

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU070122\
 Data File : VU049594.D
 Acq On : 01 Jul 2022 20:08
 Operator : SY/MD
 Sample : N3564-09 10X
 Misc : 5.98g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 WC-14B-11-2-GR

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % 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\82U062722W.M
 Title : SW846 8260

Signal : TIC: VU049594.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.826	127	151	156	rBV5	11757	29484	2.57%	0.199%
2	1.977	187	198	209	rVB	13548	18954	1.65%	0.128%
3	2.189	257	264	275	rVB2	9429	13294	1.16%	0.090%
4	3.035	513	527	531	rBV5	7350	13879	1.21%	0.094%
5	3.073	532	539	552	rVB3	16070	30885	2.69%	0.208%
6	3.179	559	572	590	rVB	17680	40487	3.53%	0.273%
7	3.353	608	626	641	rBV2	33211	73447	6.41%	0.495%
8	4.633	1009	1024	1043	rBV2	8483	25264	2.20%	0.170%
9	5.288	1210	1228	1241	rBV2	170678	360162	31.41%	2.427%
10	5.372	1241	1254	1270	rVV	306126	653441	56.99%	4.403%
11	5.459	1270	1281	1296	rVB4	20310	45137	3.94%	0.304%
12	5.588	1304	1321	1331	rBV3	30259	66997	5.84%	0.451%
13	5.700	1344	1356	1372	rVB	186983	367803	32.08%	2.478%
14	5.848	1389	1402	1413	rBV2	41301	85175	7.43%	0.574%
15	5.919	1413	1424	1434	rVV3	31722	65346	5.70%	0.440%
16	5.986	1434	1445	1462	rVB2	56573	120261	10.49%	0.810%
17	6.247	1511	1526	1548	rBV	426175	832120	72.58%	5.607%
18	6.565	1615	1625	1635	rBV9	7636	15134	1.32%	0.102%
19	6.761	1665	1686	1709	rBV2	199142	466848	40.72%	3.146%
20	6.864	1709	1718	1730	rVB3	12150	23041	2.01%	0.155%
21	6.980	1743	1754	1768	rVV4	22913	42733	3.73%	0.288%
22	7.073	1768	1783	1797	rVB3	31855	64412	5.62%	0.434%
23	7.234	1821	1833	1850	rVB4	38416	80858	7.05%	0.545%
24	7.398	1870	1884	1889	rBV3	15289	30766	2.68%	0.207%
25	7.536	1920	1927	1935	rVB3	13705	20242	1.77%	0.136%
26	7.587	1935	1943	1950	rBV5	9490	16105	1.40%	0.109%
27	7.690	1971	1975	1984	rVB4	11418	16110	1.41%	0.109%
28	7.758	1984	1996	2013	rVB4	41792	80436	7.02%	0.542%
29	7.854	2013	2026	2030	rBV	107955	178058	15.53%	1.200%
30	7.896	2030	2039	2054	rVB	663025	1146489	100.00%	7.726%
31	8.028	2068	2080	2088	rBV5	27578	68084	5.94%	0.459%
32	8.095	2096	2101	2113	rVB3	32859	51740	4.51%	0.349%
33	8.157	2113	2120	2134	rVV3	10562	21143	1.84%	0.142%
34	8.240	2134	2146	2161	rVB3	77792	158656	13.84%	1.069%
35	8.369	2175	2186	2193	rBV2	36623	66937	5.84%	0.451%
36	8.411	2193	2199	2212	rVV6	20851	41755	3.64%	0.281%

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37	8.475	2213	2219	2229	rVB7	10100	17710	1.54%	0.119%
38	8.539	2229	2239	2252	rVB5	21274	43448	3.79%	0.293%
39	8.636	2252	2269	2281	rBV2	59439	115772	10.10%	0.780%
40	8.703	2281	2290	2296	rBV5	14141	24180	2.11%	0.163%
41	8.751	2296	2305	2312	rBV6	31039	57852	5.05%	0.390%
42	8.803	2313	2321	2331	rVB6	35842	64710	5.64%	0.436%
43	8.864	2332	2340	2350	rBV	88151	136097	11.87%	0.917%
44	8.925	2350	2359	2369	rBV	106470	172435	15.04%	1.162%
45	9.070	2391	2404	2413	rBV4	38803	78096	6.81%	0.526%
46	9.121	2413	2420	2426	rVV2	31708	42433	3.70%	0.286%
47	9.166	2426	2434	2442	rBV3	55491	90445	7.89%	0.609%
48	9.330	2471	2485	2497	rBV2	49405	92898	8.10%	0.626%
49	9.417	2500	2512	2523	rBV	566577	936126	81.65%	6.308%
50	9.517	2539	2543	2550	rVB6	10537	12892	1.12%	0.087%
51	9.562	2551	2557	2571	rVB5	33310	61953	5.40%	0.417%
52	9.665	2572	2589	2596	rBV8	50623	139129	12.14%	0.938%
53	9.716	2598	2605	2612	rVB3	76126	121131	10.57%	0.816%
54	9.761	2612	2619	2644	rVB3	53412	115746	10.10%	0.780%
55	9.880	2644	2656	2671	rBV10	33014	87500	7.63%	0.590%
56	9.954	2672	2679	2688	rVB4	11852	17956	1.57%	0.121%
57	10.025	2688	2701	2713	rBV7	93652	240472	20.97%	1.620%
58	10.121	2723	2731	2743	rBV7	37633	84517	7.37%	0.570%
59	10.253	2757	2772	2786	rBV4	159971	410668	35.82%	2.767%
60	10.356	2796	2804	2811	rBV3	119820	213636	18.63%	1.440%
61	10.478	2833	2842	2852	rVB6	47357	97535	8.51%	0.657%
62	10.555	2852	2866	2876	rBV8	69383	170746	14.89%	1.151%
63	10.632	2877	2890	2903	rVV2	486500	825425	72.00%	5.562%
64	10.697	2904	2910	2924	rVV7	37649	75189	6.56%	0.507%
65	10.774	2925	2934	2935	rVV5	30786	46449	4.05%	0.313%
66	10.809	2938	2945	2962	rVB6	123824	233891	20.40%	1.576%
67	10.906	2970	2975	2980	rBV	12880	16932	1.48%	0.114%
68	10.951	2981	2989	3001	rVV5	48186	100703	8.78%	0.679%
69	11.015	3002	3009	3016	rVV6	57036	113366	9.89%	0.764%
70	11.063	3017	3024	3034	rVV4	112854	207216	18.07%	1.396%
71	11.127	3038	3044	3054	rVB6	70018	126574	11.04%	0.853%
72	11.185	3055	3062	3072	rBV8	31224	56347	4.91%	0.380%
73	11.314	3087	3102	3114	rVB8	71104	169979	14.83%	1.145%
74	11.378	3115	3122	3126	rBV6	32402	48694	4.25%	0.328%
75	11.417	3127	3134	3141	rVB	112433	146824	12.81%	0.989%
76	11.462	3142	3148	3153	rBV7	31964	48462	4.23%	0.327%

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77	11.574	3175	3183	3187	rBV6	33643	56788	4.95%	0.383%
78	11.642	3198	3204	3215	rVB5	117675	199311	17.38%	1.343%
79	11.713	3218	3226	3232	rBV	92999	144550	12.61%	0.974%
80	11.812	3247	3257	3272	rVB	587348	973201	84.89%	6.558%
81	12.099	3338	3346	3357	rBV3	72249	127129	11.09%	0.857%
82	12.182	3365	3372	3378	rBV5	107335	177403	15.47%	1.195%
83	12.571	3489	3493	3500	rVB2	94804	110877	9.67%	0.747%
84	12.838	3571	3576	3591	rVB6	127078	235129	20.51%	1.584%
85	12.931	3597	3605	3612	rBV4	100063	198793	17.34%	1.340%
86	13.053	3636	3643	3651	rVB4	106324	161491	14.09%	1.088%
87	13.201	3683	3689	3696	rBV2	46312	67532	5.89%	0.455%
88	13.452	3759	3767	3785	rVB5	198356	408182	35.60%	2.751%
89	13.838	3881	3887	3897	rBV5	69309	139374	12.16%	0.939%
90	14.121	3968	3975	3985	rBV7	93815	156153	13.62%	1.052%
91	15.275	4329	4334	4342	rVB5	60913	79180	6.91%	0.534%
92	16.262	4636	4641	4650	rVB3	65537	88717	7.74%	0.598%
93	16.407	4680	4686	4693	rBV	93719	132169	11.53%	0.891%
94	17.539	5033	5038	5050	rVB2	56528	90044	7.85%	0.607%

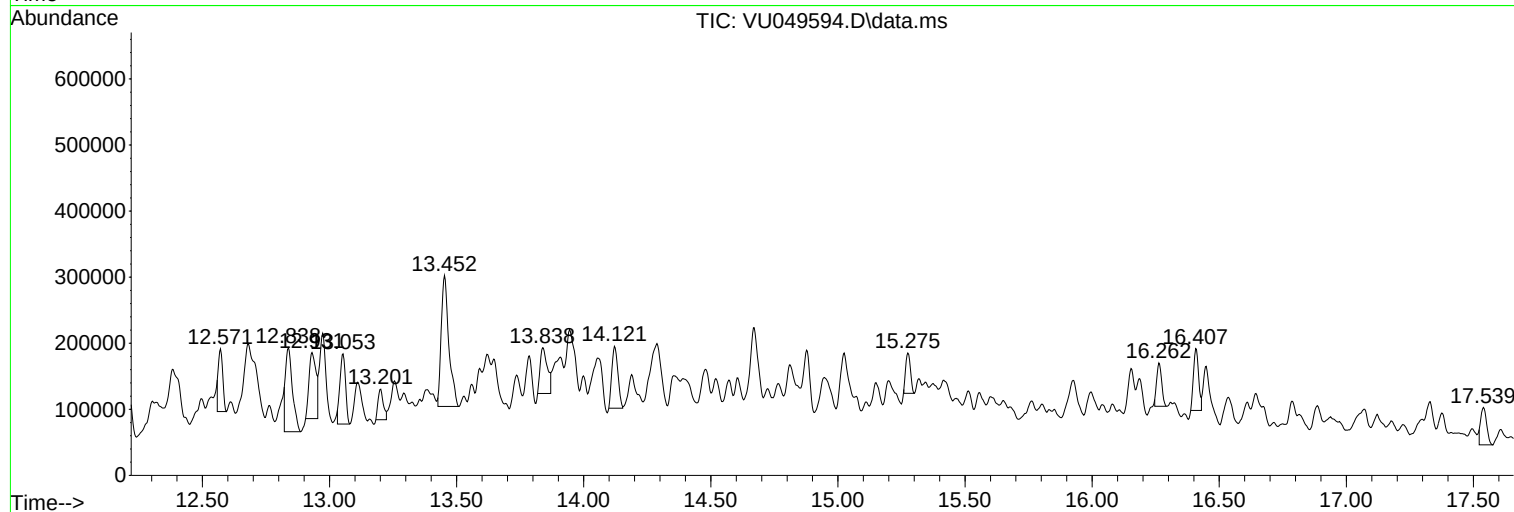
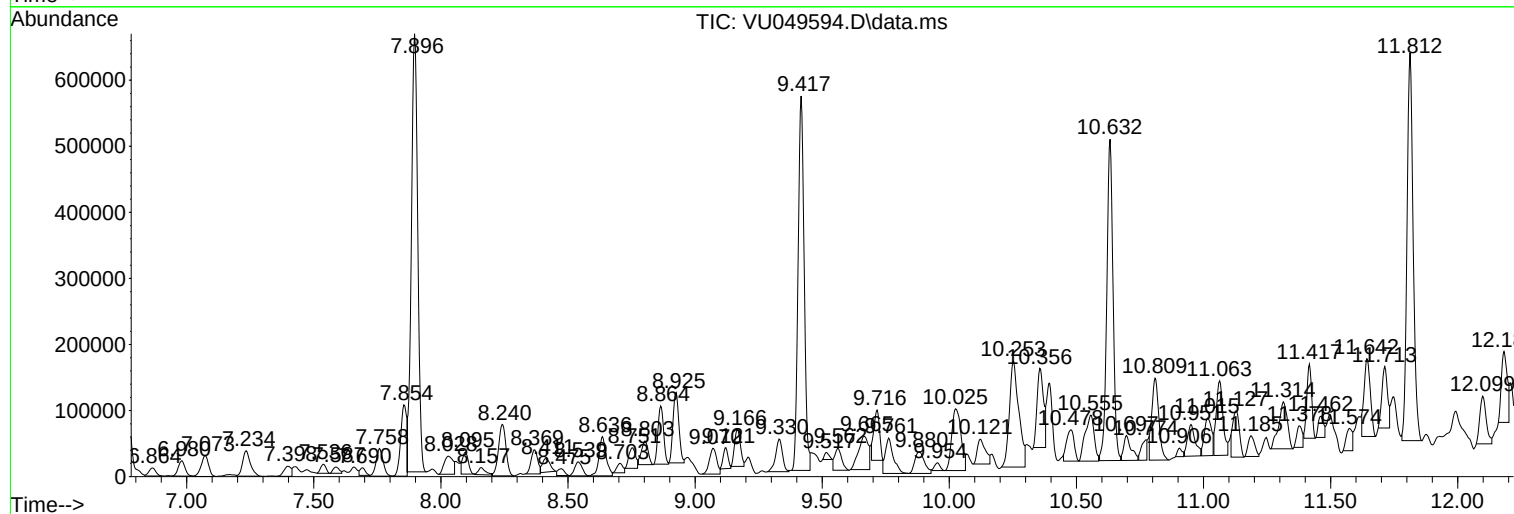
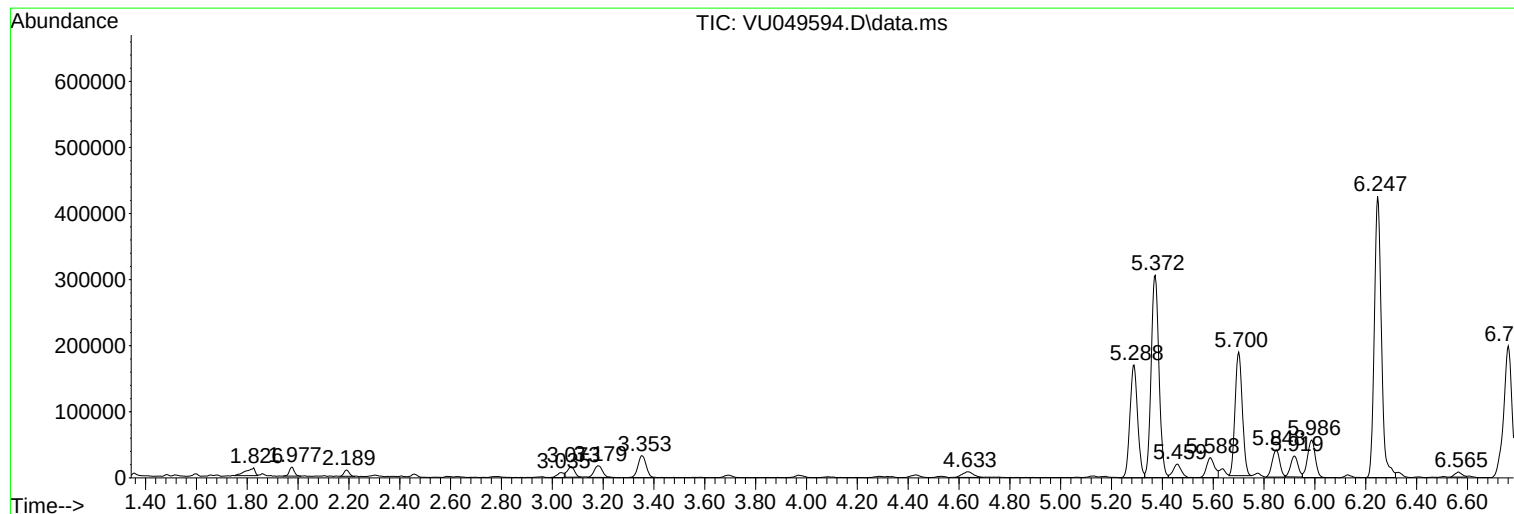
Sum of corrected areas: 14839840

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

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



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Instrument :
MSVOA_U
ClientSampleId :
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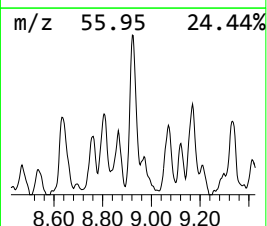
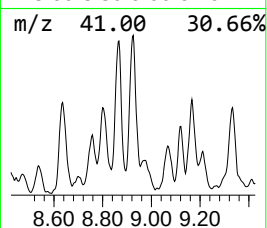
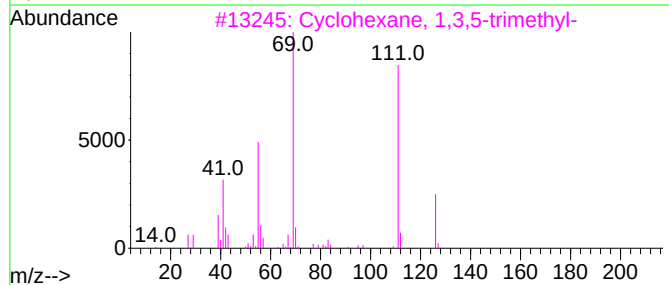
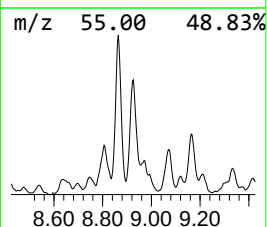
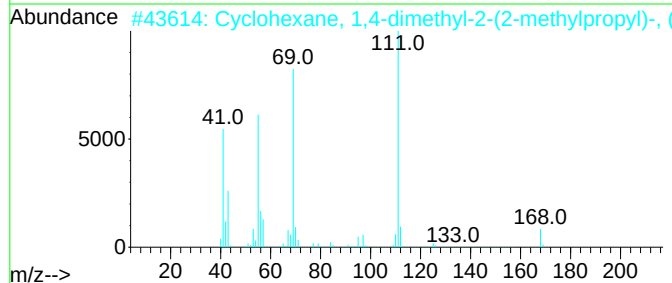
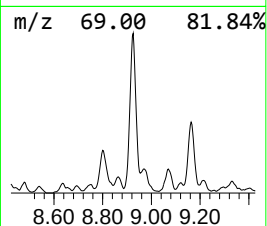
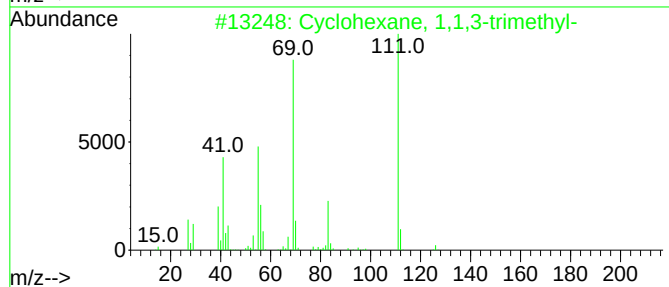
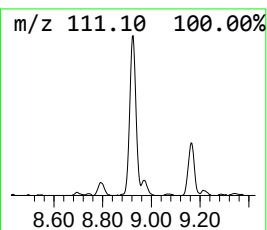
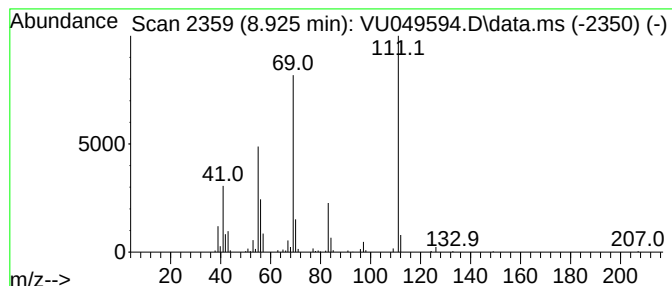
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

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

Peak Number 1 Cyclohexane, 1,1,3-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.925	9.21 ug/l	172435	Chlorobenzene-d5	9.417

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	94
2			Cyclohexane, 1,4-dimethyl-2-(2-m...	168	C12H24	054411-17-5	72
3			Cyclohexane, 1,3,5-trimethyl-	126	C9H18	001839-63-0	72
4			Cyclohexane, 1-ethyl-1,3-dimethy...	140	C10H20	062238-29-3	72
5			Cyclohexane, 1-ethyl-2,4-dimethyl-	140	C10H20	061142-69-6	72



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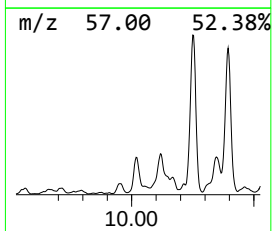
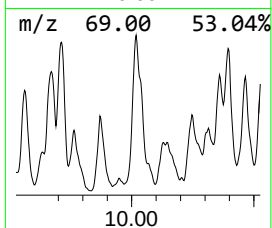
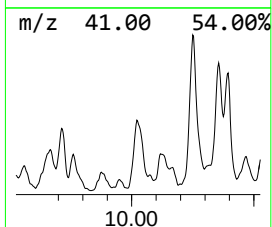
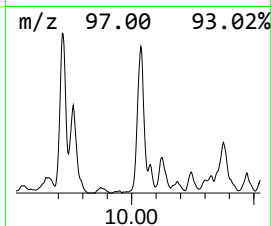
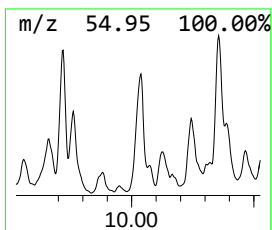
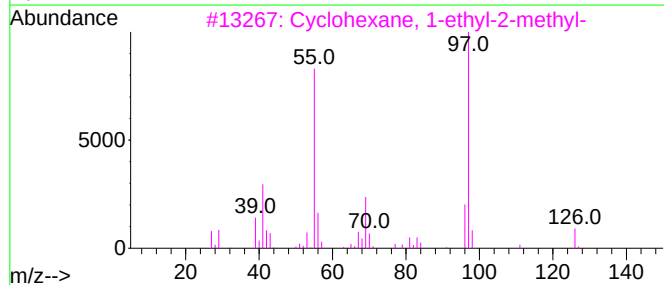
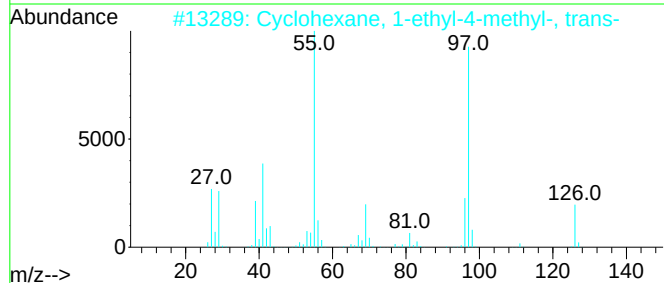
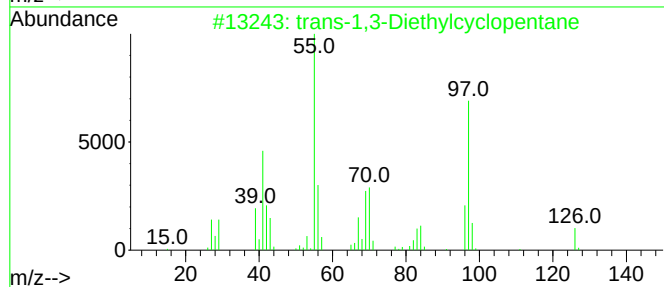
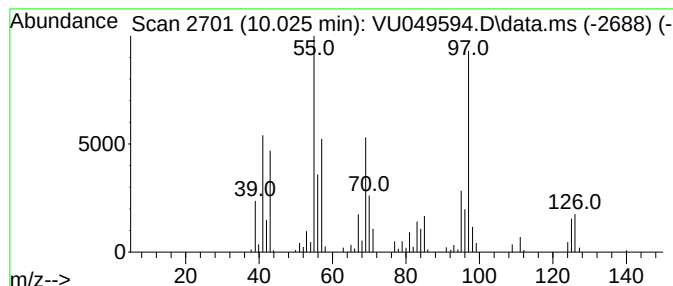
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Quant Title : SW846 8260

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

Peak Number 2 trans-1,3-Diethylcyclopentane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.025	12.84 ug/l	240472	Chlorobenzene-d5	9.417

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			trans-1,3-Diethylcyclopentane	126	C9H18	1000113-87-1	81
2			Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	006236-88-0	53
3			Cyclohexane, 1-ethyl-2-methyl-	126	C9H18	003728-54-9	52
4			1-Ethyl-3-methylcyclohexane (c,t)	126	C9H18	003728-55-0	49
5			1-Ethyl-4-methylcyclohexane	126	C9H18	003728-56-1	49



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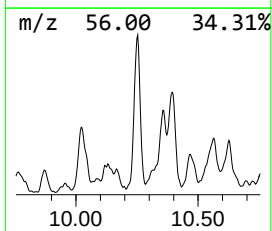
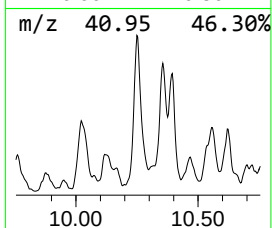
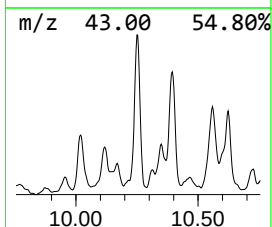
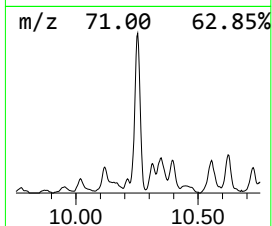
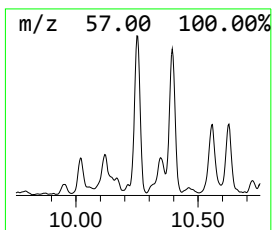
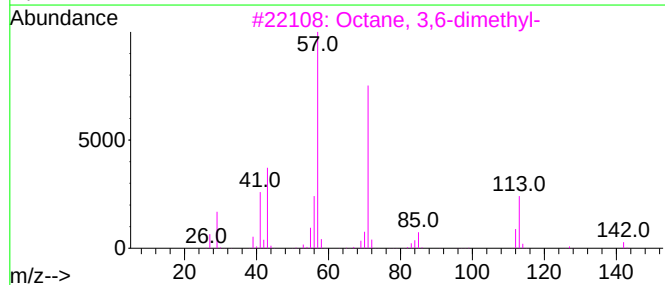
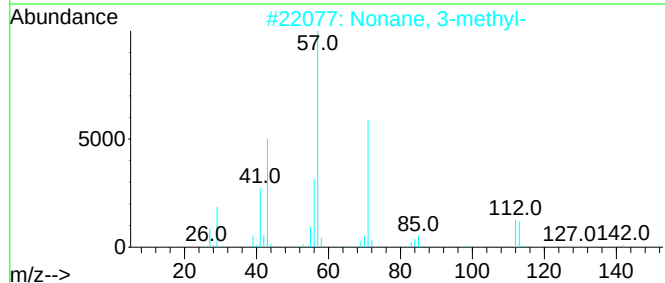
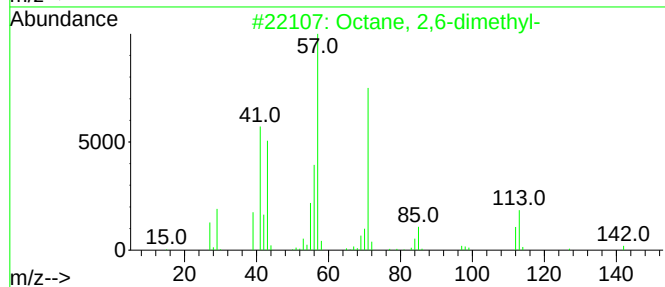
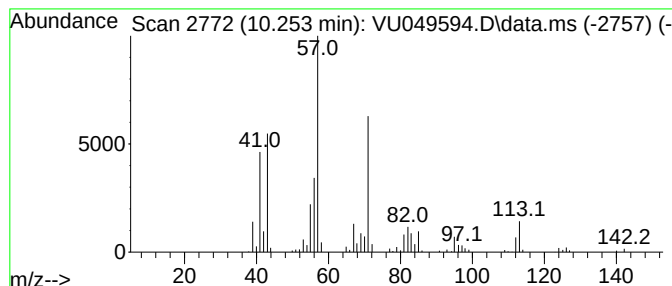
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WC-14B-11-2-GR

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

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

Peak Number 3 Octane, 2,6-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.253	21.93 ug/l	410668	Chlorobenzene-d5	9.417	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	94
2	Nonane, 3-methyl-	142	C10H22	005911-04-6	87
3	Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	87
4	Octane, 3,5-dimethyl-	142	C10H22	015869-93-9	64
5	Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU070122\
Data File : VU049594.D
Acq On : 01 Jul 2022 20:08
Operator : SY/MD
Sample : N3564-09 10X
Misc : 5.98g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
WC-14B-11-2-GR

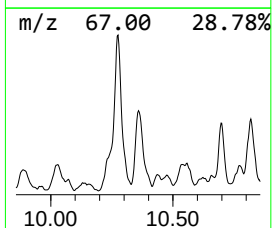
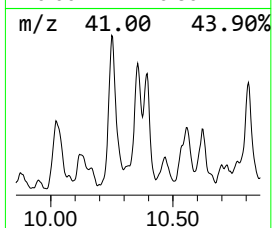
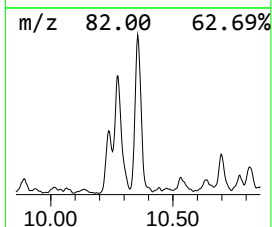
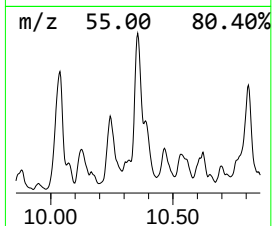
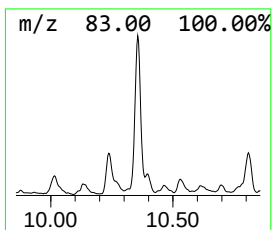
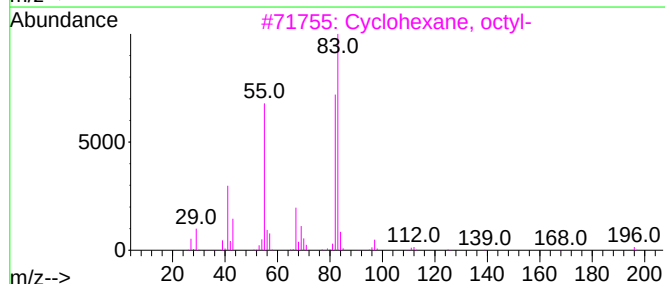
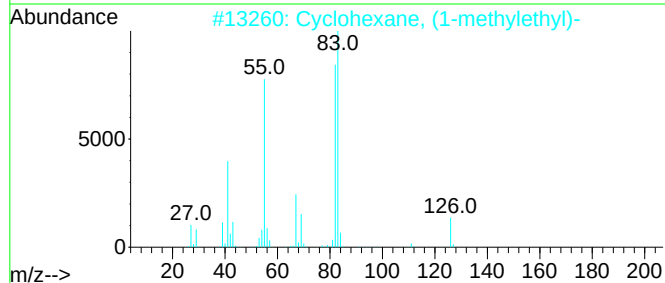
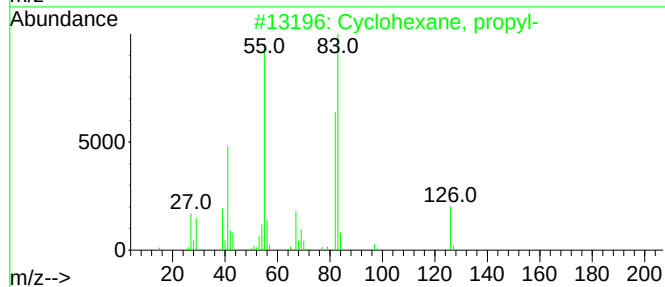
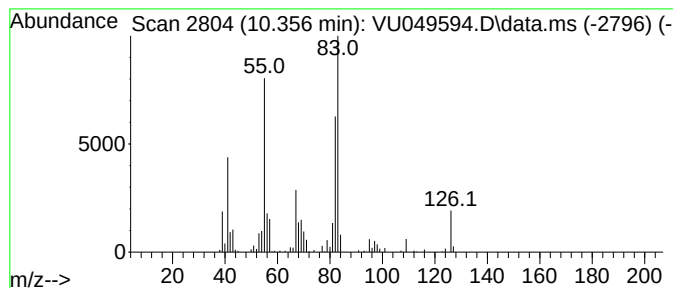
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Quant Title : SW846 8260

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

Peak Number 4 Cyclohexane, propyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.356	11.41 ug/l	213636	Chlorobenzene-d5	9.417

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, propyl-	126	C9H18	001678-92-8	87
2			Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	70
3			Cyclohexane, octyl-	196	C14H28	001795-15-9	64
4			Ethanone, 1-(1-methylcyclopentyl)-	126	C8H14O	013388-93-7	59
5			Methoxyacetic acid, cyclohexyl e...	172	C9H16O3	1000282-69-4	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU070122\
Data File : VU049594.D
Acq On : 01 Jul 2022 20:08
Operator : SY/MD
Sample : N3564-09 10X
Misc : 5.98g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
WC-14B-11-2-GR

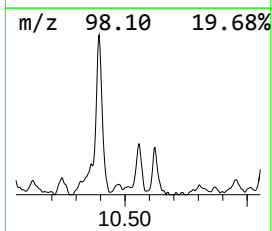
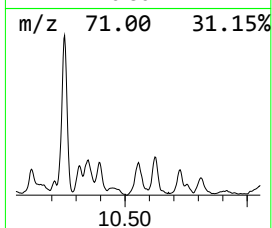
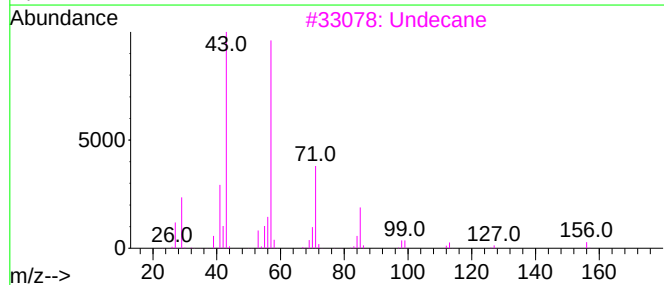
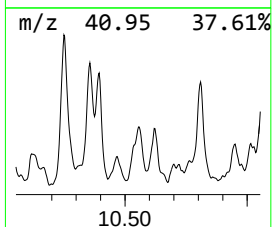
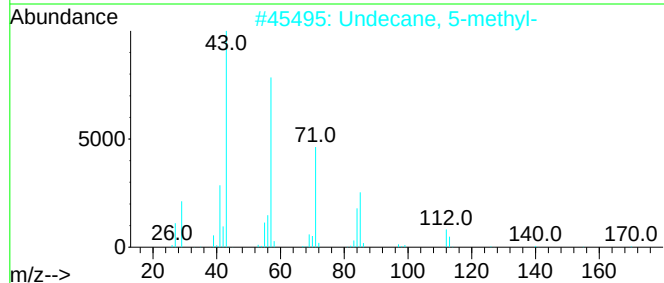
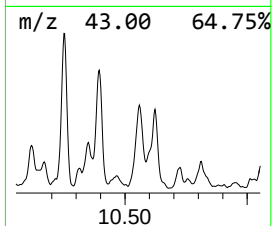
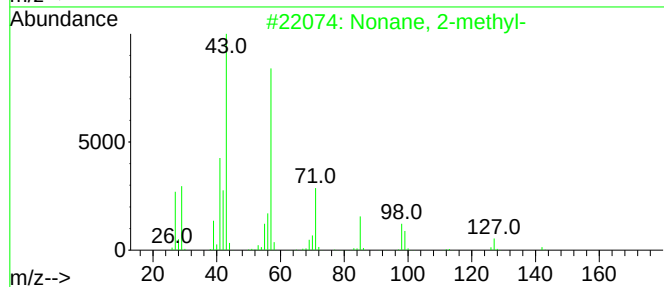
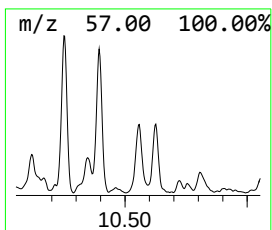
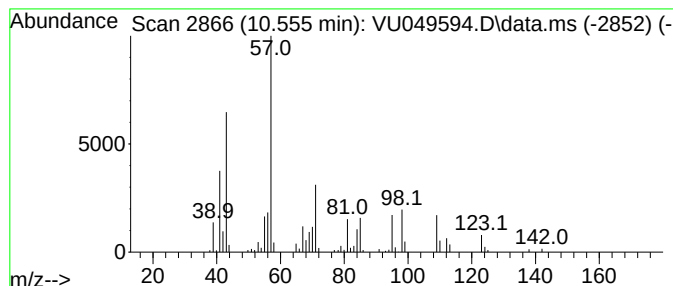
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Quant Title : SW846 8260

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

Peak Number 5 unknown10.555 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.555	9.12 ug/l	170746	Chlorobenzene-d5	9.417

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 2-methyl-	142	C10H22	000871-83-0	43
2			Undecane, 5-methyl-	170	C12H26	001632-70-8	38
3			Undecane	156	C11H24	001120-21-4	38
4			2,6-Dimethyldecane	170	C12H26	013150-81-7	38
5			Decane, 2,5,6-trimethyl-	184	C13H28	062108-23-0	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU070122\
Data File : VU049594.D
Acq On : 01 Jul 2022 20:08
Operator : SY/MD
Sample : N3564-09 10X
Misc : 5.98g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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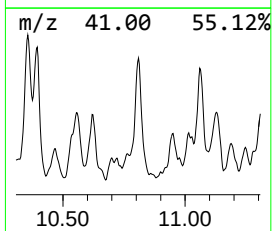
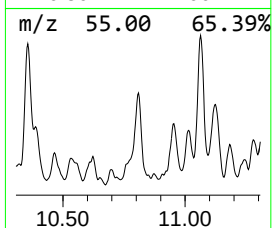
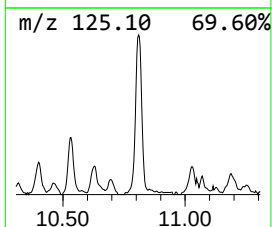
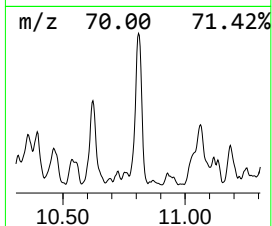
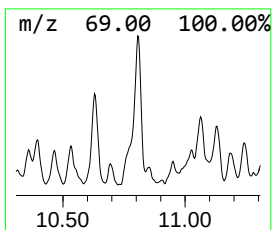
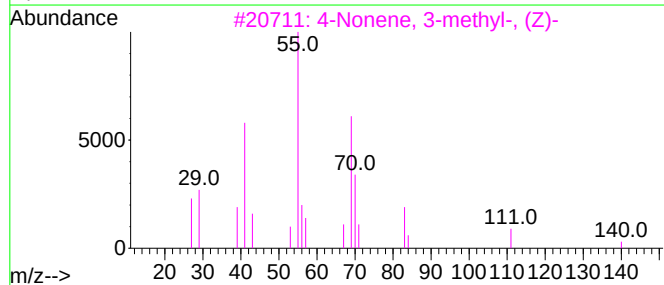
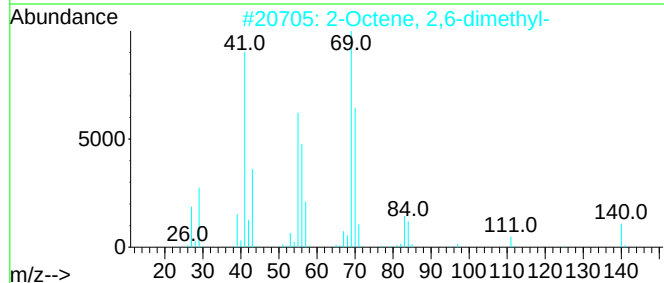
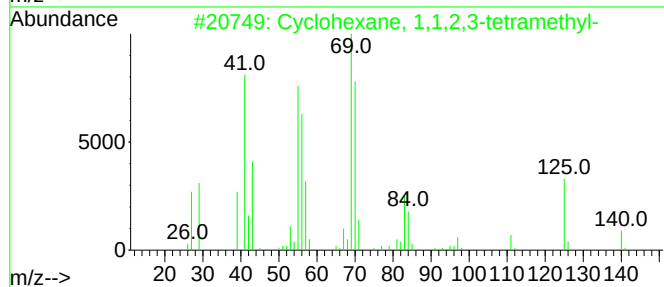
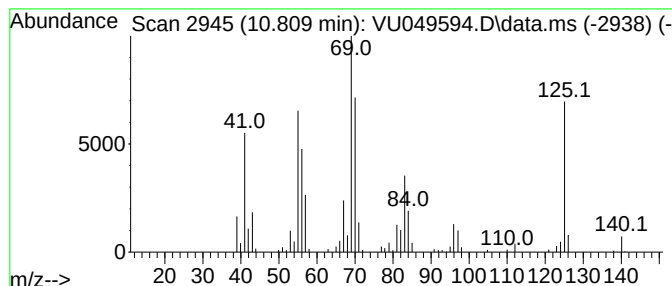
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Quant Title : SW846 8260

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

Peak Number 6 Cyclohexane, 1,1,2,3-tetram... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.809	12.02 ug/l	233891	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,1,2,3-tetramethyl-	140	C10H20	006783-92-2	68
2			2-Octene, 2,6-dimethyl-	140	C10H20	004057-42-5	62
3			4-Nonene, 3-methyl-, (Z)-	140	C10H20	063830-69-3	59
4			4-Octene, 2,6-dimethyl-, [S-(E)]-	140	C10H20	062960-76-3	58
5			1-Hexene, 3-methyl-	98	C7H14	003404-61-3	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU070122\
Data File : VU049594.D
Acq On : 01 Jul 2022 20:08
Operator : SY/MD
Sample : N3564-09 10X
Misc : 5.98g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 18 Sample Multiplier: 1

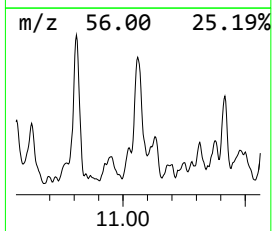
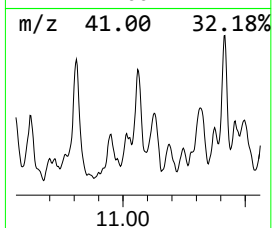
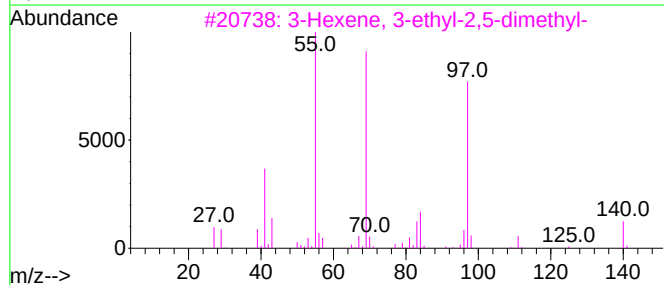
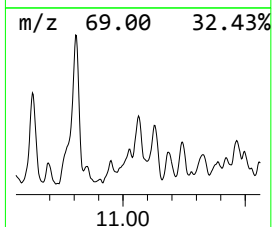
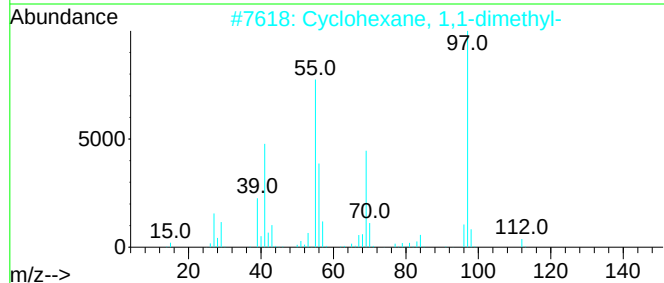
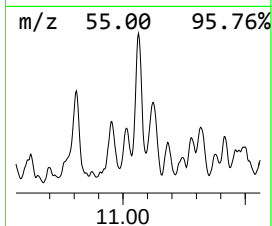
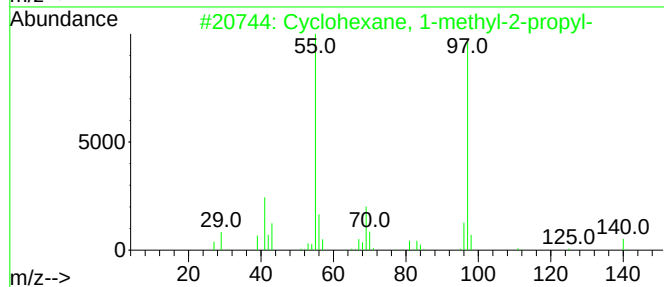
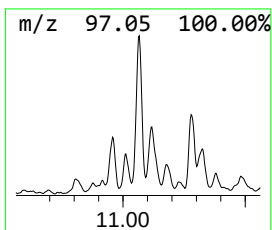
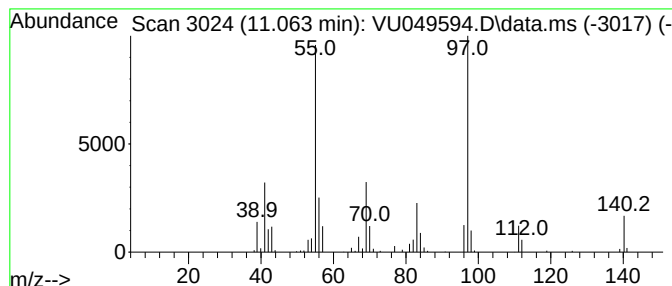
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ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

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

Peak Number 7 Cyclohexane, 1-methyl-2-pro... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD			R.T.
11.063	10.65 ug/l	207216	1,4-Dichlorobenzene-d4			11.812
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Cyclohexane, 1-methyl-2-propyl-		140	C10H20	004291-79-6	64
2	Cyclohexane, 1,1-dimethyl-		112	C8H16	000590-66-9	64
3	3-Hexene, 3-ethyl-2,5-dimethyl-		140	C10H20	062338-08-3	58
4	Cyclohexane, 1-methyl-4-(1-methy...		140	C10H20	001678-82-6	52
5	Cyclohexane, 1-methyl-4-(1-methy...		140	C10H20	006069-98-3	52



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU070122\
Data File : VU049594.D
Acq On : 01 Jul 2022 20:08
Operator : SY/MD
Sample : N3564-09 10X
Misc : 5.98g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
WC-14B-11-2-GR

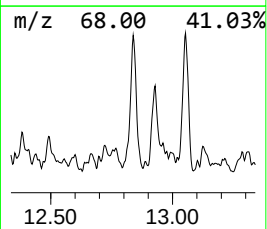
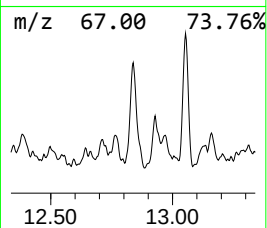
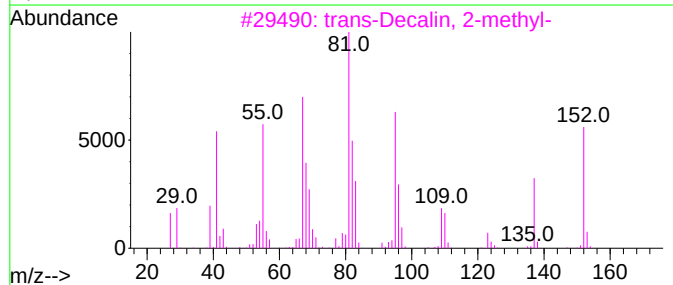
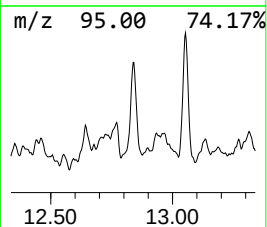
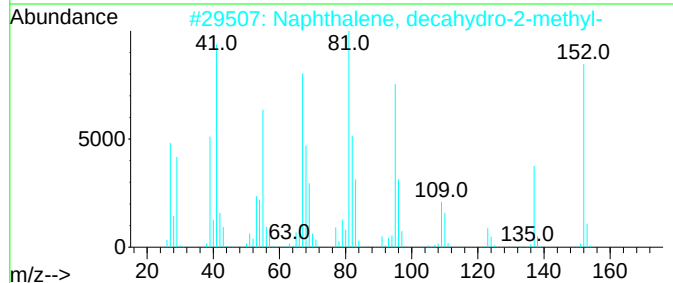
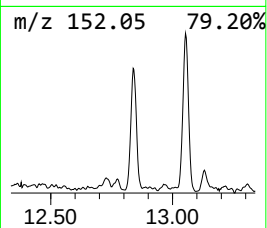
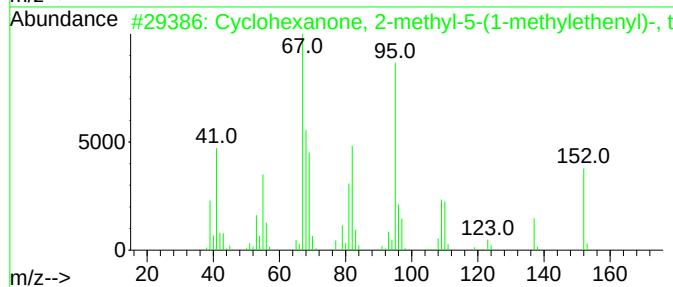
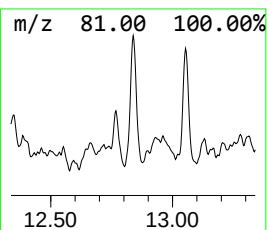
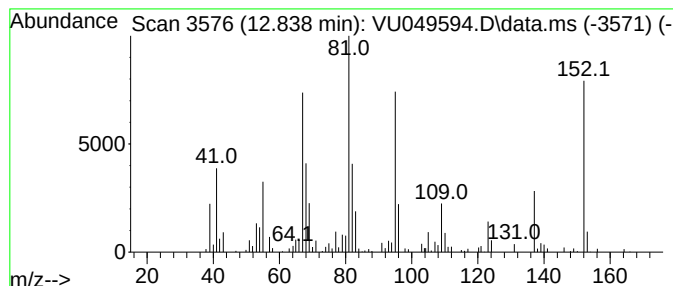
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Quant Title : SW846 8260

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

Peak Number 8 Cyclohexanone, 2-methyl-5-(... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.838	12.08 ug/l	235129	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexanone, 2-methyl-5-(1-met...	152	C10H16O	005948-04-9	90
2			Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	86
3			trans-Decalin, 2-methyl-	152	C11H20	1000152-47-3	74
4			(+)-Dihydrocarvone	152	C10H16O	005524-05-0	68
5			Cyclohexanone, 2-methyl-5-(1-met...	152	C10H16O	007764-50-3	68



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU070122\
Data File : VU049594.D
Acq On : 01 Jul 2022 20:08
Operator : SY/MD
Sample : N3564-09 10X
Misc : 5.98g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
WC-14B-11-2-GR

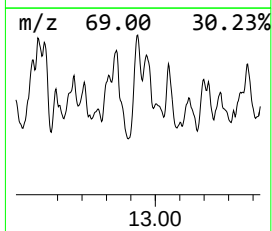
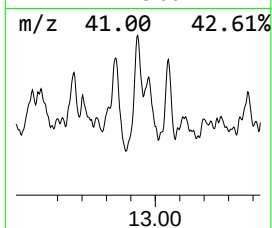
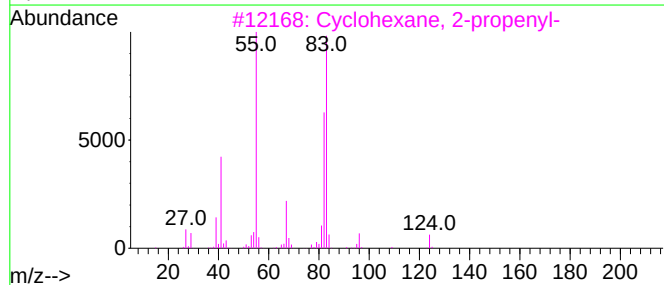
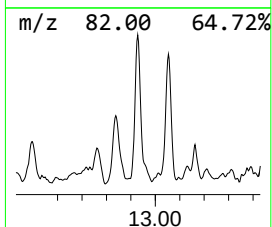
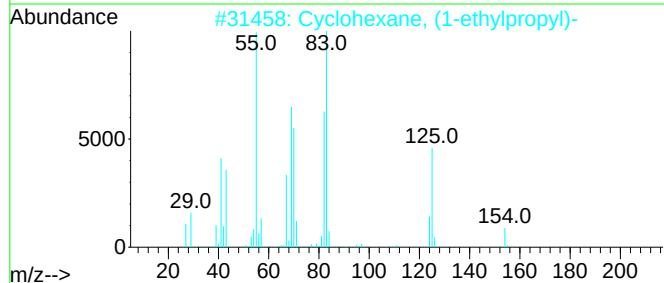
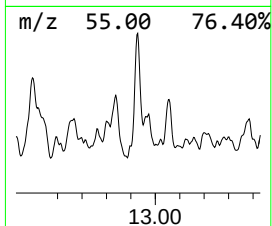
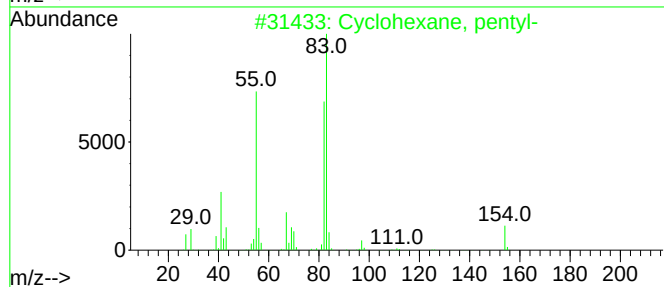
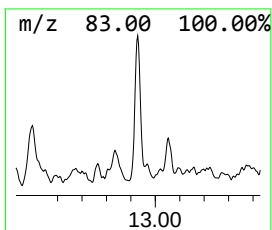
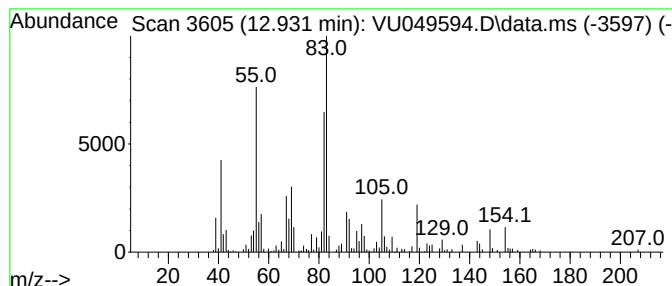
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Quant Title : SW846 8260

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

Peak Number 9 Cyclohexane, pentyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.931	10.21 ug/l	198793	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, pentyl-	154	C11H22	004292-92-6	60
2			Cyclohexane, (1-ethylpropyl)-	154	C11H22	026321-98-2	52
3			Cyclohexane, 2-propenyl-	124	C9H16	002114-42-3	49
4			Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	47
5			Cyclopentane, 1-methyl-2-(2-prop...	124	C9H16	050746-53-7	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU070122\
Data File : VU049594.D
Acq On : 01 Jul 2022 20:08
Operator : SY/MD
Sample : N3564-09 10X
Misc : 5.98g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
WC-14B-11-2-GR

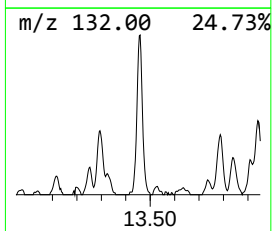
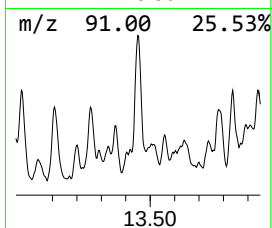
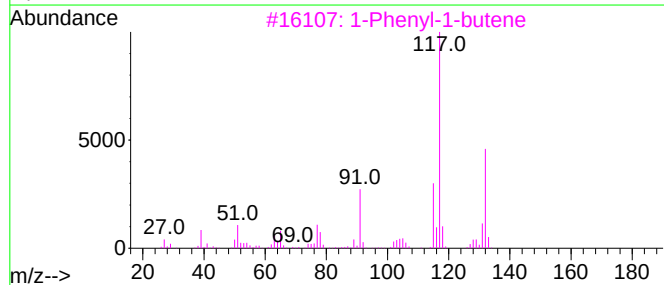
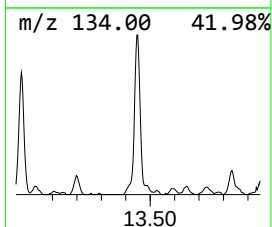
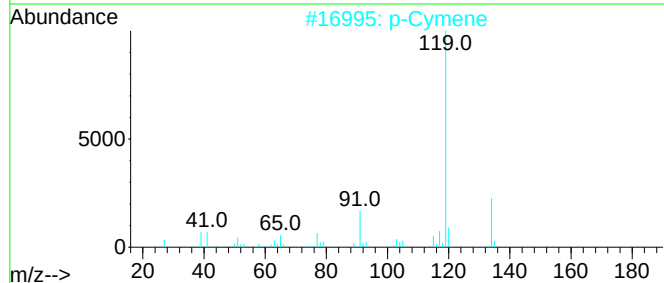
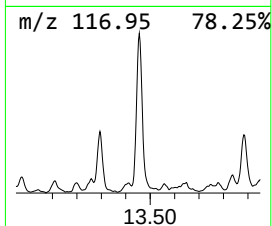
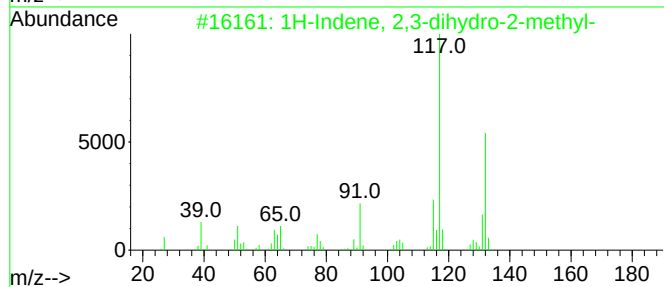
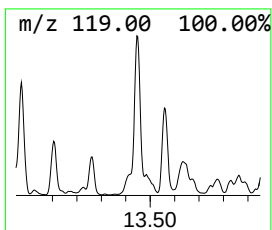
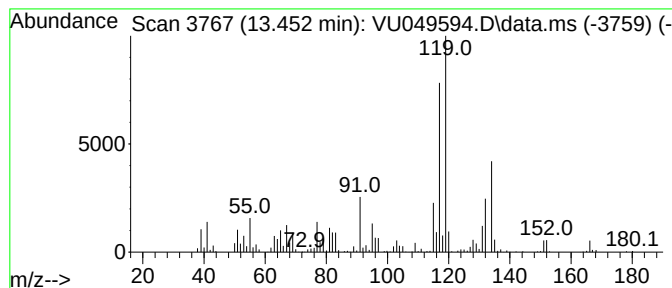
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Quant Title : SW846 8260

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

Peak Number 10 1H-Indene, 2,3-dihydro-2-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.452	20.97 ug/l	408182	1,4-Dichlorobenzene-d4	11.812

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	50
2		p-Cymene	134	C10H14	000099-87-6	50
3		1-Phenyl-1-butene	132	C10H12	000824-90-8	50
4		Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	50
5		o-Cymene	134	C10H14	000527-84-4	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU070122\
Data File : VU049594.D
Acq On : 01 Jul 2022 20:08
Operator : SY/MD
Sample : N3564-09 10X
Misc : 5.98g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
WC-14B-11-2-GR

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Cyclohexane, 1,...	8.925	9.2	ug/l	172435	3	9.417	936126	50.0
trans-1,3-Dieth...	10.025	12.8	ug/l	240472	3	9.417	936126	50.0
Octane, 2,6-dim...	10.253	21.9	ug/l	410668	3	9.417	936126	50.0
Cyclohexane, pr...	10.356	11.4	ug/l	213636	3	9.417	936126	50.0
unknown10.555	10.555	9.1	ug/l	170746	3	9.417	936126	50.0
Cyclohexane, 1,...	10.809	12.0	ug/l	233891	4	11.812	973201	50.0
Cyclohexane, 1-...	11.063	10.7	ug/l	207216	4	11.812	973201	50.0
Cyclohexanone, ...	12.838	12.1	ug/l	235129	4	11.812	973201	50.0
Cyclohexane, pe...	12.931	10.2	ug/l	198793	4	11.812	973201	50.0
1H-Indene, 2,3-...	13.452	21.0	ug/l	408182	4	11.812	973201	50.0