

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU042822\
 Data File : VU048338.D
 Acq On : 28 Apr 2022 14:04
 Operator : SY/MD
 Sample : N2587-10DL 40X
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 EXET6DL

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM042522WMA.M
 Title : VOC Analysis

Signal : TIC: VU048338.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.359	3	6	26	rVB	5248688	5162850	100.00%	24.493%
2	1.600	73	81	99	rBV	181214	237439	4.60%	1.126%
3	1.903	169	175	205	rVB	162728	253305	4.91%	1.202%
4	2.565	369	381	394	rBV	347078	571356	11.07%	2.711%
5	2.636	395	403	412	rBV	36054	59762	1.16%	0.284%
6	2.800	442	454	455	rBV5	5427	7697	0.15%	0.037%
7	3.012	515	520	521	rBV3	1338	1118	0.02%	0.005%
8	3.137	546	559	567	rBV	31703	65803	1.27%	0.312%
9	3.353	619	626	627	rBV2	899	720	0.01%	0.003%
10	3.369	627	631	640	rVB6	1516	1795	0.03%	0.009%
11	3.446	647	655	668	rVV3	1904	3621	0.07%	0.017%
12	3.690	726	731	734	rBV3	1147	1215	0.02%	0.006%
13	4.250	901	905	911	rVV4	942	1144	0.02%	0.005%
14	4.629	1010	1023	1044	rBV	194453	508465	9.85%	2.412%
15	5.063	1141	1158	1182	rBV	318024	698104	13.52%	3.312%
16	5.292	1223	1229	1231	rBV3	659	834	0.02%	0.004%
17	5.391	1251	1260	1269	rVB6	2816	5637	0.11%	0.027%
18	5.726	1340	1364	1371	rBV2	660543	1730531	33.52%	8.210%
19	6.166	1495	1501	1505	rVB4	1217	1376	0.03%	0.007%
20	6.250	1511	1527	1551	rBV	427561	815722	15.80%	3.870%
21	6.517	1599	1610	1624	rBV3	14138	29869	0.58%	0.142%
22	6.603	1633	1637	1649	rVB6	1874	2669	0.05%	0.013%
23	6.690	1649	1664	1680	rBV	410590	803119	15.56%	3.810%
24	6.928	1727	1738	1744	rBV3	1392	2371	0.05%	0.011%
25	6.996	1749	1759	1770	rVV3	13760	25295	0.49%	0.120%
26	7.144	1801	1805	1809	rVV3	833	885	0.02%	0.004%
27	7.189	1809	1819	1836	rVB6	3938	10019	0.19%	0.048%
28	7.346	1861	1868	1876	rVB5	1992	3086	0.06%	0.015%
29	7.491	1898	1913	1925	rBV	193873	351589	6.81%	1.668%
30	7.565	1925	1936	1959	rVB	234518	414501	8.03%	1.966%
31	7.658	1959	1965	1967	rBV6	972	1180	0.02%	0.006%
32	7.742	1987	1991	1997	rVB4	1036	1120	0.02%	0.005%
33	7.799	1997	2009	2025	rBV5	4338	10785	0.21%	0.051%
34	7.899	2025	2040	2052	rBV	845079	1446705	28.02%	6.863%
35	7.976	2053	2064	2080	rVB3	301290	623414	12.07%	2.958%
36	8.179	2115	2127	2143	rBV	159059	269098	5.21%	1.277%

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Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM042522WMA.M
 Title : VOC Analysis

37	8.359	2172	2183	2206	rVB2	105924	189250	3.67%	0.898%
38	8.578	2238	2251	2259	rVV	117712	203869	3.95%	0.967%
39	8.635	2259	2269	2301	rVB	618773	1101661	21.34%	5.226%
40	8.777	2307	2313	2327	rVB5	13335	24183	0.47%	0.115%
41	8.896	2347	2350	2352	rBV2	864	646	0.01%	0.003%
42	9.015	2375	2387	2407	rBV2	44025	78812	1.53%	0.374%
43	9.172	2421	2436	2447	rBV3	5112	8692	0.17%	0.041%
44	9.250	2454	2460	2463	rVV5	811	1009	0.02%	0.005%
45	9.362	2486	2495	2500	rBV4	5651	8711	0.17%	0.041%
46	9.420	2500	2513	2515	rBV	649700	903017	17.49%	4.284%
47	9.632	2572	2579	2580	rBV5	2999	2813	0.05%	0.013%
48	9.661	2584	2588	2592	rVV5	3260	4358	0.08%	0.021%
49	9.828	2633	2640	2649	rBV7	3098	5629	0.11%	0.027%
50	9.941	2667	2675	2683	rVB5	1926	3063	0.06%	0.015%
51	10.102	2714	2725	2740	rBV	8606	14934	0.29%	0.071%
52	10.224	2752	2763	2769	rVV7	1237	2018	0.04%	0.010%
53	10.253	2769	2772	2775	rVV2	656	467	0.01%	0.002%
54	10.487	2835	2845	2852	rBV6	1809	2855	0.06%	0.014%
55	10.578	2869	2873	2880	rVB5	521	494	0.01%	0.002%
56	10.651	2885	2896	2909	rBV5	20993	37921	0.73%	0.180%
57	10.709	2909	2914	2916	rBV3	866	857	0.02%	0.004%
58	10.758	2916	2929	2951	rVB	644483	1033778	20.02%	4.904%
59	10.851	2953	2958	2967	rVV5	1769	2777	0.05%	0.013%
60	10.925	2978	2981	2982	rBV	801	529	0.01%	0.003%
61	10.979	2988	2998	3010	rVB3	23101	45530	0.88%	0.216%
62	11.098	3025	3035	3049	rBV2	14297	24154	0.47%	0.115%
63	11.147	3049	3050	3056	rVB3	998	702	0.01%	0.003%
64	11.250	3075	3082	3090	rVV5	5530	8571	0.17%	0.041%
65	11.295	3091	3096	3103	rVB4	2638	3419	0.07%	0.016%
66	11.378	3115	3122	3130	rBV4	2465	3822	0.07%	0.018%
67	11.468	3139	3150	3157	rBV4	5626	9310	0.18%	0.044%
68	11.510	3157	3163	3172	rVB5	5200	7662	0.15%	0.036%
69	11.587	3176	3187	3198	rBV5	7852	13824	0.27%	0.066%
70	11.658	3205	3209	3214	rBV2	403	463	0.01%	0.002%
71	11.696	3214	3221	3224	rBV4	1403	1759	0.03%	0.008%
72	11.751	3229	3238	3245	rVV5	7386	11909	0.23%	0.056%
73	11.812	3245	3257	3271	rVV	647385	1015031	19.66%	4.815%
74	11.883	3271	3279	3291	rVV5	21572	39783	0.77%	0.189%
75	11.950	3294	3300	3316	rVB10	6323	12307	0.24%	0.058%
76	12.060	3330	3334	3337	rBV3	1079	991	0.02%	0.005%

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Integrator: RTE

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Title : VOC Analysis

77	12.195	3359	3376	3400	rBV	838960	1463477	28.35%	6.943%
78	12.475	3455	3463	3468	rBV5	1577	2259	0.04%	0.011%
79	12.574	3492	3494	3501	rVB3	509	500	0.01%	0.002%
80	12.684	3522	3528	3531	rBV4	1307	1702	0.03%	0.008%
81	12.774	3547	3556	3565	rVV	3211	4696	0.09%	0.022%
82	12.970	3603	3617	3627	rBV4	12077	20873	0.40%	0.099%
83	13.105	3655	3659	3666	rVB4	722	720	0.01%	0.003%
84	13.153	3669	3674	3677	rVB5	763	683	0.01%	0.003%
85	13.217	3687	3694	3696	rBV5	818	800	0.02%	0.004%
86	13.266	3705	3709	3712	rBV4	816	854	0.02%	0.004%
87	13.330	3723	3729	3734	rBV4	2355	3286	0.06%	0.016%
88	13.375	3734	3743	3757	rVV	34792	53041	1.03%	0.252%
89	13.446	3760	3765	3776	rVB6	1518	2306	0.04%	0.011%
90	13.732	3851	3854	3856	rBV3	888	581	0.01%	0.003%
91	13.777	3863	3868	3873	rBV6	2253	2887	0.06%	0.014%
92	14.044	3949	3951	3958	rVB3	635	646	0.01%	0.003%
93	14.089	3958	3965	3966	rBV4	2589	2136	0.04%	0.010%
94	14.333	4035	4041	4044	rBV3	887	846	0.02%	0.004%
95	14.359	4044	4049	4060	rVV5	1442	2178	0.04%	0.010%
96	14.429	4064	4071	4078	rVB5	6945	9269	0.18%	0.044%
97	14.494	4080	4091	4112	rBV2	132229	223959	4.34%	1.062%
98	14.880	4208	4211	4214	rBV4	660	460	0.01%	0.002%
99	15.545	4407	4418	4432	rBV	172744	268400	5.20%	1.273%
100	16.391	4671	4681	4690	rBV3	34633	56553	1.10%	0.268%

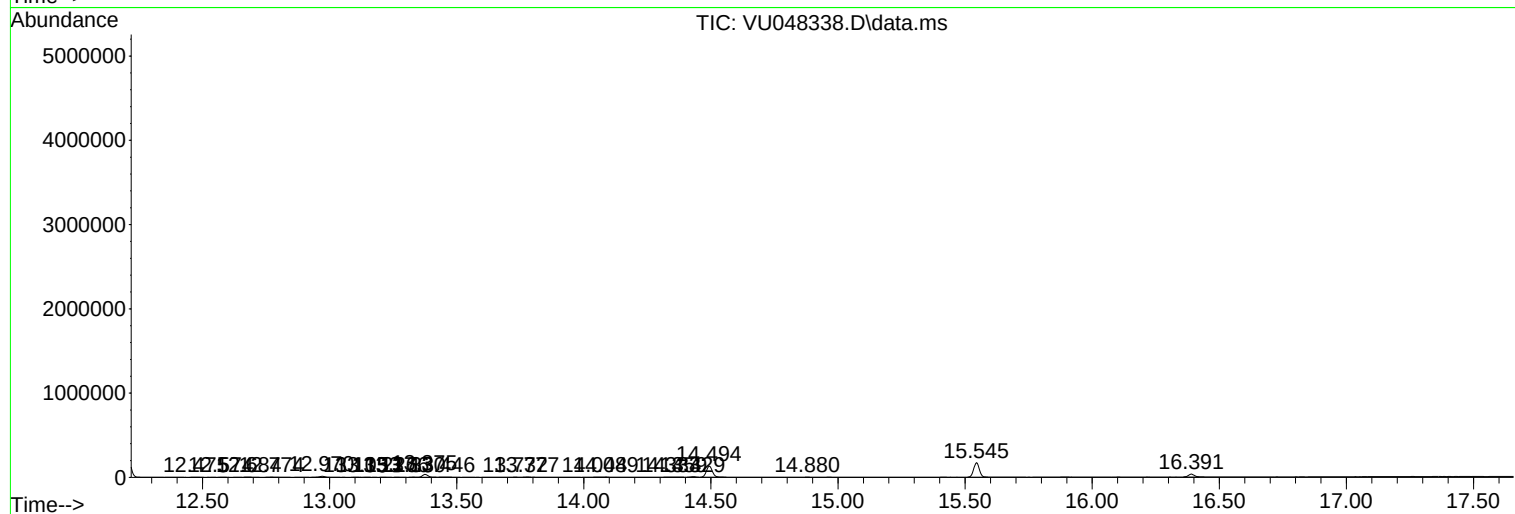
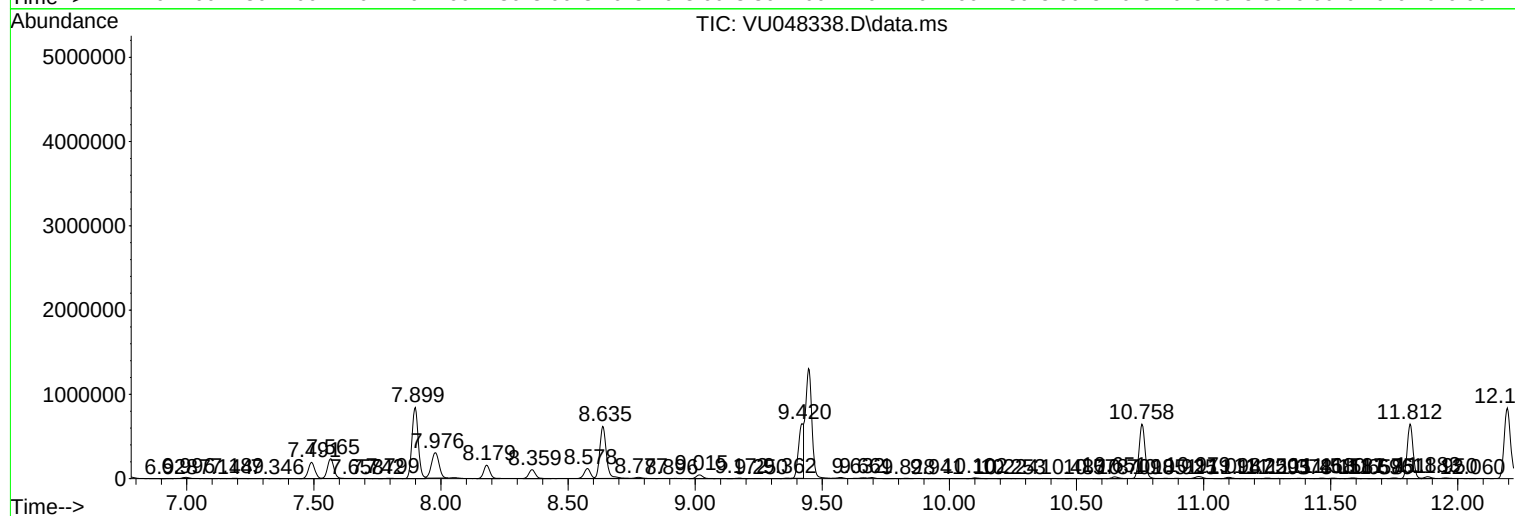
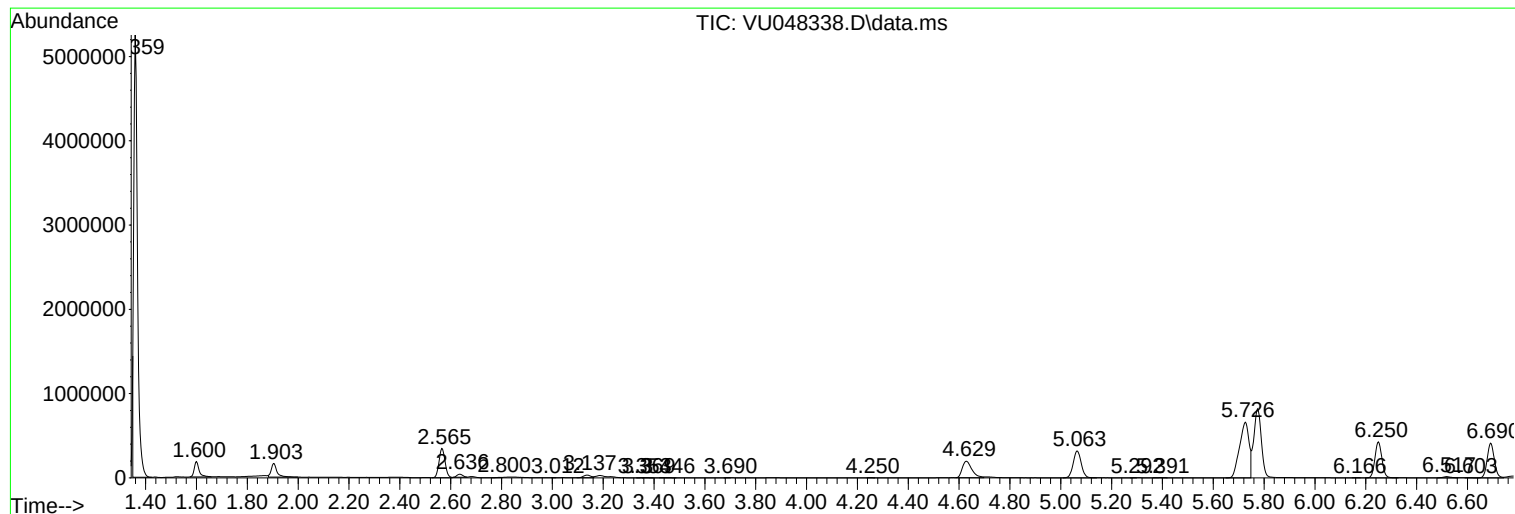
Sum of corrected areas: 21078891

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ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXET6DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM042522WMA.M
Quant Title : VOC Analysis

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



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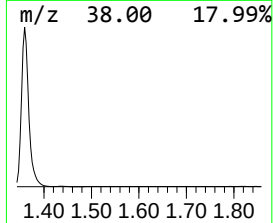
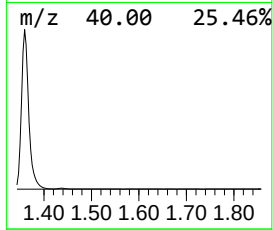
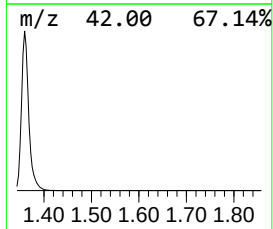
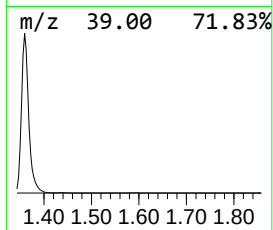
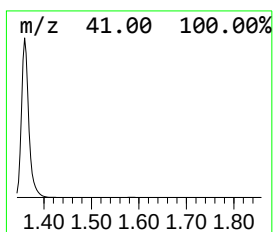
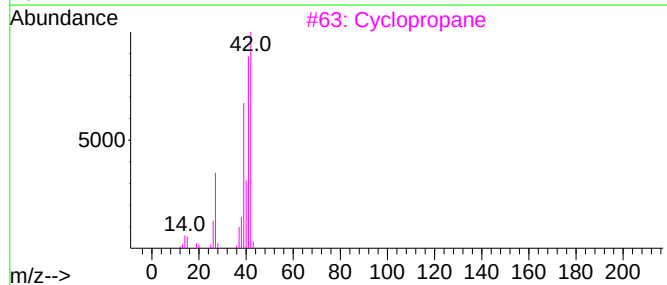
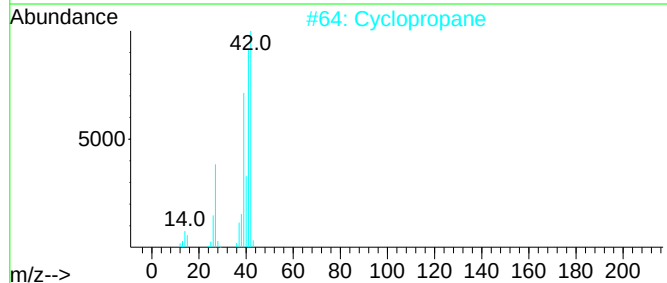
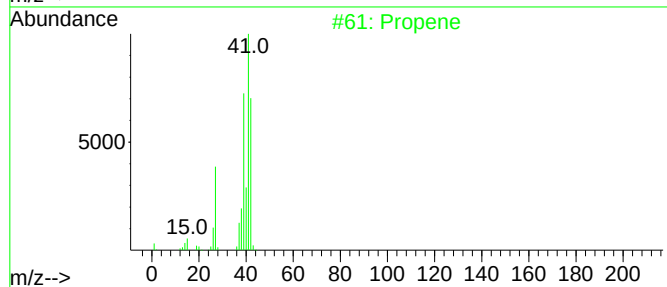
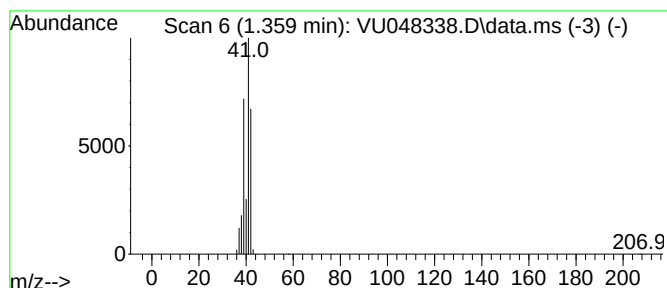
Instrument :
MSVOA_U
ClientSampleId :
EXET6DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM042522WMA.M
Quant Title : VOC Analysis

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
1.359	316.46 ug/L	5162850	1,4-Difluorobenzene	6.250	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Propene	42	C3H6	000115-07-1	90
2	Cyclopropane	42	C3H6	000075-19-4	86
3	Cyclopropane	42	C3H6	000075-19-4	78
4	Propene	42	C3H6	000115-07-1	50
5	Cyclopropene	40	C3H4	002781-85-3	16



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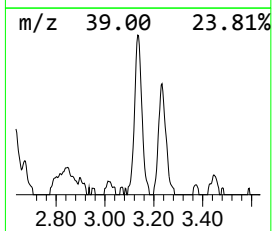
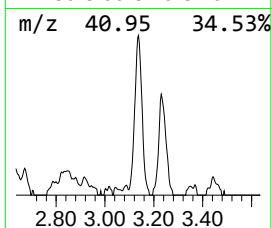
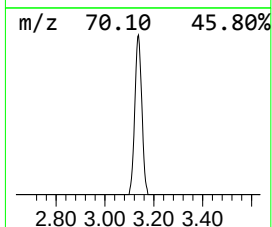
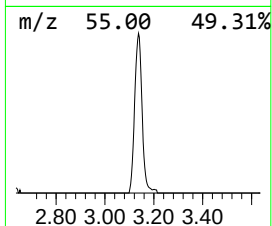
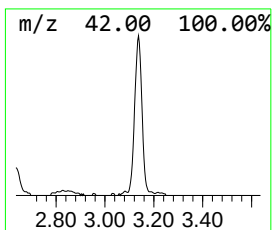
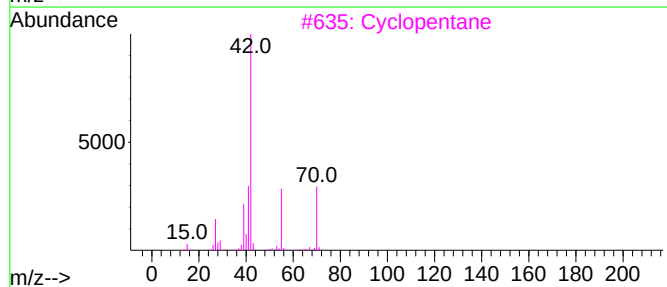
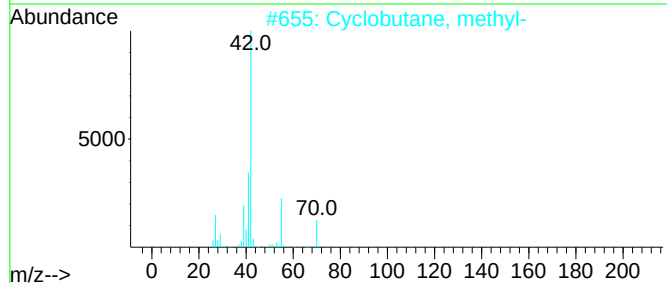
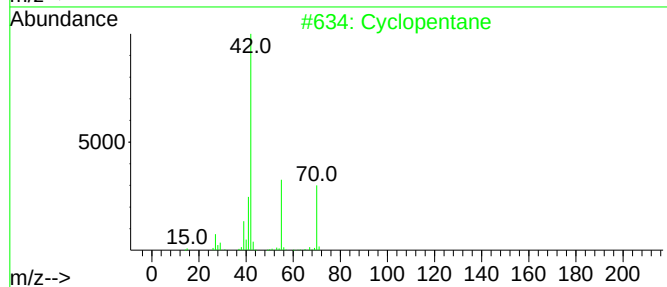
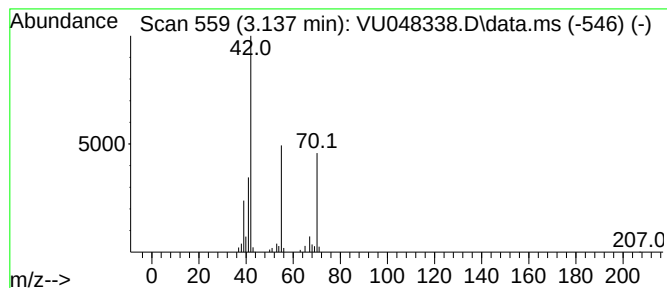
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM042522WMA.M
Quant Title : VOC Analysis

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

Peak Number 2 (DEL) Alkane: Cyclic3.137 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.137	4.03 ug/L	65803	1,4-Difluorobenzene	6.250

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane	70	C5H10	000287-92-3	86
2			Cyclobutane, methyl-	70	C5H10	000598-61-8	80
3			Cyclopentane	70	C5H10	000287-92-3	78
4			Cyclobutane, methyl-	70	C5H10	000598-61-8	72
5			1-Pentene	70	C5H10	000109-67-1	64



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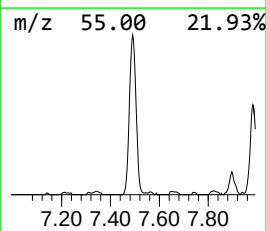
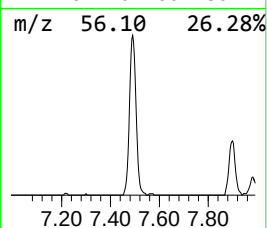
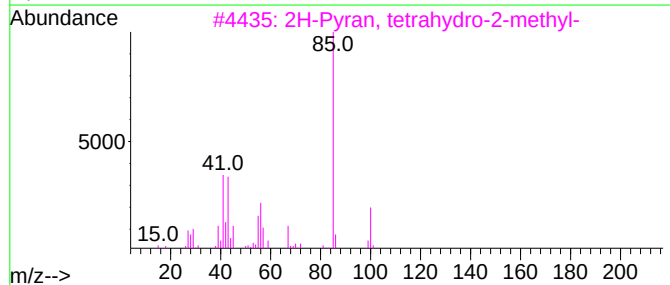
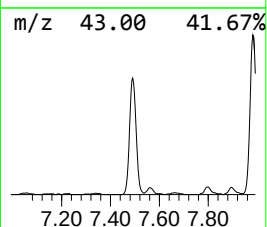
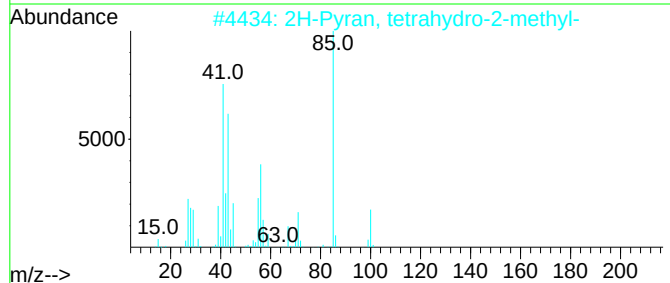
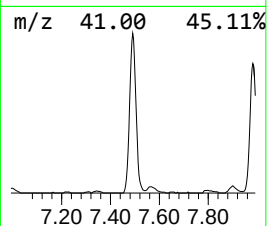
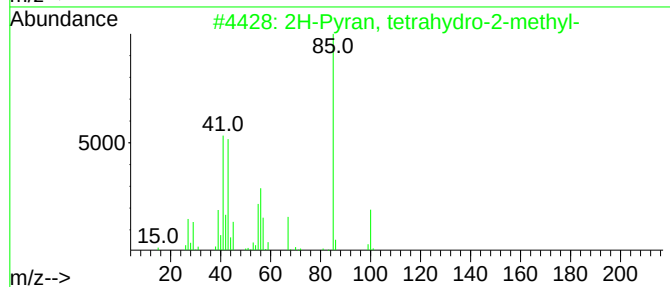
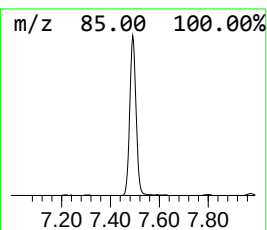
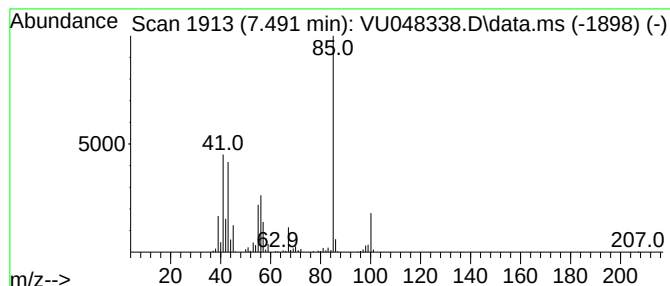
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM042522WMA.M
Quant Title : VOC Analysis

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

Peak Number 3 2H-Pyran, tetrahydro-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.491	21.55 ug/L	351589	1,4-Difluorobenzene	6.250

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2H-Pyran, tetrahydro-2-methyl-			100	C6H12O	010141-72-7	97
2	2H-Pyran, tetrahydro-2-methyl-			100	C6H12O	010141-72-7	91
3	2H-Pyran, tetrahydro-2-methyl-			100	C6H12O	010141-72-7	90
4	2H-Pyran, tetrahydro-2-methyl-			100	C6H12O	010141-72-7	83
5	2-Furanmethanol, tetrahydro-5-me...			116	C6H12O2	054774-28-6	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU042822\
Data File : VU048338.D
Acq On : 28 Apr 2022 14:04
Operator : SY/MD
Sample : N2587-10DL 40X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXET6DL

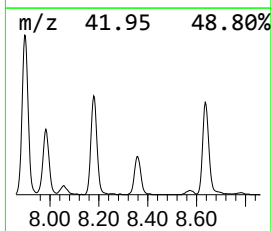
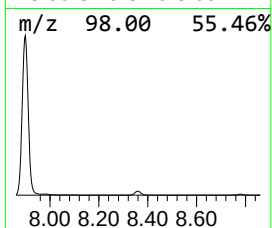
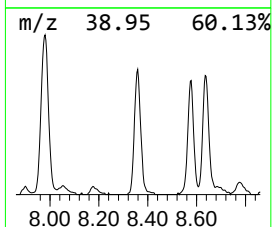
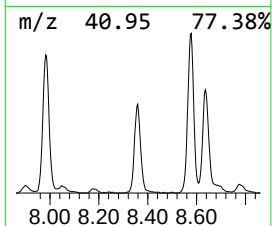
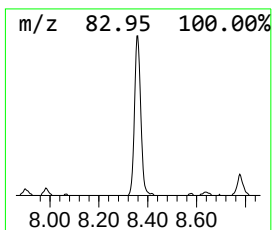
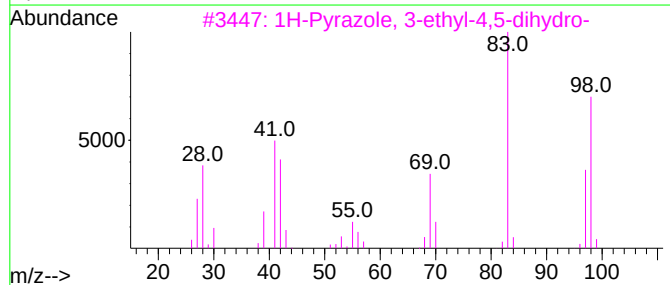
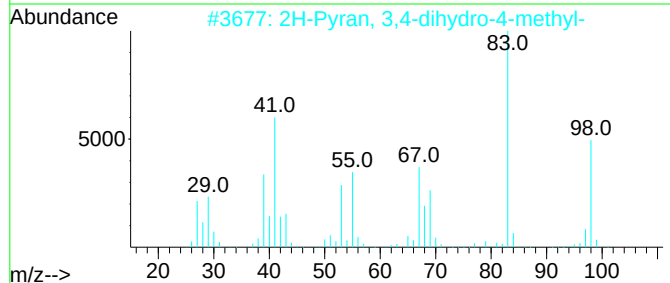
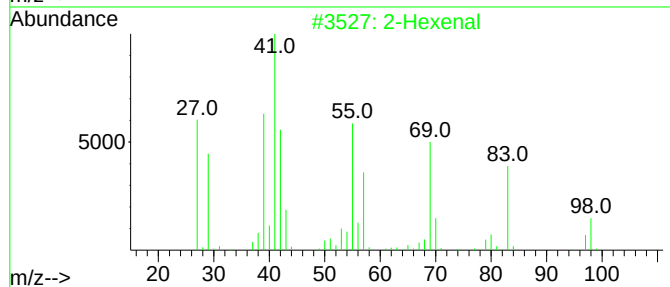
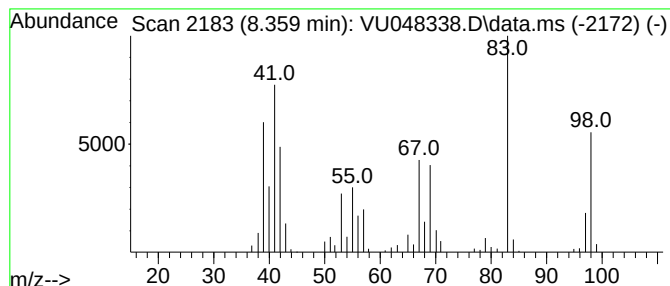
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Quant Title : VOC Analysis

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

Peak Number 5 2-Hexenal Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.359	10.48 ug/L	189250	Chlorobenzene-d5	9.417

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Hexenal	98	C6H10O	000505-57-7	59
2			2H-Pyran, 3,4-dihydro-4-methyl-	98	C6H10O	002270-61-3	58
3			1H-Pyrazole, 3-ethyl-4,5-dihydro-	98	C5H10N2	005920-29-6	49
4			1H-Pyrazole, 4,5-dihydro-4,5-dim...	98	C5H10N2	028019-94-5	47
5			1H-Pyrazole, 4,5-dihydro-5,5-dim...	98	C5H10N2	004320-85-8	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU042822\
Data File : VU048338.D
Acq On : 28 Apr 2022 14:04
Operator : SY/MD
Sample : N2587-10DL 40X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 9 Sample Multiplier: 1

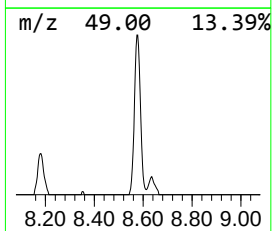
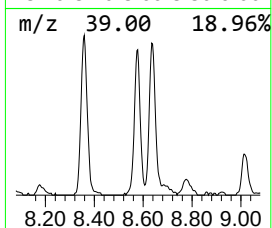
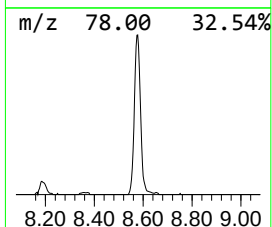
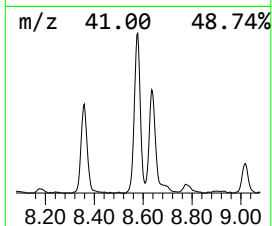
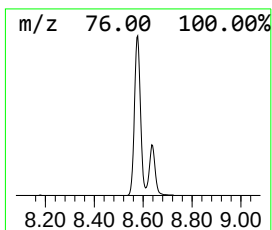
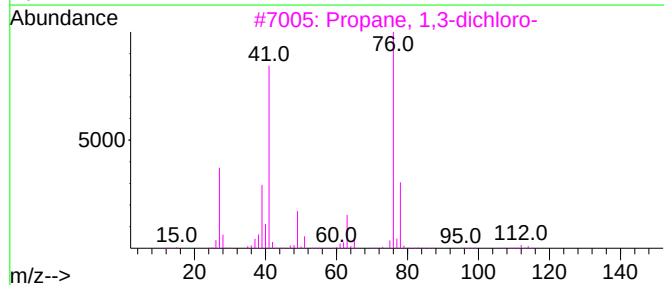
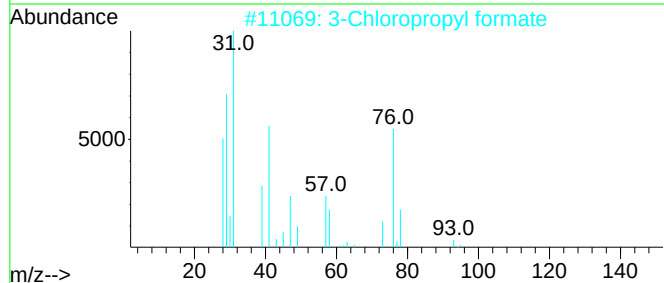
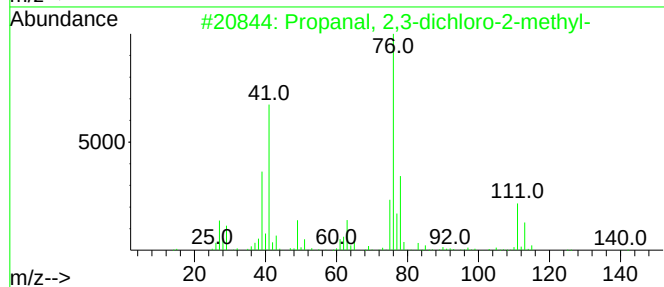
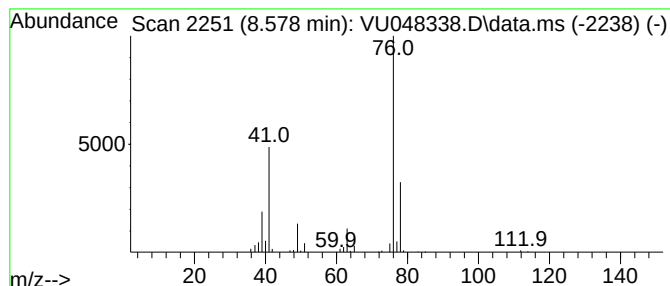
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ClientSampleId :
EXET6DL

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Quant Title : VOC Analysis

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

Peak Number 6 Propanal, 2,3-dichloro-2-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD			R.T.
8.578	11.29 ug/L	203869	Chlorobenzene-d5			9.417
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Propanal, 2,3-dichloro-2-methyl-		140	C4H6Cl2O	010141-22-7	74
2	3-Chloropropyl formate		122	C4H7ClO2	001487-44-1	43
3	Propane, 1,3-dichloro-		112	C3H6Cl2	000142-28-9	25
4	Propane, 1,3-dichloro-		112	C3H6Cl2	000142-28-9	15
5	Propane, 1,3-dichloro-		112	C3H6Cl2	000142-28-9	10



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU042822\
Data File : VU048338.D
Acq On : 28 Apr 2022 14:04
Operator : SY/MD
Sample : N2587-10DL 40X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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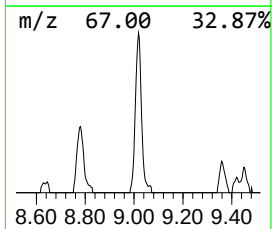
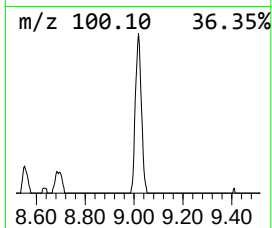
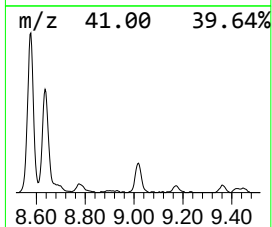
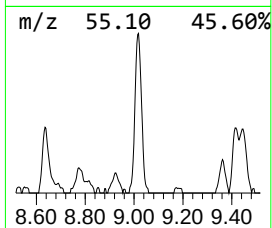
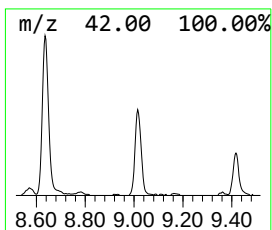
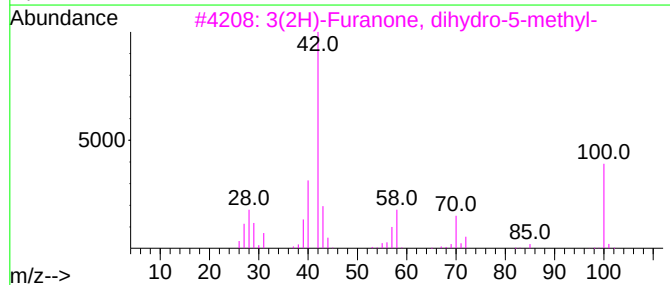
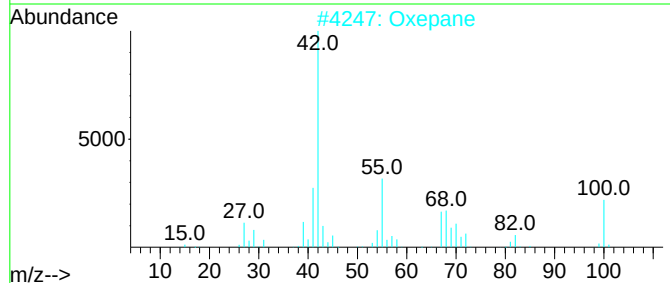
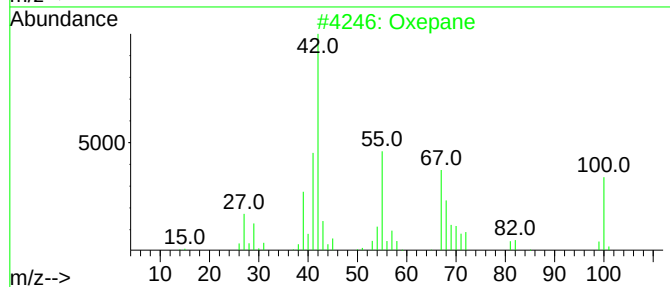
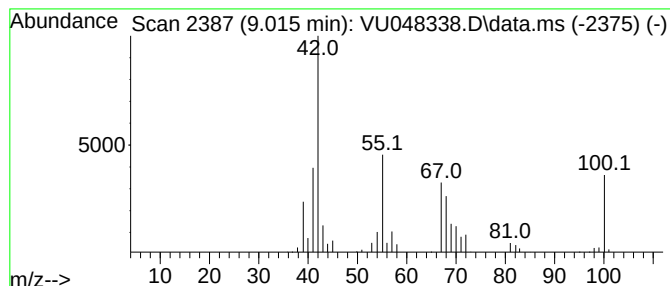
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Quant Title : VOC Analysis

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

Peak Number 7 Oxepane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.015	4.36 ug/L	78812	Chlorobenzene-d5	9.417

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxepane			100	C6H12O	000592-90-5	94
2	Oxepane			100	C6H12O	000592-90-5	74
3	3(2H)-Furanone, dihydro-5-methyl-			100	C5H8O2	034003-72-0	52
4	2,4(3H,5H)-Furandione			100	C4H4O3	004971-56-6	46
5	3-Ethyltetrahydrofuran			100	C6H12O	093716-28-0	42



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU042822\
Data File : VU048338.D
Acq On : 28 Apr 2022 14:04
Operator : SY/MD
Sample : N2587-10DL 40X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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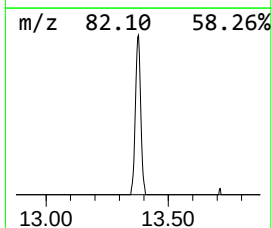
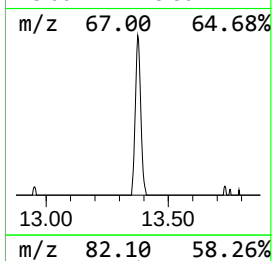
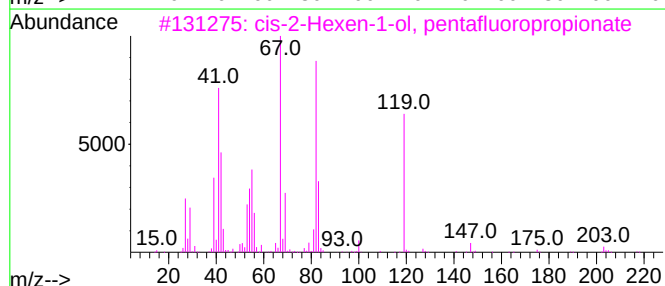
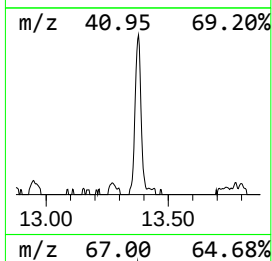
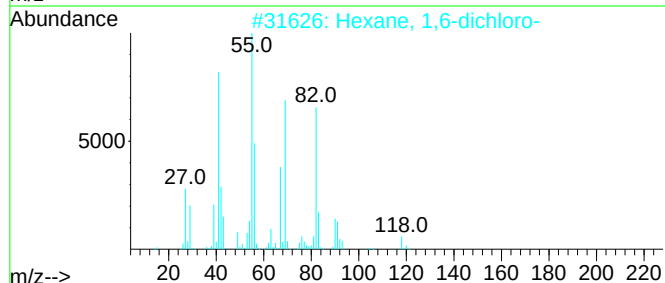
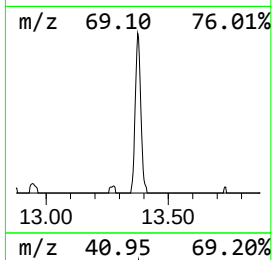
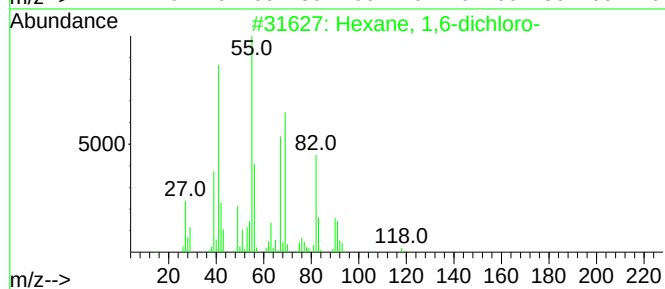
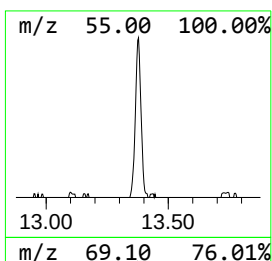
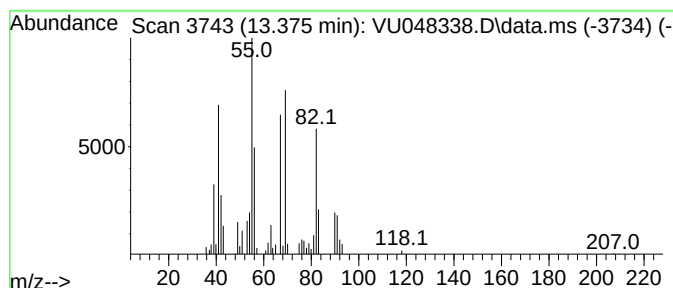
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Quant Title : VOC Analysis

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

Peak Number 8 Hexane, 1,6-dichloro- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.375	2.61 ug/L	53041	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 1,6-dichloro-	154	C6H12Cl2	002163-00-0	86
2			Hexane, 1,6-dichloro-	154	C6H12Cl2	002163-00-0	64
3			cis-2-Hexen-1-ol, pentafluoropro...	246	C9H11F5O2	1000352-74-1	38
4			cis-2-Hexen-1-ol, trifluoroacetate	196	C8H11F3O2	1000352-74-4	38
5			Cyclohexane, chloro-	118	C6H11Cl	000542-18-7	30



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU042822\
Data File : VU048338.D
Acq On : 28 Apr 2022 14:04
Operator : SY/MD
Sample : N2587-10DL 40X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXET6DL

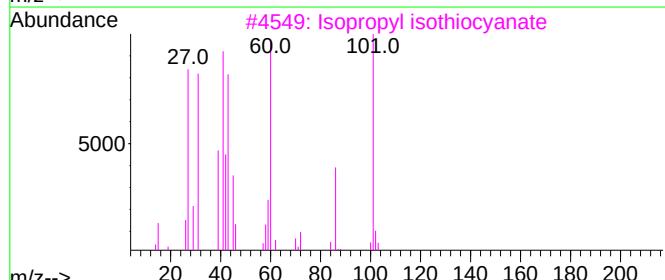
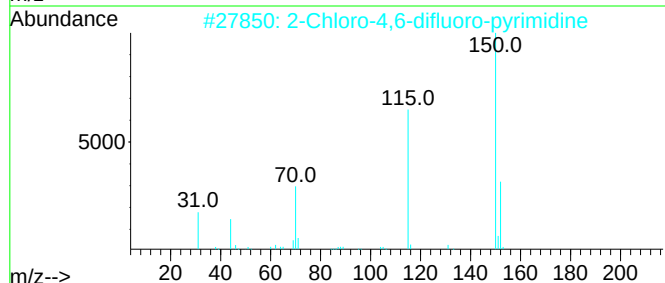
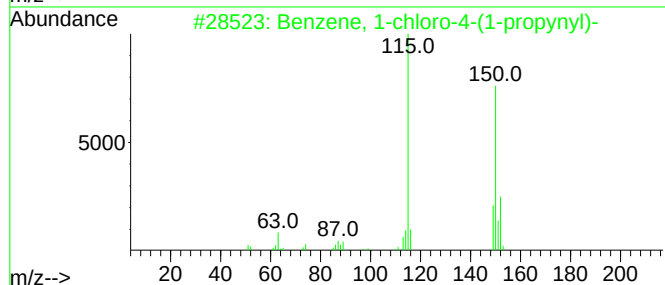
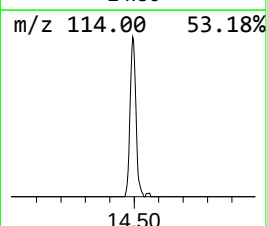
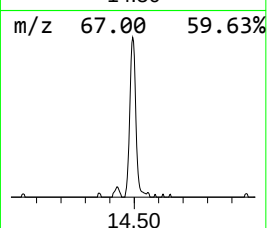
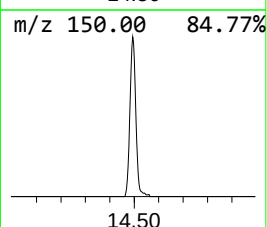
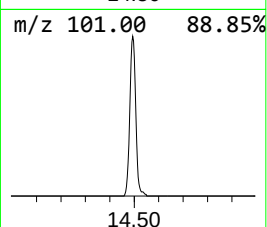
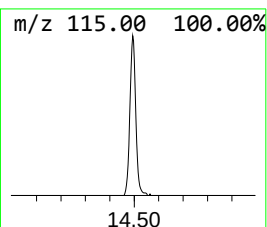
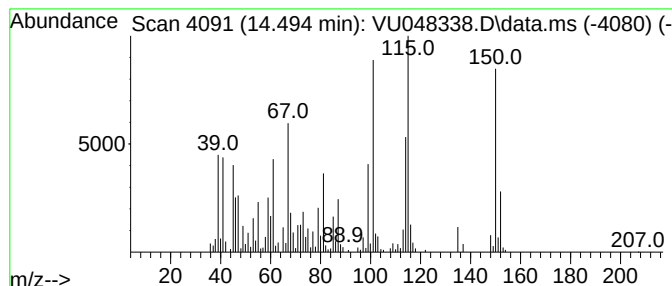
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Quant Title : VOC Analysis

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

Peak Number 9 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.494	11.03 ug/L	223959	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-chloro-4-(1-propynyl)-	150	C9H7Cl	002809-65-6	38
2			2-Chloro-4,6-difluoro-pyrimidine	150	C4HClF2N2	038953-30-9	27
3			Isopropyl isothiocyanate	101	C4H7NS	002253-73-8	15
4			Silane, ethenylethoxydimethyl-	130	C6H14OSi	005356-83-2	14
5			2-Piperidinethione	115	C5H9NS	013070-01-4	11



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU042822\
Data File : VU048338.D
Acq On : 28 Apr 2022 14:04
Operator : SY/MD
Sample : N2587-10DL 40X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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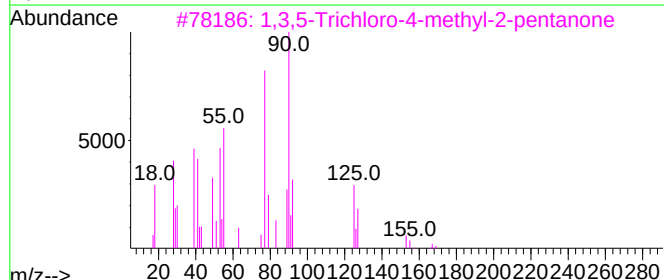
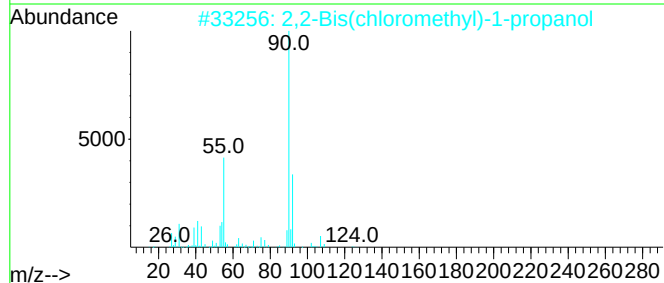
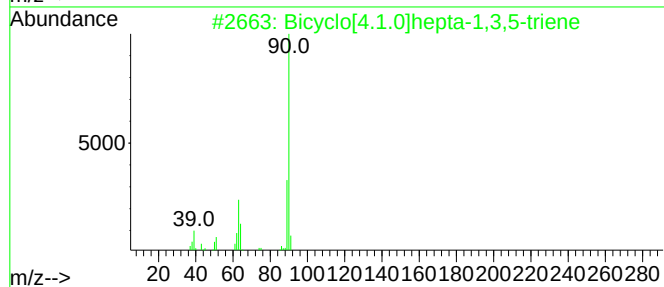
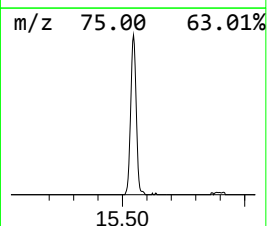
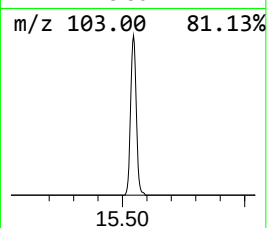
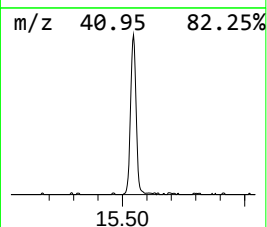
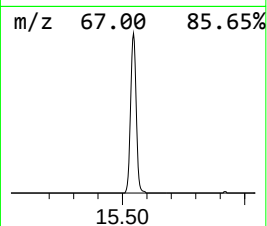
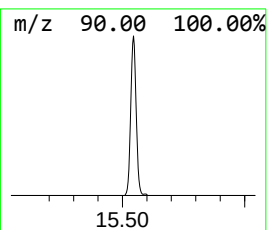
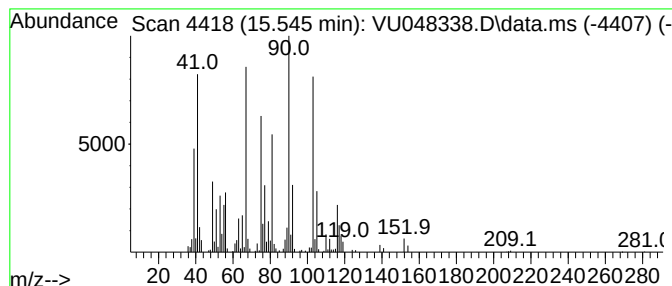
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Quant Title : VOC Analysis

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

Peak Number 10 unknown-02 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.545	13.22 ug/L	268400	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[4.1.0]hepta-1,3,5-triene	90	C7H6	004646-69-9	35
2			2,2-Bis(chloromethyl)-1-propanol	156	C5H10Cl2O	005355-54-4	32
3			1,3,5-Trichloro-4-methyl-2-penta...	202	C6H9Cl3O	1000140-03-8	28
4			4-(2-Aminoethyl)benzenesulfonamide	200	C8H12N2O2S	035303-76-5	25
5			3-Thietanol	90	C3H6OS	010304-16-2	25



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU042822\
Data File : VU048338.D
Acq On : 28 Apr 2022 14:04
Operator : SY/MD
Sample : N2587-10DL 40X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXET6DL

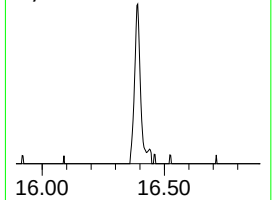
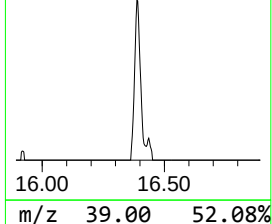
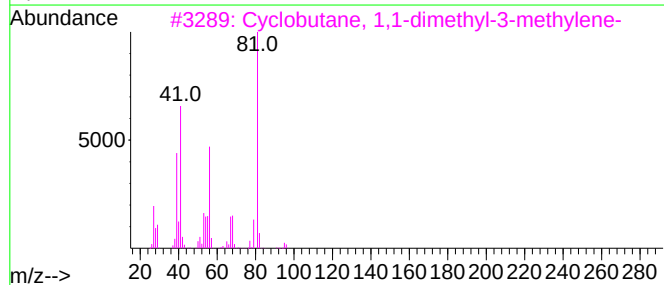
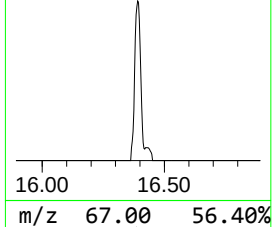
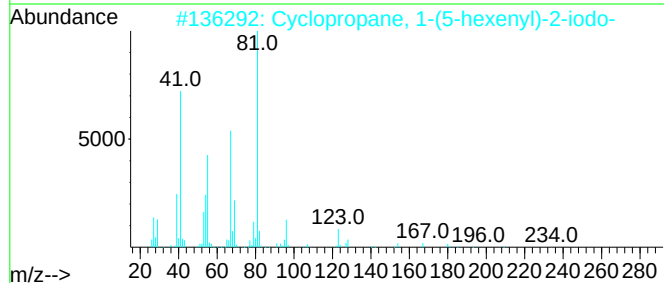
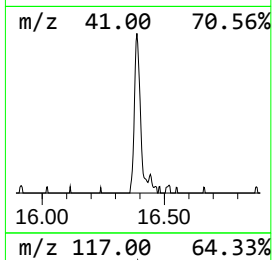
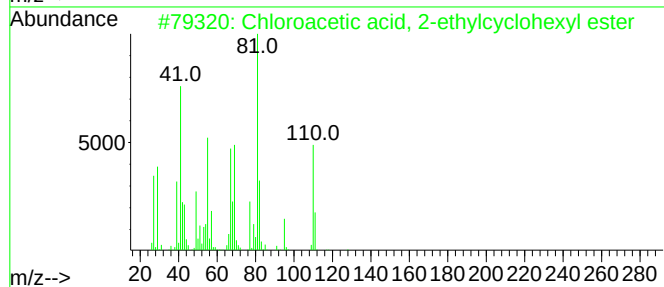
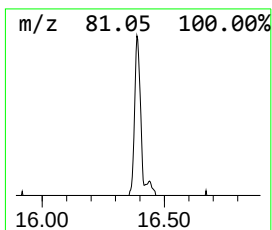
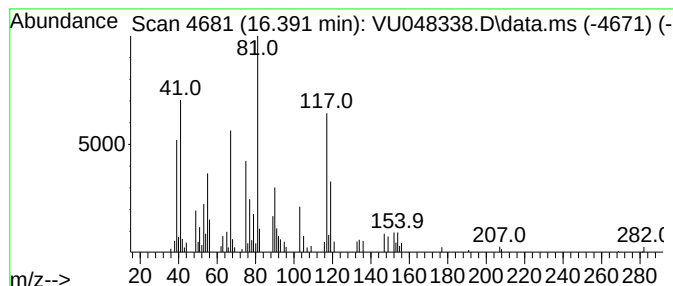
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM042522WMA.M
Quant Title : VOC Analysis

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

Peak Number 11 unknown-03 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.391	2.79 ug/L	56553	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Chloroacetic acid, 2-ethylcyclohexyl ester	204	C10H17ClO2	1000279-89-1	37
2			Cyclopropane, 1-(5-hexenyl)-2-iodo-	250	C9H15I	074685-40-8	12
3			Cyclobutane, 1,1-dimethyl-3-methylene-	96	C7H12	019037-73-1	11
4			1H-Pyrrole, 1-methyl-	81	C5H7N	000096-54-8	10
5			1,5-Hexadiene, 2-methyl-	96	C7H12	004049-81-4	10



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU042822\
 Data File : VU048338.D
 Acq On : 28 Apr 2022 14:04
 Operator : SY/MD
 Sample : N2587-10DL 40X
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 EXET6DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM042522WMA.M
 Quant Title : VOC Analysis

TIC Library : C:\Database\NIST0.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: S...	1.359	316.5	ug/L	5162850	1	6.250	815722	50.0
(DEL) Alkane: C...	3.137	4.0	ug/L	65803	1	6.250	815722	50.0
2H-Pyran, tetra...	7.491	21.6	ug/L	351589	1	6.250	815722	50.0
2-Hexenal	8.359	10.5	ug/L	189250	2	9.417	903017	50.0
Propanal, 2,3-d...	8.578	11.3	ug/L	203869	2	9.417	903017	50.0
Oxepane	9.015	4.4	ug/L	78812	2	9.417	903017	50.0
Hexane, 1,6-dic...	13.375	2.6	ug/L	53041	3	11.812	1015030	50.0
unknown-01	14.494	11.0	ug/L	223959	3	11.812	1015030	50.0
unknown-02	15.545	13.2	ug/L	268400	3	11.812	1015030	50.0
unknown-03	16.391	2.8	ug/L	56553	3	11.812	1015030	50.0