

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU042822\
 Data File : VU048336.D
 Acq On : 28 Apr 2022 13:17
 Operator : SY/MD
 Sample : N2587-08DL 40X
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 EXET4DL

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: VU048336.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	25	rVB	776021	743682	19.59%	2.827%
2	1.475	35	42	52	rBV3	39335	58283	1.53%	0.222%
3	1.600	73	81	100	rBV	185402	247591	6.52%	0.941%
4	1.896	165	173	189	rBV	128147	190308	5.01%	0.723%
5	2.565	369	381	395	rBV	334722	542494	14.29%	2.062%
6	2.784	434	449	450	rBV3	3193	5261	0.14%	0.020%
7	3.006	510	518	521	rBV4	875	1181	0.03%	0.004%
8	3.044	521	530	541	rBV7	2017	4425	0.12%	0.017%
9	3.186	553	574	600	rVB	34820	92544	2.44%	0.352%
10	3.356	624	627	630	rBV	606	475	0.01%	0.002%
11	3.446	646	655	664	rBV4	1872	3409	0.09%	0.013%
12	3.498	669	671	679	rVB3	594	557	0.01%	0.002%
13	3.687	727	730	732	rBV3	630	456	0.01%	0.002%
14	3.703	733	735	740	rVB4	891	586	0.02%	0.002%
15	3.861	780	784	786	rVV2	501	430	0.01%	0.002%
16	3.877	786	789	793	rVV2	876	748	0.02%	0.003%
17	3.896	793	795	802	rVB2	750	680	0.02%	0.003%
18	4.227	891	898	900	rBV3	613	657	0.02%	0.002%
19	4.530	984	992	996	rBV3	968	1215	0.03%	0.005%
20	4.620	1005	1020	1045	rBV	192249	467010	12.30%	1.775%
21	5.063	1140	1158	1177	rBV	300843	653335	17.21%	2.483%
22	5.282	1222	1226	1228	rBV2	547	432	0.01%	0.002%
23	5.330	1238	1241	1243	rBV2	657	468	0.01%	0.002%
24	5.378	1253	1256	1258	rBV3	777	608	0.02%	0.002%
25	5.726	1338	1364	1372	rBV2	626126	1697133	44.70%	6.451%
26	6.250	1507	1527	1548	rBV	403414	763231	20.10%	2.901%
27	6.472	1592	1596	1597	rVV	909	590	0.02%	0.002%
28	6.513	1597	1609	1624	rVB2	32341	65574	1.73%	0.249%
29	6.690	1649	1664	1679	rBV	388147	754743	19.88%	2.869%
30	6.774	1682	1690	1707	rVB2	33745	69124	1.82%	0.263%
31	6.996	1743	1759	1774	rBV	46610	92187	2.43%	0.350%
32	7.179	1805	1816	1823	rBV6	3022	6332	0.17%	0.024%
33	7.298	1846	1853	1861	rVB6	1876	2723	0.07%	0.010%
34	7.488	1897	1912	1925	rBV	332154	599278	15.78%	2.278%
35	7.565	1925	1936	1953	rVB	229194	399299	10.52%	1.518%
36	7.690	1967	1975	1985	rVB7	1649	2806	0.07%	0.011%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

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Sampling : 1

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Stop Thrs : 0

Peak Location: TOP

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Peak separation: 5

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

37	7.787	2000	2005	2007	rBV3	1570	1231	0.03%	0.005%
38	7.896	2025	2039	2053	rBV	802179	1396320	36.77%	5.307%
39	7.980	2054	2065	2086	rVB	480768	859115	22.63%	3.265%
40	8.179	2114	2127	2144	rBV	155557	260218	6.85%	0.989%
41	8.356	2175	2182	2190	rBV6	3484	6055	0.16%	0.023%
42	8.472	2213	2218	2222	rBV2	384	449	0.01%	0.002%
43	8.510	2225	2230	2232	rBV3	1303	1071	0.03%	0.004%
44	8.574	2235	2250	2260	rVV	2281660	3797021	100.00%	14.432%
45	8.632	2260	2268	2297	rVB	639234	1155762	30.44%	4.393%
46	8.980	2372	2376	2379	rVB2	571	533	0.01%	0.002%
47	9.018	2379	2388	2398	rBV6	2976	4695	0.12%	0.018%
48	9.169	2422	2435	2448	rBV3	12208	22673	0.60%	0.086%
49	9.362	2484	2495	2502	rBV2	46329	74143	1.95%	0.282%
50	9.446	2502	2521	2545	rVB2	1521263	3464911	91.25%	13.170%
51	9.578	2554	2562	2571	rVB3	21130	32554	0.86%	0.124%
52	9.632	2571	2579	2590	rVB6	7783	13545	0.36%	0.051%
53	9.693	2590	2598	2608	rVB4	2055	3633	0.10%	0.014%
54	9.845	2634	2645	2655	rVV5	3297	6320	0.17%	0.024%
55	9.941	2664	2675	2686	rVV	34240	57453	1.51%	0.218%
56	10.108	2715	2727	2737	rBV7	1978	3619	0.10%	0.014%
57	10.250	2764	2771	2781	rBV5	955	1462	0.04%	0.006%
58	10.340	2796	2799	2804	rBV3	506	428	0.01%	0.002%
59	10.484	2839	2844	2846	rVV	818	742	0.02%	0.003%
60	10.578	2867	2873	2874	rBV2	410	444	0.01%	0.002%
61	10.652	2887	2896	2909	rBV5	12356	23186	0.61%	0.088%
62	10.706	2909	2913	2915	rVV2	1390	1271	0.03%	0.005%
63	10.758	2916	2929	2954	rVB2	656550	1219857	32.13%	4.636%
64	10.893	2965	2971	2977	rBV4	1947	2414	0.06%	0.009%
65	10.973	2985	2996	3006	rBV3	10286	16174	0.43%	0.061%
66	11.195	3056	3065	3072	rBV5	2018	2661	0.07%	0.010%
67	11.250	3074	3082	3091	rVB5	5052	7166	0.19%	0.027%
68	11.375	3114	3121	3131	rBV6	2438	3581	0.09%	0.014%
69	11.462	3139	3148	3156	rBV10	4060	7479	0.20%	0.028%
70	11.510	3156	3163	3173	rVB4	7398	11901	0.31%	0.045%
71	11.584	3178	3186	3189	rBV5	1319	1506	0.04%	0.006%
72	11.700	3213	3222	3226	rVV6	1448	2403	0.06%	0.009%
73	11.751	3230	3238	3245	rVV4	4623	7268	0.19%	0.028%
74	11.812	3245	3257	3270	rVV	655389	1046265	27.55%	3.977%
75	11.880	3270	3278	3290	rVV2	43709	76815	2.02%	0.292%
76	11.954	3293	3301	3315	rVV6	11504	21758	0.57%	0.083%

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

77	12.021	3315	3322	3330	rVB3	6315	8948	0.24%	0.034%
78	12.098	3344	3346	3350	rBV4	797	654	0.02%	0.002%
79	12.195	3360	3376	3394	rBV	856153	1449435	38.17%	5.509%
80	12.330	3413	3418	3421	rBV4	945	953	0.03%	0.004%
81	12.401	3435	3440	3443	rBV3	623	429	0.01%	0.002%
82	12.471	3449	3462	3475	rVV2	30734	45717	1.20%	0.174%
83	12.706	3525	3535	3547	rVB	38720	59055	1.56%	0.224%
84	12.774	3547	3556	3566	rVB3	2320	3836	0.10%	0.015%
85	12.970	3605	3617	3629	rBV	50283	82103	2.16%	0.312%
86	13.160	3665	3676	3683	rBV7	1094	2184	0.06%	0.008%
87	13.221	3690	3695	3698	rBV5	961	1082	0.03%	0.004%
88	13.269	3705	3710	3712	rBV3	953	1080	0.03%	0.004%
89	13.333	3721	3730	3732	rBV4	4636	5799	0.15%	0.022%
90	13.375	3732	3743	3764	rVB	750346	1122318	29.56%	4.266%
91	13.716	3843	3849	3858	rBV4	1057	1402	0.04%	0.005%
92	13.780	3862	3869	3872	rBV5	1644	2089	0.06%	0.008%
93	14.092	3961	3966	3974	rBV2	1247	1354	0.04%	0.005%
94	14.359	4036	4049	4060	rBV4	9792	15607	0.41%	0.059%
95	14.430	4060	4071	4081	rBV2	107238	166711	4.39%	0.634%
96	14.494	4081	4091	4109	rVB	217093	358678	9.45%	1.363%
97	14.950	4225	4233	4248	rBV7	13363	25906	0.68%	0.098%
98	15.278	4331	4335	4340	rBV7	1555	1839	0.05%	0.007%
99	15.545	4405	4418	4436	rBV	1099008	1684946	44.38%	6.404%
100	16.391	4670	4681	4691	rBV2	109040	181788	4.79%	0.691%

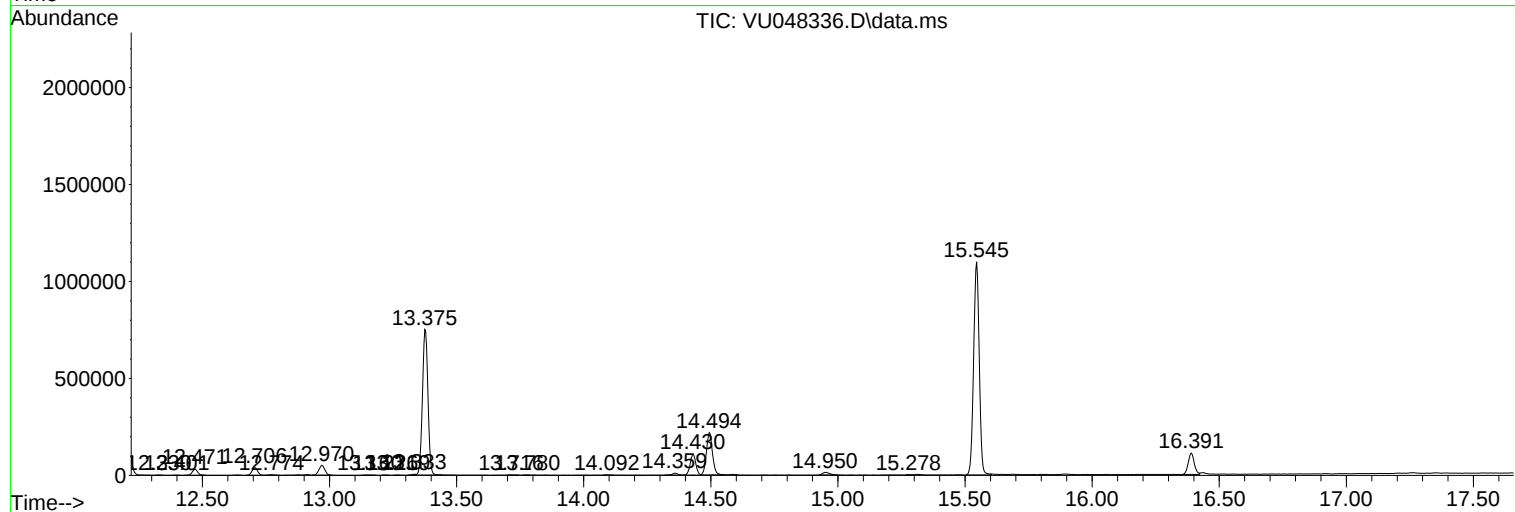
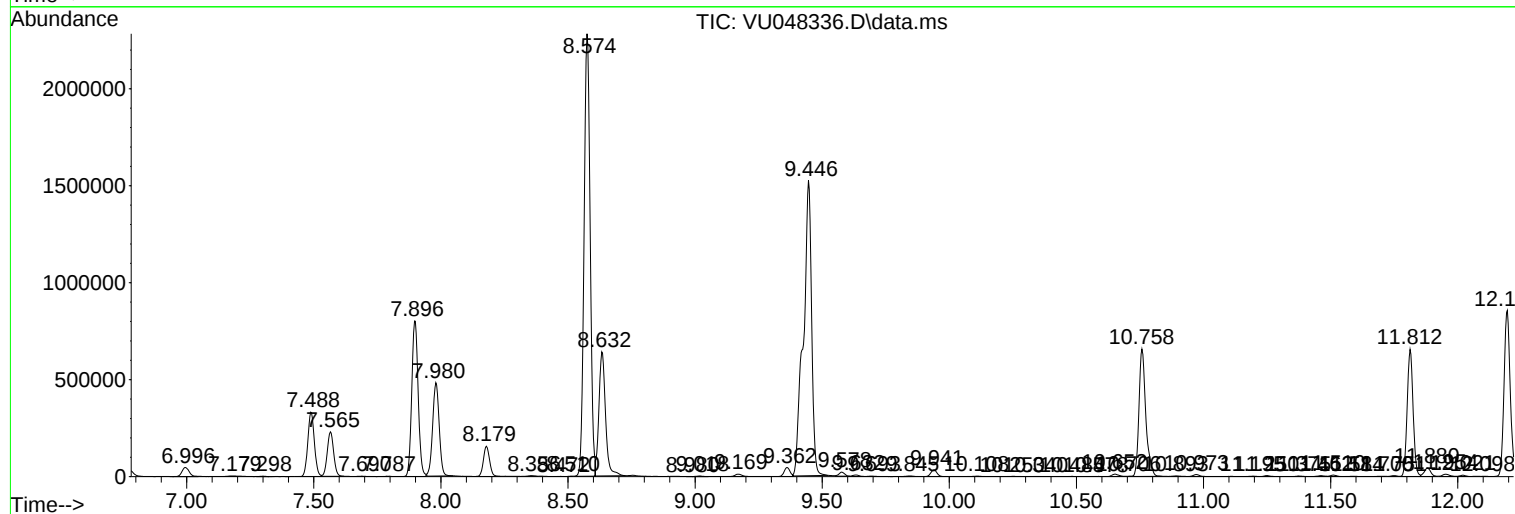
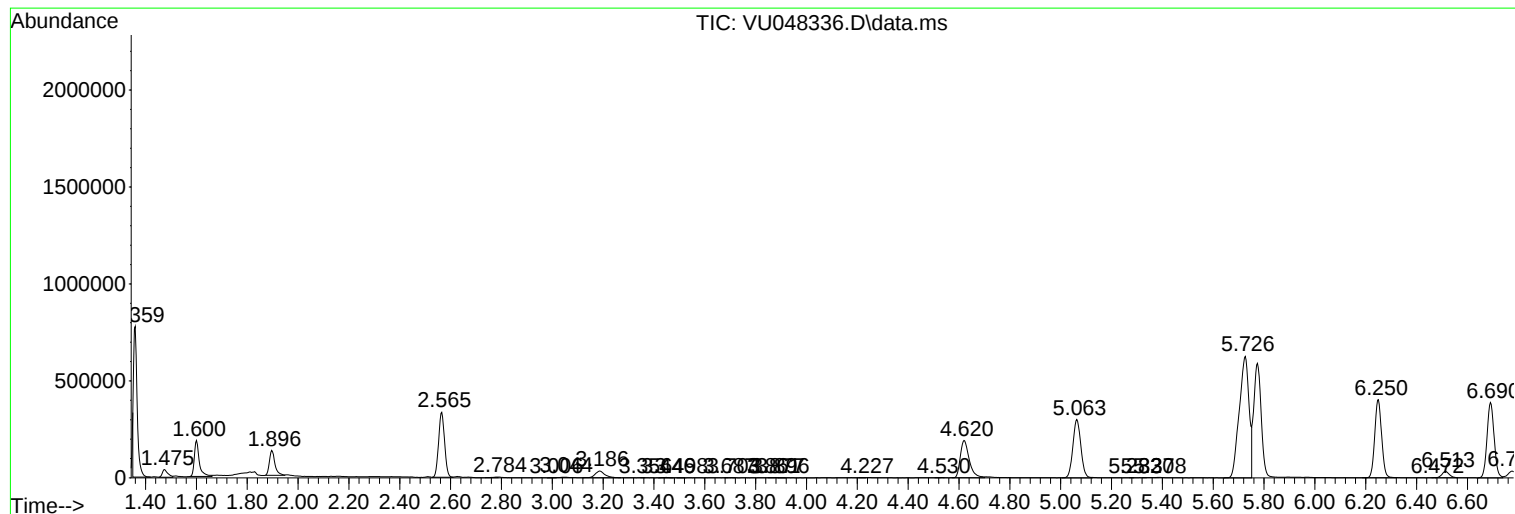
Sum of corrected areas: 26309870

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ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXET4DL

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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Misc : 5.0mL/MSVOA_U/WATER
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Instrument :
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ClientSampleId :
EXET4DL

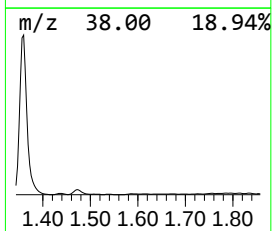
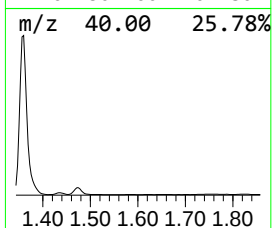
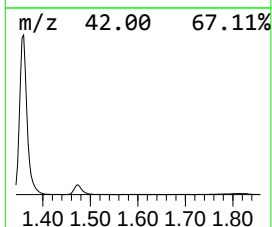
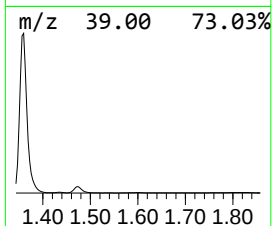
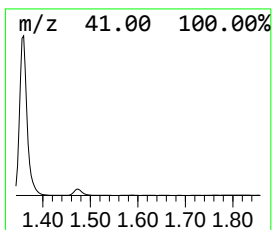
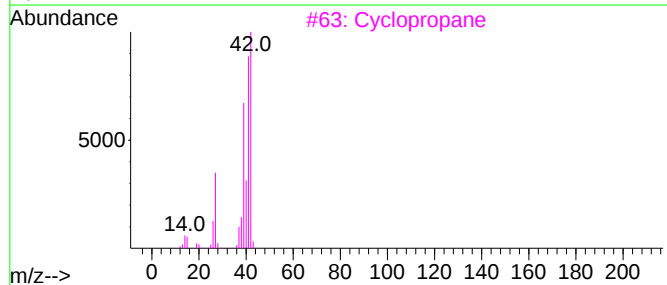
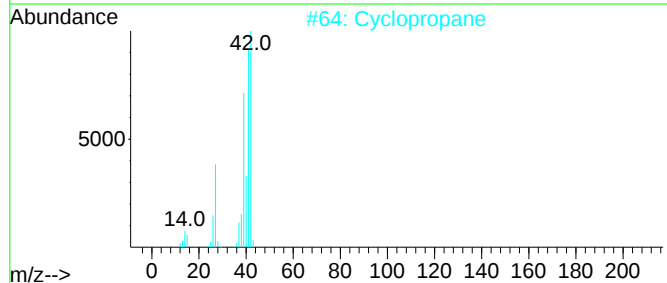
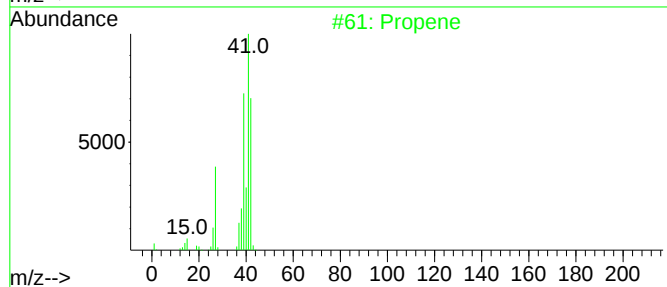
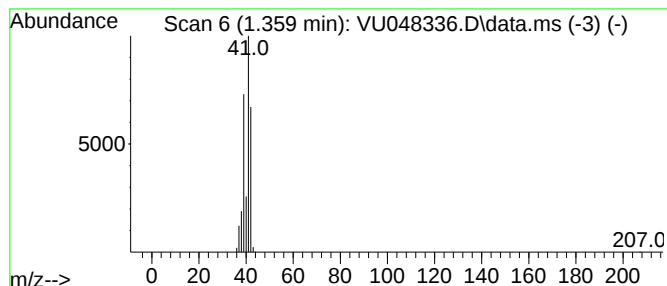
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM042522WMA.M
Quant Title : VOC Analysis

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.359	48.72 ug/L	743682	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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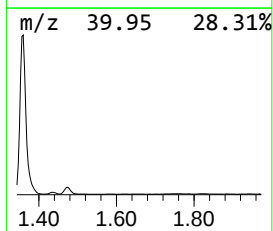
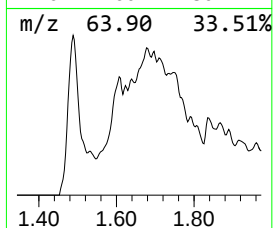
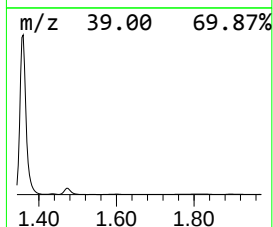
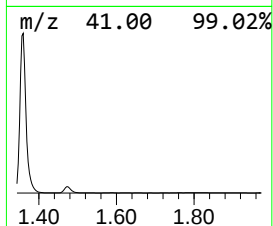
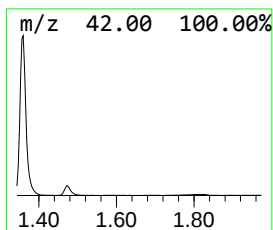
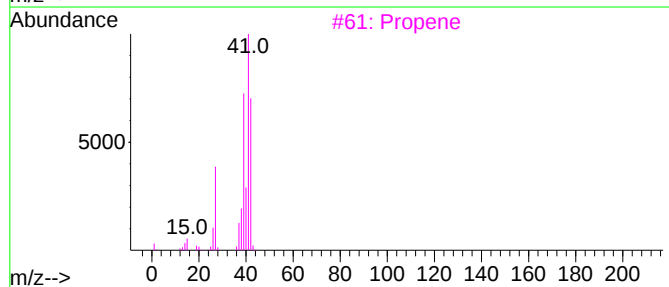
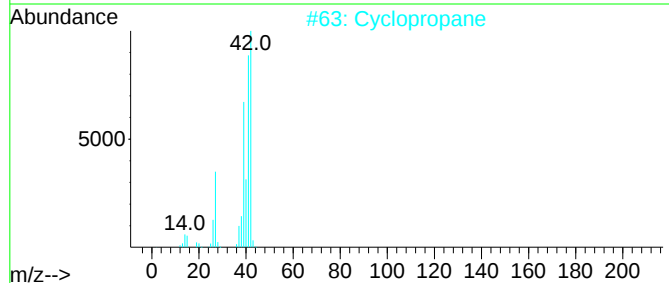
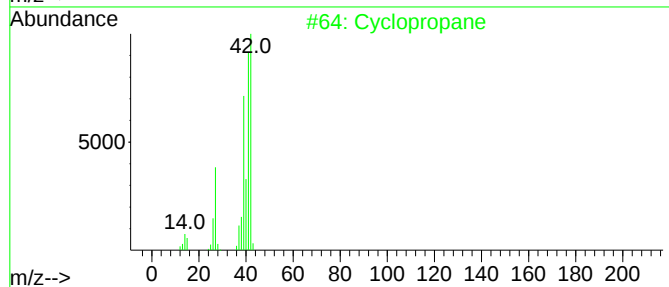
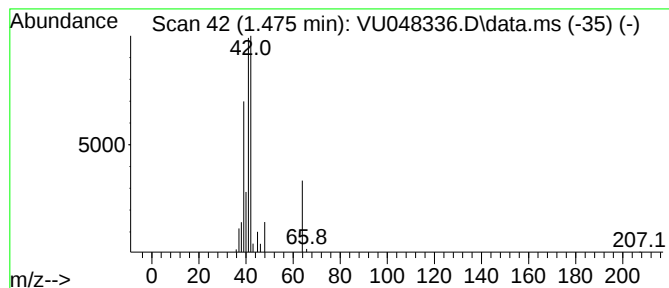
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM042522WMA.M
Quant Title : VOC Analysis

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

Peak Number 2 (DEL) Alkane: Cyclic1.475 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.475	3.82 ug/L	58283	1,4-Difluorobenzene	6.250

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopropane	42	C3H6	000075-19-4	72
2			Cyclopropane	42	C3H6	000075-19-4	72
3			Propene	42	C3H6	000115-07-1	64
4			2-Oxetanone, 4-methylene-	84	C4H4O2	000674-82-8	47
5			Cyclopropane	42	C3H6	000075-19-4	47



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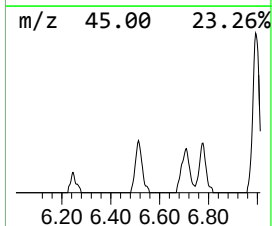
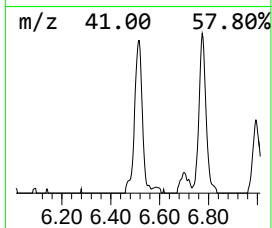
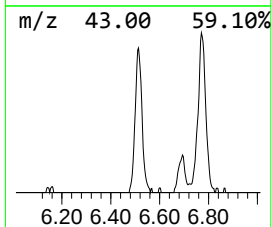
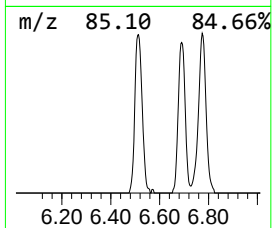
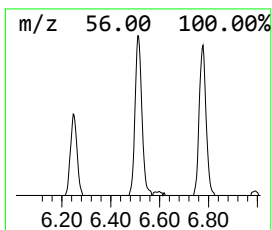
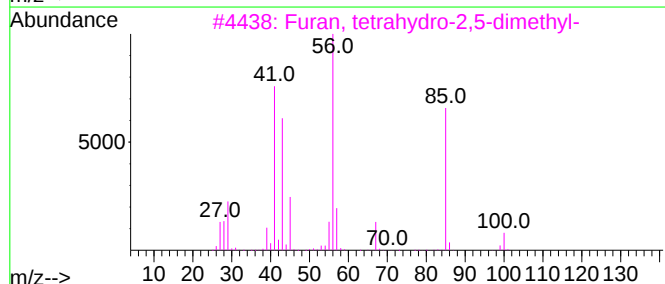
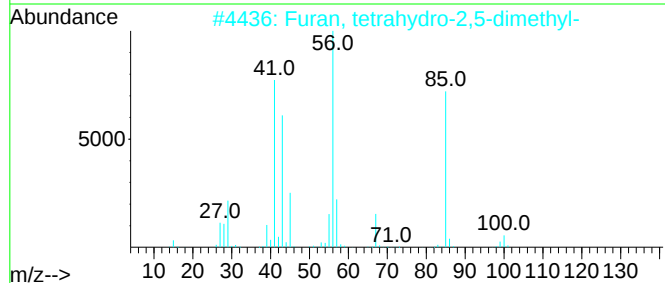
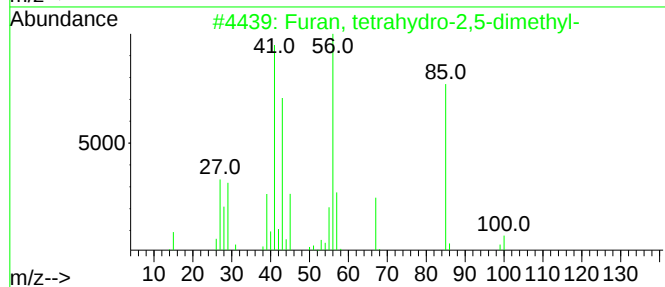
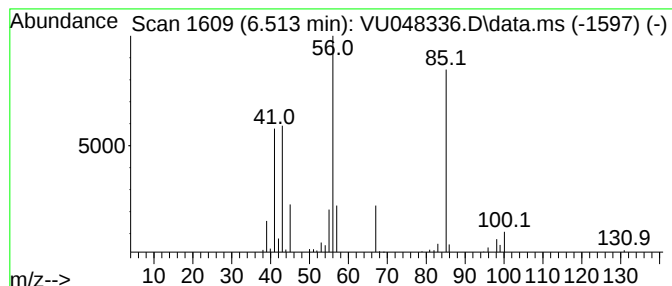
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM042522WMA.M
Quant Title : VOC Analysis

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

Peak Number 4 Furan, tetrahydro-2,5-dimet... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.513	4.30 ug/L	65574	1,4-Difluorobenzene	6.250

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Furan, tetrahydro-2,5-dimethyl-	100	C6H12O	001003-38-9	91
2			Furan, tetrahydro-2,5-dimethyl-	100	C6H12O	001003-38-9	91
3			Furan, tetrahydro-2,5-dimethyl-	100	C6H12O	001003-38-9	87
4			2(3H)-Furanone, dihydro-5-methyl-	100	C5H8O2	000108-29-2	72
5			Furan, tetrahydro-2,5-dimethyl-,...	100	C6H12O	038484-59-2	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU042822\
Data File : VU048336.D
Acq On : 28 Apr 2022 13:17
Operator : SY/MD
Sample : N2587-08DL 40X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXET4DL

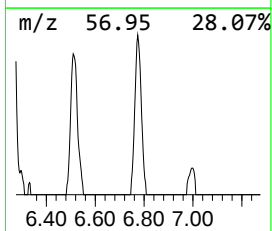
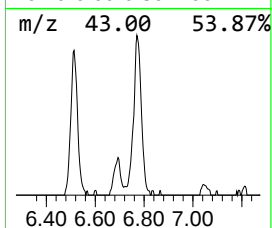
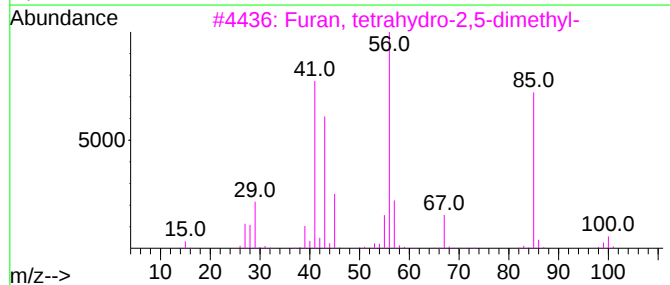
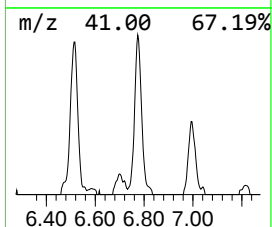
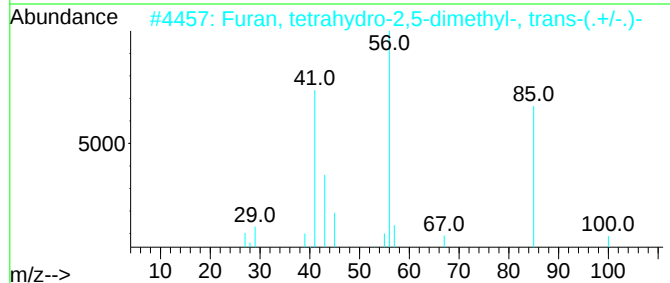
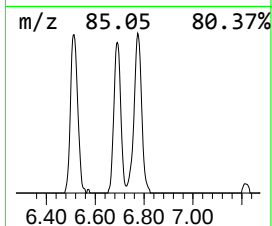
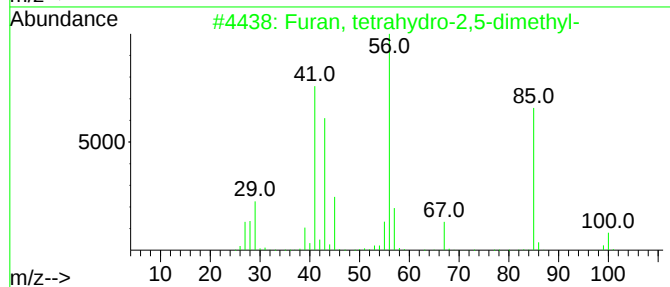
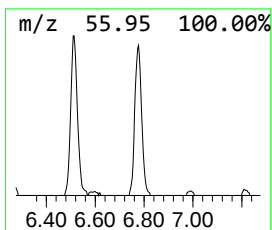
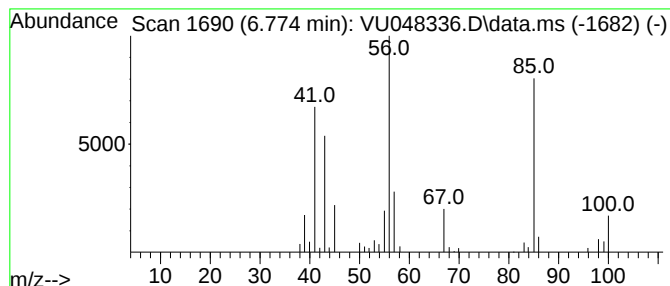
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Quant Title : VOC Analysis

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

Peak Number 5 Furan, tetrahydro-2,5-dimet... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.774	4.53 ug/L	69124	1,4-Difluorobenzene	6.250

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Furan, tetrahydro-2,5-dimethyl-	100	C6H12O	001003-38-9	83
2			Furan, tetrahydro-2,5-dimethyl-,...	100	C6H12O	038484-59-2	80
3			Furan, tetrahydro-2,5-dimethyl-	100	C6H12O	001003-38-9	78
4			Furan, tetrahydro-2,5-dimethyl-	100	C6H12O	001003-38-9	78
5			2(3H)-Furanone, dihydro-5-methyl-	100	C5H8O2	000108-29-2	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU042822\
Data File : VU048336.D
Acq On : 28 Apr 2022 13:17
Operator : SY/MD
Sample : N2587-08DL 40X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXET4DL

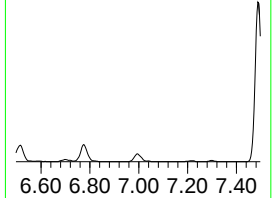
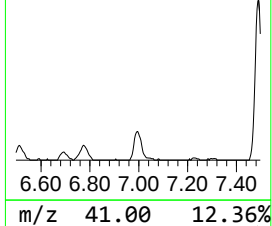
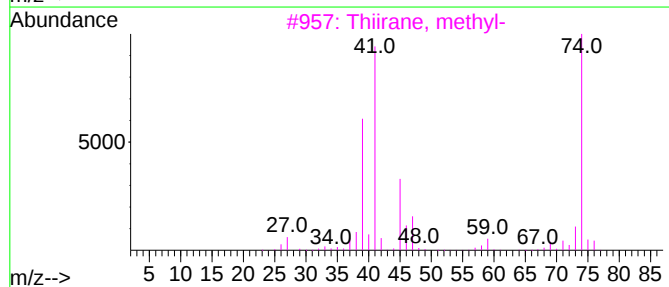
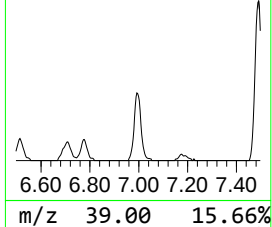
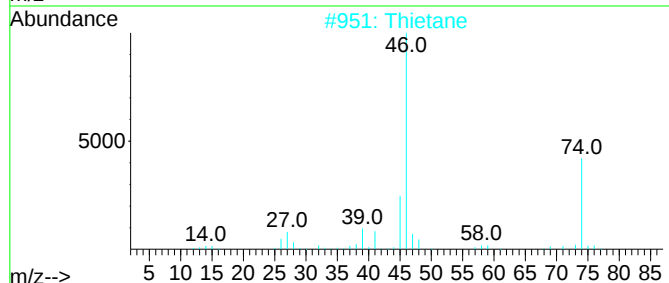
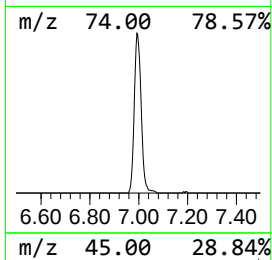
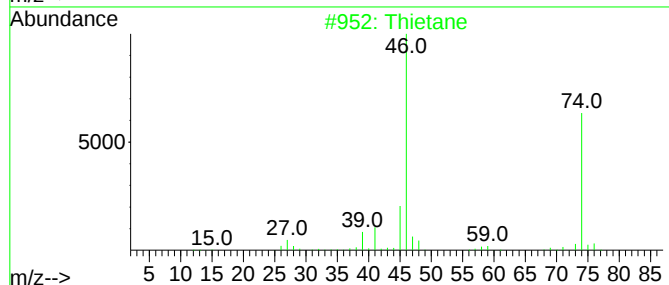
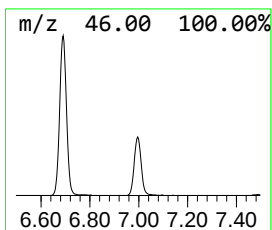
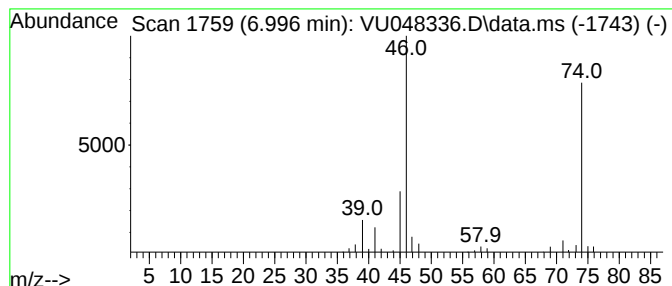
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Quant Title : VOC Analysis

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

Peak Number 6 Thietane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.996	6.04 ug/L	92187	1,4-Difluorobenzene	6.250

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Thietane	74	C3H6S	000287-27-4	91
2			Thietane	74	C3H6S	000287-27-4	90
3			Thiirane, methyl-	74	C3H6S	001072-43-1	25
4			Allyl mercaptan	74	C3H6S	000870-23-5	9
5			Thiirane, methyl-	74	C3H6S	001072-43-1	9



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU042822\
Data File : VU048336.D
Acq On : 28 Apr 2022 13:17
Operator : SY/MD
Sample : N2587-08DL 40X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXET4DL

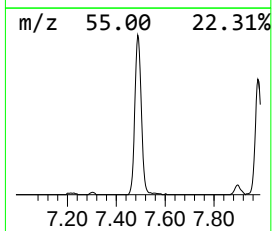
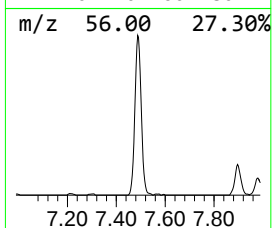
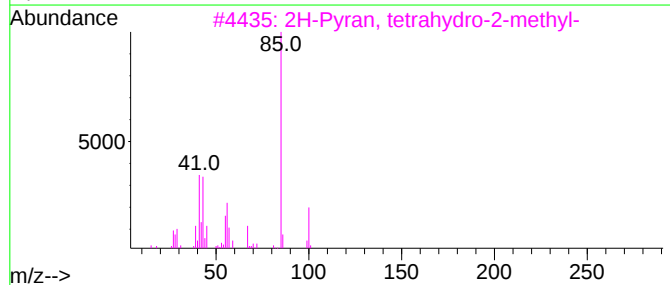
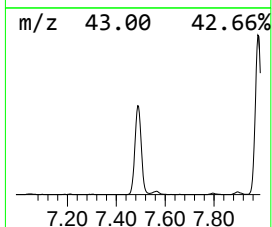
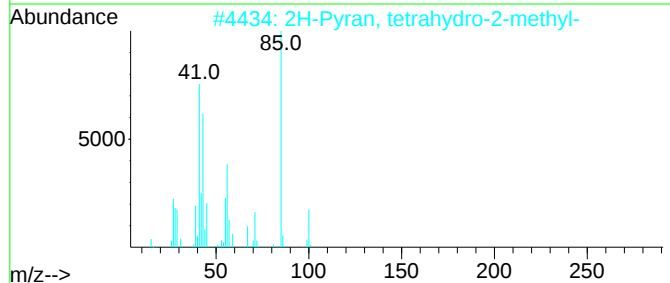
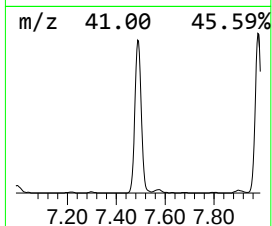
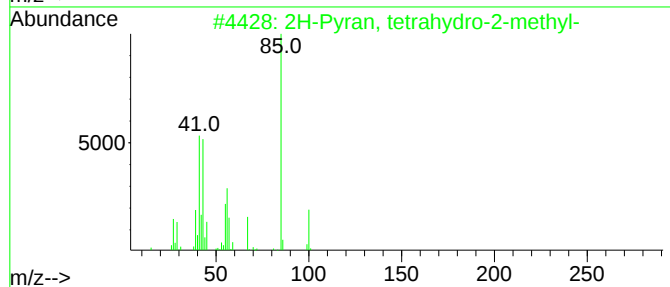
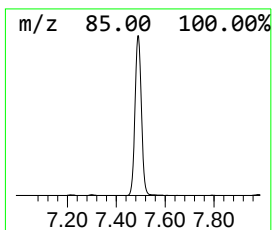
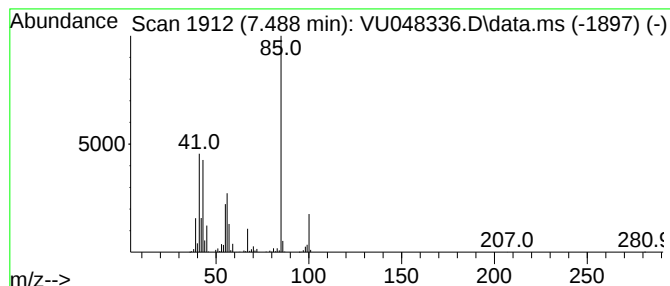
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Quant Title : VOC Analysis

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

Peak Number 7 2H-Pyran, tetrahydro-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.488	39.26 ug/L	599278	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	2-Furanmethanol, tetrahydro-5-me...	116	C6H12O2	054774-28-6	59		
5	2H-Pyran, tetrahydro-2-methyl-	100	C6H12O	010141-72-7	58		



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU042822\
Data File : VU048336.D
Acq On : 28 Apr 2022 13:17
Operator : SY/MD
Sample : N2587-08DL 40X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXET4DL

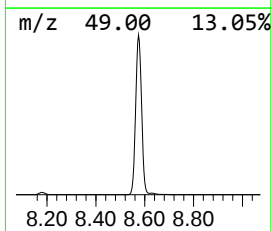
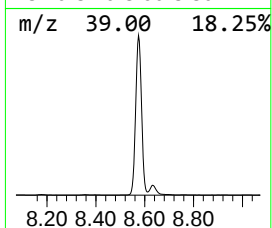
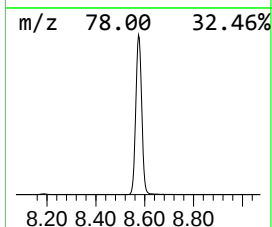
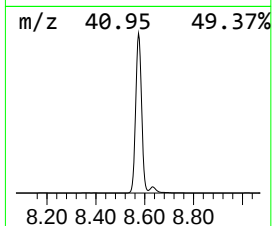
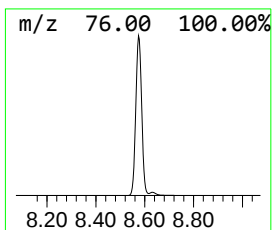
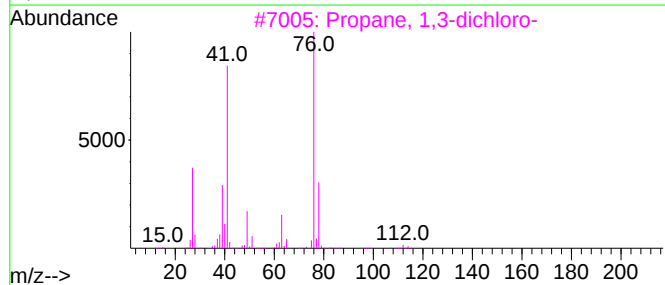
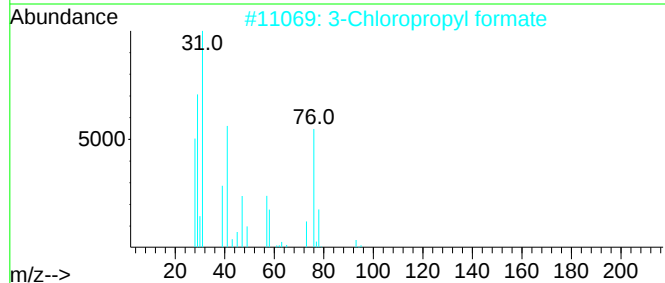
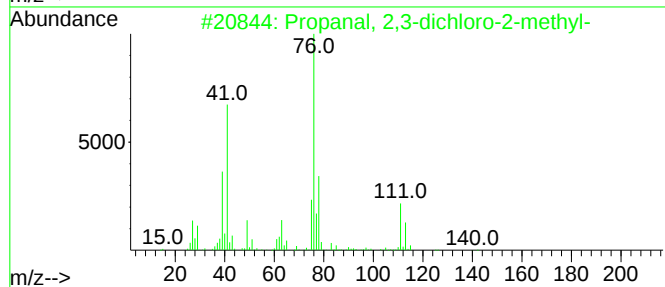
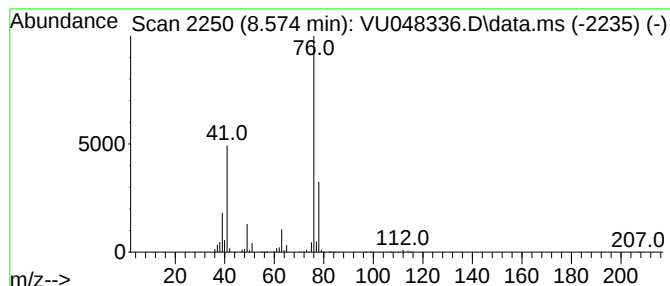
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Quant Title : VOC Analysis

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

Peak Number 9 Propanal, 2,3-dichloro-2-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.574	54.79 ug/L	3797020	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 : VU048336.D
Acq On : 28 Apr 2022 13:17
Operator : SY/MD
Sample : N2587-08DL 40X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXET4DL

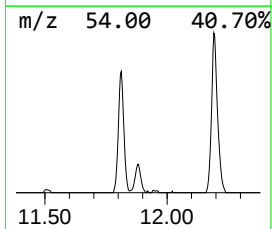
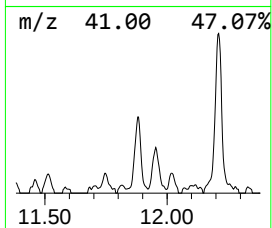
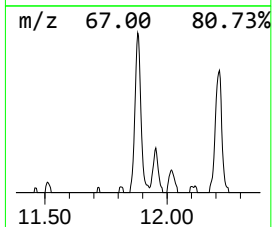
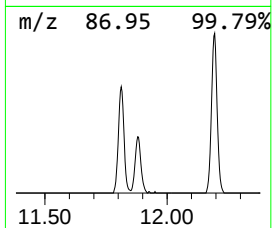
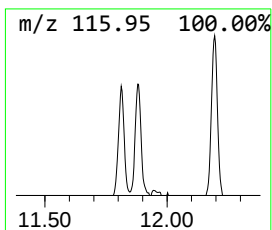
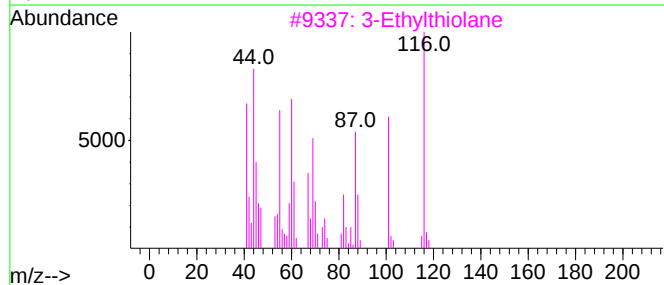
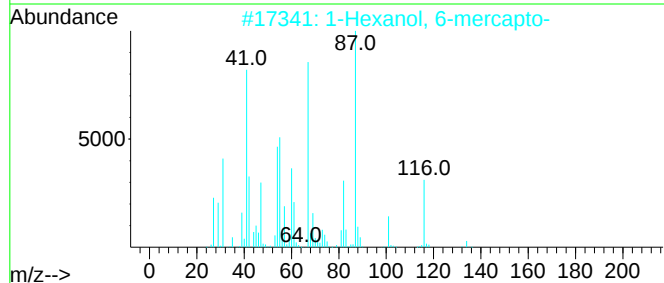
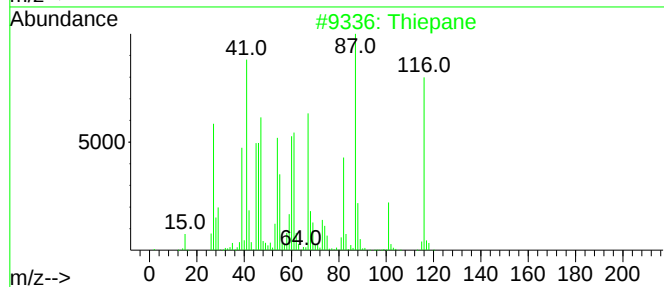
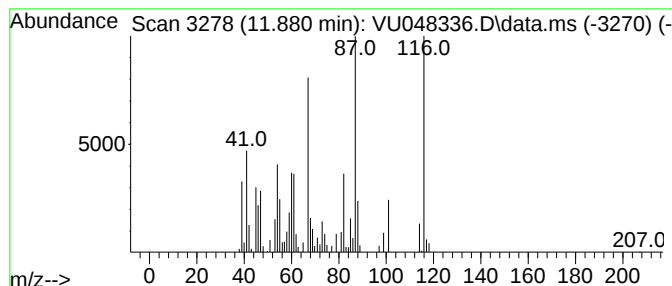
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Quant Title : VOC Analysis

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

Peak Number 10 Thiepane Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.880	3.67 ug/L	76815	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Thiepane	116	C6H12S	004753-80-4	94
2			1-Hexanol, 6-mercapto-	134	C6H14OS	001633-78-9	50
3			3-Ethylthiolane	116	C6H12S	062184-67-2	38
4			Hydrazine, 1-ethyl-1-(1-methylpr...	116	C6H16N2	020325-97-7	38
5			Cyclopentanethiol, 2-methyl-, tr...	116	C6H12S	001638-93-3	37



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU042822\
Data File : VU048336.D
Acq On : 28 Apr 2022 13:17
Operator : SY/MD
Sample : N2587-08DL 40X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXET4DL

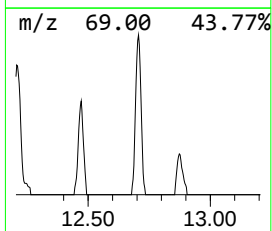
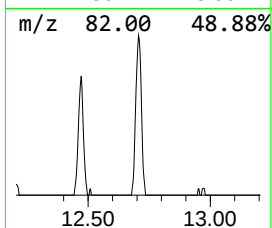
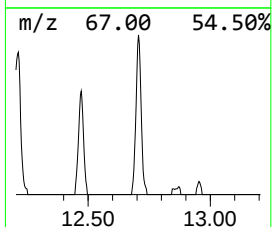
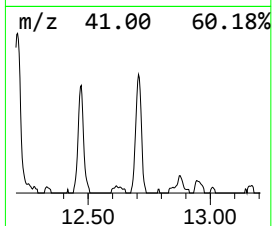
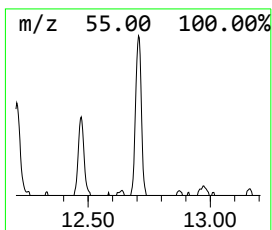
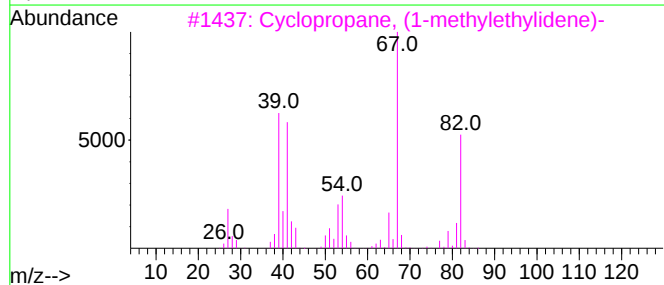
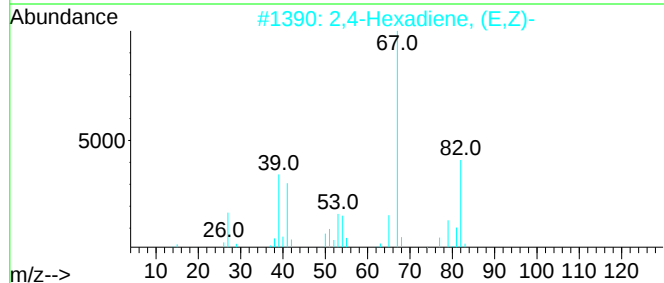
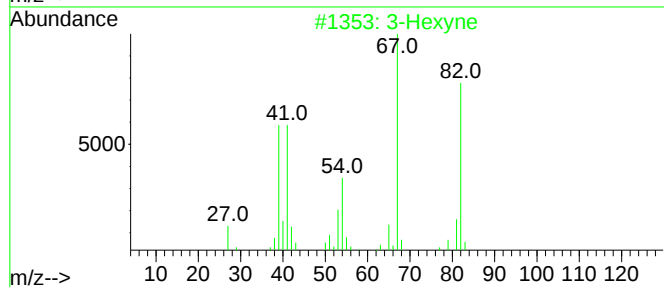
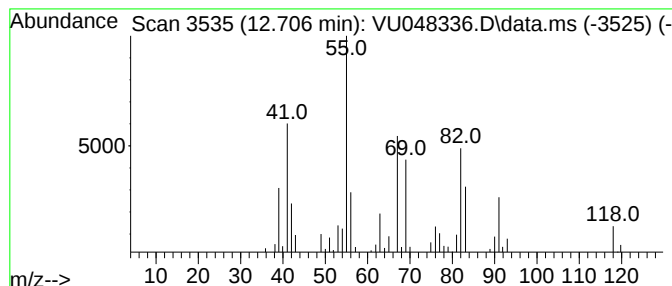
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Quant Title : VOC Analysis

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

Peak Number 11 unknown-01 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.706	2.82 ug/L	59055	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Hexyne	82	C6H10	000928-49-4	38
2			2,4-Hexadiene, (E,Z)-	82	C6H10	005194-50-3	38
3			Cyclopropane, (1-methylethylidene)-	82	C6H10	004741-86-0	38
4			1,4-Hexadiene	82	C6H10	000592-45-0	30
5			4-Methyl-1,3-pentadiene	82	C6H10	000926-56-7	30



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU042822\
Data File : VU048336.D
Acq On : 28 Apr 2022 13:17
Operator : SY/MD
Sample : N2587-08DL 40X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXET4DL

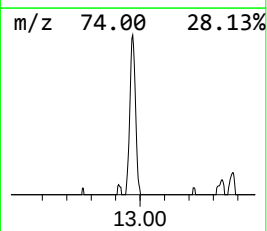
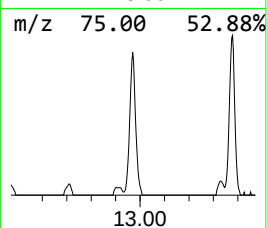
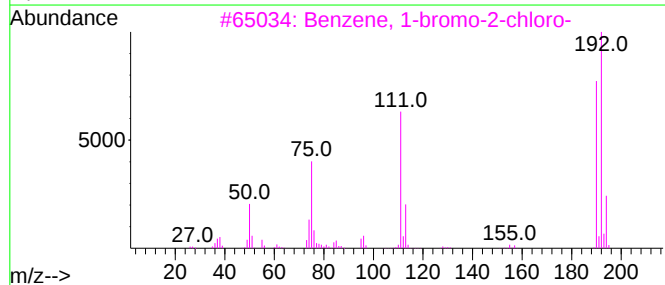
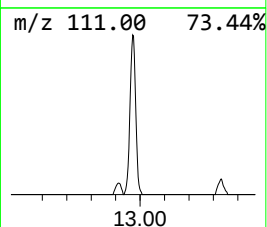
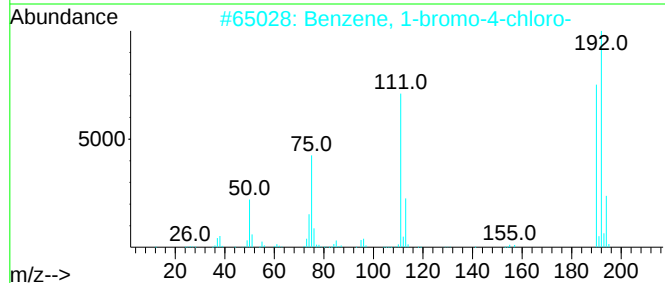
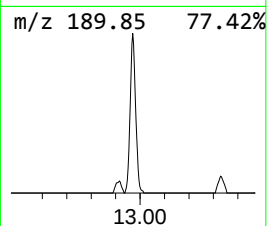
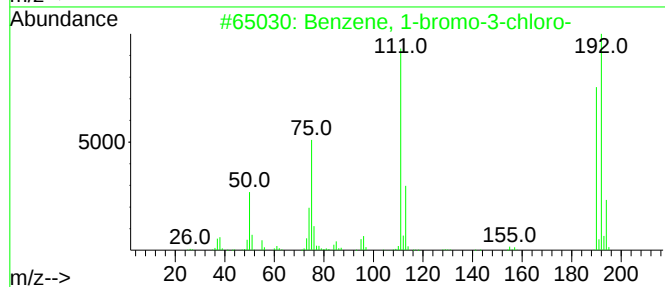
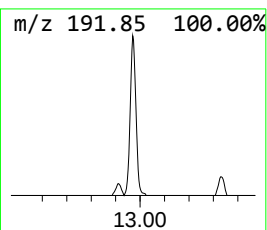
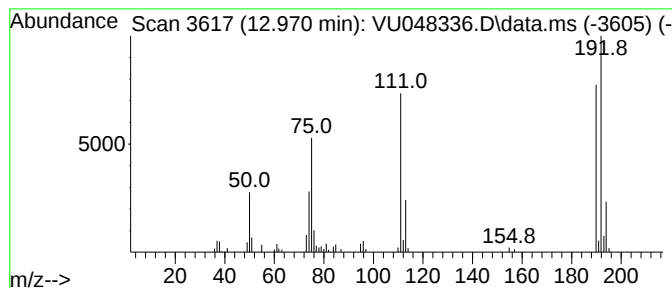
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Quant Title : VOC Analysis

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

Peak Number 12 Benzene, 1-bromo-3-chloro- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.970	3.92 ug/L	82103	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	98
2			Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	98
3			Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	98
4			Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	98
5			Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	97



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU042822\
Data File : VU048336.D
Acq On : 28 Apr 2022 13:17
Operator : SY/MD
Sample : N2587-08DL 40X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXET4DL

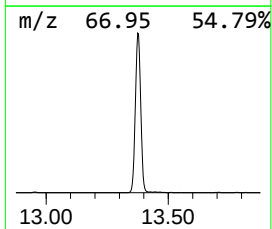
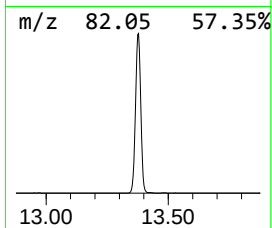
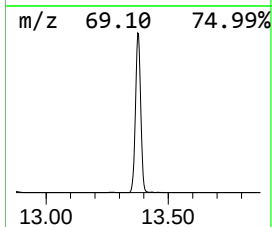
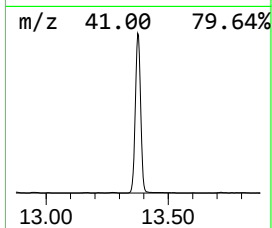
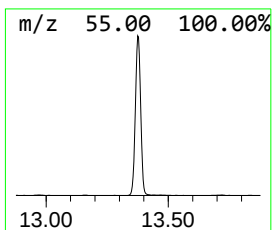
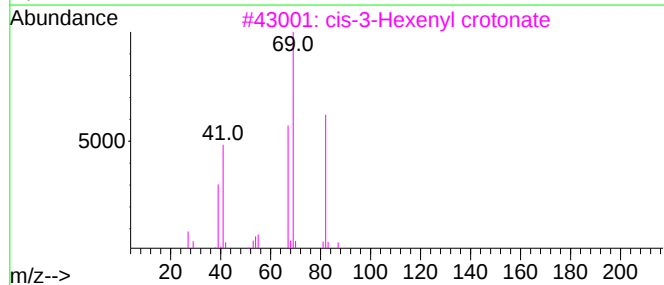
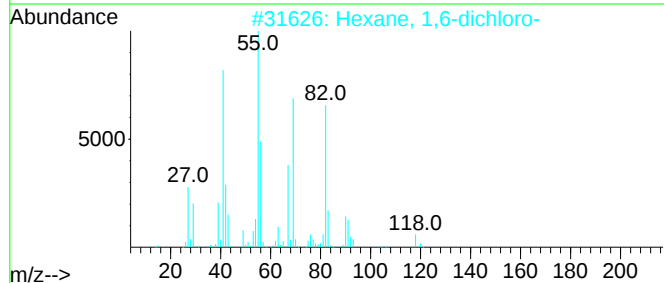
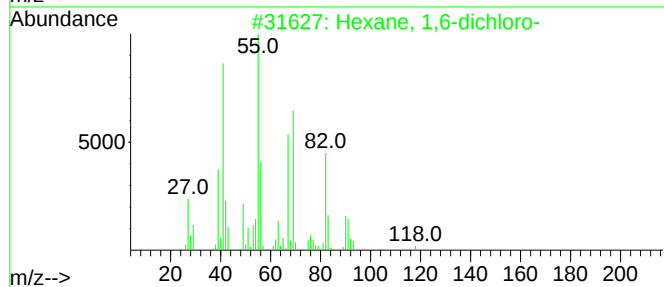
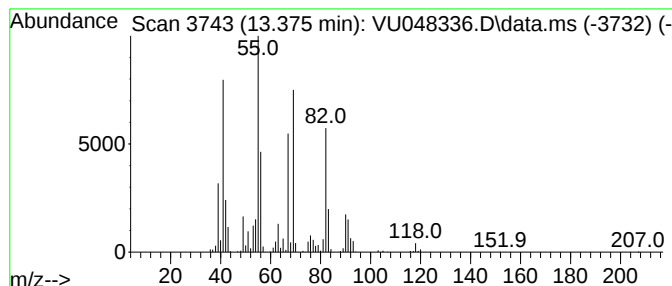
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Quant Title : VOC Analysis

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

Peak Number 13 Hexane, 1,6-dichloro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.375	53.63 ug/L	1122320	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	90
2			Hexane, 1,6-dichloro-	154	C6H12Cl2	002163-00-0	47
3			cis-3-Hexenyl crotonate	168	C10H16O2	1000429-17-4	47
4			1-Hexene, 6-chloro-	118	C6H11Cl	000928-89-2	43
5			Cyclohexane, chloro-	118	C6H11Cl	000542-18-7	42



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU042822\
Data File : VU048336.D
Acq On : 28 Apr 2022 13:17
Operator : SY/MD
Sample : N2587-08DL 40X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXET4DL

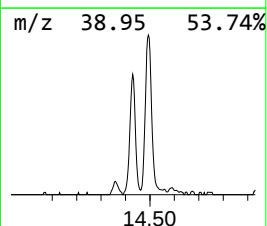
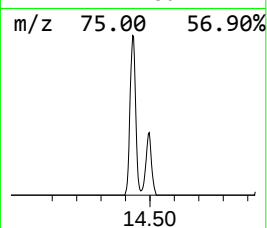
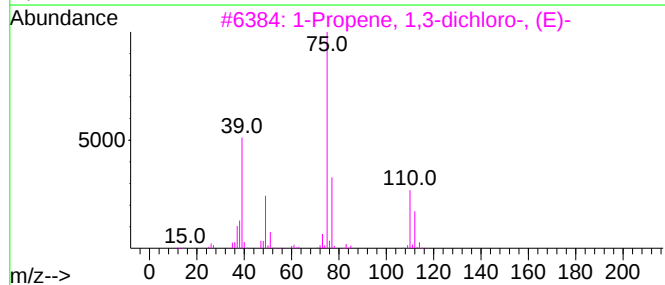
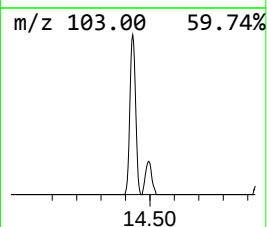
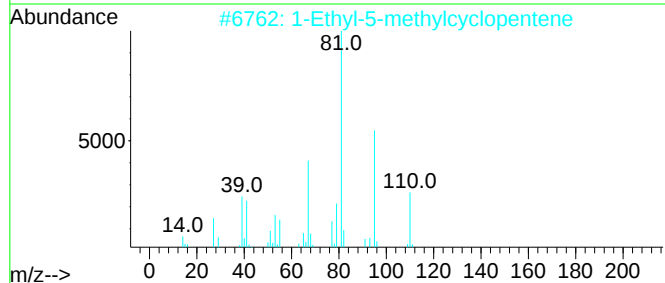
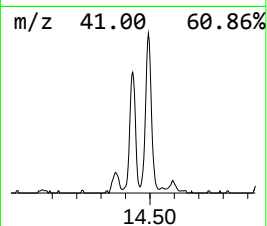
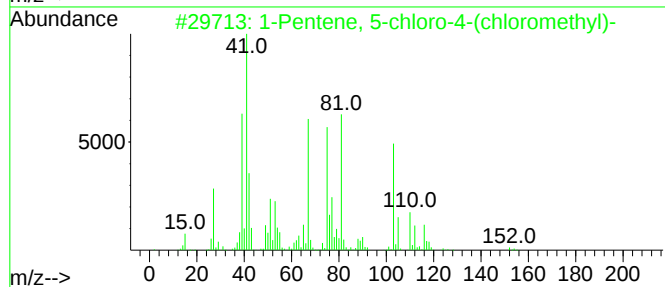
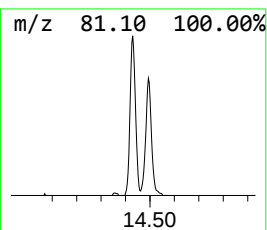
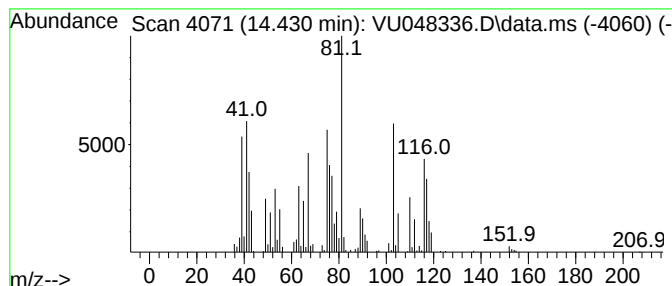
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Quant Title : VOC Analysis

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

Peak Number 14 unknown-02 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.430	7.97 ug/L	166711	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Pentene, 5-chloro-4-(chloromet...	152	C6H10Cl2	027990-74-5	42
2			1-Ethyl-5-methylcyclopentene	110	C8H14	097797-57-4	18
3			1-Propene, 1,3-dichloro-, (E)-	110	C3H4Cl2	010061-02-6	18
4			Cyclohexane, ethylidene-	110	C8H14	001003-64-1	18
5			Cyclohexene, 1-ethyl-	110	C8H14	001453-24-3	14



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU042822\
Data File : VU048336.D
Acq On : 28 Apr 2022 13:17
Operator : SY/MD
Sample : N2587-08DL 40X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXET4DL

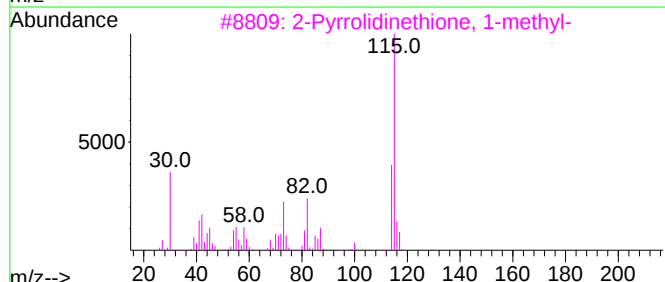
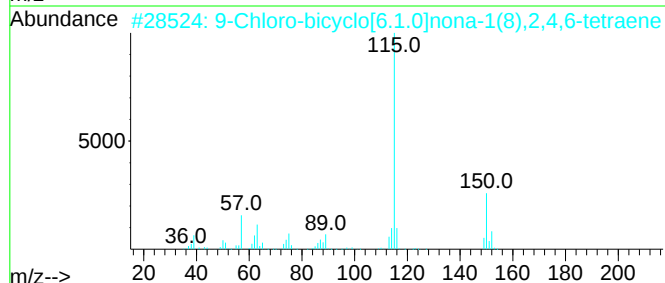
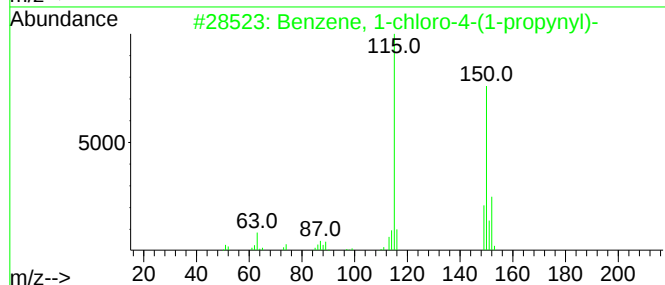
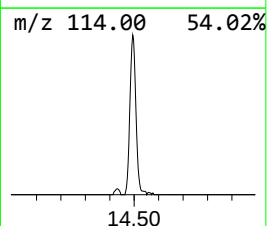
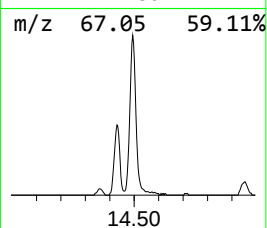
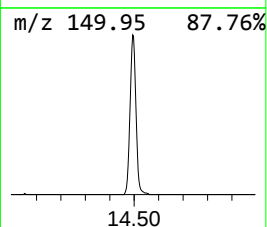
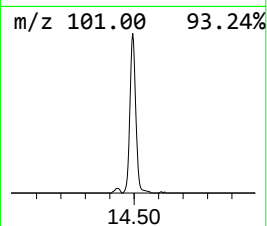
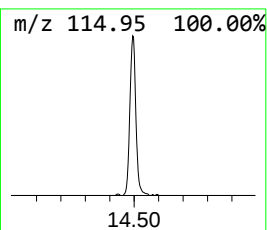
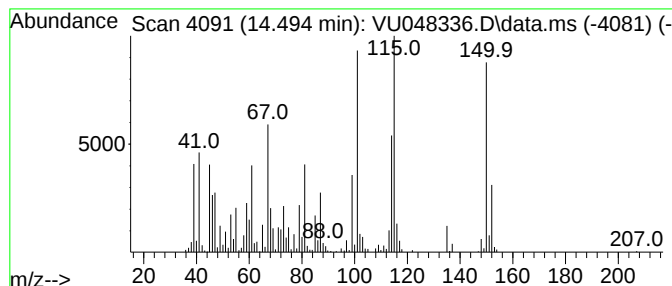
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Quant Title : VOC Analysis

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

Peak Number 15 unknown-03 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.494	17.14 ug/L	358678	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	45
2			9-Chloro-bicyclo[6.1.0]nona-1(8)...	150	C9H7Cl	1010295-96-4	35
3			2-Pyrrolidinethione, 1-methyl-	115	C5H9NS	010441-57-3	18
4			2-Benzothiazolamine	150	C7H6N2S	000136-95-8	11
5			2-Pyrrolidinethione	101	C4H7NS	002295-35-4	11



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU042822\
Data File : VU048336.D
Acq On : 28 Apr 2022 13:17
Operator : SY/MD
Sample : N2587-08DL 40X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXET4DL

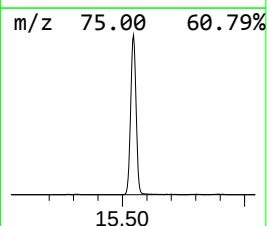
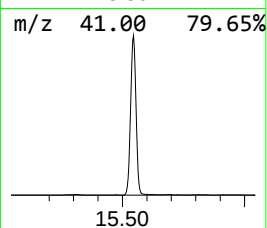
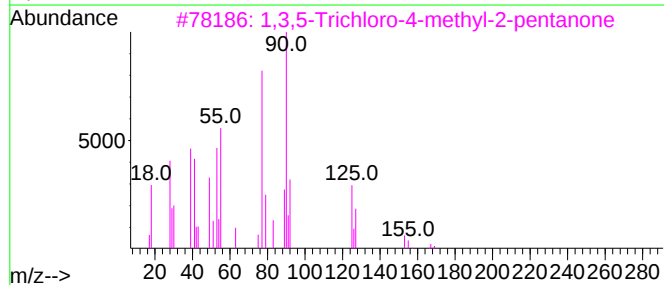
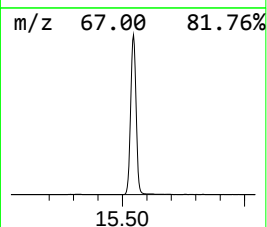
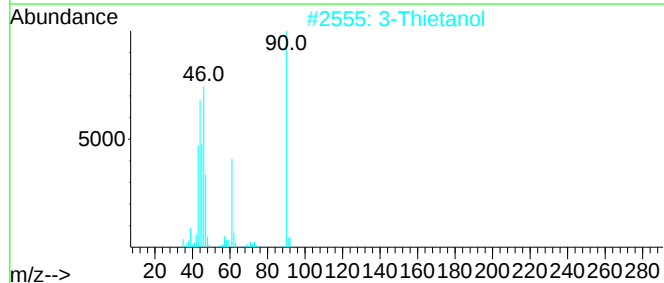
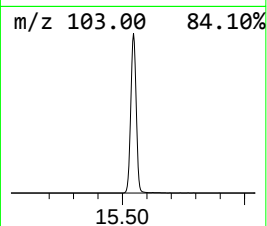
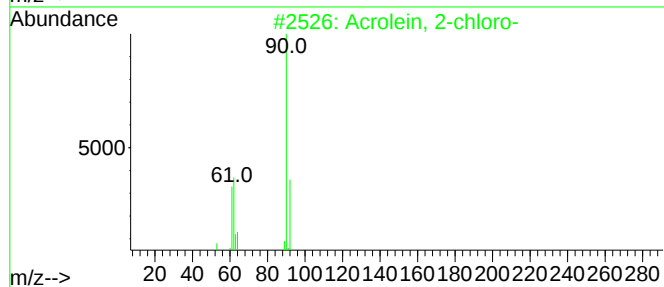
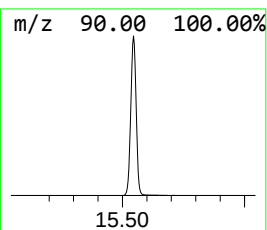
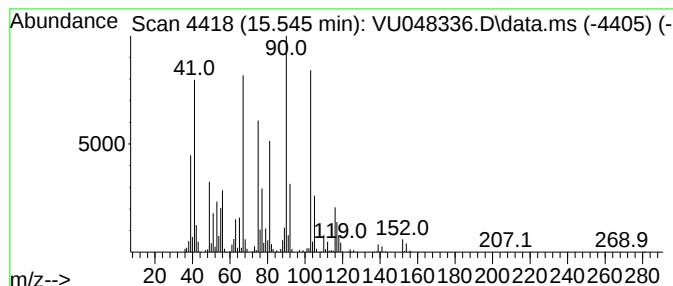
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Quant Title : VOC Analysis

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

Peak Number 16 unknown-04 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.545	80.52 ug/L	1684950	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acrolein, 2-chloro-	90	C3H3ClO	000683-51-2	43
2			3-Thietanol	90	C3H6OS	010304-16-2	25
3			1,3,5-Trichloro-4-methyl-2-penta...	202	C6H9Cl3O	1000140-03-8	25
4			Propane, 2-(methylthio)-	90	C4H10S	001551-21-9	25
5			2,2-Bis(chloromethyl)-1-propanol	156	C5H10Cl2O	005355-54-4	23



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU042822\
Data File : VU048336.D
Acq On : 28 Apr 2022 13:17
Operator : SY/MD
Sample : N2587-08DL 40X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
EXET4DL

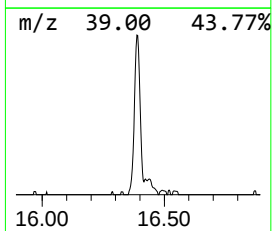
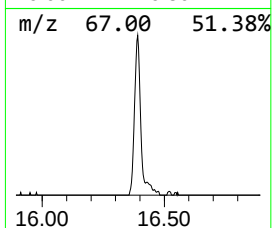
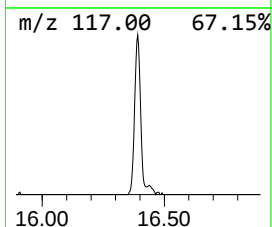
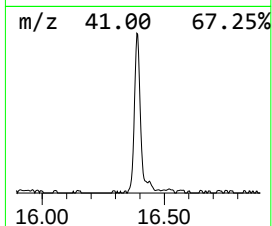
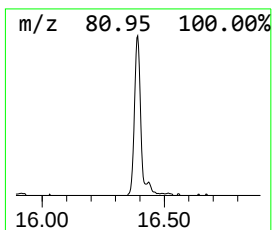
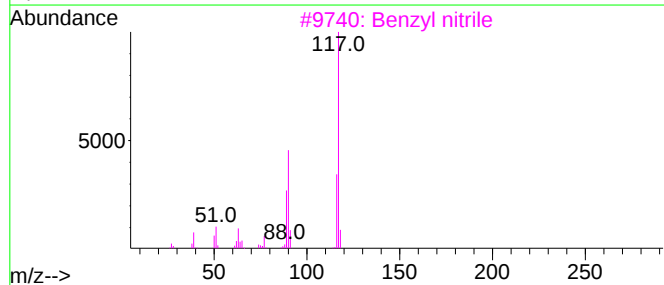
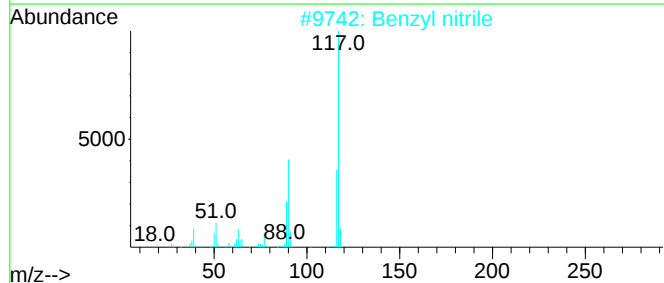
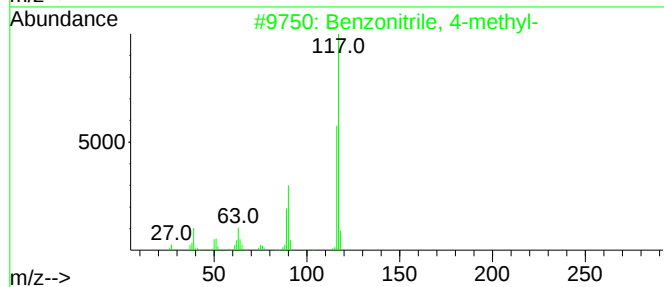
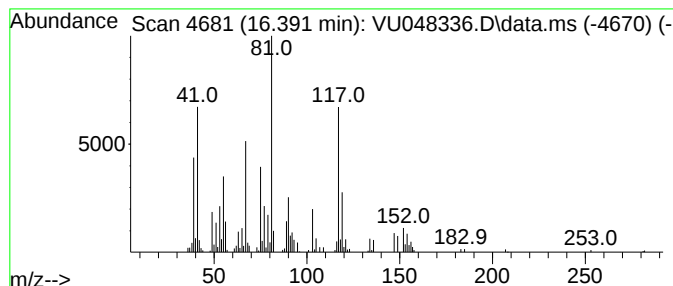
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Quant Title : VOC Analysis

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

Peak Number 17 unknown-05 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.391	8.69 ug/L	181788	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzonitrile, 4-methyl-	117	C8H7N	000104-85-8	25
2			Benzyl nitrile	117	C8H7N	000140-29-4	25
3			Benzyl nitrile	117	C8H7N	000140-29-4	25
4			Benzonitrile, 2-methyl-	117	C8H7N	000529-19-1	15
5			Benzyl nitrile	117	C8H7N	000140-29-4	15



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU042822\
 Data File : VU048336.D
 Acq On : 28 Apr 2022 13:17
 Operator : SY/MD
 Sample : N2587-08DL 40X
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 EXET4DL

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	48.7	ug/L	743682	1	6.250	763231	50.0
(DEL) Alkane: C...	1.475	3.8	ug/L	58283	1	6.250	763231	50.0
Furan, tetrahyd...	6.513	4.3	ug/L	65574	1	6.250	763231	50.0
Furan, tetrahyd...	6.774	4.5	ug/L	69124	1	6.250	763231	50.0
Thietane	6.996	6.0	ug/L	92187	1	6.250	763231	50.0
2H-Pyran, tetra...	7.488	39.3	ug/L	599278	1	6.250	763231	50.0
Propanal, 2,3-d...	8.574	54.8	ug/L	3797020	2	9.417	3464910	50.0
Thiepane	11.880	3.7	ug/L	76815	3	11.812	1046270	50.0
unknown-01	12.706	2.8	ug/L	59055	3	11.812	1046270	50.0
Benzene, 1-brom...	12.970	3.9	ug/L	82103	3	11.812	1046270	50.0
Hexane, 1,6-dic...	13.375	53.6	ug/L	1122320	3	11.812	1046270	50.0
unknown-02	14.430	8.0	ug/L	166711	3	11.812	1046270	50.0
unknown-03	14.494	17.1	ug/L	358678	3	11.812	1046270	50.0
unknown-04	15.545	80.5	ug/L	1684950	3	11.812	1046270	50.0
unknown-05	16.391	8.7	ug/L	181788	3	11.812	1046270	50.0