methyl-	158	C10H22O	000544-01-4	72
3		Pentane, 3-ethyl-	100	C7H16	000617-78-7	72
4		Heptane, 4-methyl-	114	C8H18	000589-53-7	58
5		Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	56



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

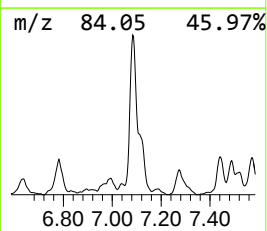
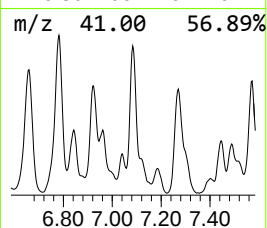
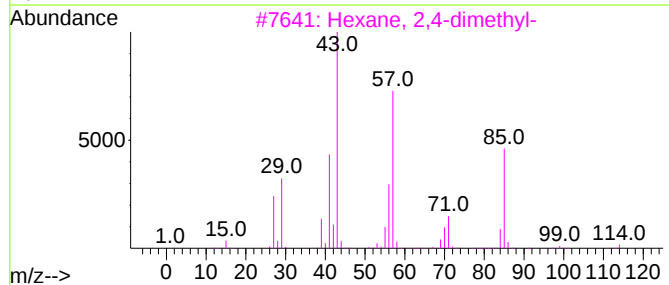
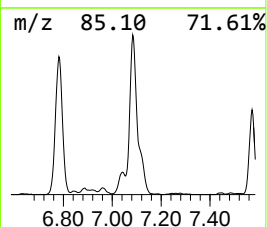
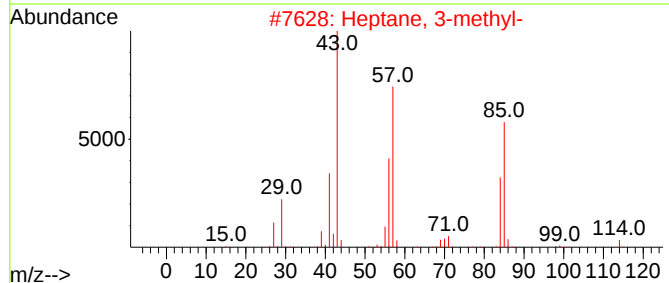
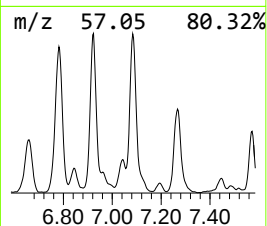
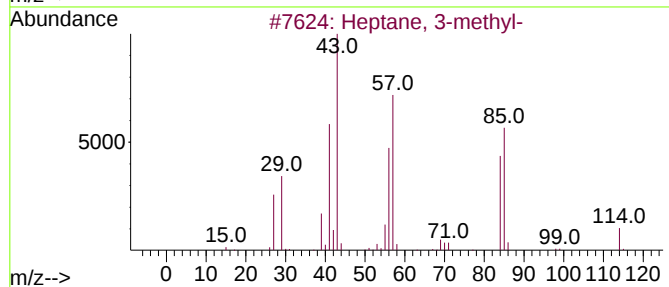
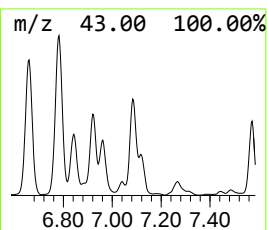
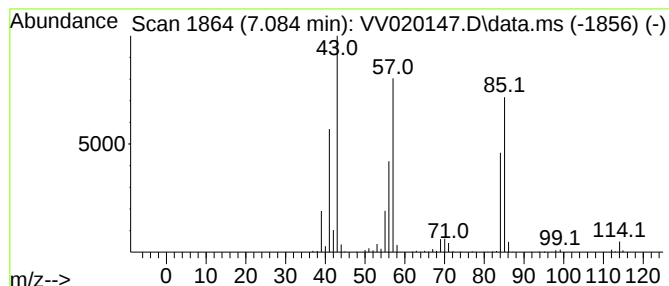
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 25 (DEL) Alkane: Straight-Chai... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.084	22.39 ug/L	449872	1,4-Difluorobenzene	5.621

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptane, 3-methyl-	114	C8H18	000589-81-1	91
2		Heptane, 3-methyl-	114	C8H18	000589-81-1	91
3		Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	64
4		Nonane, 5-methyl-	142	C10H22	015869-85-9	64
5		Hexane, 1-(hexyloxy)-2-methyl-	200	C13H28O	074421-17-3	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

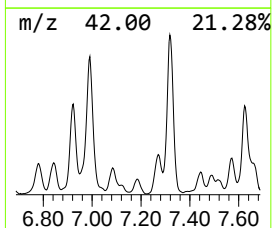
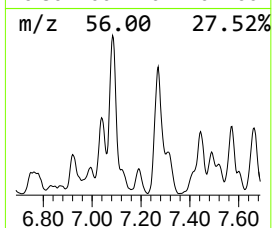
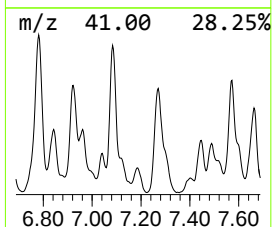
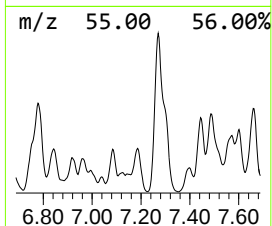
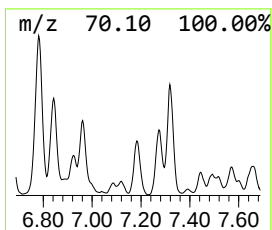
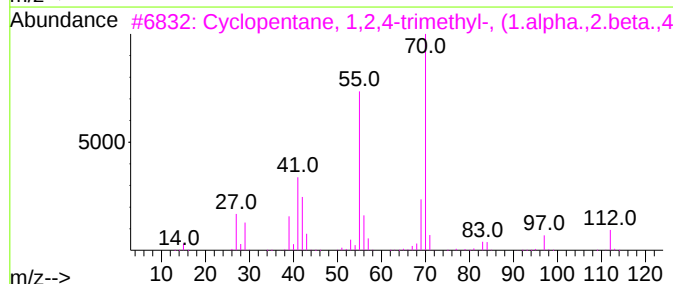
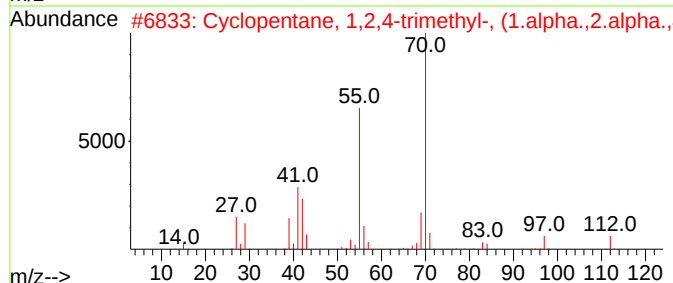
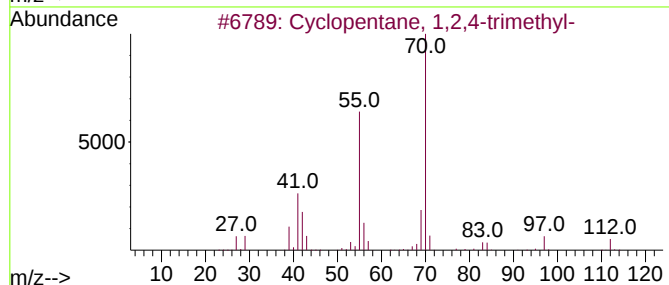
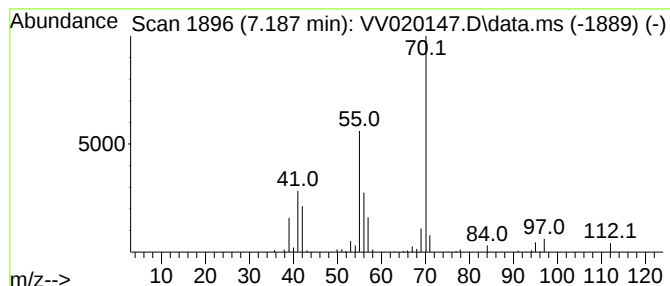
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 26 (DEL) Alkane: Cyclic7.187 Concentration Rank 66

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.187	6.83 ug/L	137286	1,4-Difluorobenzene	5.621

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, 1,2,4-trimethyl-	112	C8H16	002815-58-9	72
2			Cyclopentane, 1,2,4-trimethyl-, ...	112	C8H16	004850-28-6	64
3			Cyclopentane, 1,2,4-trimethyl-, ...	112	C8H16	016883-48-0	59
4			Cyclopentane, 1,2,3-trimethyl-, ...	112	C8H16	002613-69-6	53
5			Heptane, 3-methylene-	112	C8H16	001632-16-2	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

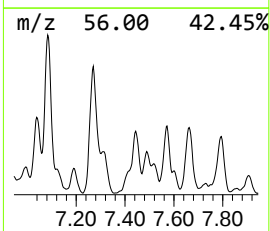
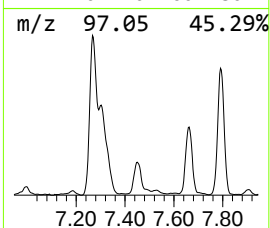
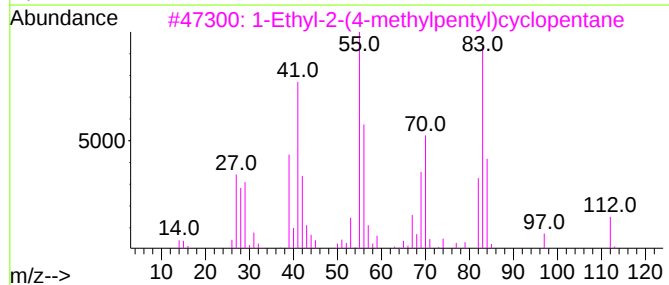
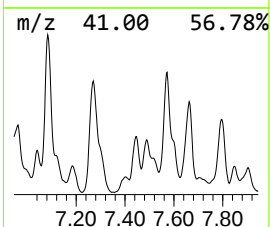
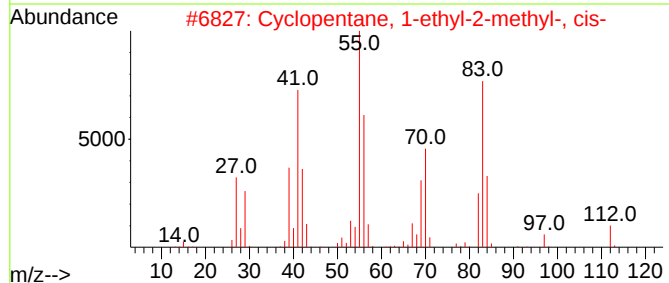
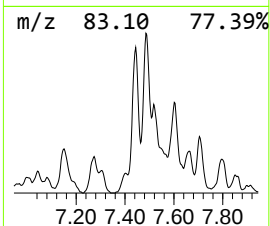
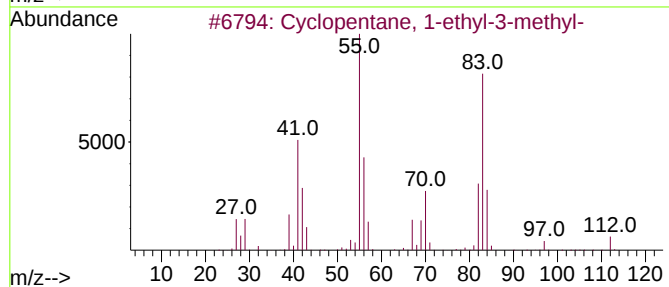
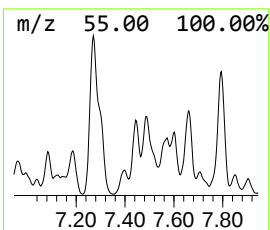
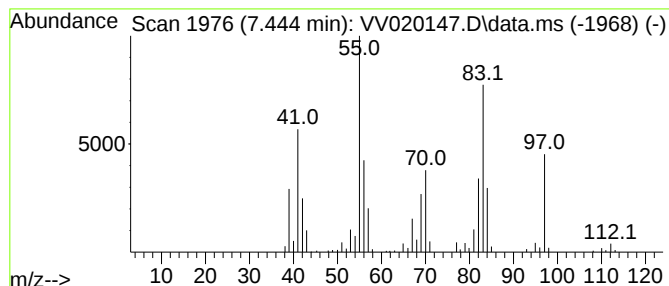
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 27 (DEL) Alkane: Cyclic7.444 Concentration Rank 57

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.444	8.24 ug/L	225164	Chlorobenzene-d5	8.858

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopentane, 1-ethyl-3-methyl-	112	C8H16	003726-47-4	87
2		Cyclopentane, 1-ethyl-2-methyl-,...	112	C8H16	000930-89-2	81
3		1-Ethyl-2-(4-methylpentyl)cyclop...	182	C13H26	219726-60-0	72
4		Cyclopentane, 1-ethyl-2-methyl-,...	112	C8H16	000930-89-2	47
5		Hexane, 3-methyl-4-methylene-	112	C8H16	003404-67-9	46



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

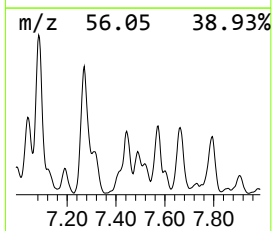
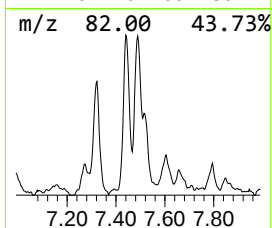
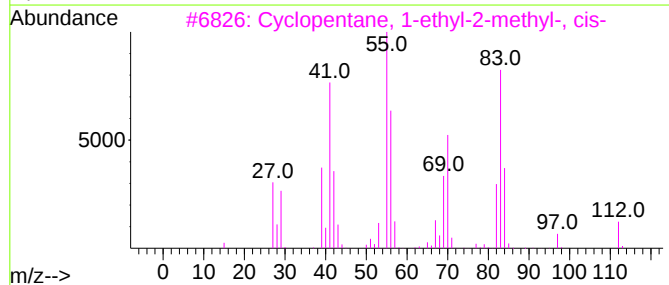
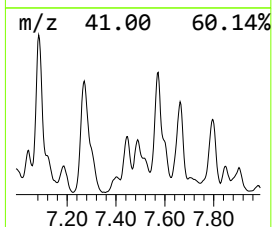
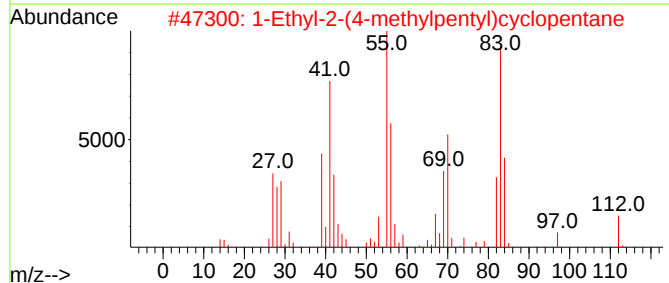
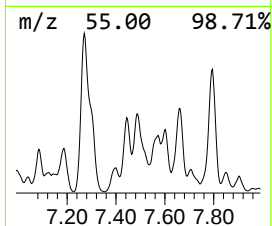
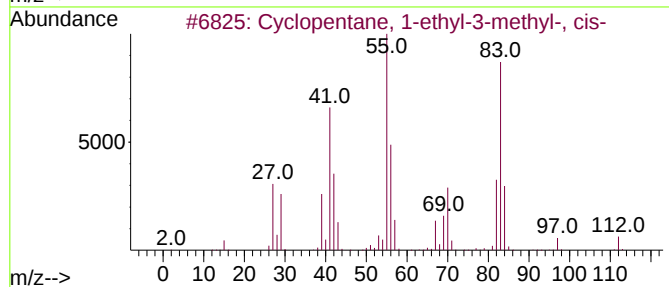
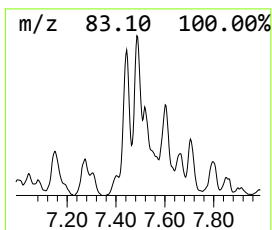
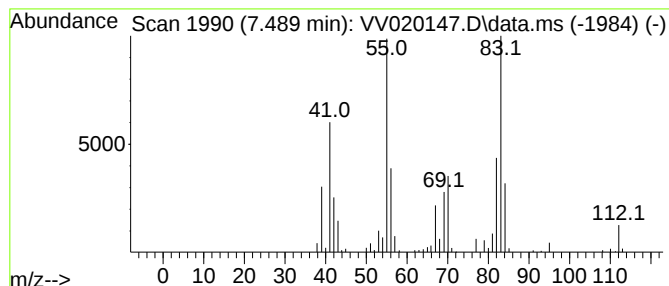
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 28 (DEL) Alkane: Cyclic7.489 Concentration Rank 50

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.489	10.36 ug/L	283227	Chlorobenzene-d5	8.858

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, 1-ethyl-3-methyl-,...	112	C8H16	002613-66-3	58
2			1-Ethyl-2-(4-methylpentyl)cyclop...	182	C13H26	219726-60-0	58
3			Cyclopentane, 1-ethyl-2-methyl-,...	112	C8H16	000930-89-2	53
4			Cyclopentane, 1-ethyl-2-methyl-,...	112	C8H16	000930-89-2	49
5			Cyclopentane, 1-ethyl-2-methyl-,...	112	C8H16	000930-89-2	49



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

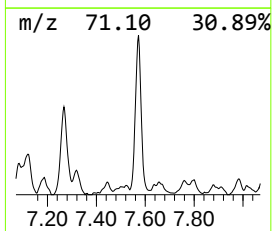
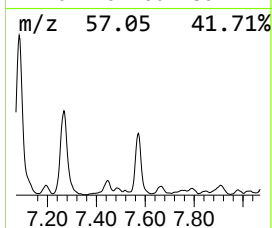
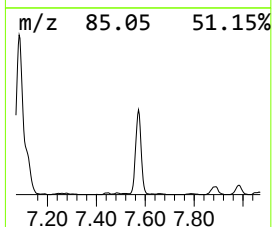
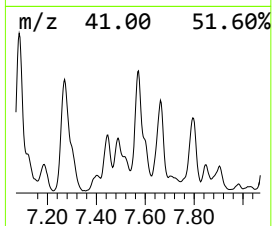
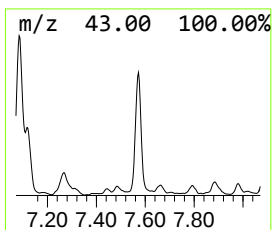
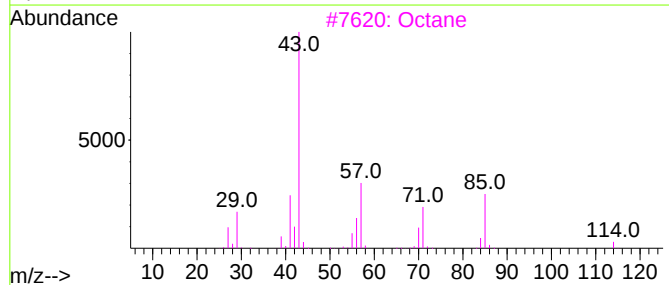
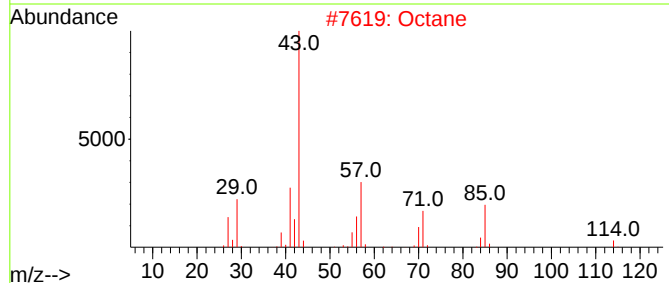
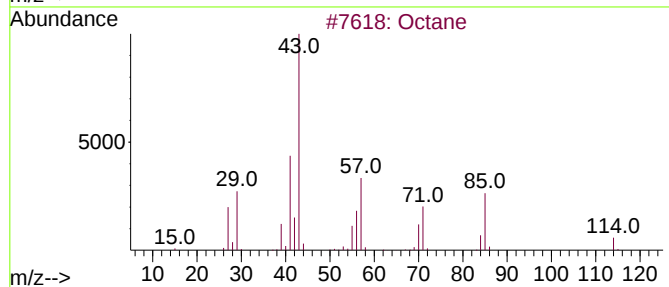
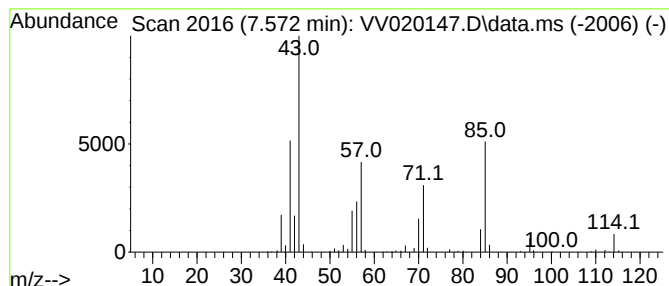
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 29 (DEL) Alkane: Straight-Chai... Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.572	18.54 ug/L	506890	Chlorobenzene-d5	8.858

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane			114	C8H18	000111-65-9	91
2	Octane			114	C8H18	000111-65-9	91
3	Octane			114	C8H18	000111-65-9	91
4	Heptane, 2,4-dimethyl-			128	C9H20	002213-23-2	72
5	Hexane, 2,4-dimethyl-			114	C8H18	000589-43-5	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

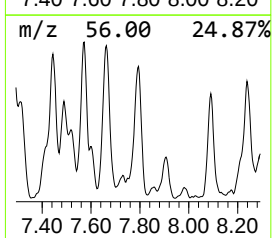
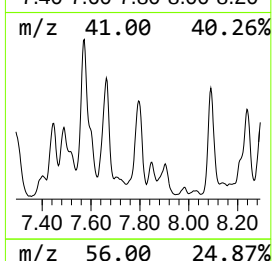
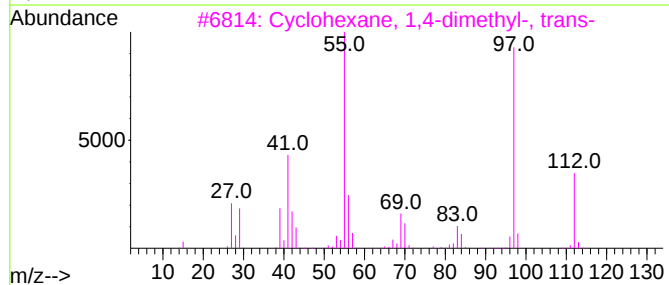
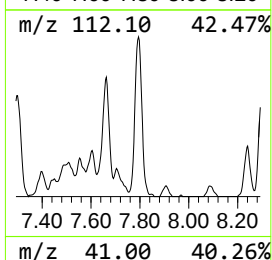
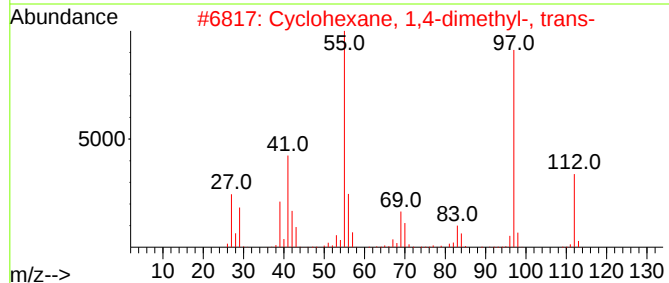
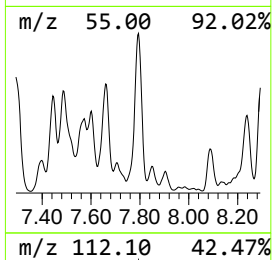
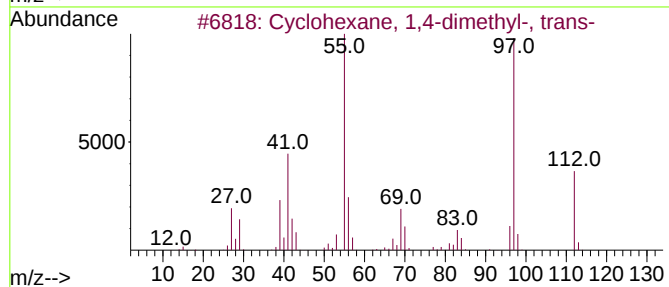
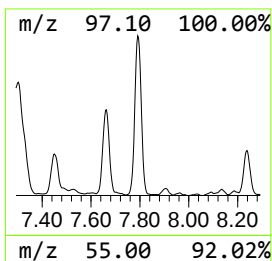
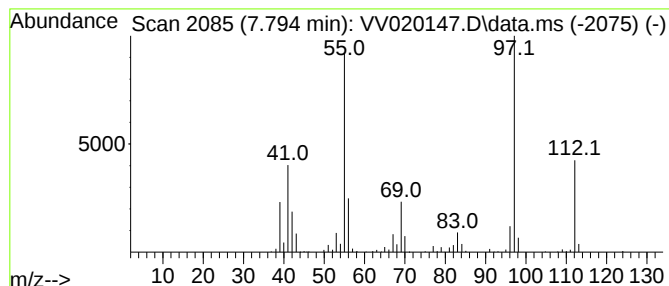
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 30 (DEL) Alkane: Cyclic7.794 Concentration Rank 45

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.794	11.39 ug/L	311274	Chlorobenzene-d5	8.858

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,4-dimethyl-, trans-	112	C8H16	002207-04-7	91
2			Cyclohexane, 1,4-dimethyl-, trans-	112	C8H16	002207-04-7	91
3			Cyclohexane, 1,4-dimethyl-, trans-	112	C8H16	002207-04-7	91
4			Cyclohexane, 1,3-dimethyl-, trans-	112	C8H16	002207-03-6	91
5			Cyclohexane, 1,3-dimethyl-, cis-	112	C8H16	000638-04-0	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

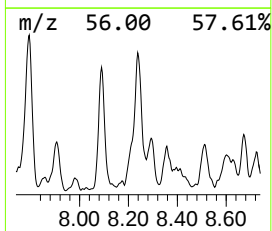
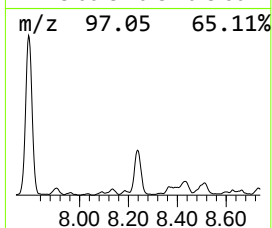
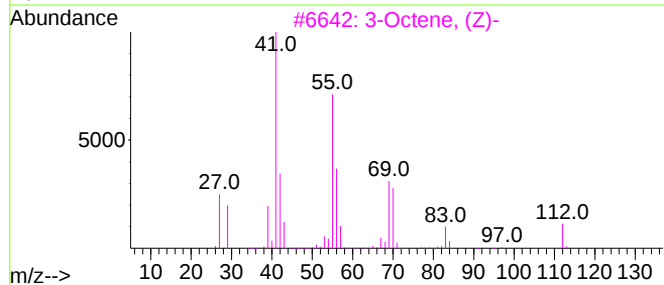
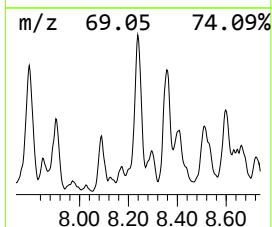
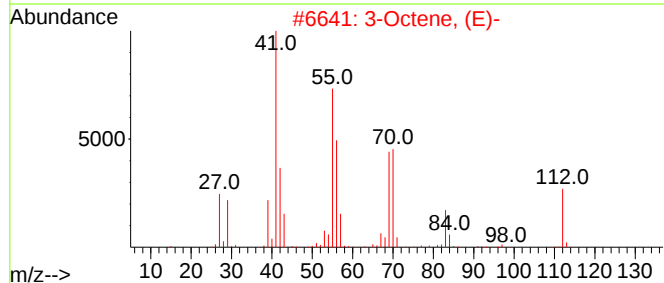
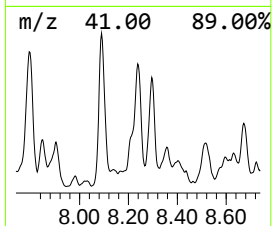
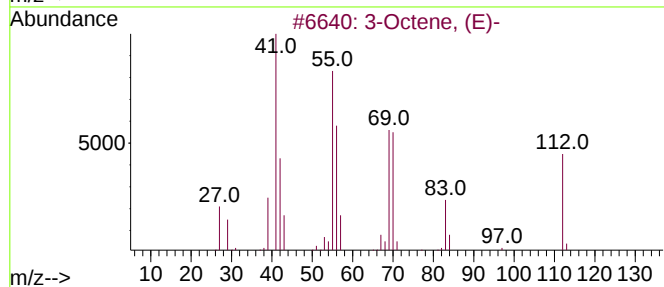
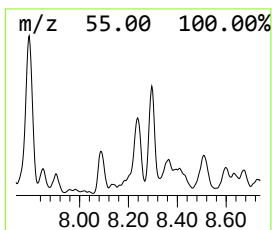
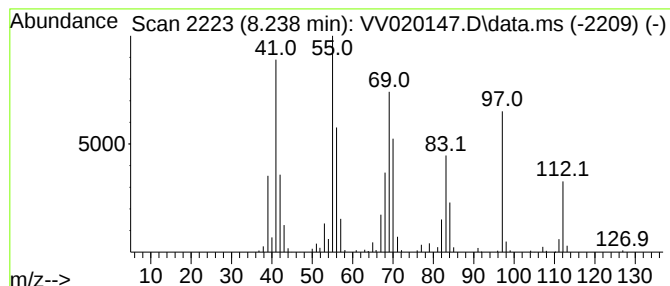
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 32 (DEL) Alkane: Straight-Chai... Concentration Rank 51

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.238	9.84 ug/L	268970	Chlorobenzene-d5	8.858

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Octene, (E)-		112	C8H16	014919-01-8	76
2	3-Octene, (E)-		112	C8H16	014919-01-8	76
3	3-Octene, (Z)-		112	C8H16	014850-22-7	70
4	3-Octene, (E)-		112	C8H16	014919-01-8	70
5	3-Octene, (Z)-		112	C8H16	014850-22-7	70



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

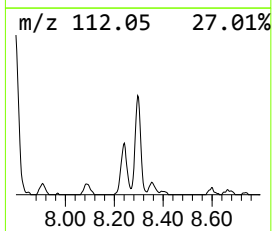
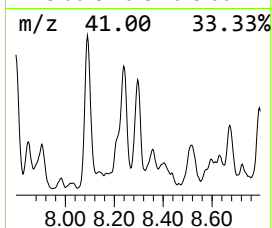
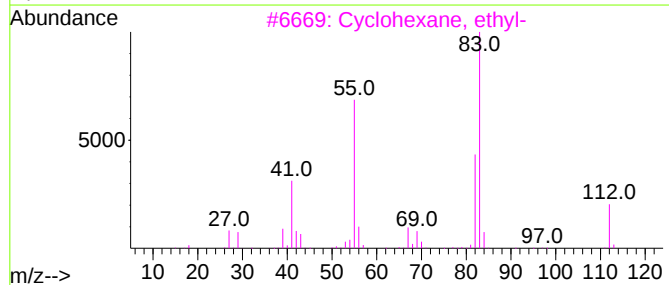
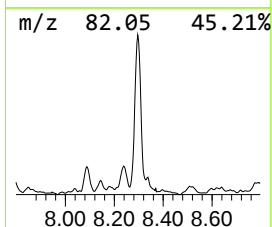
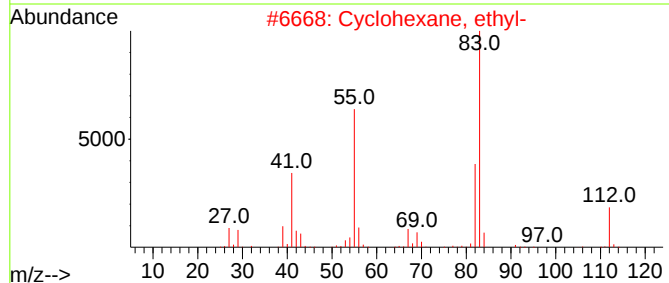
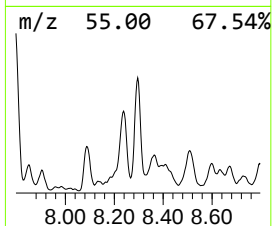
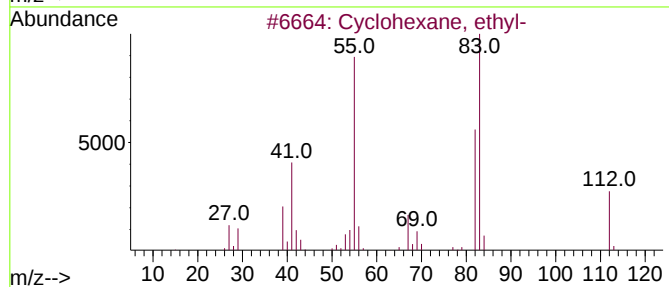
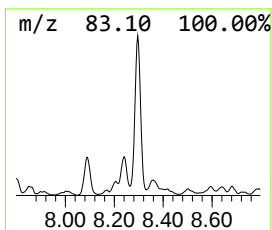
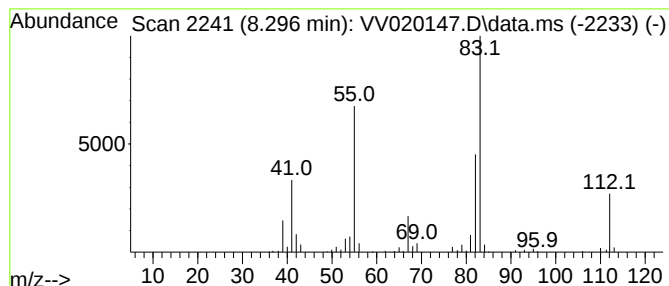
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 33 (DEL) Alkane: Cyclic8.296 Concentration Rank 68

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.296	6.64 ug/L	181567	Chlorobenzene-d5	8.858

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane, ethyl-	112	C8H16	001678-91-7	87
2		Cyclohexane, ethyl-	112	C8H16	001678-91-7	86
3		Cyclohexane, ethyl-	112	C8H16	001678-91-7	86
4		Cyclohexane, ethyl-	112	C8H16	001678-91-7	78
5		Cyclohexane, ethyl-	112	C8H16	001678-91-7	78



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

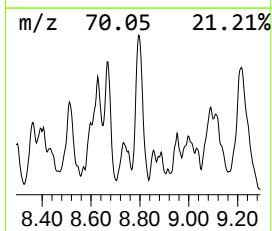
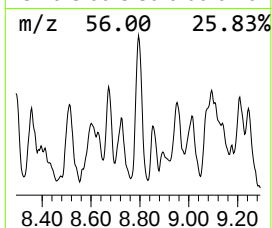
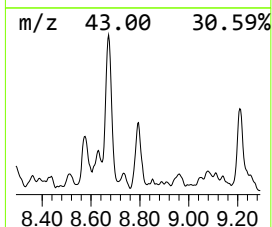
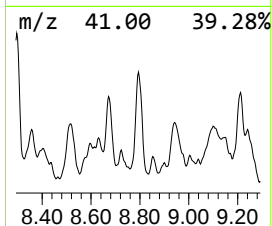
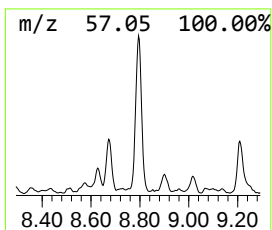
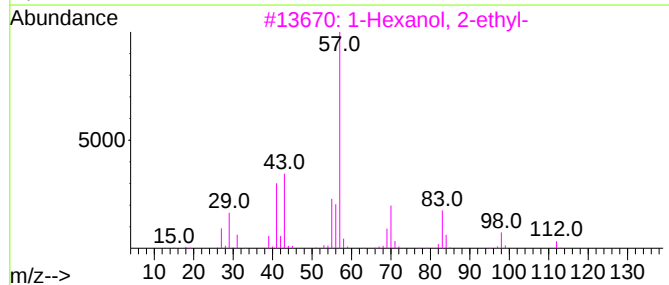
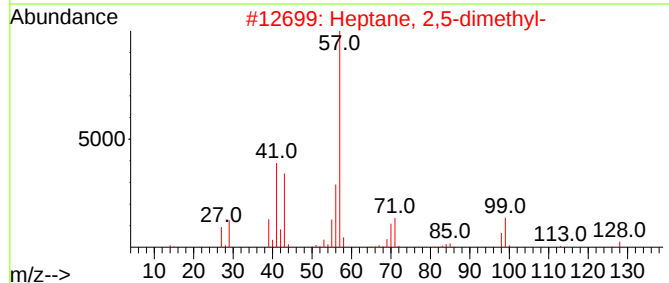
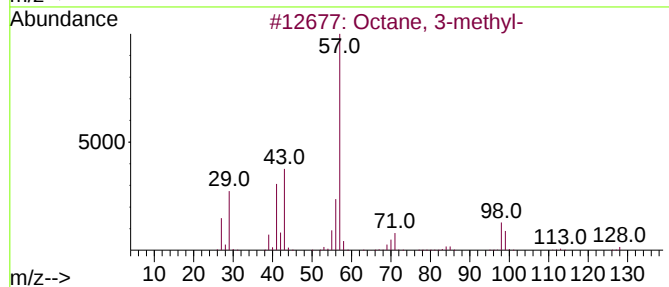
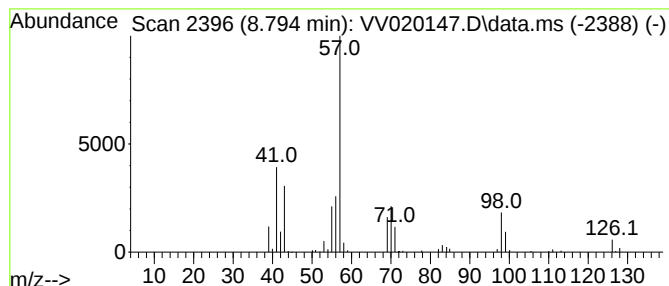
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 36 (DEL) Alkane: Straight-Chai... Concentration Rank 76

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.794	4.70 ug/L	128591	Chlorobenzene-d5	8.858

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octane, 3-methyl-	128	C9H20	002216-33-3	64
2		Heptane, 2,5-dimethyl-	128	C9H20	002216-30-0	62
3		1-Hexanol, 2-ethyl-	130	C8H18O	000104-76-7	59
4		Heptane, 1,1'-oxybis-	214	C14H30O	000629-64-1	59
5		Heptane, 3-ethyl-2-methyl-	142	C10H22	014676-29-0	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

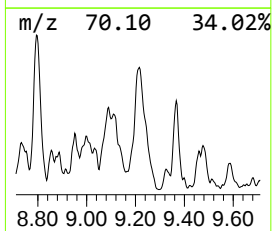
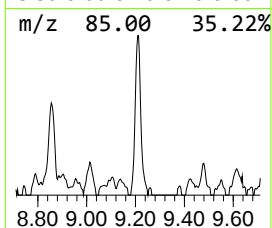
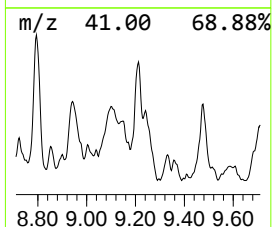
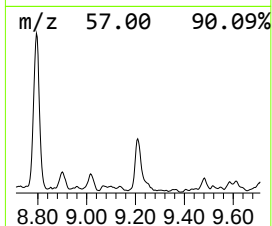
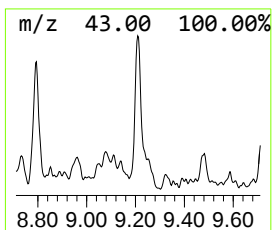
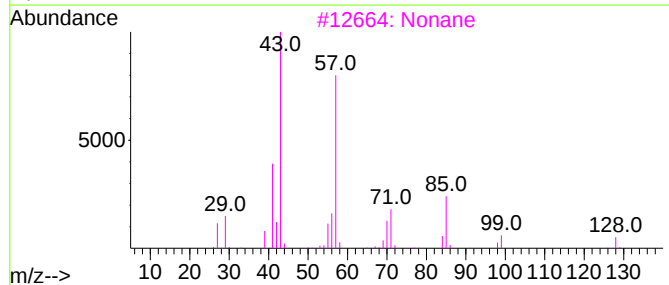
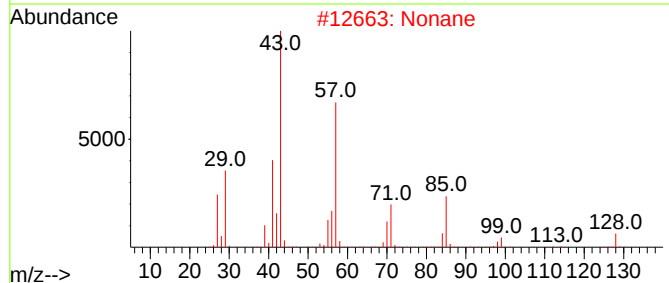
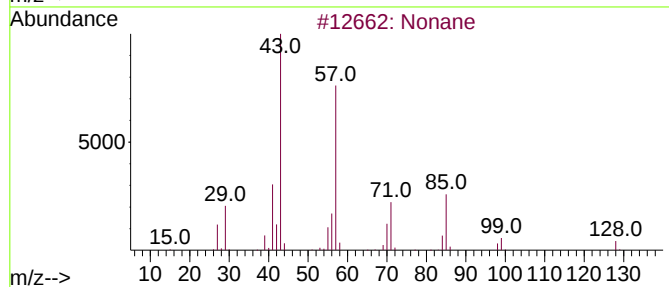
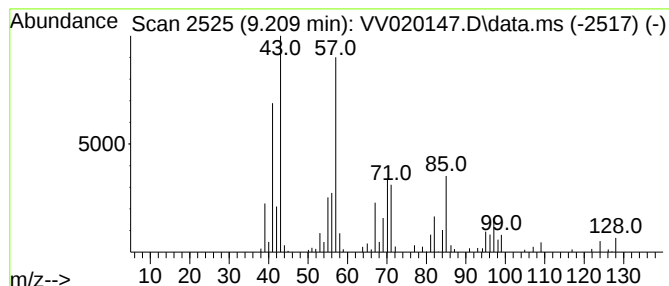
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 38 (DEL) Alkane: Straight-Chai... Concentration Rank 56

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.209	8.32 ug/L	227449	Chlorobenzene-d5	8.858

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane	128	C9H20	000111-84-2	70
2			Nonane	128	C9H20	000111-84-2	70
3			Nonane	128	C9H20	000111-84-2	49
4			Ether, heptyl hexyl	200	C13H28O	007289-40-9	43
5			Heptane, 2,4,6-trimethyl-	142	C10H22	002613-61-8	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

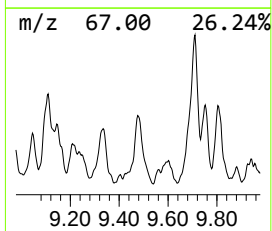
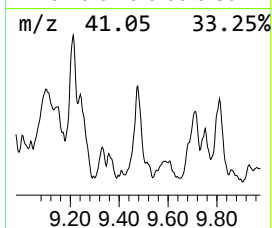
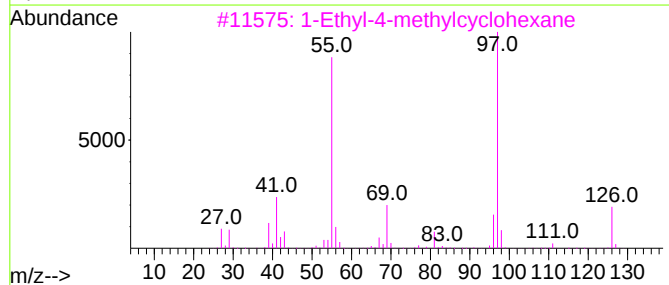
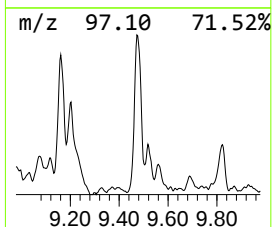
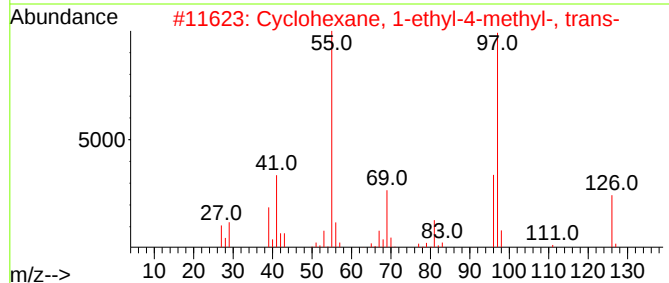
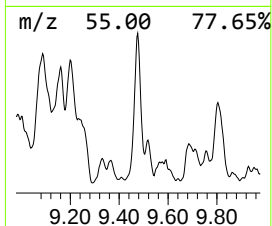
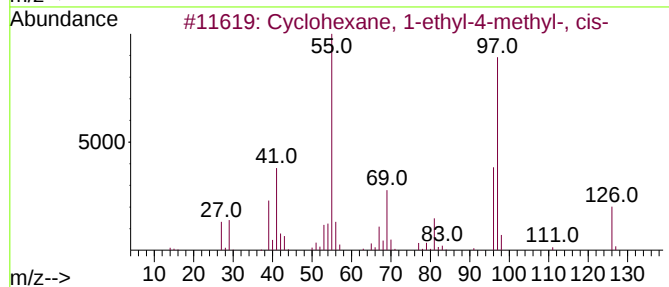
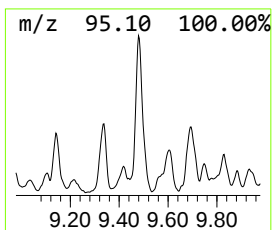
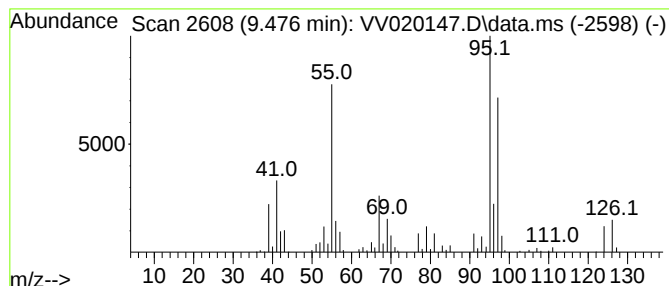
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 39 (DEL) Alkane: Cyclic9.476 Concentration Rank 62

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.476	7.32 ug/L	199951	Chlorobenzene-d5	8.858

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	004926-78-7	46
2			Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	006236-88-0	46
3			1-Ethyl-4-methylcyclohexane	126	C9H18	003728-56-1	46
4			Furan, 4-methyl-2-propyl-	124	C8H12O	006148-37-4	43
5			Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-78-8	42



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

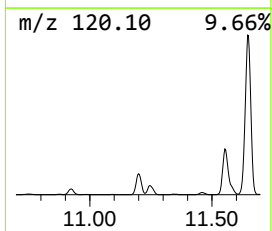
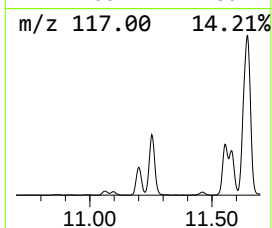
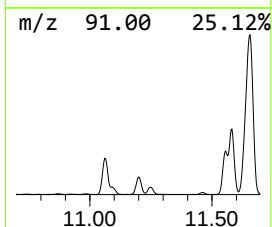
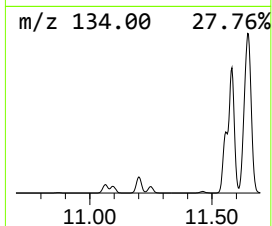
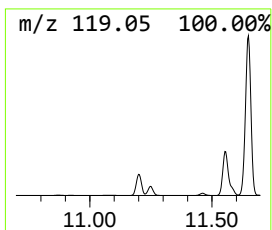
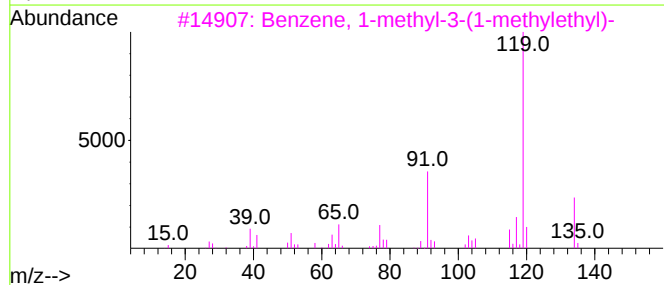
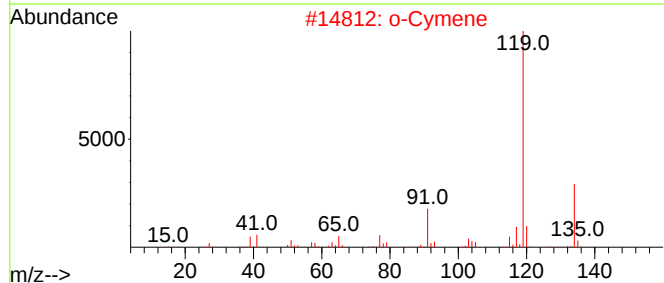
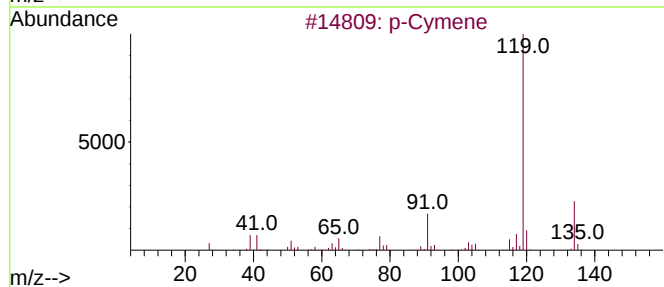
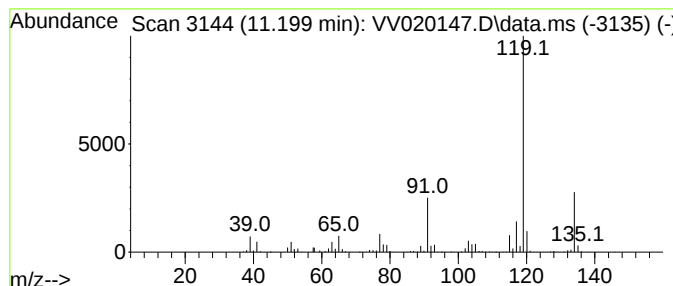
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 44 p-Cymene Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.199	14.81 ug/L	484858	1,4-Dichlorobenzene-d4	11.254

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			p-Cymene	134	C10H14	000099-87-6	97
2			o-Cymene	134	C10H14	000527-84-4	95
3			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
4			p-Cymene	134	C10H14	000099-87-6	95
5			o-Cymene	134	C10H14	000527-84-4	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

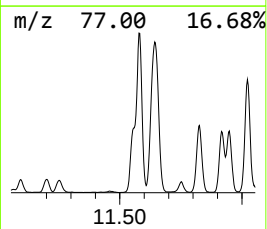
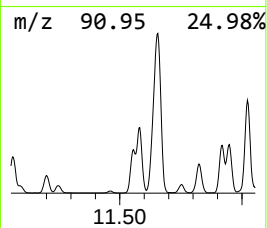
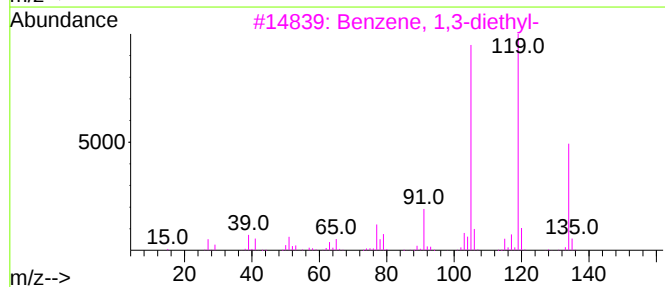
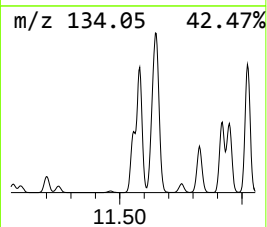
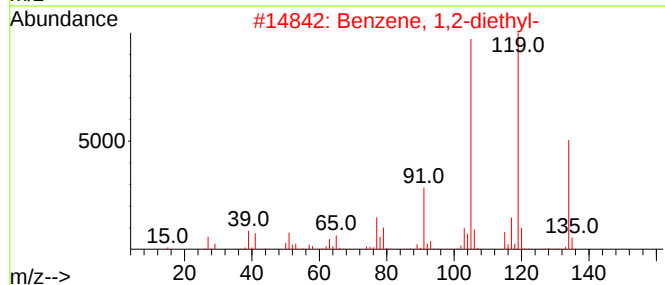
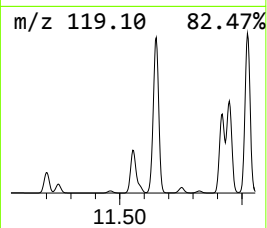
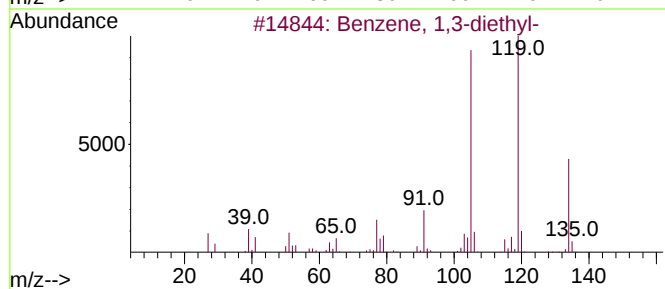
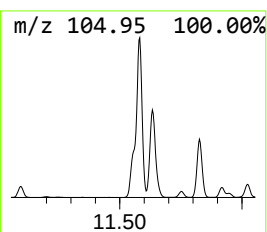
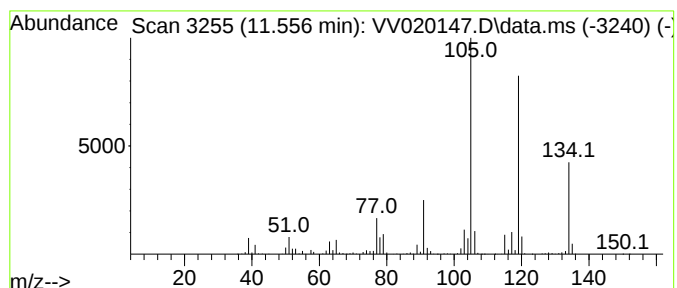
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 45 Benzene, 1,3-diethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.556	49.05 ug/L	1605360	1,4-Dichlorobenzene-d4	11.254

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	96
2			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
3			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	95
4			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	94
5			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

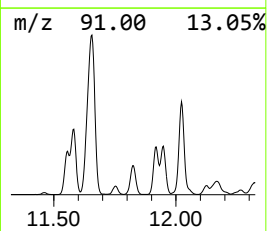
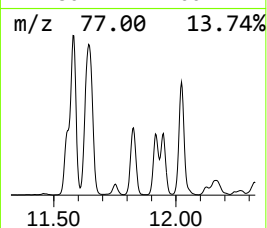
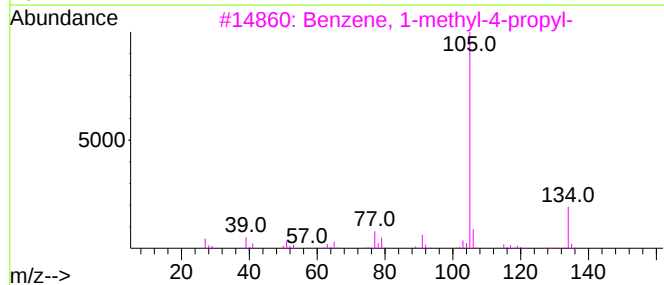
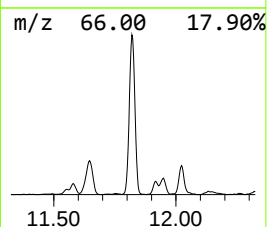
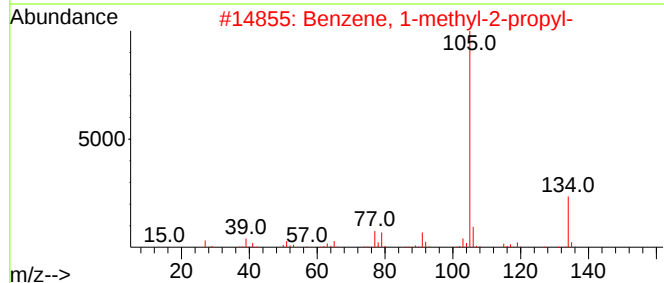
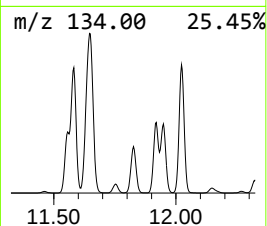
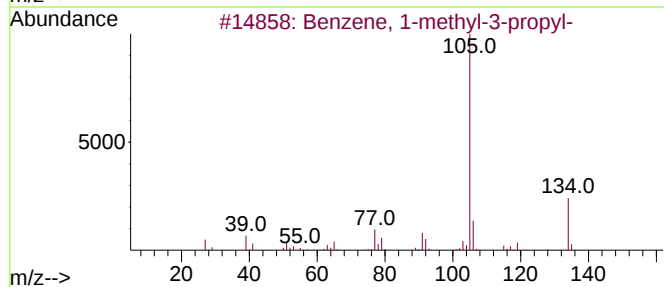
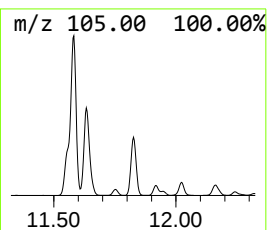
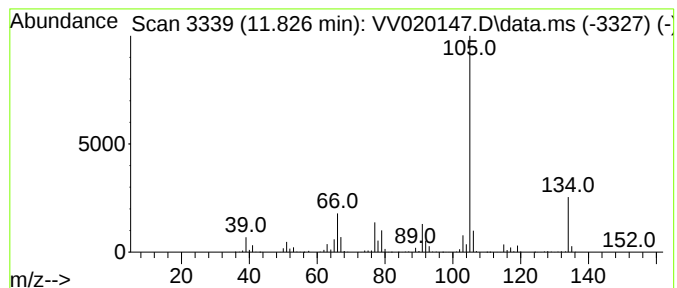
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 47 Benzene, 1-methyl-3-propyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.826	48.96 ug/L	1602390	1,4-Dichlorobenzene-d4	11.254

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	90
2			Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	90
3			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	87
4			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	87
5			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

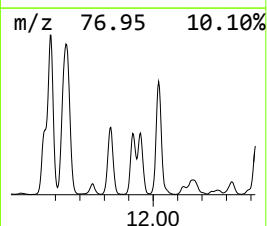
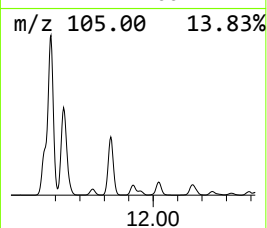
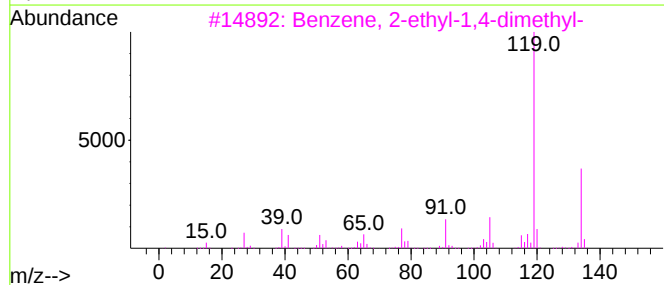
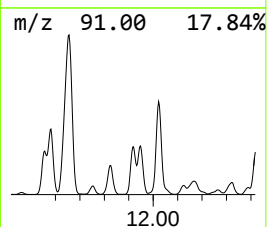
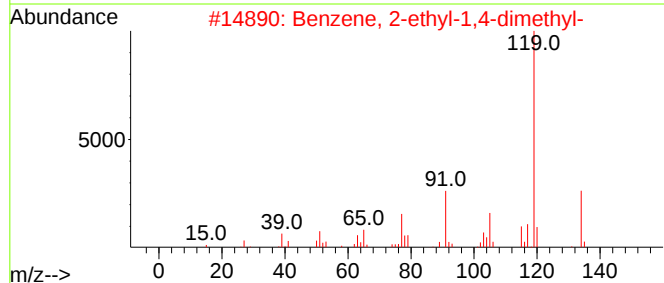
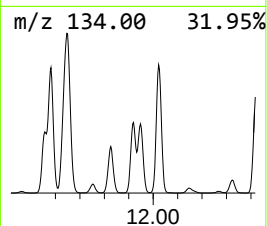
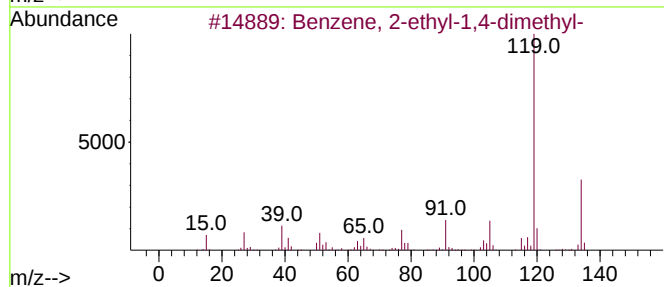
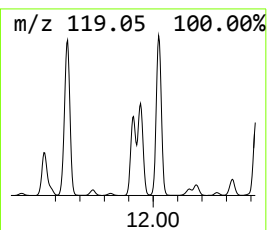
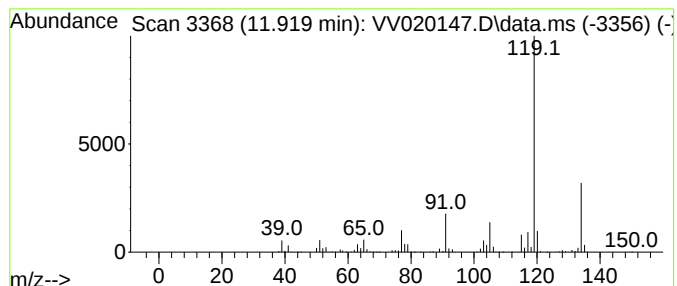
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 48 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.919	56.95 ug/L	1863900	1,4-Dichlorobenzene-d4	11.254

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	97
2			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
3			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
4			o-Cymene	134	C10H14	000527-84-4	95
5			o-Cymene	134	C10H14	000527-84-4	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

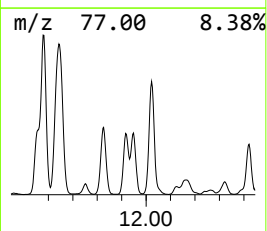
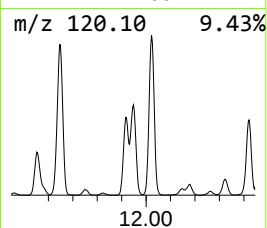
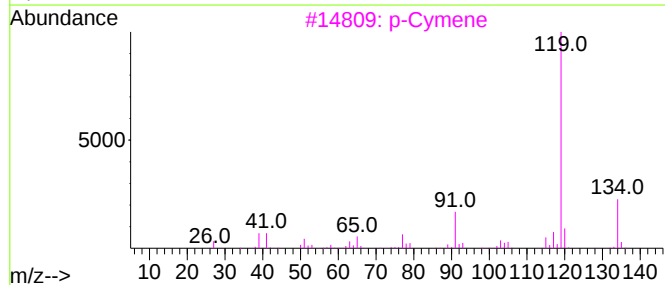
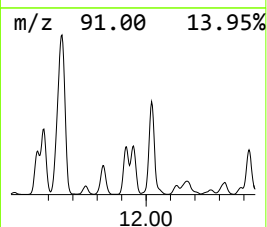
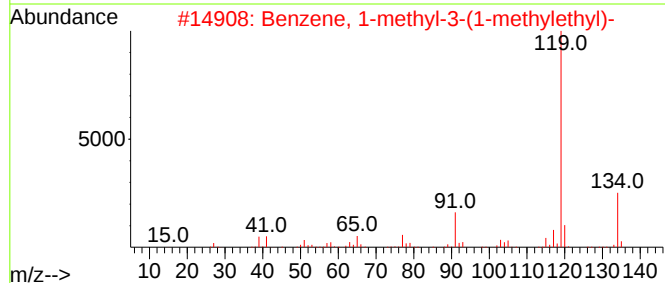
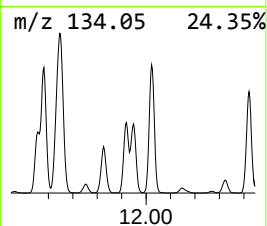
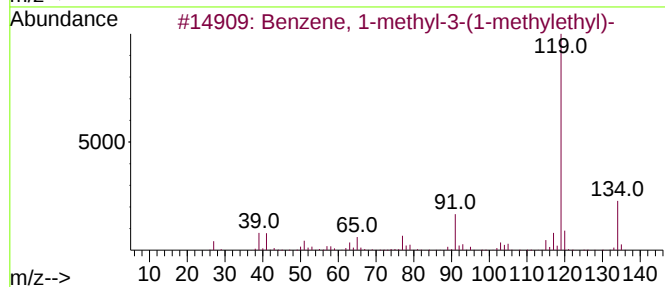
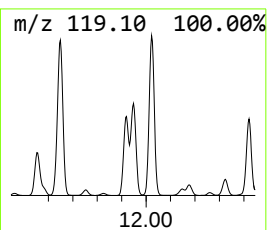
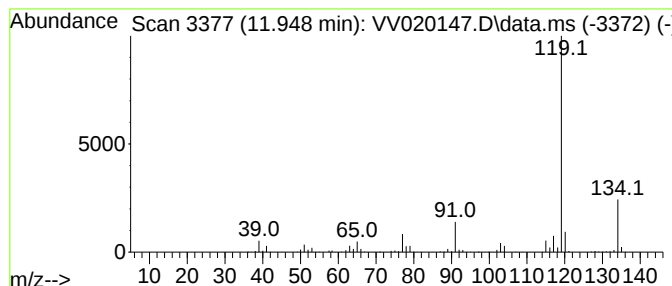
Instrument :
MSVOA_V
ClientSampleId :
BG501

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM012121WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 49 Benzene, 1-methyl-3-(1-meth... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD			R.T.
11.948	55.67 ug/L	1822060	1,4-Dichlorobenzene-d4			11.254
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-3-(1-methyleth...		134	C10H14	000535-77-3	94
2	Benzene, 1-methyl-3-(1-methyleth...		134	C10H14	000535-77-3	94
3	p-Cymene		134	C10H14	000099-87-6	94
4	Benzene, 2-ethyl-1,3-dimethyl-		134	C10H14	002870-04-4	91
5	1,3-Cyclopentadiene, 1,2,3,4-tet...		134	C10H14	076089-59-3	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

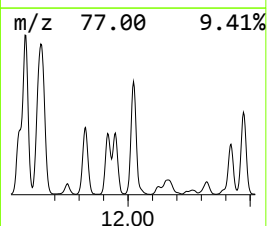
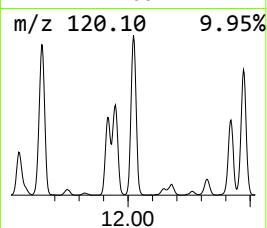
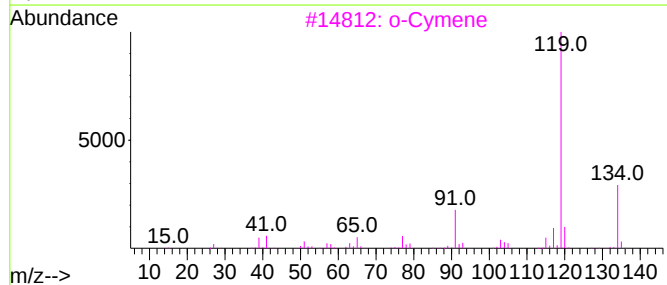
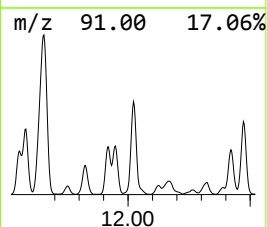
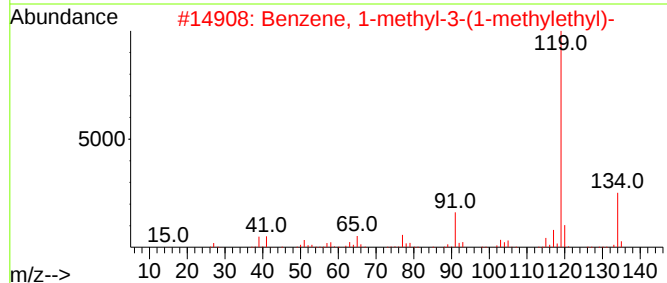
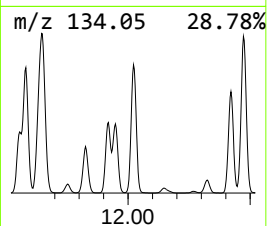
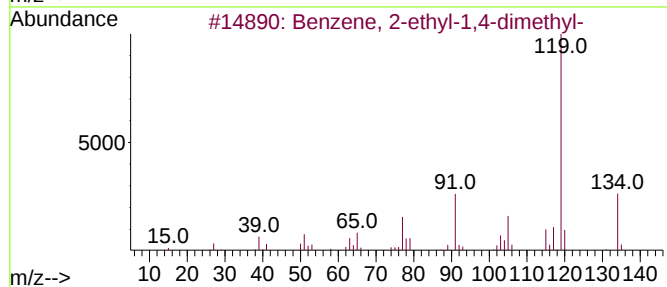
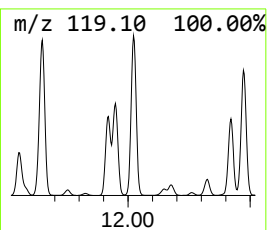
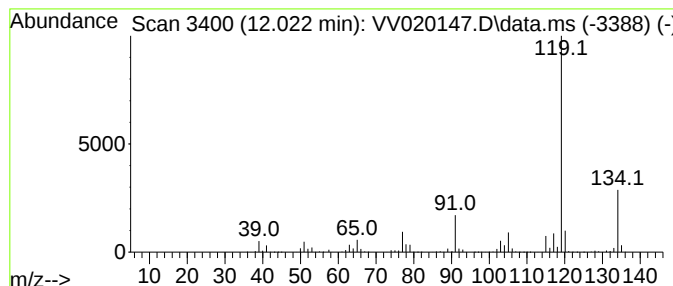
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 50 o-Cymene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.022	113.90 ug/L	3727580	1,4-Dichlorobenzene-d4	11.254

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2			Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	95
3			o-Cymene	134	C10H14	000527-84-4	95
4			o-Cymene	134	C10H14	000527-84-4	95
5			o-Cymene	134	C10H14	000527-84-4	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

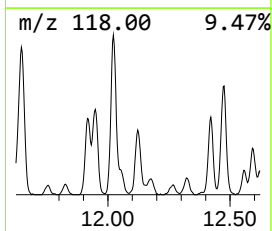
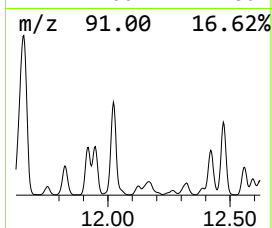
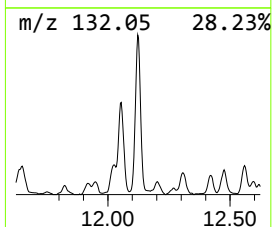
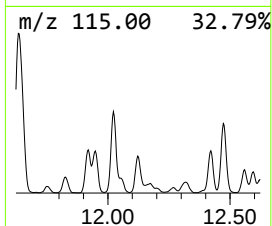
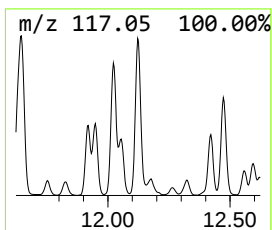
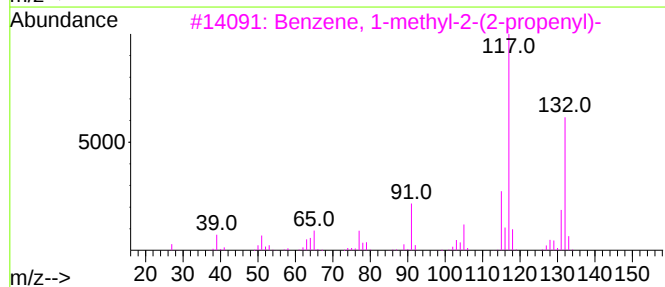
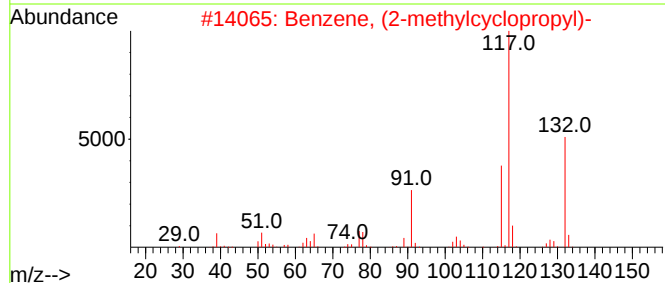
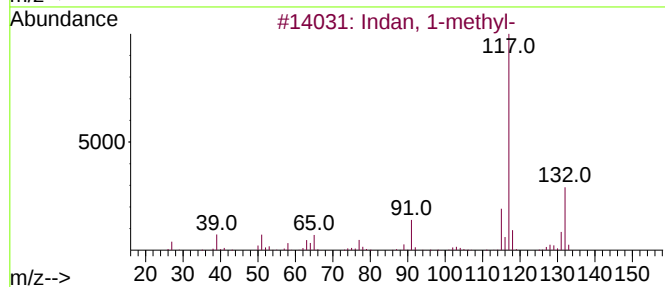
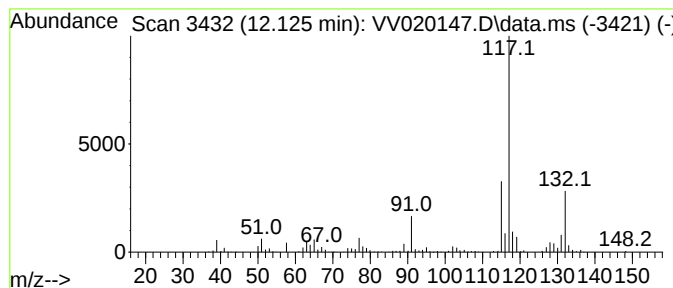
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 51 Indan, 1-methyl- Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.125	14.51 ug/L	475030	1,4-Dichlorobenzene-d4	11.254

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indan, 1-methyl-	132	C10H12	000767-58-8	87
2			Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	87
3			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	87
4			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
5			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

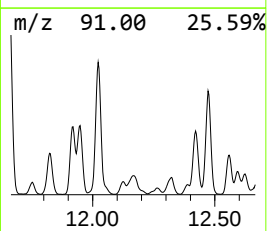
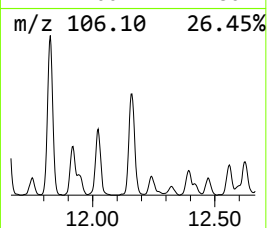
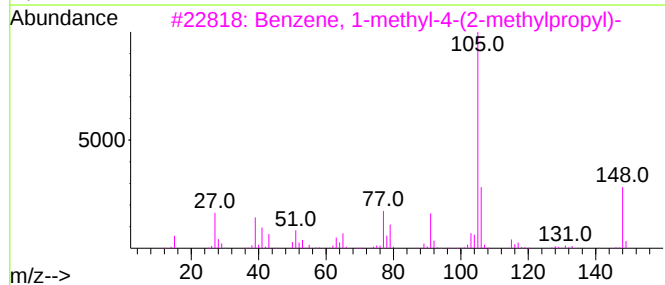
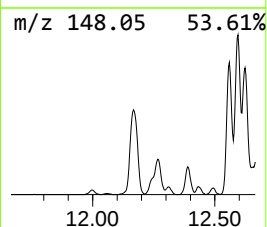
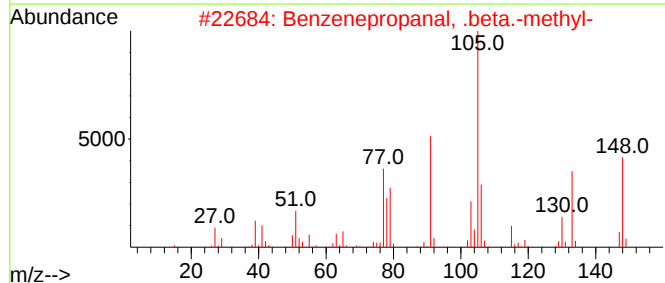
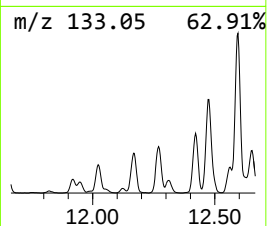
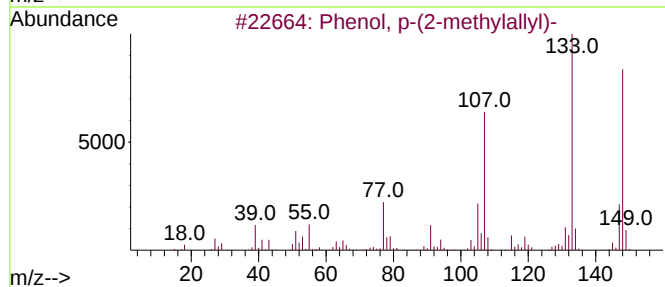
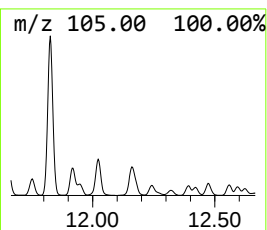
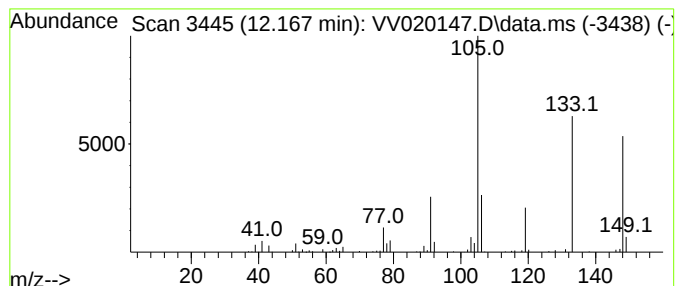
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 52 Phenol, p-(2-methylallyl)- Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.167	20.89 ug/L	683698	1,4-Dichlorobenzene-d4	11.254

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, p-(2-methylallyl)-	148	C10H12O	033641-78-0	50
2			Benzenepropanal, .beta.-methyl-	148	C10H12O	016251-77-7	50
3			Benzene, 1-methyl-4-(2-methylpro...	148	C11H16	005161-04-6	50
4			6-Methyl-4-indanol	148	C10H12O	020294-32-0	50
5			2-Methylindan-2-ol	148	C10H12O	033223-84-6	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

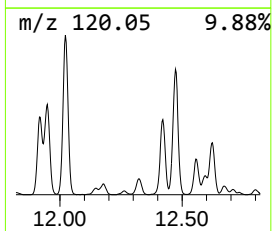
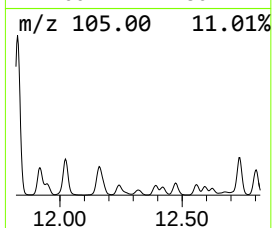
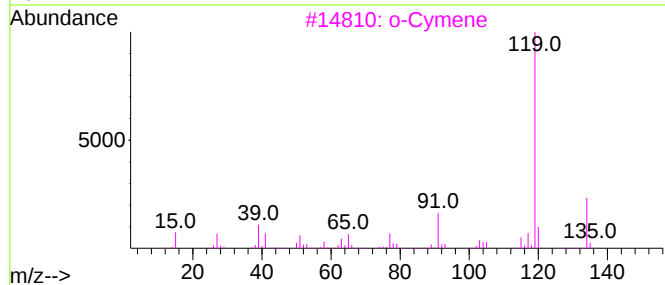
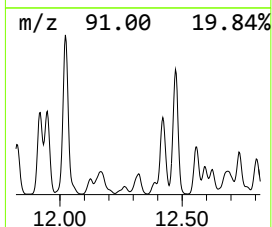
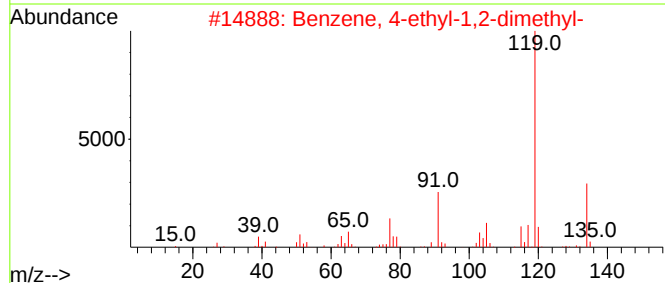
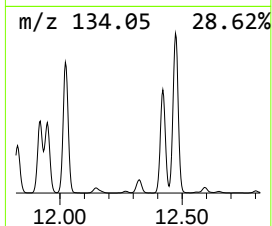
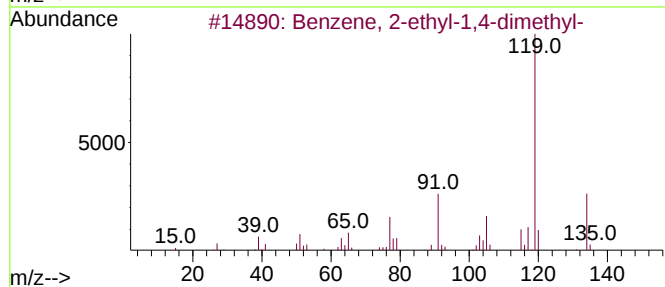
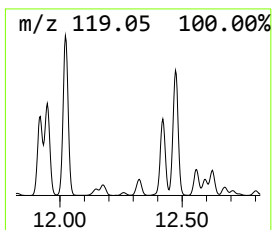
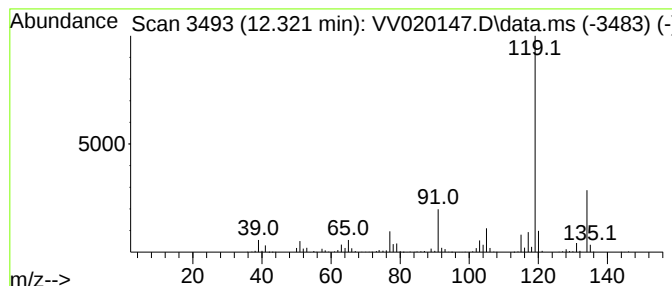
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 54 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.321	15.91 ug/L	520829	1,4-Dichlorobenzene-d4	11.254

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			o-Cymene	134	C10H14	000527-84-4	95
4			o-Cymene	134	C10H14	000527-84-4	95
5			p-Cymene	134	C10H14	000099-87-6	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

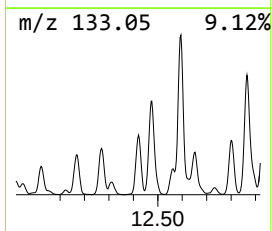
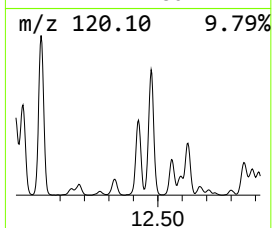
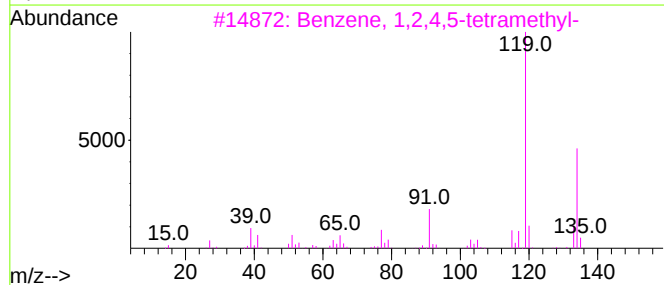
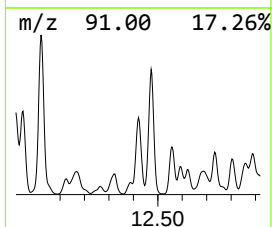
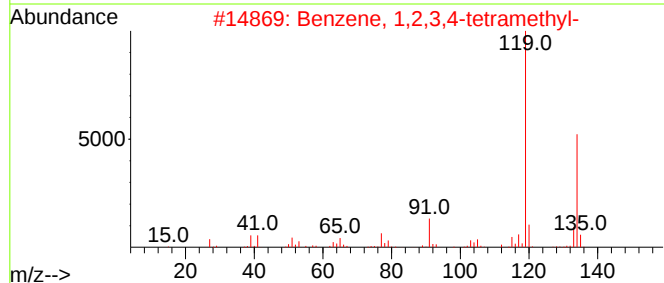
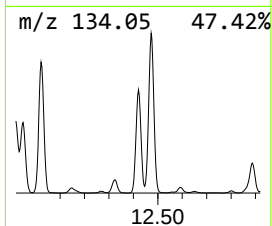
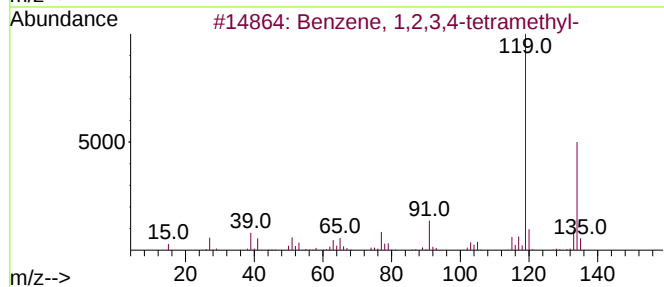
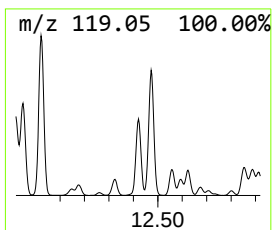
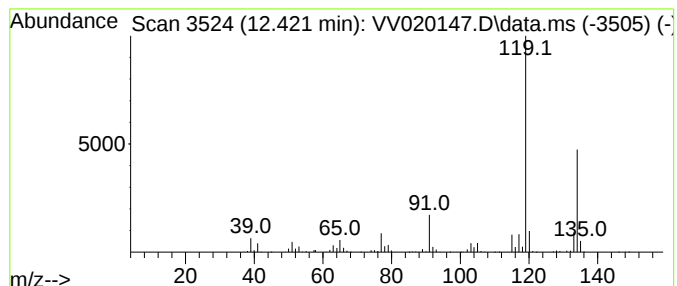
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 55 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.421	60.40 ug/L	1976900	1,4-Dichlorobenzene-d4	11.254

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

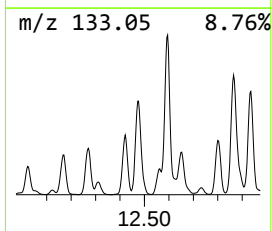
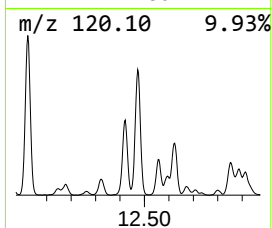
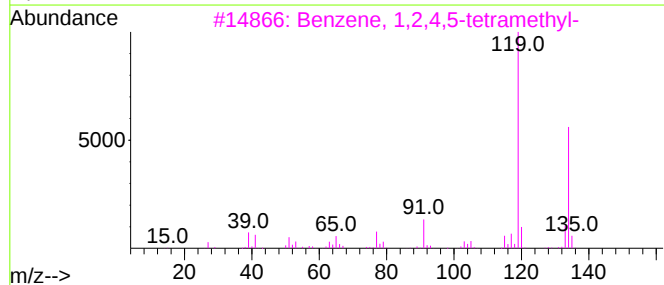
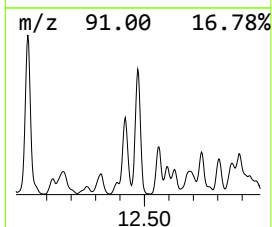
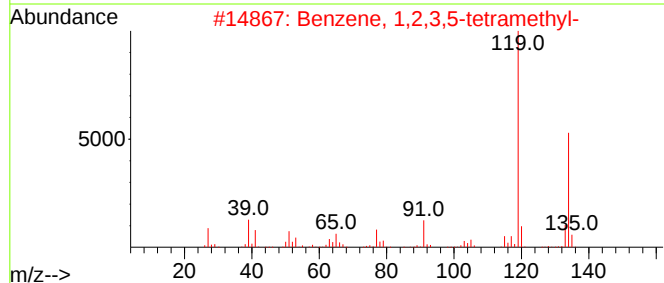
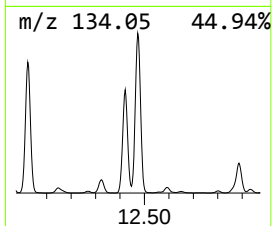
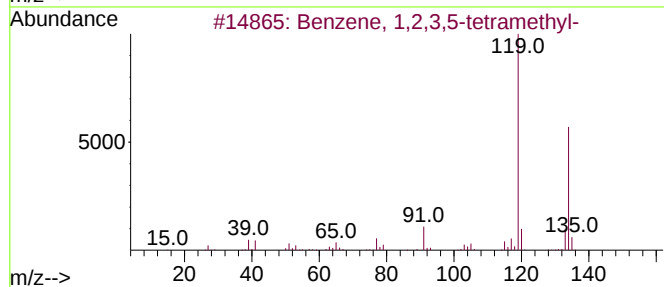
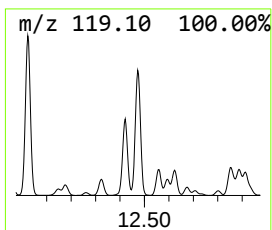
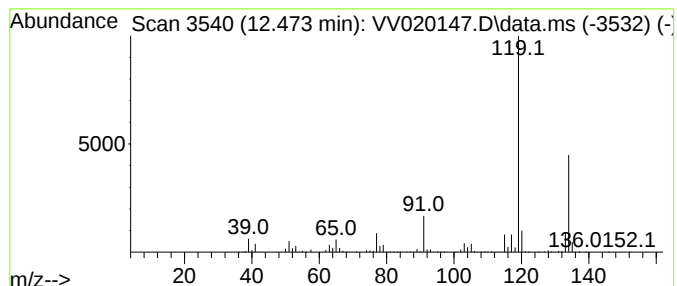
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 56 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.473	92.35 ug/L	3022520	1,4-Dichlorobenzene-d4	11.254

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
2			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
3			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
4			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
5			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

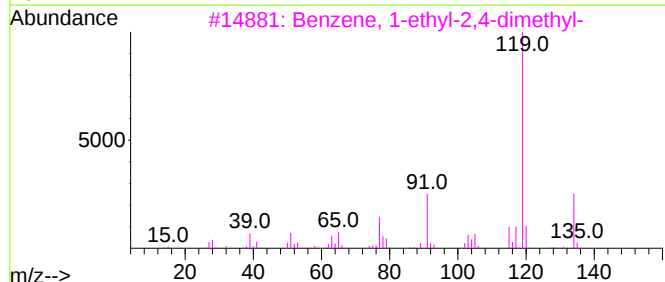
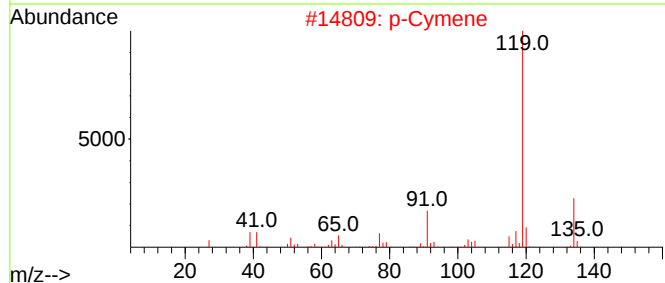
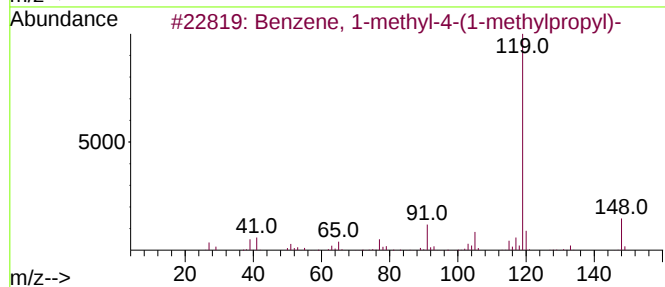
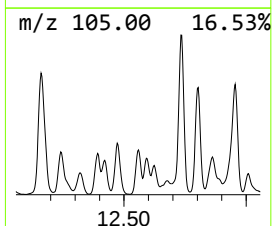
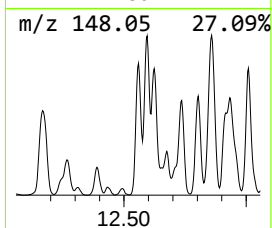
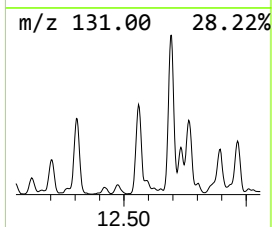
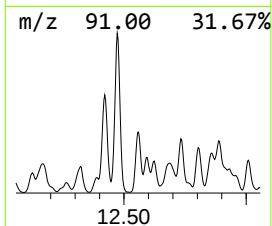
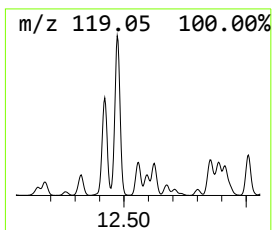
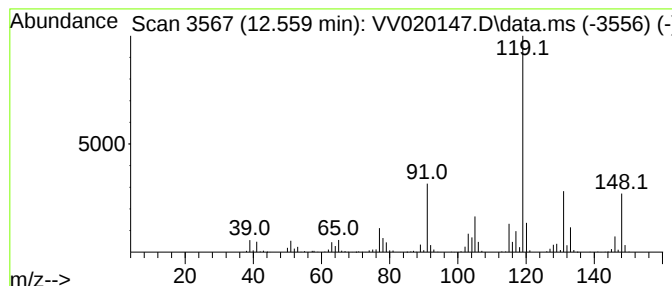
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 57 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.559	25.55 ug/L	836043	1,4-Dichlorobenzene-d4	11.254

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	70
2			p-Cymene	134	C10H14	000099-87-6	60
3			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	60
4			o-Cymene	134	C10H14	000527-84-4	60
5			o-Cymene	134	C10H14	000527-84-4	60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

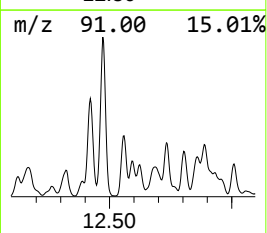
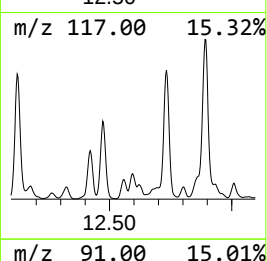
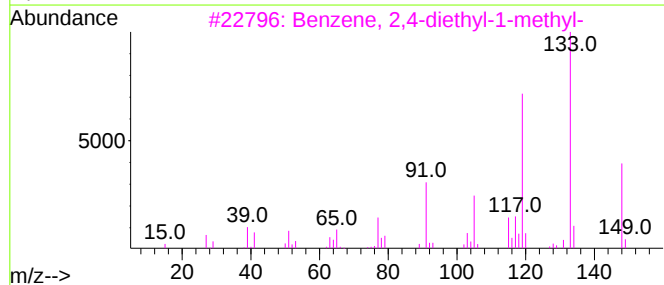
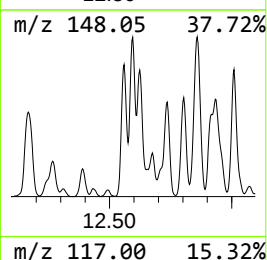
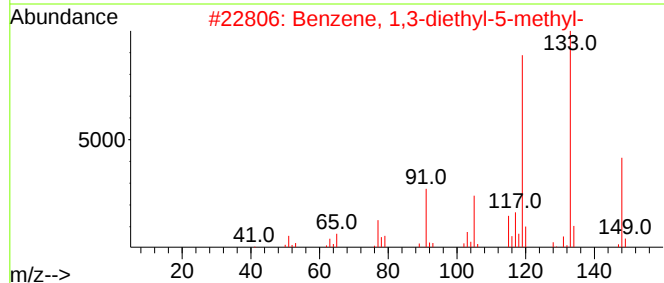
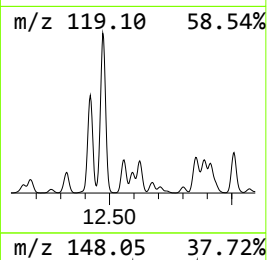
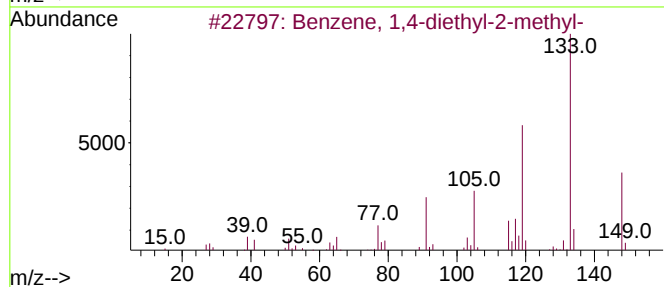
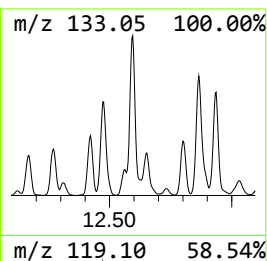
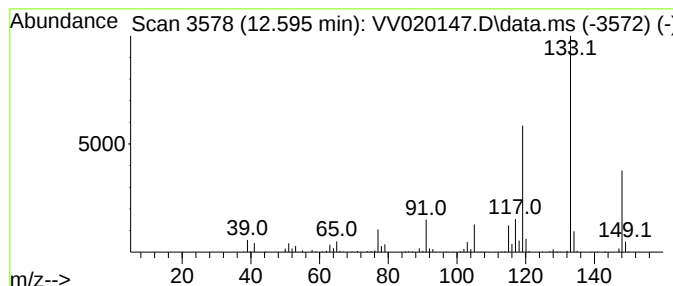
Instrument :
MSVOA_V
ClientSampleId :
BG501

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM012121WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 58 Benzene, 1,4-diethyl-2-methyl- Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD			R.T.
12.595	15.83 ug/L	518230	1,4-Dichlorobenzene-d4			11.254
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Benzene, 1,4-diethyl-2-methyl-		148	C11H16	013632-94-5	94
2	Benzene, 1,3-diethyl-5-methyl-		148	C11H16	002050-24-0	94
3	Benzene, 2,4-diethyl-1-methyl-		148	C11H16	001758-85-6	94
4	Benzene, pentamethyl-		148	C11H16	000700-12-9	76
5	Benzene, 1,3-diethyl-5-methyl-		148	C11H16	002050-24-0	76



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

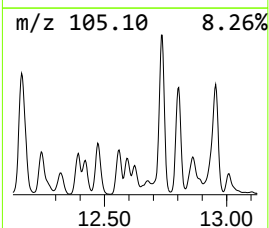
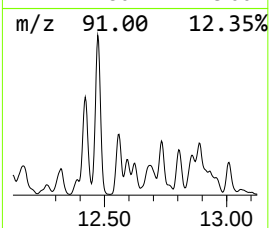
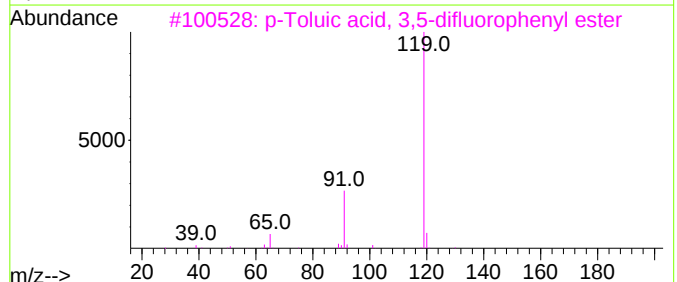
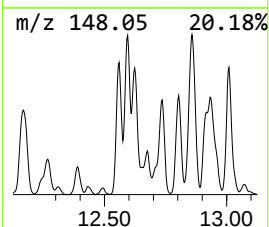
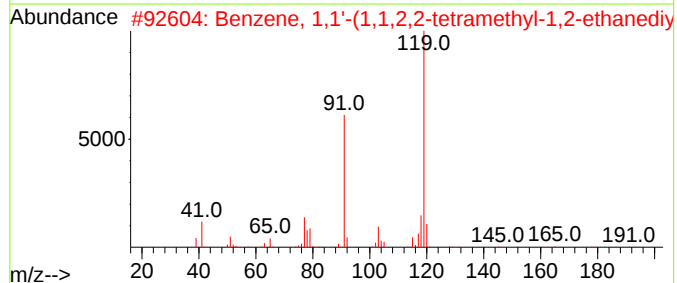
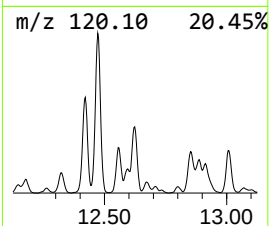
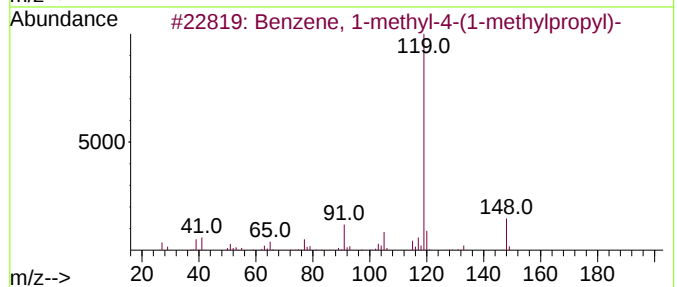
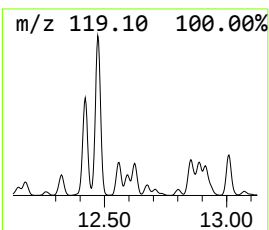
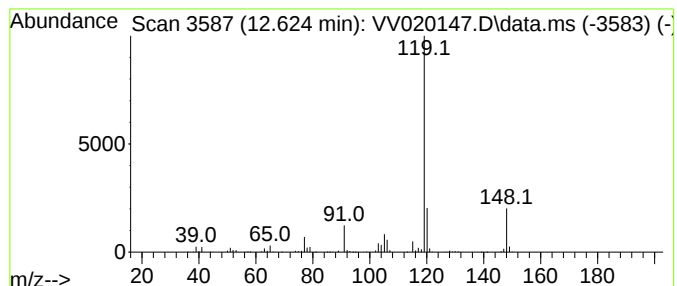
Instrument :
MSVOA_V
ClientSampleId :
BG501

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM012121WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 59 Benzene, 1-methyl-4-(1-meth... Concentration Rank 47

R.T.	EstConc	Area	Relative to ISTD			R.T.
12.624	10.91 ug/L	356937	1,4-Dichlorobenzene-d4			11.254
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-4-(1-methylpro...		148	C11H16	001595-16-0	72
2	Benzene, 1,1'-(1,1,2,2-tetrameth...		238	C18H22	001889-67-4	53
3	p-Toluic acid, 3,5-difluoropheny...		248	C14H10F2O2	1000292-64-0	47
4	Benzene, 1-ethyl-3,5-dimethyl-		134	C10H14	000934-74-7	42
5	Benzene, 1-ethyl-3,5-dimethyl-		134	C10H14	000934-74-7	42



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

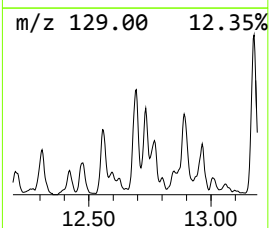
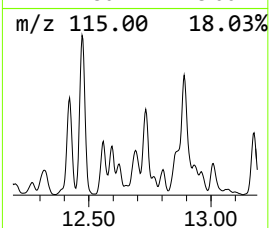
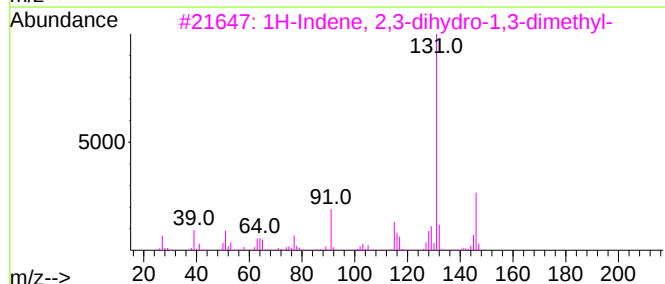
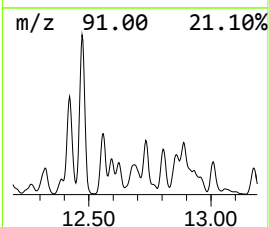
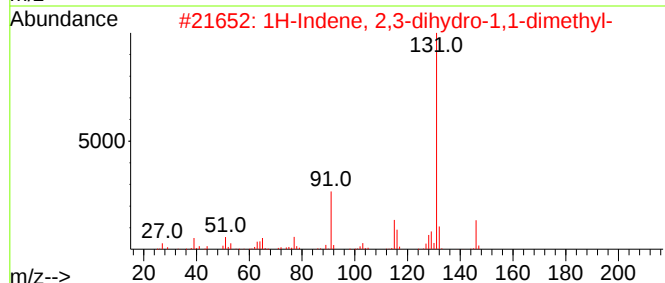
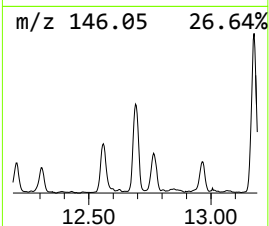
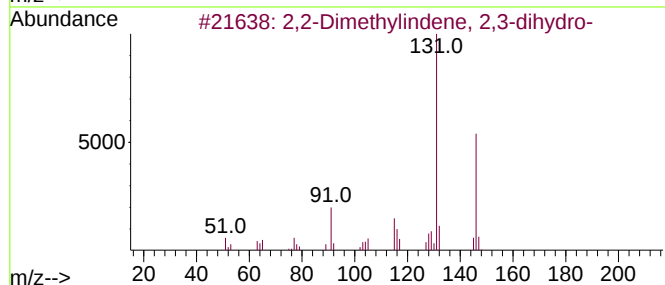
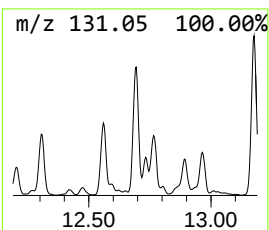
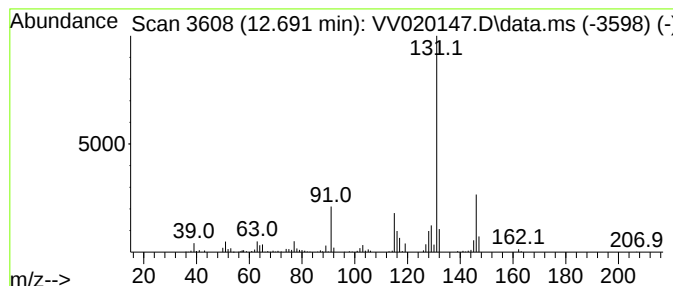
Instrument :
MSVOA_V
ClientSampleId :
BG501

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM012121WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 60 2,2-Dimethylindene, 2,3-dih... Concentration Rank 44

R.T.	EstConc	Area	Relative to ISTD			R.T.
12.691	11.60 ug/L	379668	1,4-Dichlorobenzene-d4			11.254
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	2,2-Dimethylindene, 2,3-dihydro-		146	C11H14	020836-11-7	94
2	1H-Indene, 2,3-dihydro-1,1-dimet...		146	C11H14	004912-92-9	91
3	1H-Indene, 2,3-dihydro-1,3-dimet...		146	C11H14	004175-53-5	91
4	Benzene, (1,2-dimethyl-1-propenyl)-		146	C11H14	000769-57-3	91
5	Bicyclo[4.2.0]octa-1,3,5-triene,...		146	C11H14	027087-54-3	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

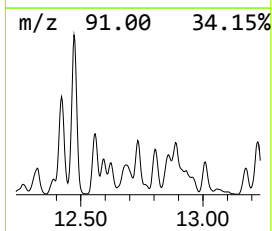
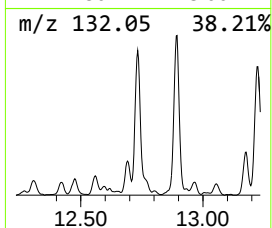
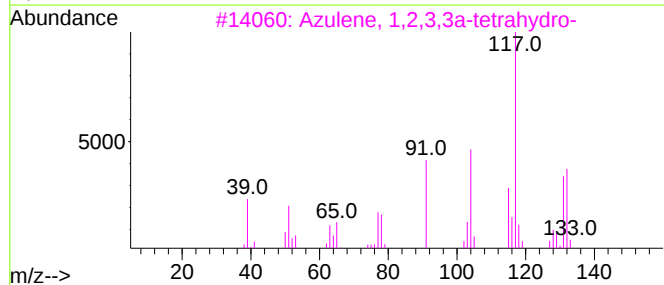
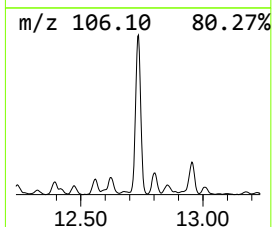
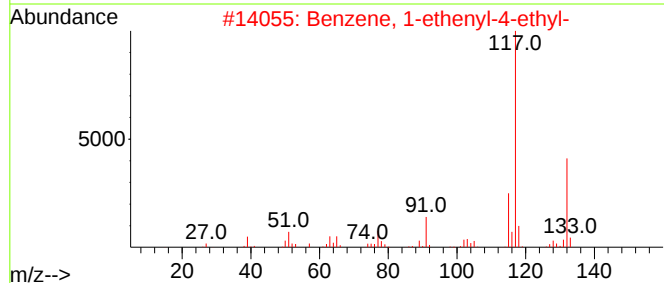
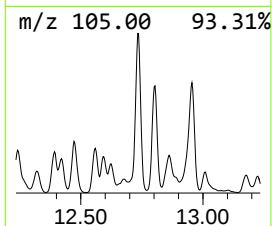
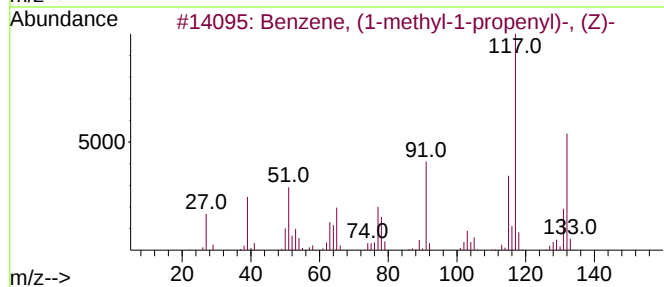
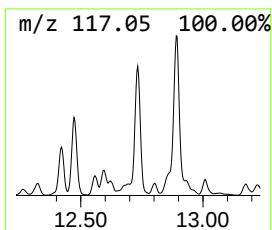
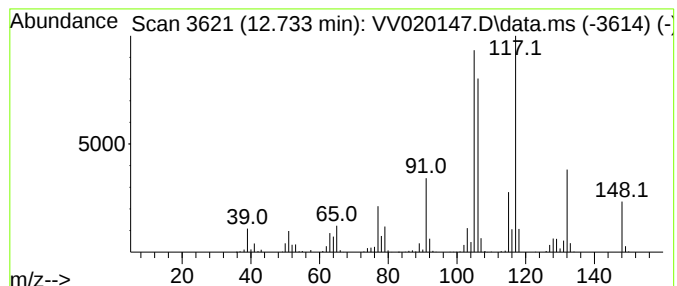
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 61 Benzene, (1-methyl-1-propen... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.733	21.76 ug/L	712276	1,4-Dichlorobenzene-d4	11.254

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	60
2			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	55
3			Azulene, 1,2,3,3a-tetrahydro-	132	C10H12	033877-87-1	55
4			1-Phenyl-1-butene	132	C10H12	000824-90-8	55
5			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	55



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

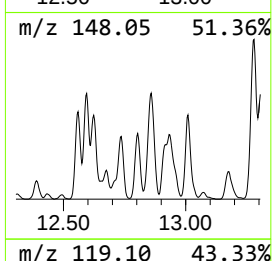
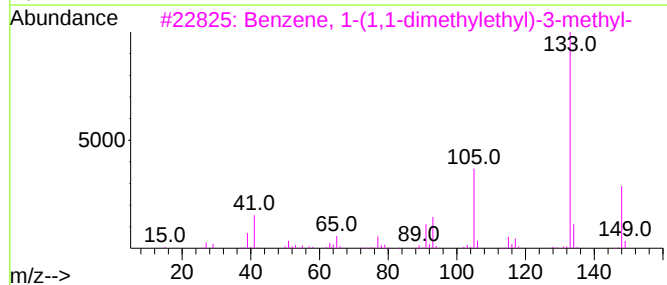
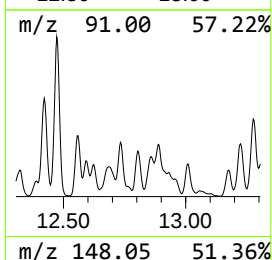
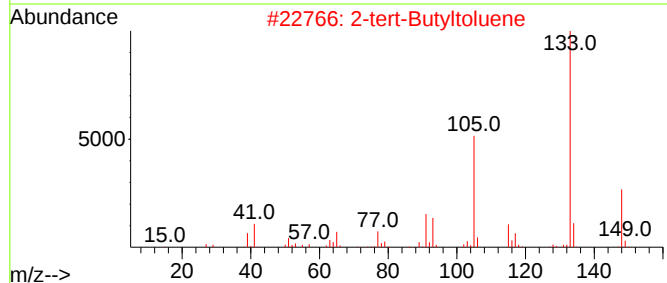
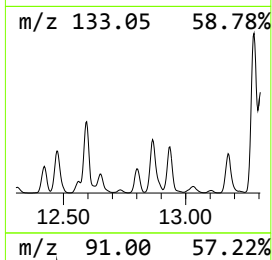
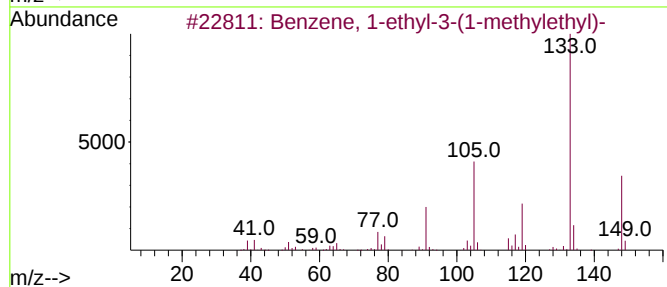
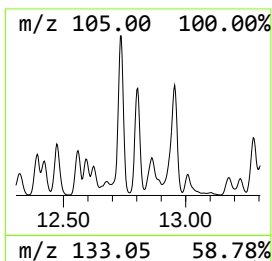
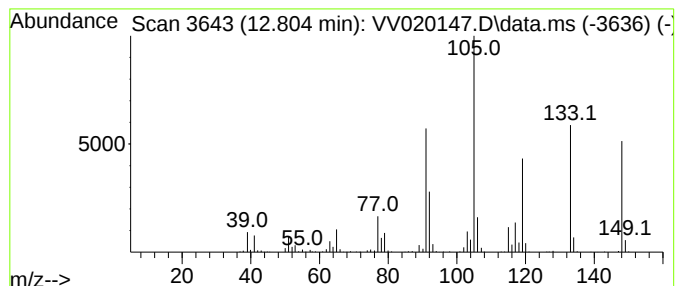
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM012121WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 62 2-tert-Butyltoluene Concentration Rank 46

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.804	11.21 ug/L	366795	1,4-Dichlorobenzene-d4	11.254

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3-(1-methylethyl)-	148	C11H16	004920-99-4	62
2			2-tert-Butyltoluene	148	C11H16	001074-92-6	58
3			Benzene, 1-(1,1-dimethylethyl)-3...	148	C11H16	001075-38-3	58
4			1-(2,3-Dimethylphenyl)ethanone	148	C10H12O	002142-71-4	55
5			Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	52



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

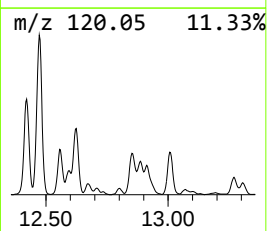
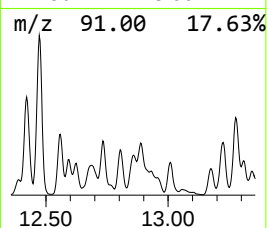
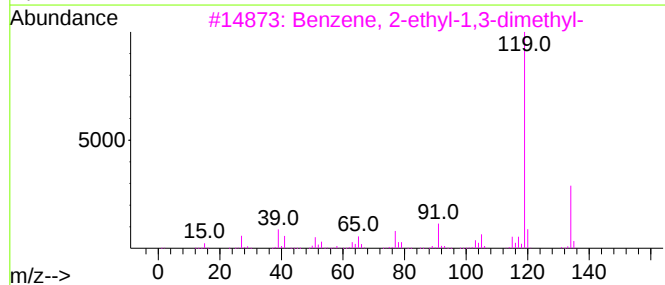
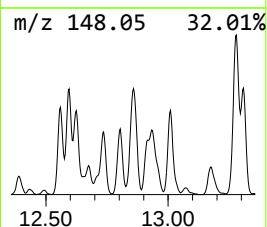
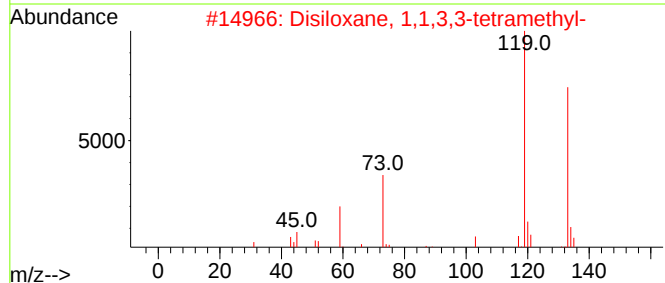
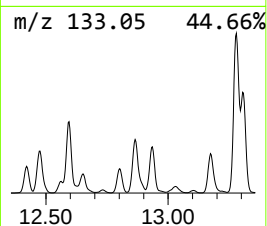
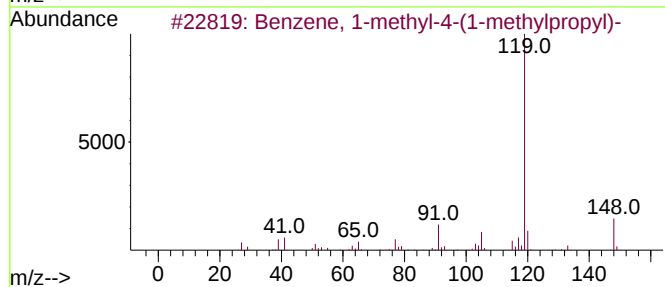
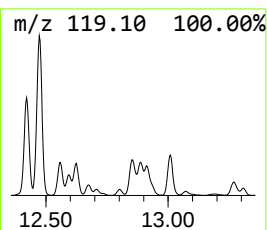
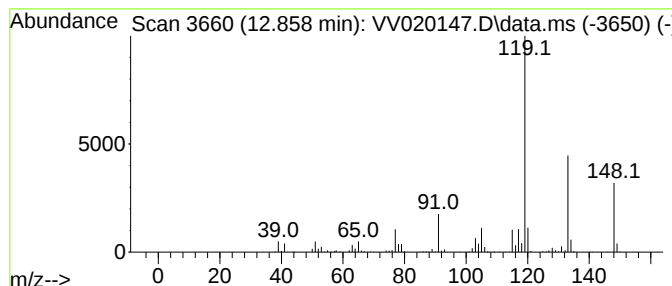
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 63 Benzene, 2-ethyl-1,3-dimethyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.858	24.19 ug/L	791714	1,4-Dichlorobenzene-d4	11.254

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	68
2			Disiloxane, 1,1,3,3-tetramethyl-	134	C4H14OSi2	003277-26-7	64
3			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	64
4			o-Cymene	134	C10H14	000527-84-4	60
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

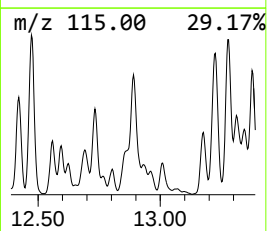
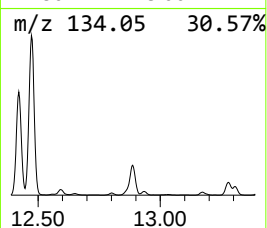
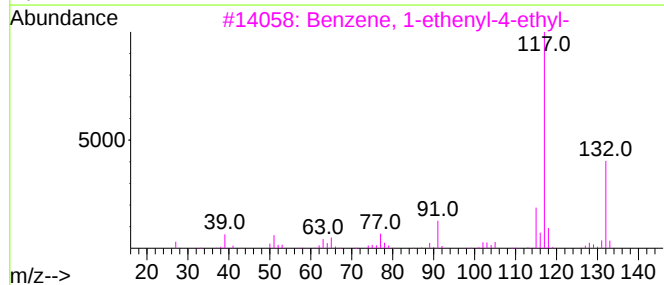
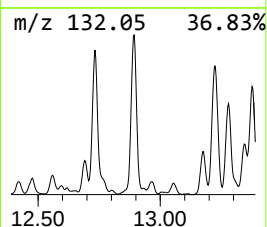
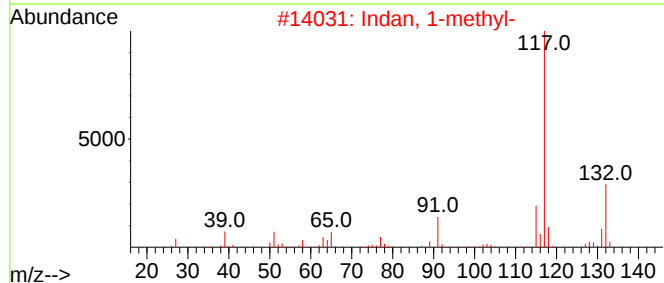
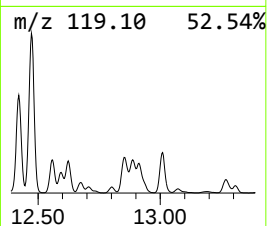
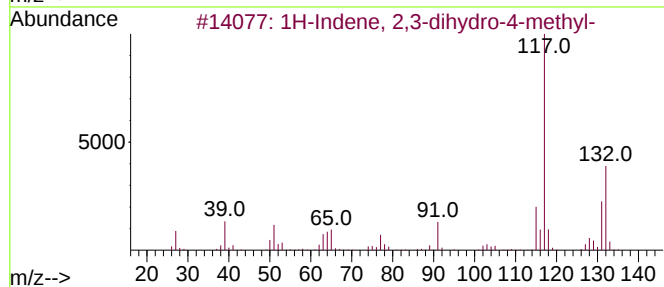
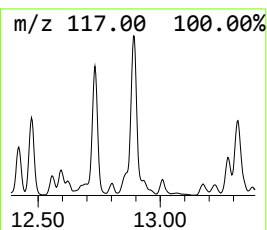
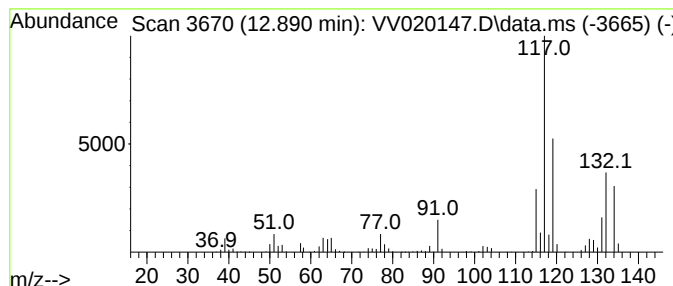
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 64 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.890	21.99 ug/L	719529	1,4-Dichlorobenzene-d4	11.254

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
2			Indan, 1-methyl-	132	C10H12	000767-58-8	60
3			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	55
4			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	53
5			Indan, 1-methyl-	132	C10H12	000767-58-8	52



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

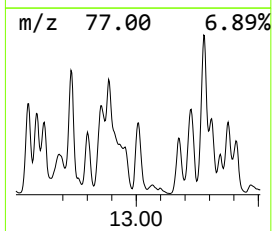
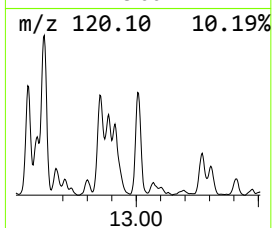
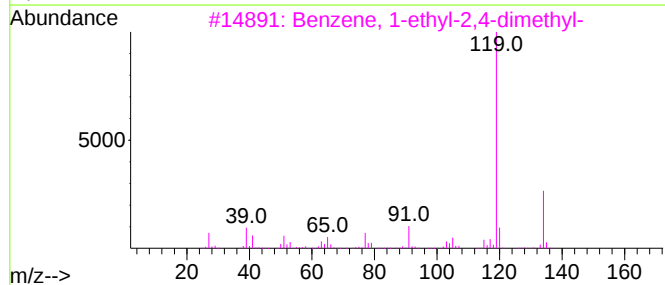
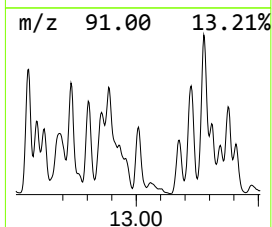
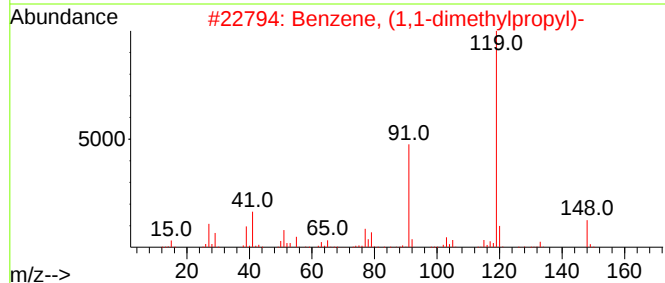
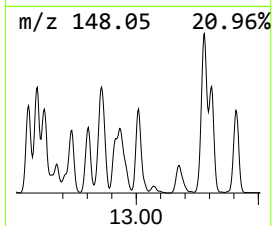
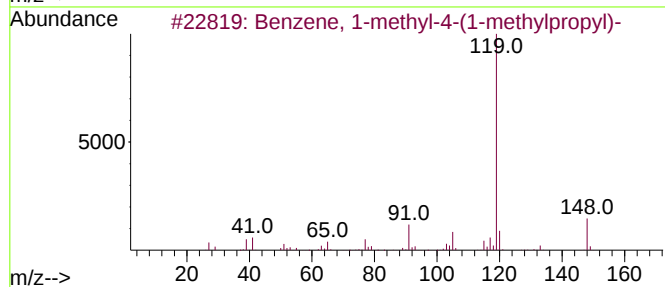
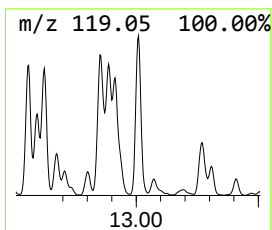
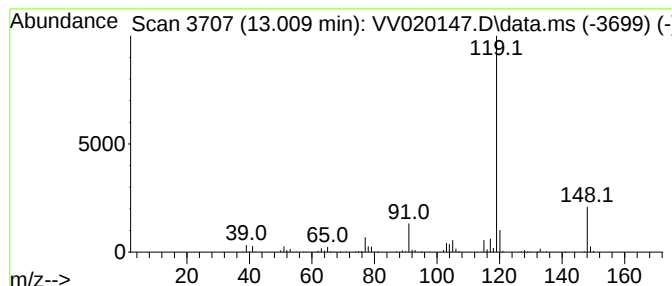
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 65 Benzene, (1,1-dimethylpropyl)- Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.009	16.68 ug/L	545747	1,4-Dichlorobenzene-d4	11.254

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	90
2			Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	78
3			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	72
4			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	72
5			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

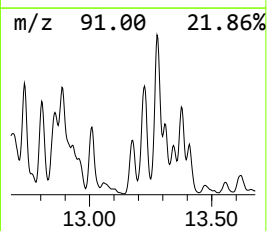
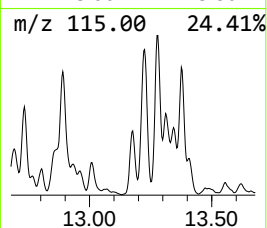
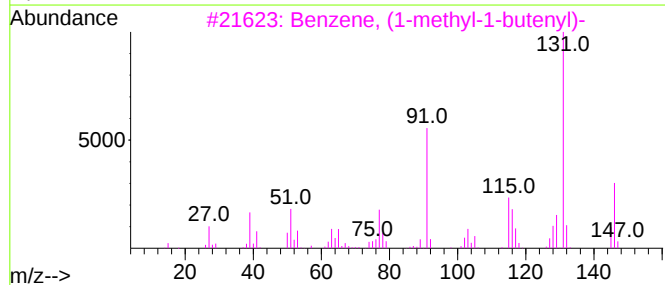
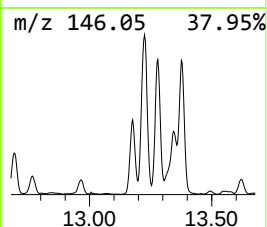
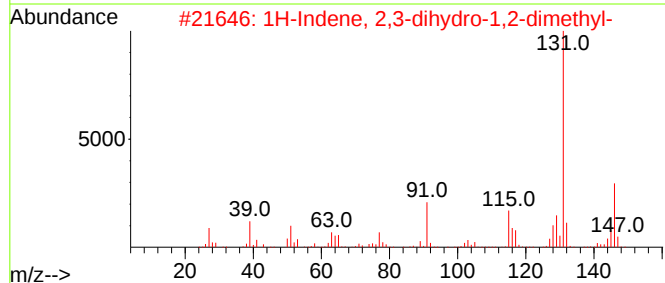
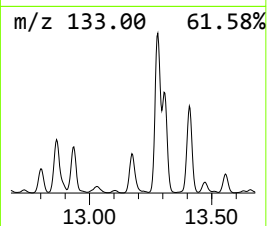
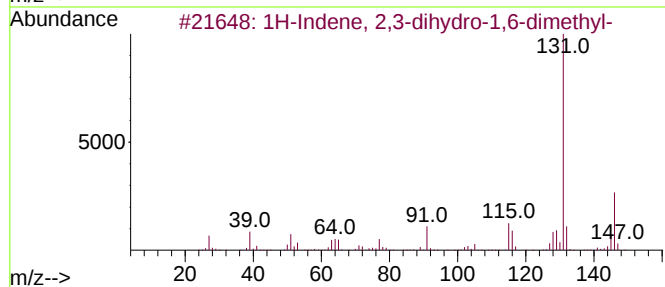
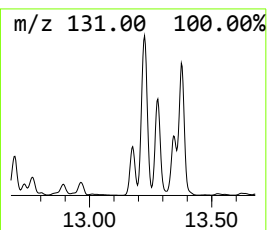
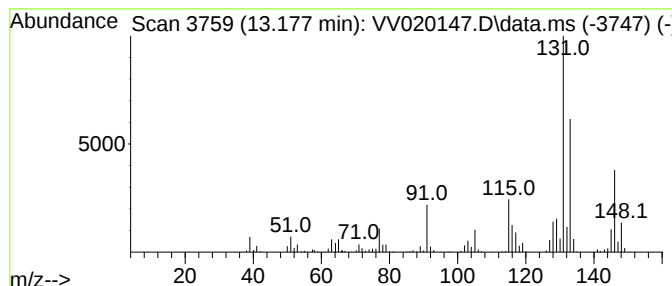
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 66 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.177	20.82 ug/L	681255	1,4-Dichlorobenzene-d4	11.254

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	92
2			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	89
3			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	89
4			Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	86
5			Benzene, 1-methyl-3-(1-methyl-2-...	146	C11H14	052161-57-6	86



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

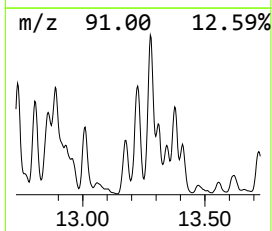
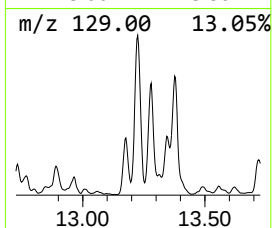
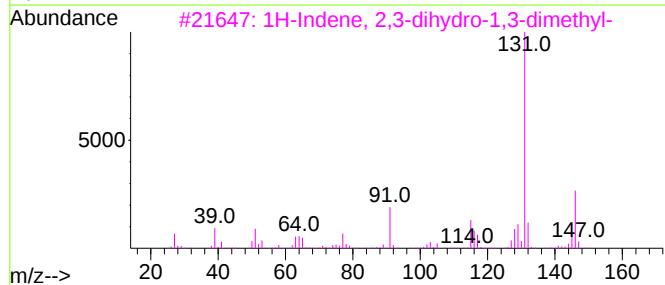
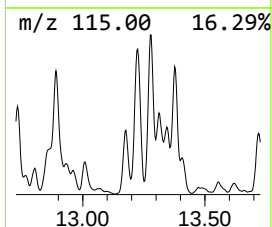
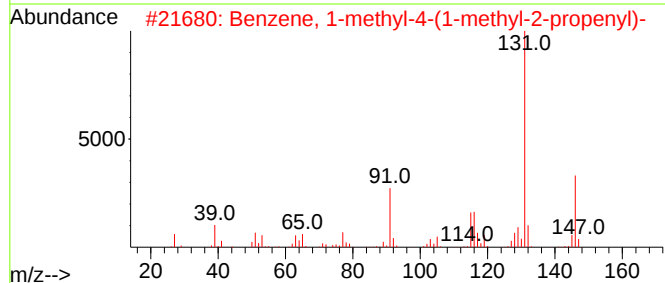
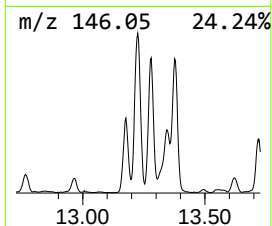
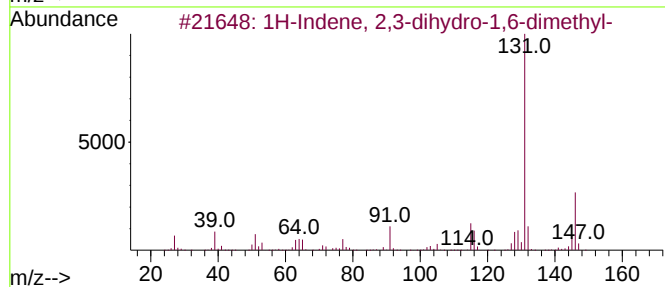
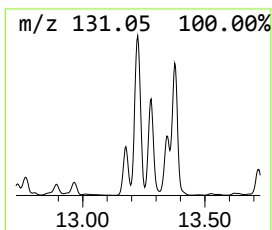
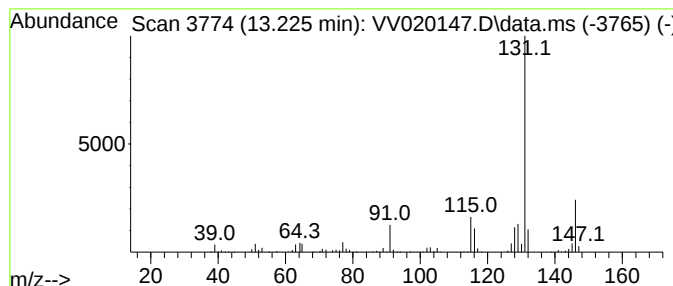
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 67 Benzene, 1-methyl-4-(1-meth... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.225	34.16 ug/L	1117920	1,4-Dichlorobenzene-d4	11.254

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	94
2			Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	91
3			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	91
4			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	91
5			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

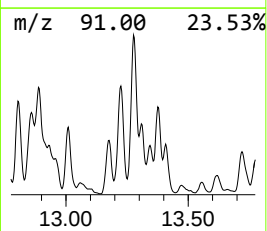
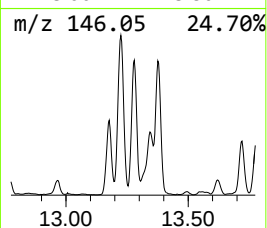
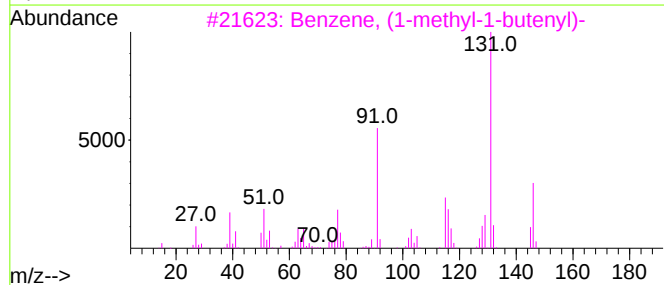
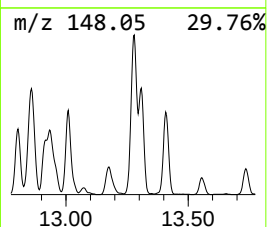
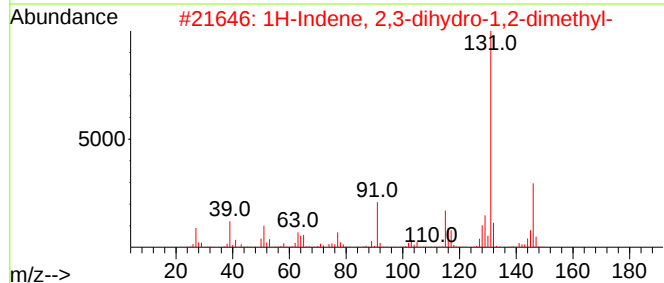
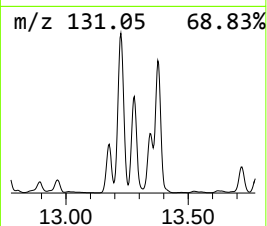
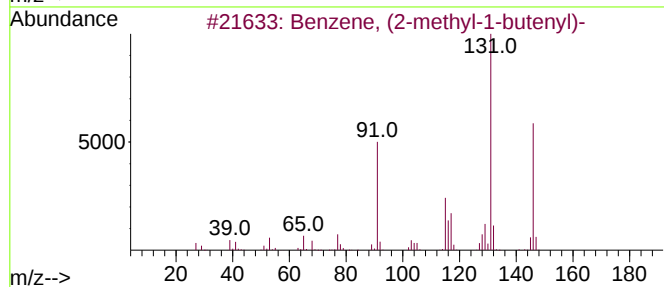
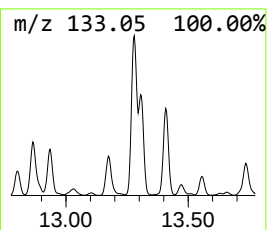
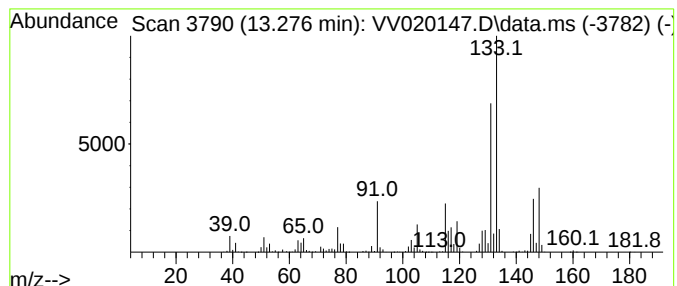
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 68 Benzene, (2-methyl-1-butenyl)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.276	53.97 ug/L	1766290	1,4-Dichlorobenzene-d4	11.254

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	86
2			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	62
3			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	49
4			Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	49
5			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	46



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

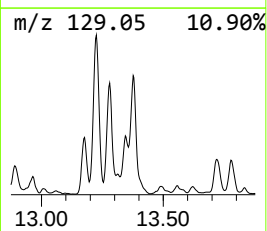
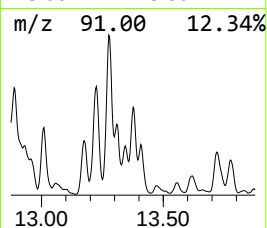
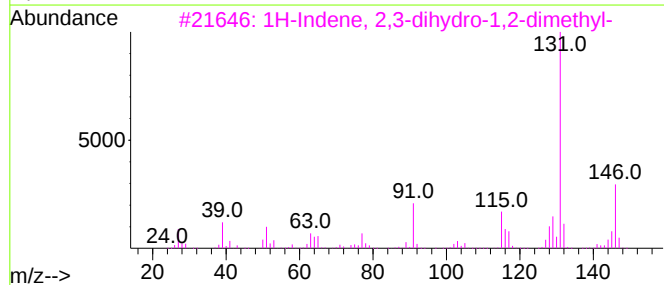
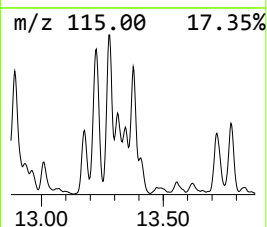
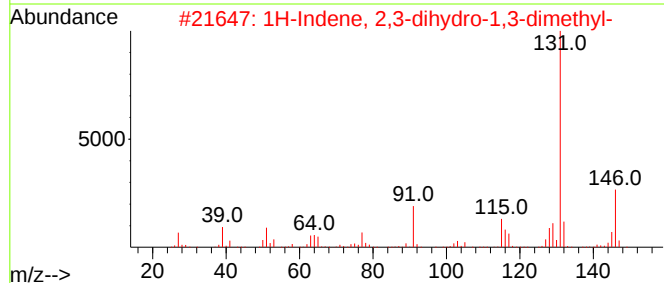
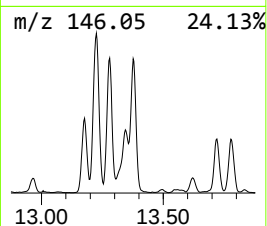
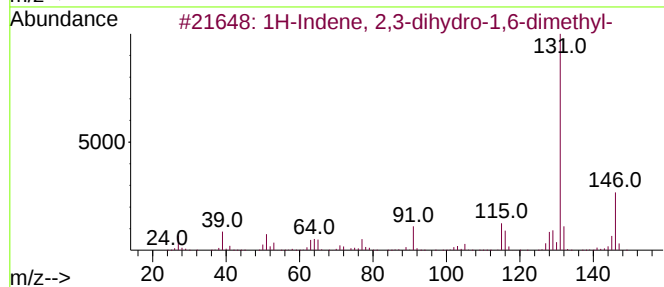
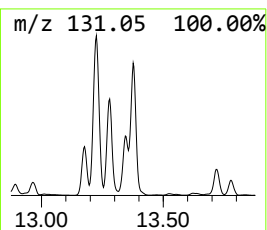
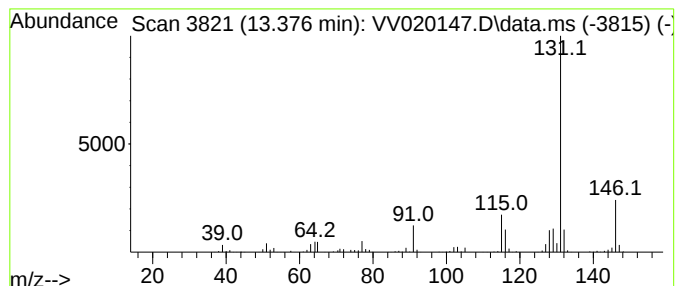
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 70 1H-Indene, 2,3-dihydro-1,3-... Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.376	17.72 ug/L	579861	1,4-Dichlorobenzene-d4	11.254

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	94
2			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	91
3			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	91
4			2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	91
5			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

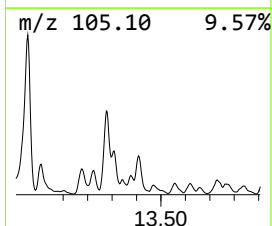
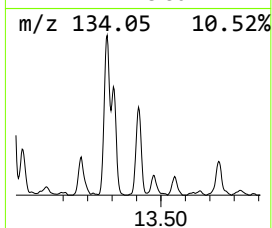
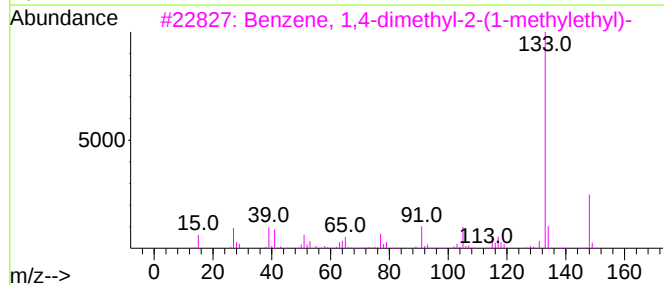
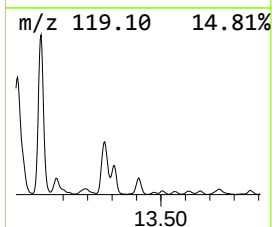
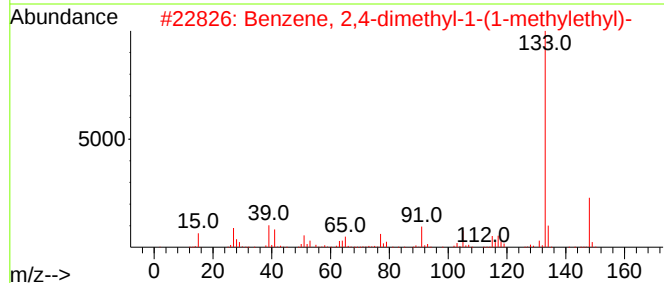
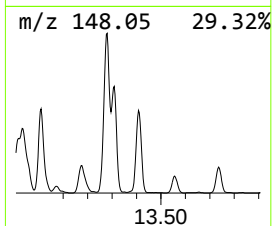
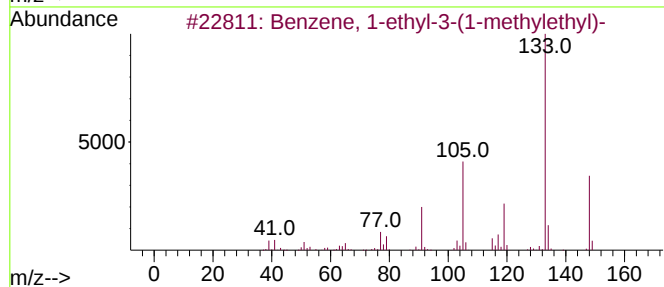
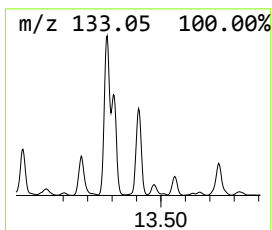
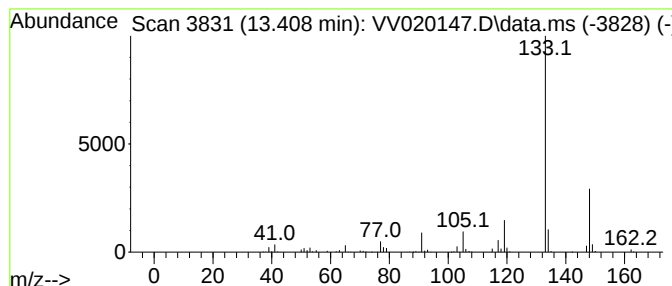
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 71 Benzene, 1-ethyl-3-(1-methy... Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.408	13.91 ug/L	455299	1,4-Dichlorobenzene-d4	11.254

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3-(1-methylethyl)-	148	C11H16	004920-99-4	87
2			Benzene, 2,4-dimethyl-1-(1-methy...	148	C11H16	004706-89-2	86
3			Benzene, 1,4-dimethyl-2-(1-methy...	148	C11H16	004132-72-3	86
4			Benzene, pentamethyl-	148	C11H16	000700-12-9	80
5			Benzene, 1,3-dimethyl-5-(1-methy...	148	C11H16	004706-90-5	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

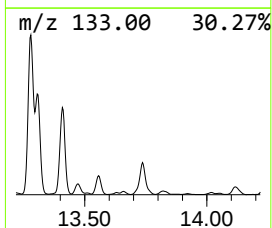
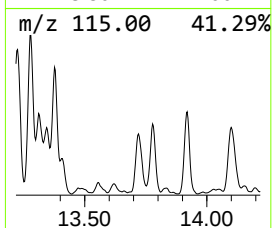
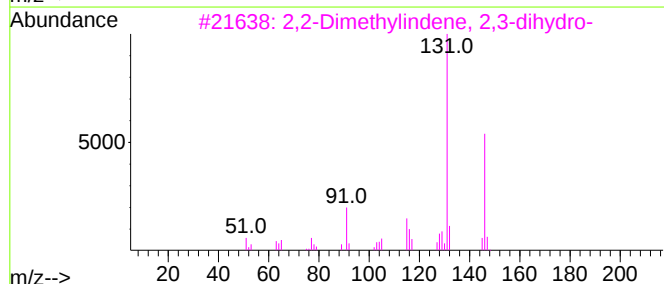
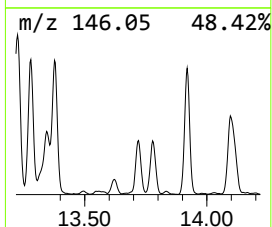
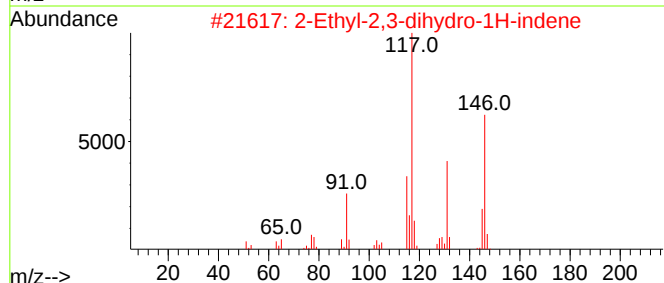
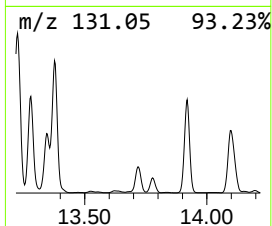
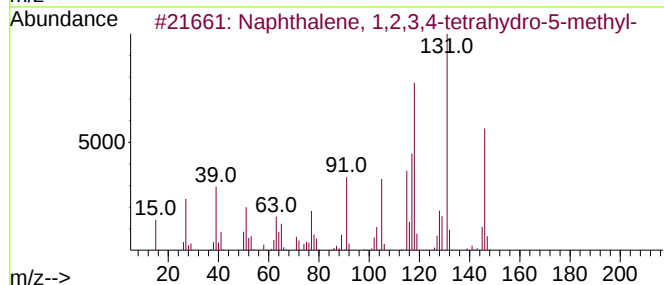
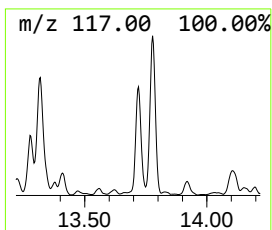
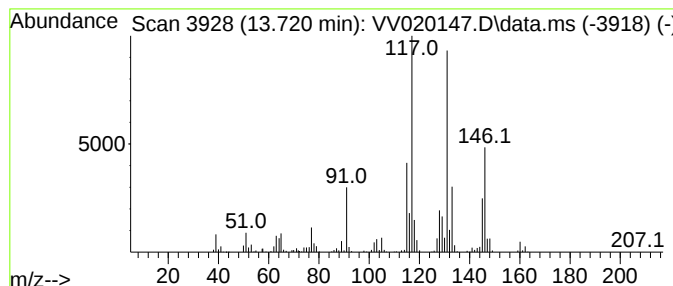
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 74 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.720	18.91 ug/L	618919	1,4-Dichlorobenzene-d4	11.254

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	64
2			2-Ethyl-2,3-dihydro-1H-indene	146	C11H14	056147-63-8	64
3			2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	60
4			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	58
5			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

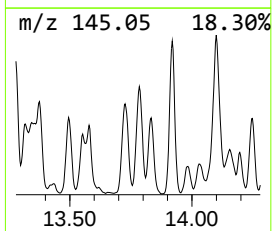
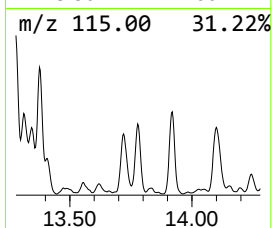
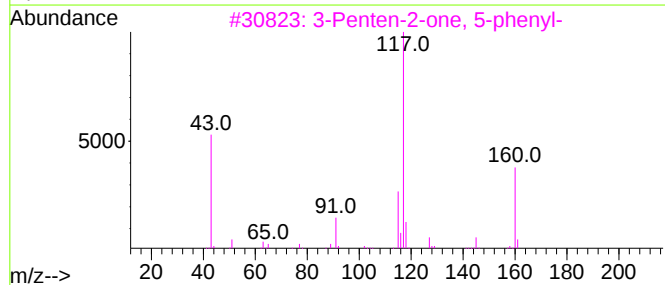
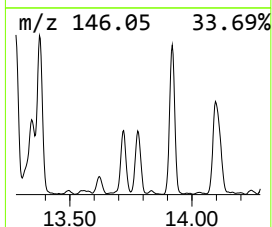
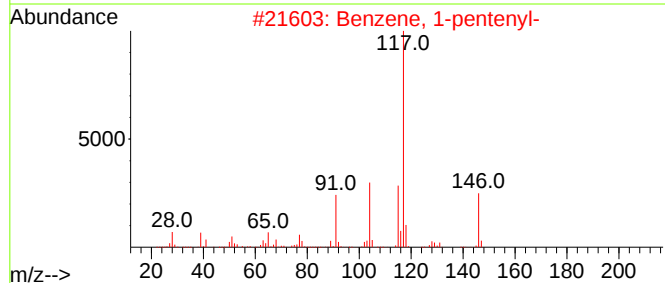
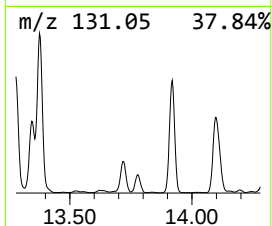
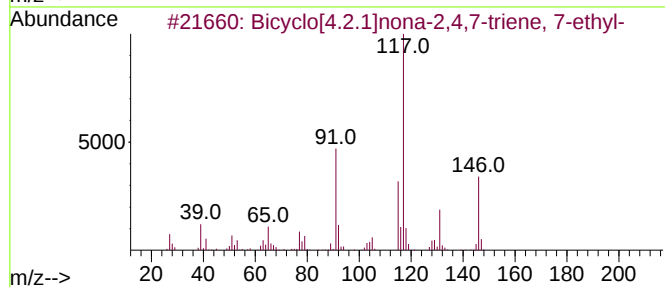
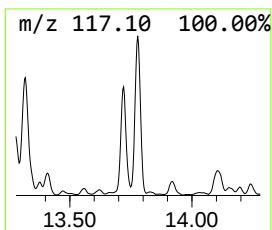
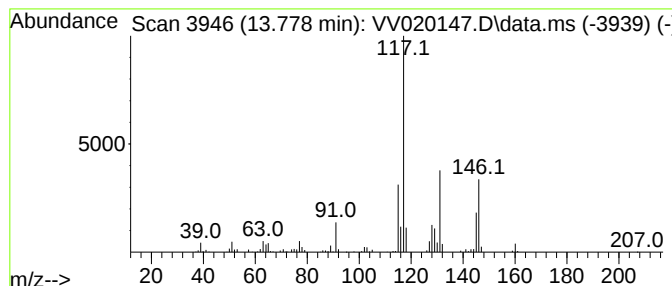
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 75 Bicyclo[4.2.1]nona-2,4,7-tr... Concentration Rank 40

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.778	13.35 ug/L	436773	1,4-Dichlorobenzene-d4	11.254

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[4.2.1]nona-2,4,7-triene,...	146	C11H14	1000164-42-6	64
2			Benzene, 1-pentenyl-	146	C11H14	000826-18-6	59
3			3-Penten-2-one, 5-phenyl-	160	C11H12O	010521-97-8	58
4			2-Ethyl-2,3-dihydro-1H-indene	146	C11H14	056147-63-8	53
5			Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

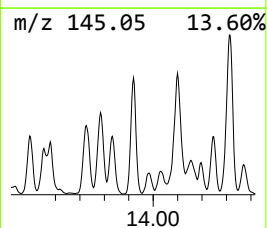
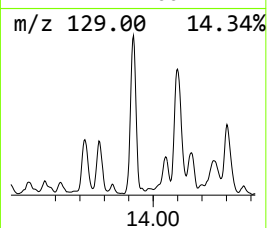
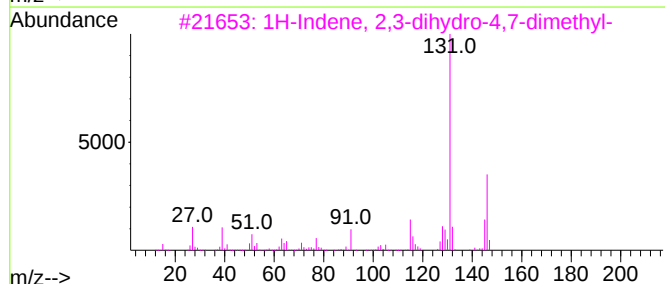
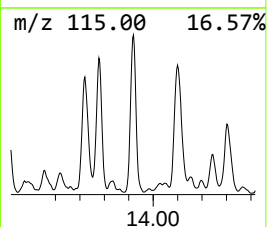
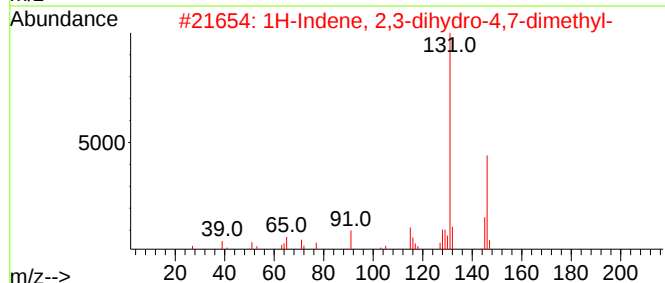
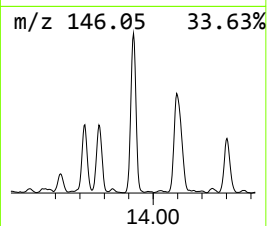
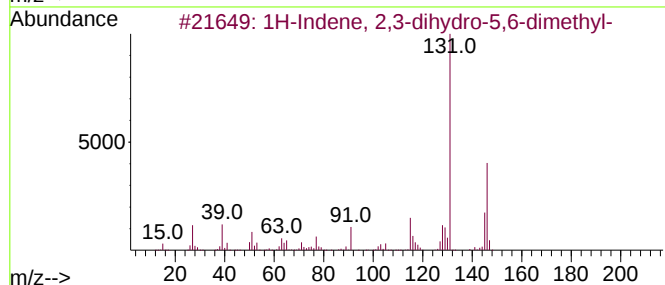
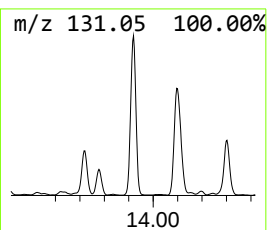
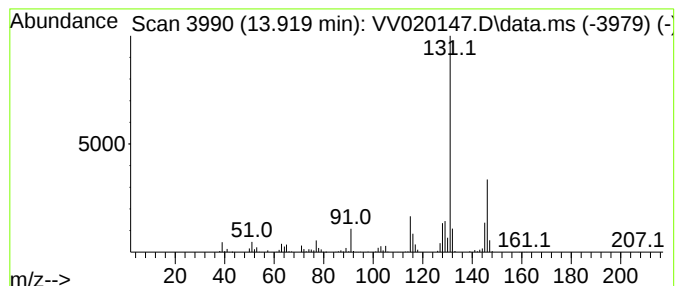
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 76 1H-Indene, 2,3-dihydro-5,6-... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.919	24.94 ug/L	816215	1,4-Dichlorobenzene-d4	11.254

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	96
2			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	94
3			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	94
4			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	94
5			1H-Indene, 2,3-dihydro-4,6-dimet...	146	C11H14	001685-82-1	93



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

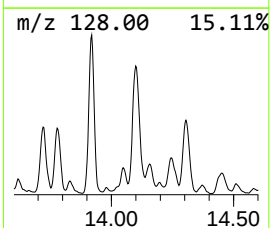
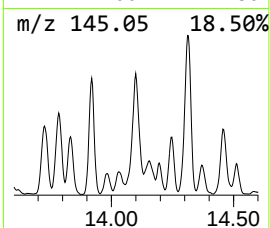
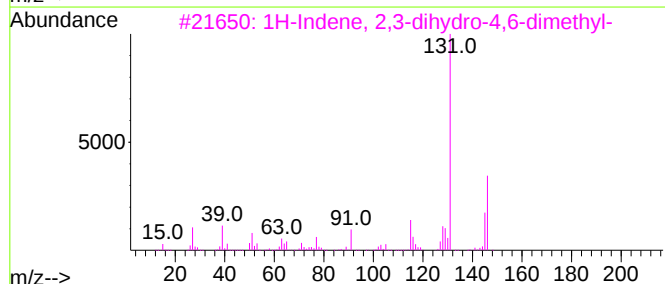
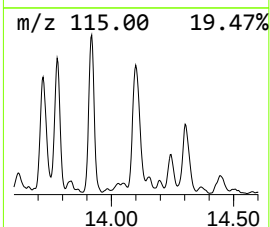
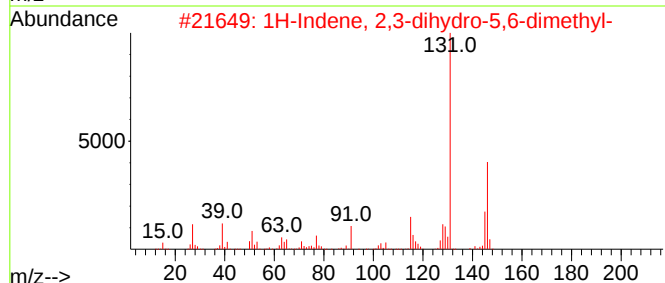
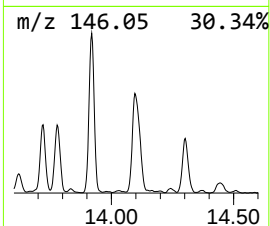
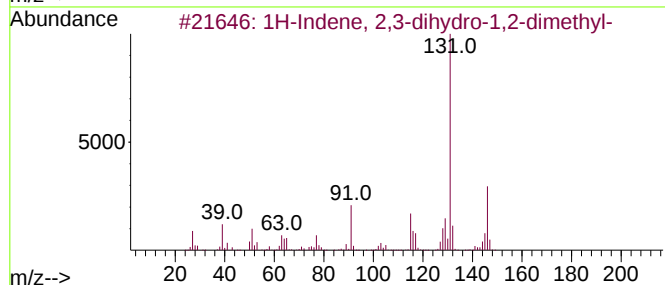
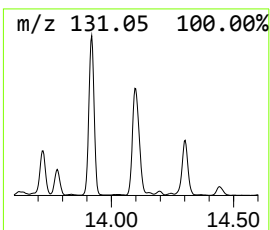
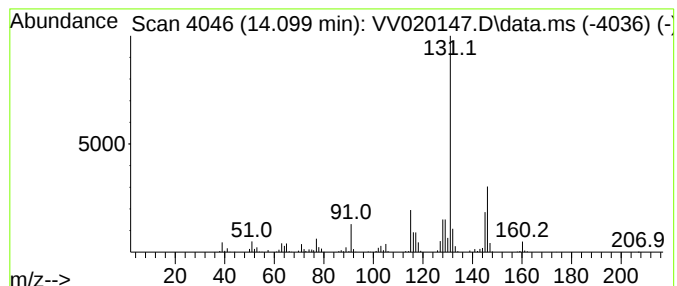
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Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 77 1H-Indene, 2,3-dihydro-4,6-... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.100	25.33 ug/L	829068	1,4-Dichlorobenzene-d4	11.254

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	94
2			1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	93
3			1H-Indene, 2,3-dihydro-4,6-dimet...	146	C11H14	001685-82-1	91
4			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	90
5			Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV012221\
Data File : VV020147.D
Acq On : 22 Jan 2021 19:47
Operator : SY/MD
Sample : M1201-06
Misc : 5.0mL/MSVOA_V/WATER
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
BG501

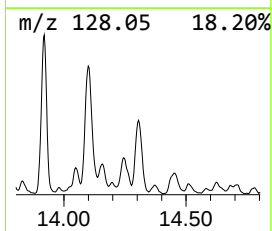
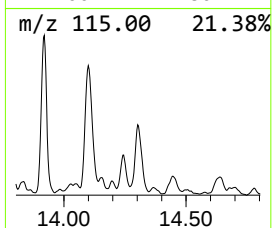
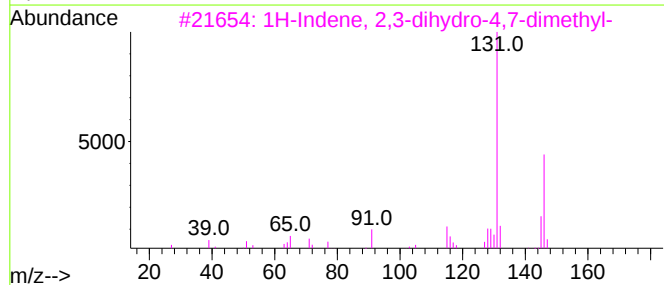
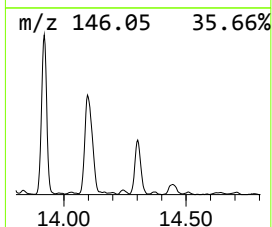
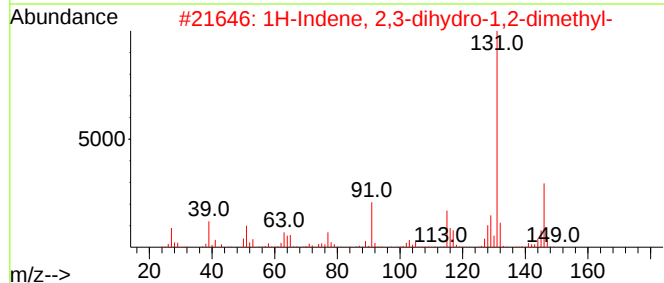
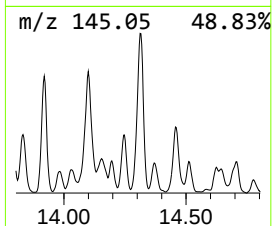
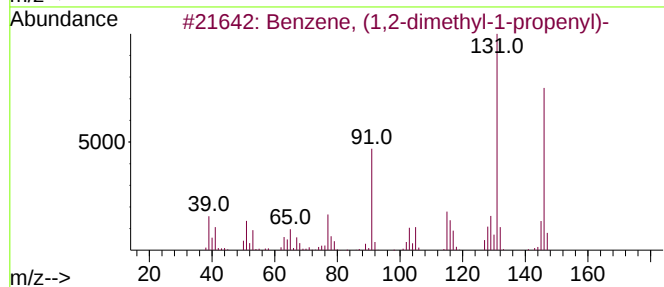
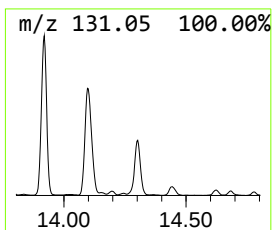
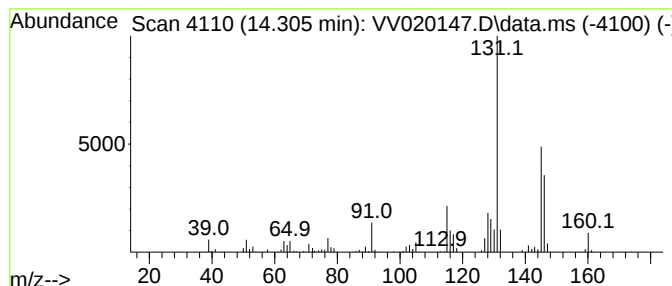
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM012121WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 79 Benzene, (1,2-dimethyl-1-propenyl)- Concentration Rank 39

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.305	13.55 ug/L	443311	1,4-Dichlorobenzene-d4	11.254

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	94
2			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	72
3			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	70
4			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	70
5			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	70



Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VW012221\
 Data File : VW020147.D
 Acq On : 22 Jan 2021 19:47
 Operator : SY/MD
 Sample : M1201-06
 Misc : 5.0mL/MSVOA_V/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 BG501

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SOMVLM012121WMA.M
 Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
 TIC Integration Parameters: LSCINT.P

					--Internal Standard--			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
(DEL) Alkane: S...	1.110	7.0	ug/L	140129	1	5.621	1004440	50.0
(DEL) Alkane: S...	2.534	10.5	ug/L	211576	1	5.621	1004440	50.0
(DEL) Alkane: S...	2.765	14.7	ug/L	295674	1	5.621	1004440	50.0
(DEL) Alkane: S...	3.045	41.4	ug/L	830603	1	5.621	1004440	50.0
(DEL) Alkane: S...	3.672	9.6	ug/L	193473	1	5.621	1004440	50.0
(DEL) Alkane: C...	3.769	25.7	ug/L	515548	1	5.621	1004440	50.0
(DEL) Alkane: S...	4.769	31.9	ug/L	640666	1	5.621	1004440	50.0
(DEL) Alkane: S...	4.913	44.1	ug/L	885471	1	5.621	1004440	50.0
(DEL) Alkane: C...	5.183	12.5	ug/L	251612	1	5.621	1004440	50.0
(DEL) Alkane: S...	5.238	42.7	ug/L	858347	1	5.621	1004440	50.0
(DEL) Alkane: C...	5.331	6.9	ug/L	137881	1	5.621	1004440	50.0
(DEL) Alkane: S...	5.495	28.9	ug/L	581635	1	5.621	1004440	50.0
(DEL) Alkane: S...	6.203	9.2	ug/L	184609	1	5.621	1004440	50.0
(DEL) Alkane: S...	6.261	12.2	ug/L	245741	1	5.621	1004440	50.0
(DEL) Alkane: C...	6.370	6.8	ug/L	136781	1	5.621	1004440	50.0
(DEL) Alkane: C...	6.466	12.9	ug/L	259150	1	5.621	1004440	50.0
(DEL) Alkane: S...	6.659	39.4	ug/L	791388	1	5.621	1004440	50.0
(DEL) Alkane: S...	6.781	50.4	ug/L	1011710	1	5.621	1004440	50.0
(DEL) Alkane: S...	6.843	10.4	ug/L	209774	1	5.621	1004440	50.0
(DEL) Alkane: S...	6.923	17.3	ug/L	347605	1	5.621	1004440	50.0
(DEL) Alkane: S...	6.961	5.6	ug/L	112309	1	5.621	1004440	50.0
(DEL) Alkane: S...	7.084	22.4	ug/L	449872	1	5.621	1004440	50.0
(DEL) Alkane: C...	7.187	6.8	ug/L	137286	1	5.621	1004440	50.0
(DEL) Alkane: C...	7.444	8.2	ug/L	225164	2	8.858	1366720	50.0
(DEL) Alkane: C...	7.489	10.4	ug/L	283227	2	8.858	1366720	50.0
(DEL) Alkane: S...	7.572	18.5	ug/L	506890	2	8.858	1366720	50.0
(DEL) Alkane: C...	7.794	11.4	ug/L	311274	2	8.858	1366720	50.0
(DEL) Alkane: S...	8.238	9.8	ug/L	268970	2	8.858	1366720	50.0
(DEL) Alkane: C...	8.296	6.6	ug/L	181567	2	8.858	1366720	50.0
(DEL) Alkane: S...	8.794	4.7	ug/L	128591	2	8.858	1366720	50.0
(DEL) Alkane: S...	9.209	8.3	ug/L	227449	2	8.858	1366720	50.0
(DEL) Alkane: C...	9.476	7.3	ug/L	199951	2	8.858	1366720	50.0
p-Cymene	11.199	14.8	ug/L	484858	3	11.254	1636400	50.0
Benzene, 1,3-di...	11.556	49.0	ug/L	1605360	3	11.254	1636400	50.0
Benzene, 1-meth...	11.826	49.0	ug/L	1602390	3	11.254	1636400	50.0
Benzene, 2-ethy...	11.919	57.0	ug/L	1863900	3	11.254	1636400	50.0
Benzene, 1-meth...	11.948	55.7	ug/L	1822060	3	11.254	1636400	50.0
o-Cymene	12.022	113.9	ug/L	3727580	3	11.254	1636400	50.0
Indan, 1-methyl-	12.125	14.5	ug/L	475030	3	11.254	1636400	50.0
Phenol, p-(2-me...	12.167	20.9	ug/L	683698	3	11.254	1636400	50.0
Benzene, 4-ethy...	12.321	15.9	ug/L	520829	3	11.254	1636400	50.0
Benzene, 1,2,3,...	12.421	60.4	ug/L	1976900	3	11.254	1636400	50.0
Benzene, 1,2,3,...	12.473	92.3	ug/L	3022520	3	11.254	1636400	50.0
Benzene, 1-ethy...	12.559	25.6	ug/L	836043	3	11.254	1636400	50.0
Benzene, 1,4-di...	12.595	15.8	ug/L	518230	3	11.254	1636400	50.0
Benzene, 1-meth...	12.624	10.9	ug/L	356937	3	11.254	1636400	50.0
2,2-Dimethylind...	12.691	11.6	ug/L	379668	3	11.254	1636400	50.0
Benzene, (1-met...	12.733	21.8	ug/L	712276	3	11.254	1636400	50.0
2-tert-Butyltol...	12.804	11.2	ug/L	366795	3	11.254	1636400	50.0
Benzene, 2-ethy...	12.858	24.2	ug/L	791714	3	11.254	1636400	50.0
1H-Indene, 2,3-...	12.890	22.0	ug/L	719529	3	11.254	1636400	50.0
Benzene, (1,1-d...	13.009	16.7	ug/L	545747	3	11.254	1636400	50.0

1H-Indene, 2,3-...	13.177	20.8	ug/L	681255	3	11.254	1636400	50.0
Benzene, 1-meth...	13.225	34.2	ug/L	1117920	3	11.254	1636400	50.0
Benzene, (2-met...	13.276	54.0	ug/L	1766290	3	11.254	1636400	50.0
1H-Indene, 2,3-...	13.376	17.7	ug/L	579861	3	11.254	1636400	50.0
Benzene, 1-ethy...	13.408	13.9	ug/L	455299	3	11.254	1636400	50.0
Naphthalene, 1,...	13.720	18.9	ug/L	618919	3	11.254	1636400	50.0
Bicyclo[4.2.1]n...	13.778	13.4	ug/L	436773	3	11.254	1636400	50.0
1H-Indene, 2,3-...	13.919	24.9	ug/L	816215	3	11.254	1636400	50.0
1H-Indene, 2,3-...	14.100	25.3	ug/L	829068	3	11.254	1636400	50.0
Benzene, (1,2-d...	14.305	13.6	ug/L	443311	3	11.254	1636400	50.0

Instrument :
MSVOA_V
ClientSampleId :
BG501