

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU053122\
 Data File : VU048771.D
 Acq On : 01 Jun 2022 04:57
 Operator : SY/MD
 Sample : N3075-03 20X
 Misc : 6.11g/10.0mL/100uL/5.0mL/MSVOA_U/MEOH
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 SS-3

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\82U052622W.M
 Title : SW846 8260

Signal : TIC: VU048771.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.125	236	244	264	rVB	525824	714524	4.53%	0.167%
2	3.012	494	520	532	rBV	1026044	2510342	15.91%	0.585%
3	3.298	592	609	629	rVB	665502	1507462	9.55%	0.352%
4	3.639	694	715	740	rBV	1612283	3624938	22.97%	0.845%
5	4.289	887	917	931	rBV5	146651	552868	3.50%	0.129%
6	4.488	962	979	996	rVV	964295	2400149	15.21%	0.560%
7	5.096	1145	1168	1177	rBV2	274020	711419	4.51%	0.166%
8	5.350	1232	1247	1263	rVV7	2352455	6910224	43.79%	1.612%
9	5.433	1264	1273	1297	rVB3	658299	1851816	11.74%	0.432%
10	5.565	1297	1314	1343	rBV	2022840	4902306	31.07%	1.143%
11	5.826	1383	1395	1407	rVV	885259	1862591	11.80%	0.434%
12	5.899	1407	1418	1426	rVV2	672041	1448724	9.18%	0.338%
13	5.967	1427	1439	1453	rVB3	1159923	2780743	17.62%	0.649%
14	6.115	1469	1485	1510	rBV	5382231	10202021	64.66%	2.379%
15	6.279	1510	1536	1559	rBV3	1101934	4023243	25.50%	0.938%
16	6.546	1607	1619	1646	rVB6	644329	2000667	12.68%	0.467%
17	6.748	1659	1682	1694	rBV	6940789	15778908	100.00%	3.680%
18	6.854	1706	1715	1728	rVB	790897	1495311	9.48%	0.349%
19	6.967	1738	1750	1765	rBV	993663	2106522	13.35%	0.491%
20	7.060	1767	1779	1791	rVB3	950943	2002632	12.69%	0.467%
21	7.150	1791	1807	1820	rBV4	456997	1219889	7.73%	0.284%
22	7.227	1821	1831	1851	rVB4	1055192	2778541	17.61%	0.648%
23	7.327	1851	1862	1870	rBV4	328801	653318	4.14%	0.152%
24	7.388	1870	1881	1886	rBV2	901142	1691387	10.72%	0.394%
25	7.491	1902	1913	1921	rBV	5590570	10013443	63.46%	2.335%
26	7.530	1922	1925	1935	rVB2	1972643	2596778	16.46%	0.606%
27	7.652	1953	1963	1982	rVB	3800805	6955617	44.08%	1.622%
28	7.755	1987	1995	2009	rVB6	1028505	2081924	13.19%	0.486%
29	7.851	2009	2025	2032	rBV2	4832452	10062475	63.77%	2.347%
30	8.022	2067	2078	2085	rBV4	1202857	2864964	18.16%	0.668%
31	8.128	2095	2111	2130	rVB	7480603	14338779	90.87%	3.344%
32	8.240	2132	2146	2161	rVB6	2878102	7119547	45.12%	1.660%
33	8.362	2173	2184	2191	rBV3	1828650	3379502	21.42%	0.788%
34	8.407	2192	2198	2211	rVB2	1018004	1847546	11.71%	0.431%
35	8.533	2228	2237	2248	rVB2	1055325	1734734	10.99%	0.405%
36	8.633	2250	2268	2281	rBV	3374383	6292540	39.88%	1.468%

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37	8.703	2281	2290	2295	rBV5	492813	858280	5.44%	0.200%
38	8.755	2296	2306	2313	rBV	2101320	4009695	25.41%	0.935%
39	8.864	2330	2340	2349	rBV	3982254	6536966	41.43%	1.525%
40	8.922	2350	2358	2368	rBV	3837108	6306371	39.97%	1.471%
41	9.063	2390	2402	2411	rBV7	1555851	3534840	22.40%	0.824%
42	9.118	2412	2419	2426	rVV	2108982	3349956	21.23%	0.781%
43	9.166	2426	2434	2440	rVV4	2304363	4389728	27.82%	1.024%
44	9.211	2440	2448	2459	rVB2	7162084	12042580	76.32%	2.809%
45	9.333	2476	2486	2499	rVB	6674510	10850605	68.77%	2.531%
46	9.568	2550	2559	2573	rBV2	1581116	3236162	20.51%	0.755%
47	9.716	2598	2605	2610	rBV3	1903629	2854769	18.09%	0.666%
48	9.886	2646	2658	2670	rBV5	792866	1919092	12.16%	0.448%
49	9.954	2671	2679	2689	rBV7	461141	923067	5.85%	0.215%
50	10.025	2689	2701	2712	rBV7	2778315	6804942	43.13%	1.587%
51	10.118	2723	2730	2742	rBV4	1462511	2714712	17.20%	0.633%
52	10.253	2754	2772	2786	rBV4	5404201	12685241	80.39%	2.958%
53	10.356	2795	2804	2812	rBV4	3791329	6988214	44.29%	1.630%
54	10.485	2833	2844	2853	rBV3	1145448	2501236	15.85%	0.583%
55	10.558	2854	2867	2875	rVV4	2003152	4678380	29.65%	1.091%
56	10.626	2876	2888	2892	rVV2	3815433	7490675	47.47%	1.747%
57	10.652	2893	2896	2905	rVV	3706621	4705046	29.82%	1.097%
58	10.697	2906	2910	2919	rVV4	638868	1089448	6.90%	0.254%
59	10.758	2921	2929	2938	rVV	3165379	5215445	33.05%	1.216%
60	10.809	2939	2945	2963	rVB6	1792861	3448651	21.86%	0.804%
61	10.909	2968	2976	2984	rBV	1684290	2760197	17.49%	0.644%
62	11.025	3004	3012	3018	rBV3	1293902	2097593	13.29%	0.489%
63	11.066	3019	3025	3028	rBV2	1284734	1628831	10.32%	0.380%
64	11.304	3087	3099	3113	rBV5	1973582	4731050	29.98%	1.103%
65	11.375	3115	3121	3127	rVV3	978771	1417891	8.99%	0.331%
66	11.417	3127	3134	3142	rVV	3672933	5097286	32.30%	1.189%
67	11.472	3142	3151	3174	rVB3	2791355	6924198	43.88%	1.615%
68	11.575	3177	3183	3189	rBV2	750168	1149299	7.28%	0.268%
69	11.642	3199	3204	3215	rVB4	1796201	2902331	18.39%	0.677%
70	11.713	3217	3226	3233	rBV2	2227120	3783072	23.98%	0.882%
71	11.902	3271	3285	3297	rVB3	6298395	13325648	84.45%	3.108%
72	12.002	3311	3316	3335	rVB4	2313037	4040711	25.61%	0.942%
73	12.102	3336	3347	3363	rBV2	4588113	9596280	60.82%	2.238%
74	12.192	3364	3375	3390	rVB5	5836189	12336730	78.18%	2.877%
75	12.382	3427	3434	3449	rVB	2610309	4776347	30.27%	1.114%
76	12.468	3453	3461	3467	rBV	1876068	3139037	19.89%	0.732%

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77	12.574	3486	3494	3502	rVB	4880731	6761056	42.85%	1.577%
78	12.684	3517	3528	3549	rVB5	3251080	8931920	56.61%	2.083%
79	12.877	3583	3588	3596	rVB3	1468191	2069432	13.12%	0.483%
80	12.931	3597	3605	3609	rBV3	1257302	2042145	12.94%	0.476%
81	12.973	3612	3618	3626	rVB	2702307	3996022	25.33%	0.932%
82	13.031	3630	3636	3651	rVB3	2304700	5399767	34.22%	1.259%
83	13.111	3652	3661	3663	rBV	1373464	1873129	11.87%	0.437%
84	13.295	3710	3718	3728	rVB3	2632502	4220633	26.75%	0.984%
85	13.356	3731	3737	3745	rVB3	962603	1342163	8.51%	0.313%
86	13.455	3746	3768	3783	rBV5	5981896	14779260	93.66%	3.447%
87	13.562	3796	3801	3805	rBV	684262	791615	5.02%	0.185%
88	13.629	3816	3822	3845	rVB9	1577518	3902571	24.73%	0.910%
89	13.735	3846	3855	3862	rBV2	885767	1561076	9.89%	0.364%
90	13.787	3863	3871	3879	rVV2	1515862	2396179	15.19%	0.559%
91	13.838	3879	3887	3902	rVB5	1567042	3154639	19.99%	0.736%
92	13.944	3914	3920	3934	rVB2	1213108	2488323	15.77%	0.580%
93	14.089	3956	3965	3984	rVB	3686931	6733827	42.68%	1.570%
94	14.288	4020	4027	4039	rVB5	593052	1189477	7.54%	0.277%
95	14.484	4082	4088	4108	rVB2	866572	1516685	9.61%	0.354%
96	14.668	4136	4145	4157	rBV3	857504	1675013	10.62%	0.391%
97	14.812	4182	4190	4203	rBV4	327890	717428	4.55%	0.167%
98	14.877	4203	4210	4220	rVB2	557330	910305	5.77%	0.212%
99	15.221	4307	4317	4331	rVB	2766607	4432204	28.09%	1.034%
100	15.410	4365	4376	4387	rBV	1229407	2044850	12.96%	0.477%

Sum of corrected areas: 428788715

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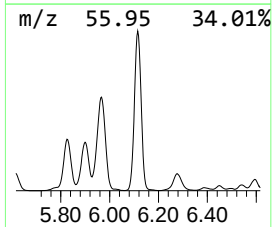
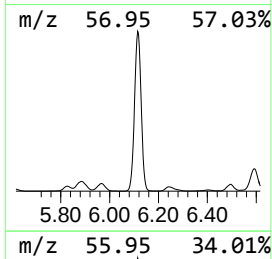
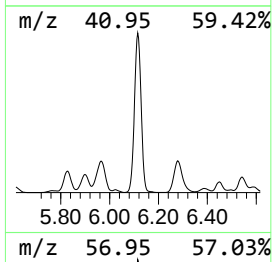
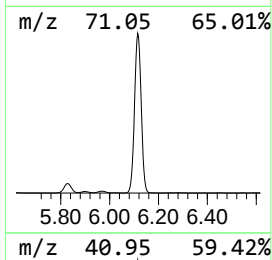
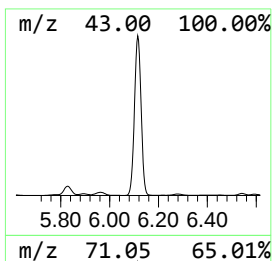
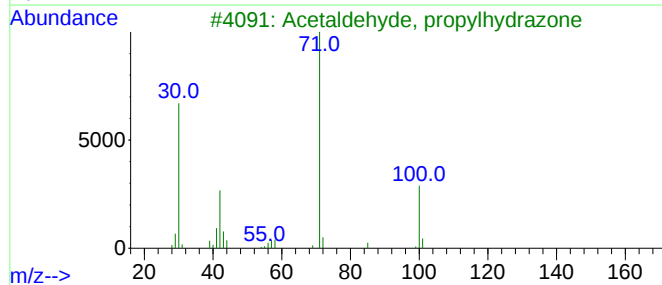
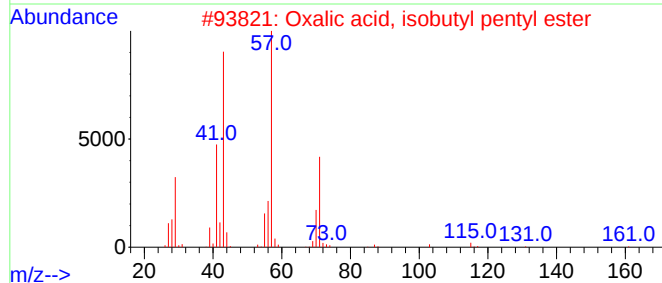
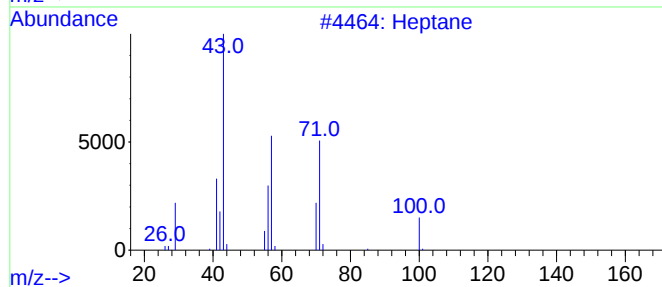
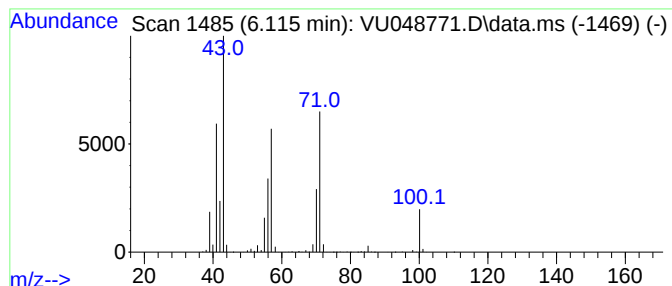
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U052622W.M
Quant Title : SW846 8260

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

Peak Number 1 Heptane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.115	126.79 ug/l	10202000	1,4-Difluorobenzene	6.244

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane	100	C7H16	000142-82-5	91
2			Oxalic acid, isobutyl pentyl ester	216	C11H20O4	1000309-37-0	50
3			Acetaldehyde, propylhydrazone	100	C5H12N2	007422-88-0	47
4			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	40
5			Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	38



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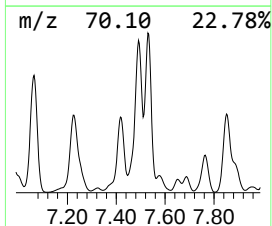
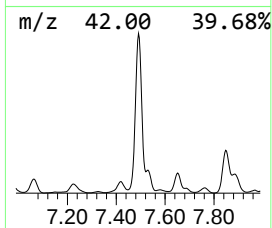
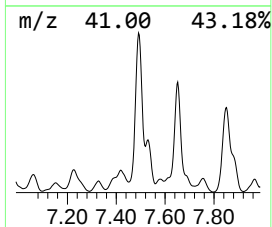
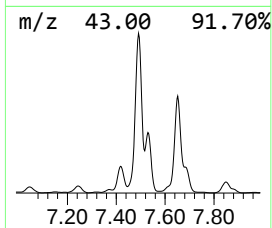
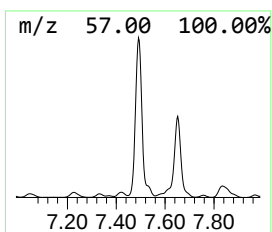
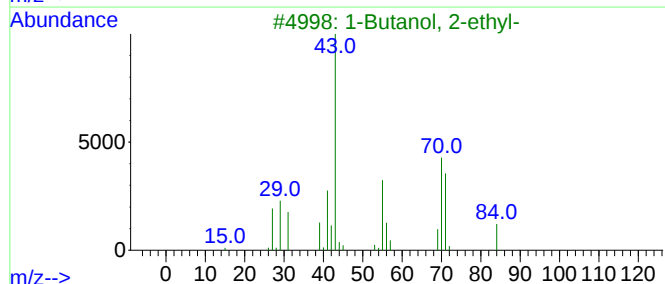
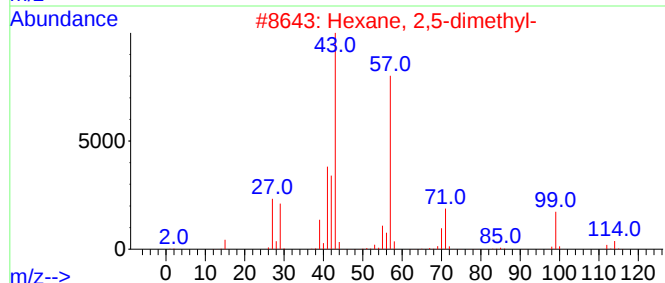
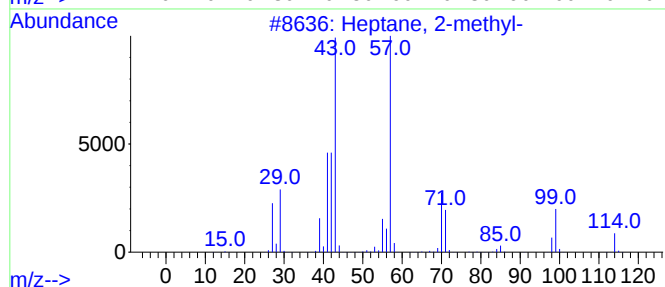
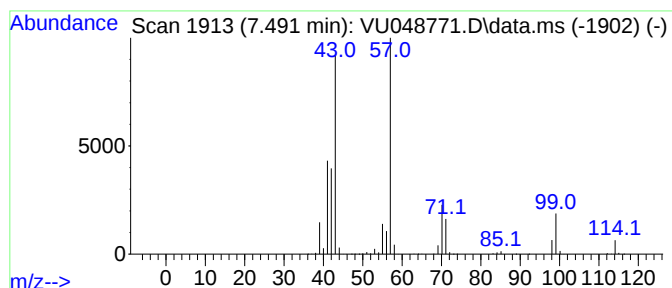
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U052622W.M
Quant Title : SW846 8260

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

Peak Number 2 Heptane, 2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.491	124.44 ug/l	10013400	1,4-Difluorobenzene	6.244

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane, 2-methyl-	114	C8H18	000592-27-8	94
2			Hexane, 2,5-dimethyl-	114	C8H18	000592-13-2	87
3			1-Butanol, 2-ethyl-	102	C6H14O	000097-95-0	35
4			Hexane, 2,4,4-trimethyl-	128	C9H20	016747-30-1	32
5			Cyclohexanol, 4-methyl-, cis-	114	C7H14O	007731-28-4	32



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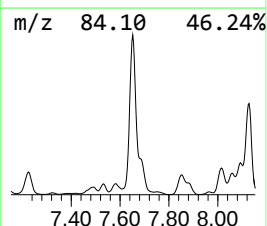
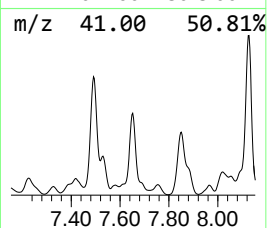
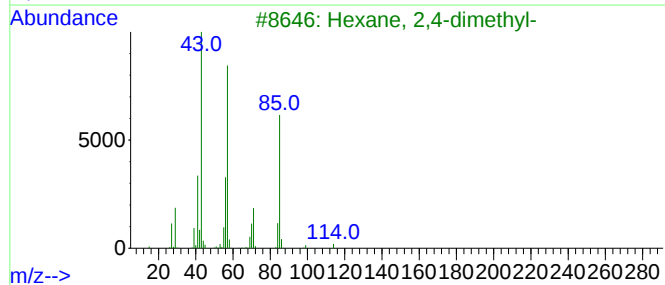
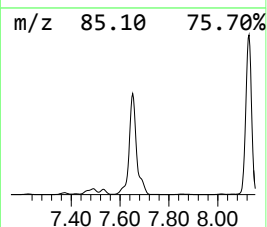
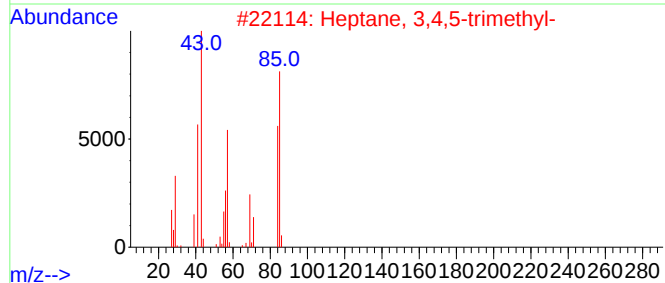
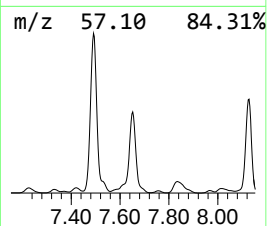
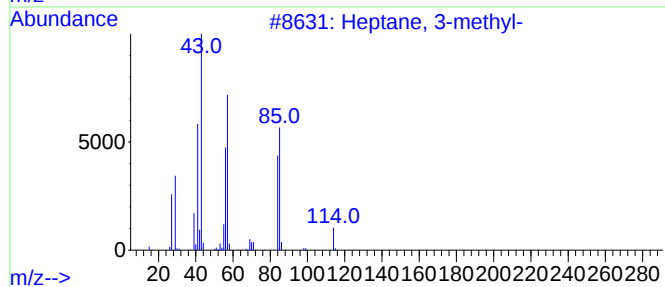
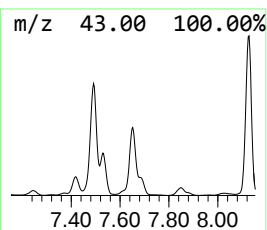
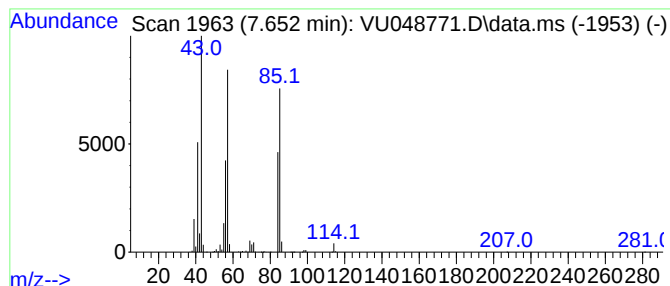
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U052622W.M
Quant Title : SW846 8260

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

Peak Number 3 Heptane, 3-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.652	86.44 ug/l	6955620	1,4-Difluorobenzene	6.244

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane, 3-methyl-	114	C8H18	000589-81-1	94
2			Heptane, 3,4,5-trimethyl-	142	C10H22	020278-89-1	86
3			Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	74
4			Pentane, 3-ethyl-2,4-dimethyl-	128	C9H20	001068-87-7	59
5			2-Methylpentyl isovalerate	186	C11H22O2	1000236-38-7	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU053122\
Data File : VU048771.D
Acq On : 01 Jun 2022 04:57
Operator : SY/MD
Sample : N3075-03 20X
Misc : 6.11g/10.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 47 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
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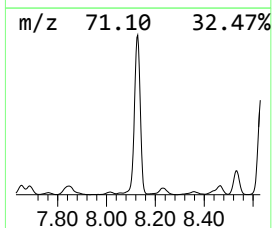
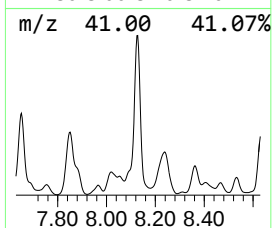
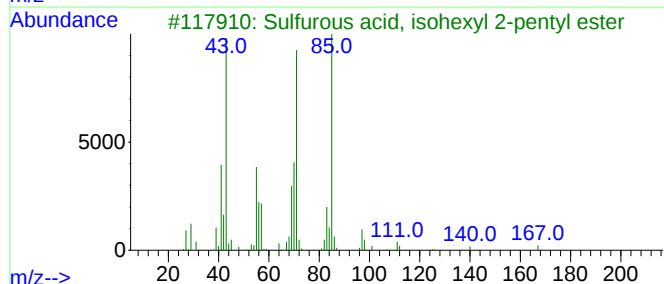
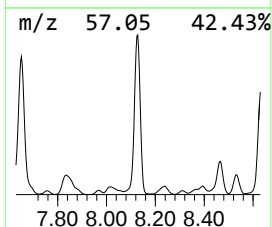
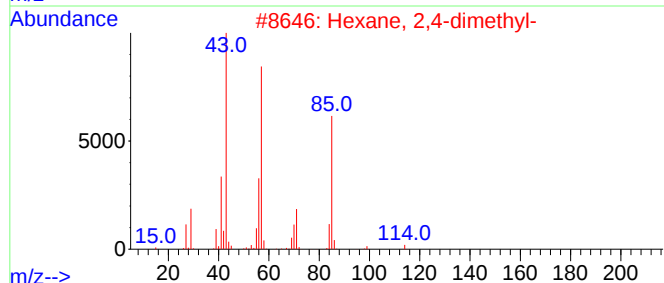
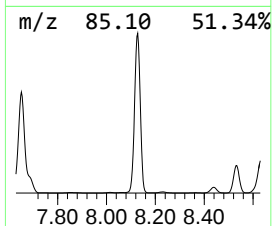
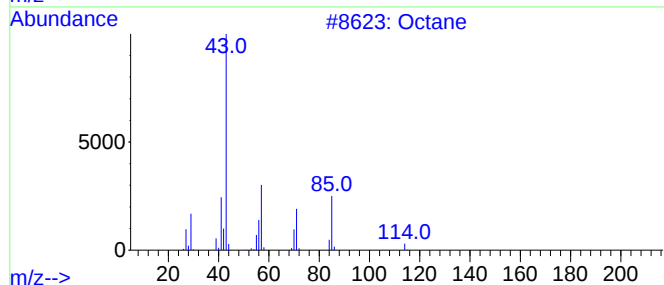
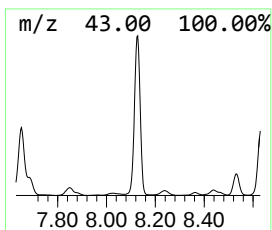
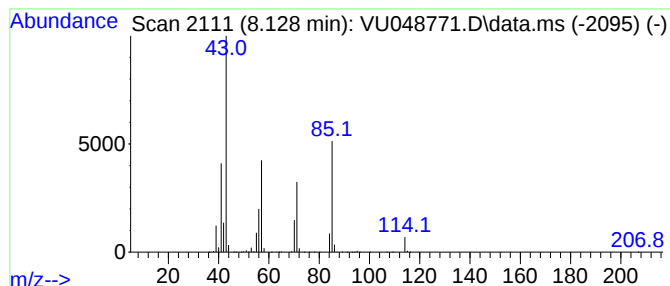
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Quant Title : SW846 8260

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

Peak Number 4 Octane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.128	66.07 ug/l	14338800	Chlorobenzene-d5	9.420

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane			114	C8H18	000111-65-9	91
2	Hexane, 2,4-dimethyl-			114	C8H18	000589-43-5	64
3	Sulfurous acid, isohexyl 2-penty...			236	C11H24O3S	1000309-15-5	64
4	Undecane, 2,4-dimethyl-			184	C13H28	017312-80-0	64
5	Ethanol, 2-(hexyloxy)-			146	C8H18O2	000112-25-4	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU053122\
Data File : VU048771.D
Acq On : 01 Jun 2022 04:57
Operator : SY/MD
Sample : N3075-03 20X
Misc : 6.11g/10.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 47 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
SS-3

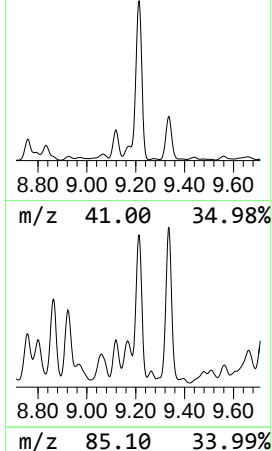
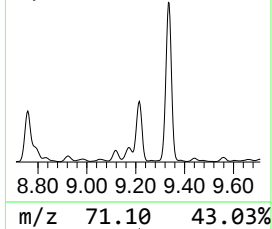
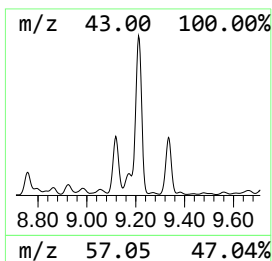
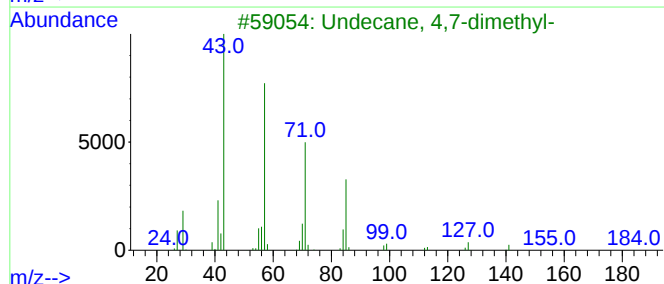
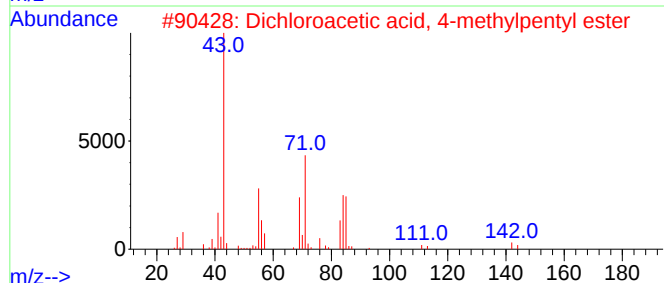
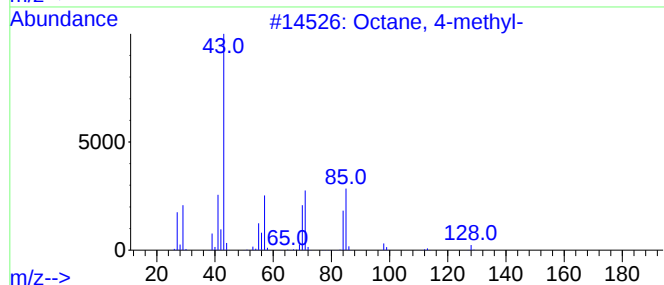
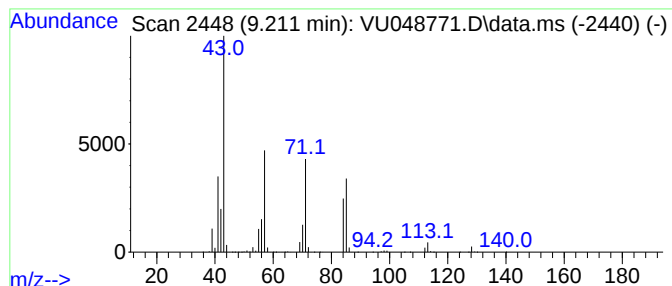
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Quant Title : SW846 8260

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

Peak Number 5 Octane, 4-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.211	55.49 ug/l	12042600	Chlorobenzene-d5	9.420

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 4-methyl-	128	C9H20	002216-34-4	64
2			Dichloroacetic acid, 4-methylpen...	212	C8H14Cl2O2	1000282-43-7	64
3			Undecane, 4,7-dimethyl-	184	C13H28	017301-32-5	53
4			Sulfurous acid, 2-propyl tetradec...	320	C17H36O3S	1010309-12-5	53
5			Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU053122\
Data File : VU048771.D
Acq On : 01 Jun 2022 04:57
Operator : SY/MD
Sample : N3075-03 20X
Misc : 6.11g/10.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 47 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
SS-3

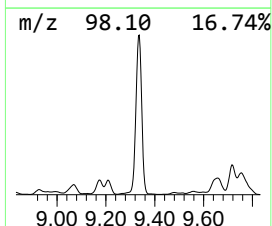
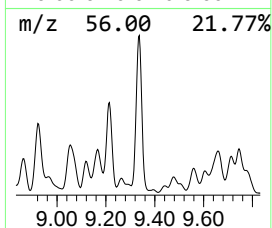
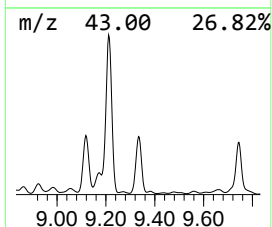
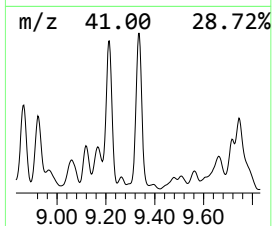
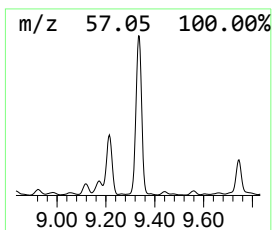
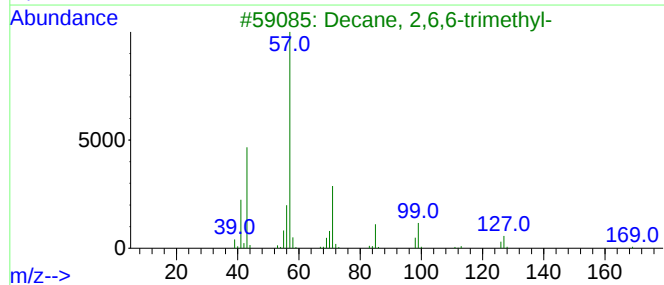
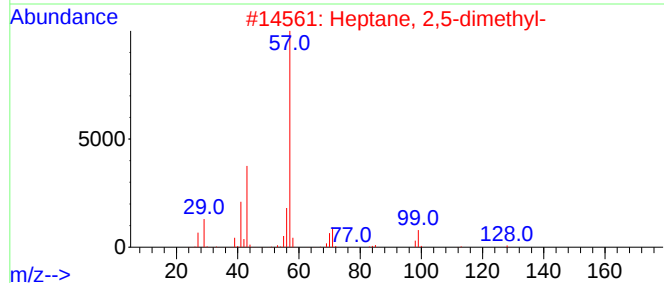
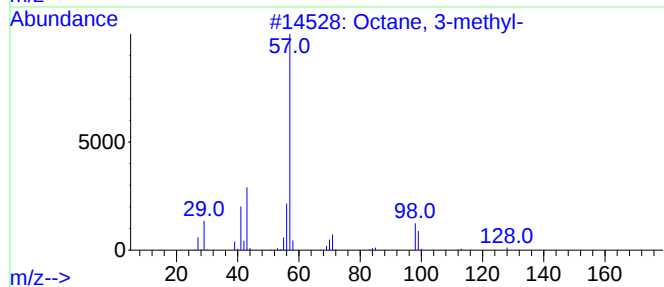
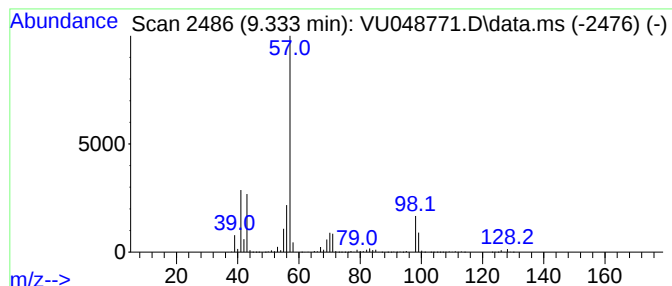
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Quant Title : SW846 8260

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

Peak Number 6 Octane, 3-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.333	50.00 ug/l	10850600	Chlorobenzene-d5	9.420

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 3-methyl-	128	C9H20	002216-33-3	91
2			Heptane, 2,5-dimethyl-	128	C9H20	002216-30-0	80
3			Decane, 2,6,6-trimethyl-	184	C13H28	062108-24-1	64
4			Undecane, 6-methyl-	170	C12H26	017302-33-9	59
5			Heptane, 4-propyl-	142	C10H22	003178-29-8	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU053122\
Data File : VU048771.D
Acq On : 01 Jun 2022 04:57
Operator : SY/MD
Sample : N3075-03 20X
Misc : 6.11g/10.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 47 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
SS-3

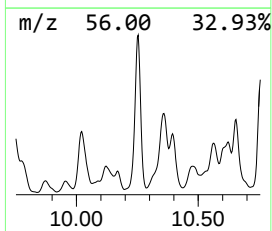
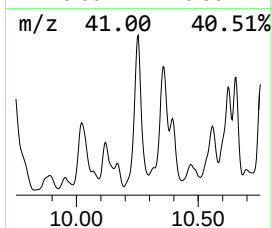
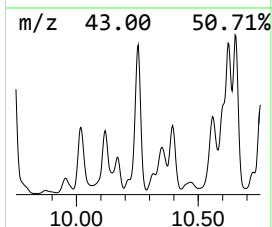
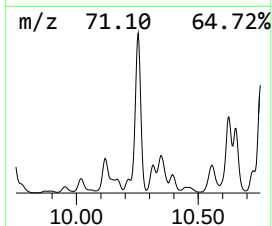
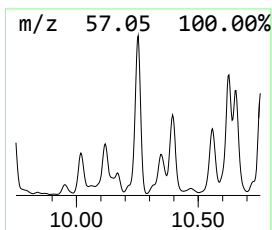
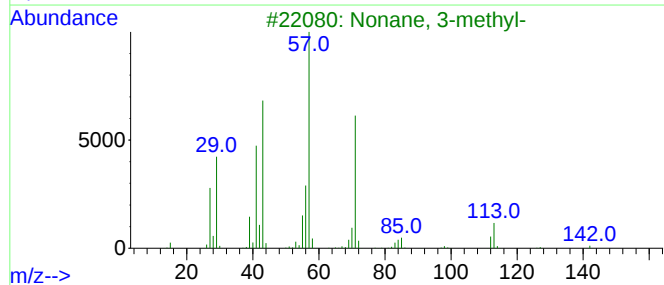
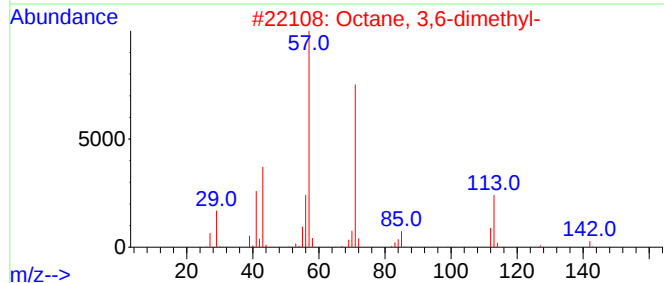
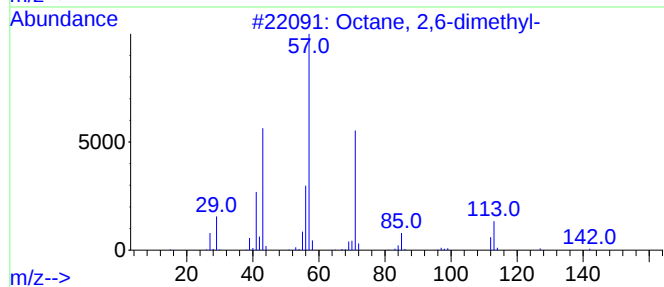
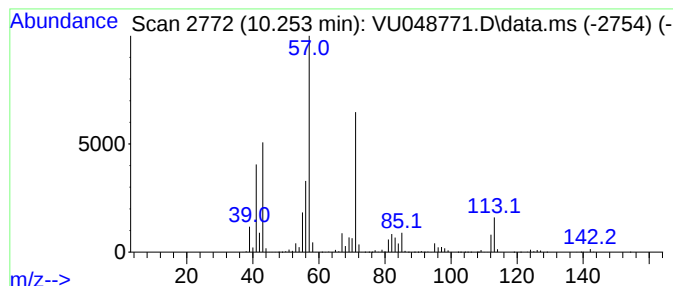
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Quant Title : SW846 8260

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

Peak Number 7 Octane, 2,6-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.253	58.45 ug/l	12685200	Chlorobenzene-d5	9.420

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	95
2			Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	91
3			Nonane, 3-methyl-	142	C10H22	005911-04-6	90
4			Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	64
5			Sulfurous acid, 2-ethylhexyl pen...	264	C13H28O3S	1000309-18-9	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU053122\
Data File : VU048771.D
Acq On : 01 Jun 2022 04:57
Operator : SY/MD
Sample : N3075-03 20X
Misc : 6.11g/10.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 47 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
SS-3

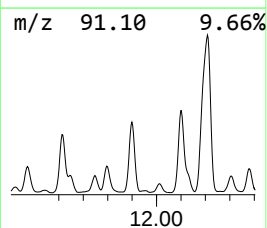
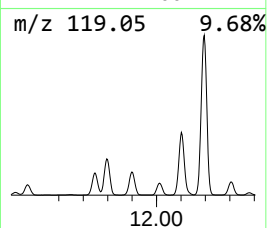
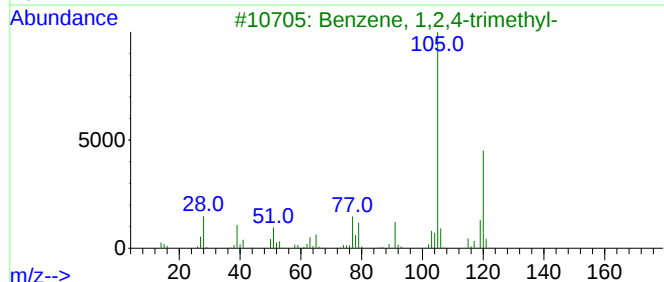
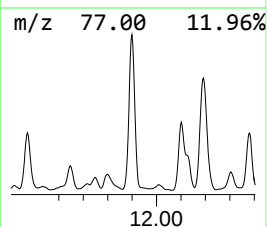
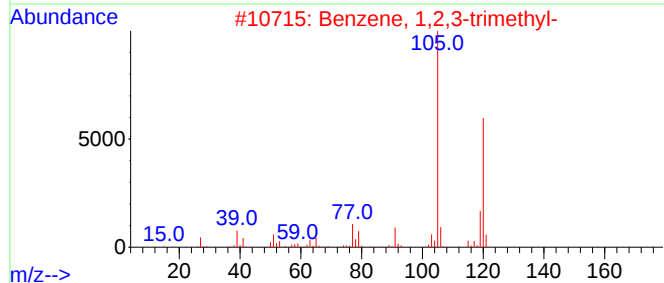
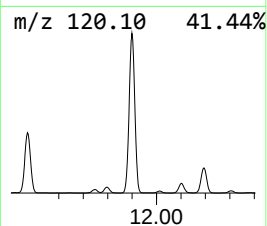
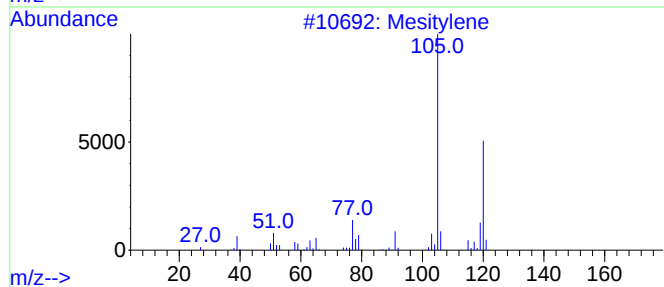
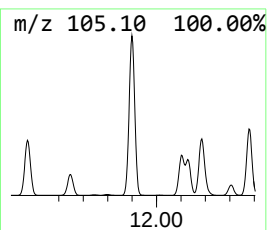
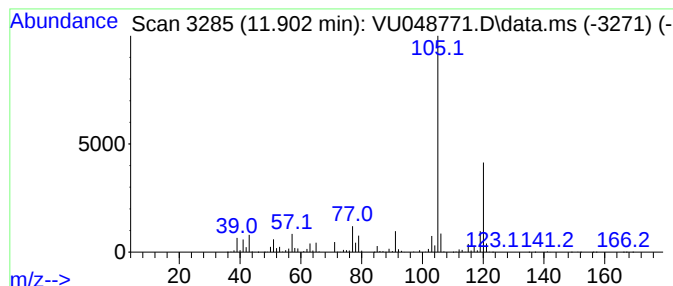
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,3-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.903	50.00 ug/l	13325600	1,4-Dichlorobenzene-d4	11.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Mesitylene	120	C9H12	000108-67-8	97
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU053122\
Data File : VU048771.D
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Operator : SY/MD
Sample : N3075-03 20X
Misc : 6.11g/10.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 47 Sample Multiplier: 1

Instrument :
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ClientSampleId :
SS-3

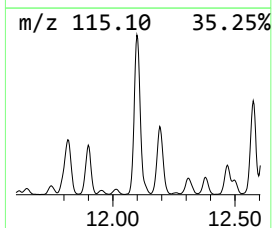
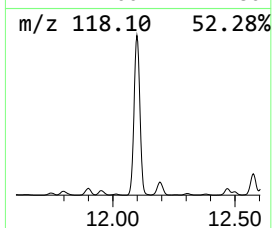
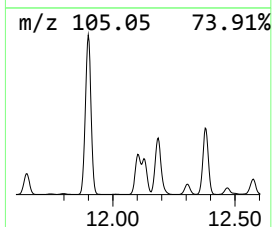
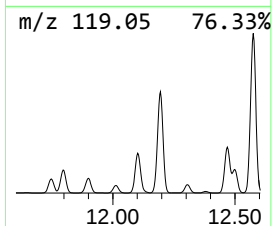
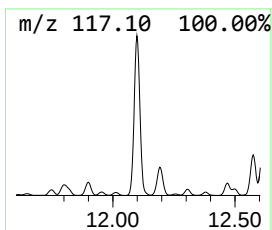
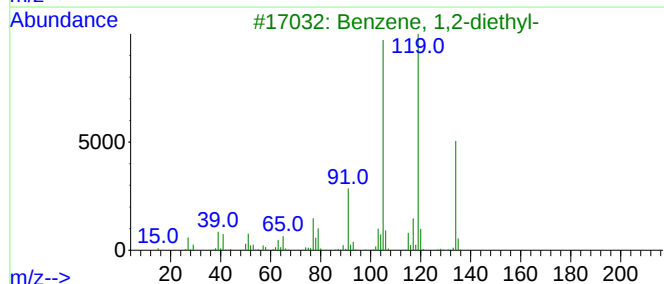
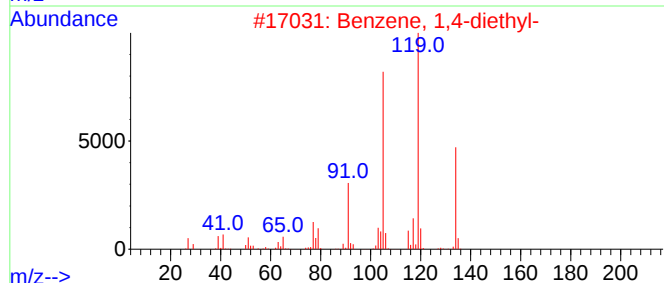
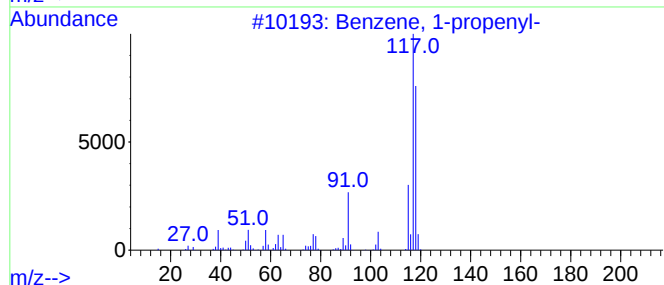
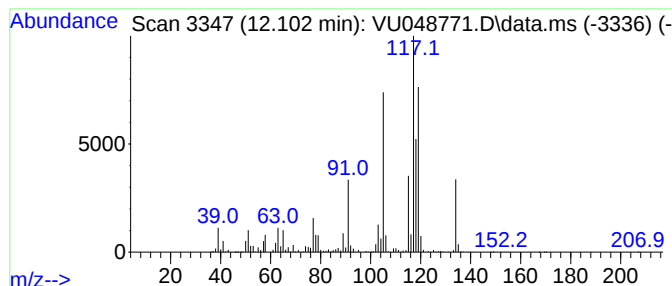
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1-propenyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.102	36.01 ug/l	9596280	1,4-Dichlorobenzene-d4	11.819

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-propenyl-	118	C9H10	000637-50-3	64
2			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	55
3			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	55
4			Benzene, cyclopropyl-	118	C9H10	000873-49-4	50
5			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU053122\
Data File : VU048771.D
Acq On : 01 Jun 2022 04:57
Operator : SY/MD
Sample : N3075-03 20X
Misc : 6.11g/10.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 47 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
SS-3

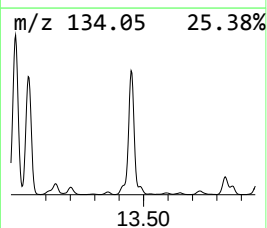
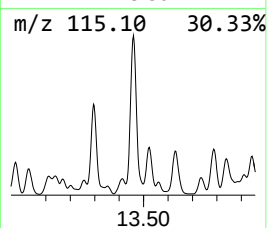
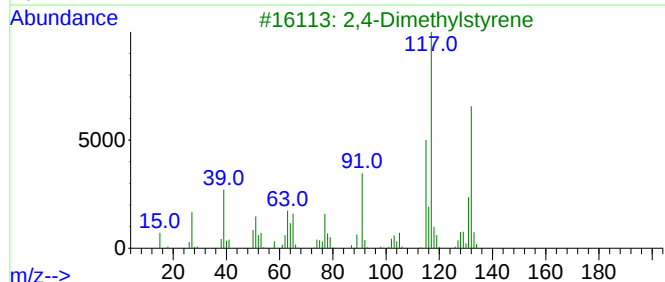
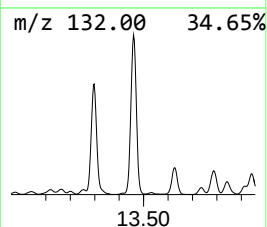
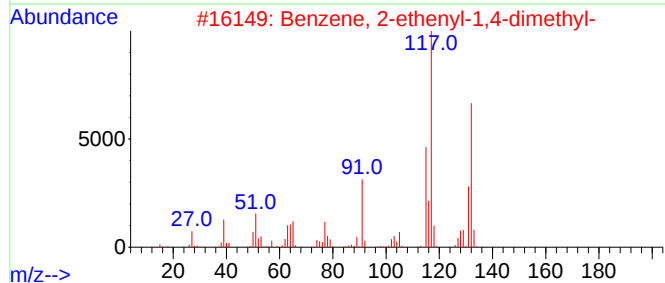
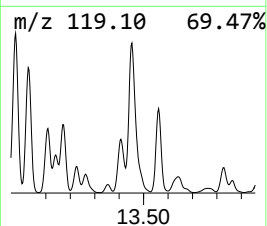
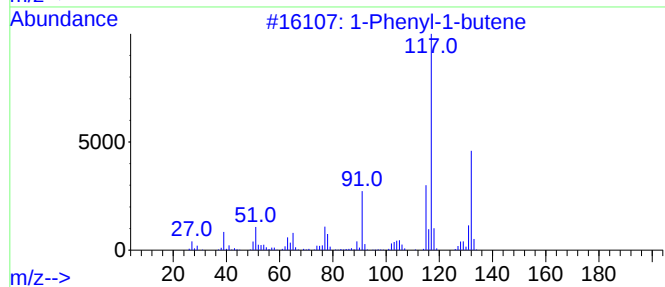
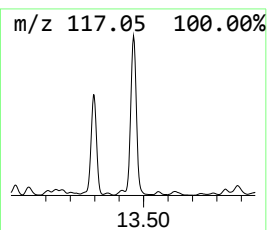
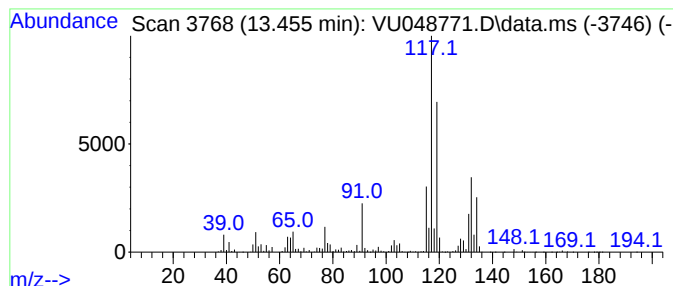
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U052622W.M
Quant Title : SW846 8260

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

Peak Number 10 1-Phenyl-1-butene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.456	55.45 ug/l	14779300	1,4-Dichlorobenzene-d4	11.819

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Phenyl-1-butene	132	C10H12	000824-90-8	96
2		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
3		2,4-Dimethylstyrene	132	C10H12	002234-20-0	90
4		Benzene, 2-butenyl-	132	C10H12	001560-06-1	70
5		Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	70



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU053122\
Data File : VU048771.D
Acq On : 01 Jun 2022 04:57
Operator : SY/MD
Sample : N3075-03 20X
Misc : 6.11g/10.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 47 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
SS-3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U052622W.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
Heptane	6.115	126.8	ug/l	10202000	2	6.244	4023240	50.0
Heptane, 2-methyl-	7.491	124.4	ug/l	10013400	2	6.244	4023240	50.0
Heptane, 3-methyl-	7.652	86.4	ug/l	6955620	2	6.244	4023240	50.0
Octane	8.128	66.1	ug/l	14338800	3	9.420	10850600	50.0
Octane, 4-methyl-	9.211	55.5	ug/l	12042600	3	9.420	10850600	50.0
Octane, 3-methyl-	9.333	50.0	ug/l	10850600	3	9.420	10850600	50.0
Octane, 2,6-dim...	10.253	58.5	ug/l	12685200	3	9.420	10850600	50.0
Benzene, 1,2,3-...	11.903	50.0	ug/l	13325600	4	11.819	13325600	50.0
Benzene, 1-prop...	12.102	36.0	ug/l	9596280	4	11.819	13325600	50.0
1-Phenyl-1-butene	13.456	55.5	ug/l	14779300	4	11.819	13325600	50.0