

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060421\
 Data File : VX022419.D
 Acq On : 04 Jun 2021 17:48
 Operator : JC/MD
 Sample : M2583-04 10X
 Misc : 4.99g/10mL/100uL/5.0mL/MSVOA_X/MEOH
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 40955

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_X\Method\82X051921W.M
 Title : SW846 8260

Signal : TIC: VX022419.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.362	41	45	49	rBV2	9501	14509	1.14%	0.123%
2	1.563	70	78	79	rBV8	5843	12784	1.00%	0.109%
3	2.392	207	214	228	rBV	700505	1122204	87.98%	9.536%
4	2.947	296	305	315	rBV	6089	15823	1.24%	0.134%
5	5.391	695	706	716	rBV2	48069	140177	10.99%	1.191%
6	5.562	716	734	744	rBV2	76271	252654	19.81%	2.147%
7	5.824	768	777	791	rBV3	19324	55597	4.36%	0.472%
8	5.964	791	800	811	rBV	60279	162958	12.78%	1.385%
9	6.159	820	832	842	rBV8	3774	18218	1.43%	0.155%
10	6.617	898	907	916	rVB2	24947	60171	4.72%	0.511%
11	6.769	922	932	945	rBV	142049	338785	26.56%	2.879%
12	7.379	1022	1032	1042	rBV4	10332	27341	2.14%	0.232%
13	8.653	1235	1241	1247	rBV	292015	464420	36.41%	3.947%
14	8.720	1247	1252	1266	rVB	810511	1275488	100.00%	10.839%
15	8.915	1277	1284	1289	rBV3	9157	13940	1.09%	0.118%
16	9.909	1443	1447	1452	rVB	12198	17241	1.35%	0.147%
17	10.012	1457	1464	1466	rBV	11091	14909	1.17%	0.127%
18	10.055	1466	1471	1484	rVB	289787	403247	31.62%	3.427%
19	10.195	1489	1494	1500	rVV	42577	59475	4.66%	0.505%
20	10.305	1500	1512	1518	rVV	173708	250729	19.66%	2.131%
21	10.366	1518	1522	1532	rVB	36015	52160	4.09%	0.443%
22	10.592	1552	1559	1563	rBV4	9363	14110	1.11%	0.120%
23	10.646	1563	1568	1573	rBV	91119	120155	9.42%	1.021%
24	10.787	1583	1591	1595	rBV3	16902	27446	2.15%	0.233%
25	10.860	1595	1603	1607	rBV3	11542	20850	1.63%	0.177%
26	10.963	1614	1620	1624	rBV2	13613	22547	1.77%	0.192%
27	11.085	1634	1640	1644	rVV	255678	336024	26.34%	2.856%
28	11.116	1644	1645	1651	rVB2	34681	37514	2.94%	0.319%
29	11.195	1651	1658	1661	rBV	24412	35005	2.74%	0.297%
30	11.305	1666	1676	1684	rVV	58675	87572	6.87%	0.744%
31	11.384	1684	1689	1696	rVV	223166	407727	31.97%	3.465%
32	11.457	1696	1701	1703	rVV	126092	182586	14.31%	1.552%
33	11.482	1703	1705	1713	rVB	126769	133611	10.48%	1.135%
34	11.616	1723	1727	1734	rVB	96042	124789	9.78%	1.060%
35	11.689	1734	1739	1741	rBV2	14290	18260	1.43%	0.155%
36	11.719	1741	1744	1746	rVV	18548	20805	1.63%	0.177%

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060421\
 Data File : VX022419.D
 Acq On : 04 Jun 2021 17:48
 Operator : JC/MD
 Sample : M2583-04 10X
 Misc : 4.99g/10mL/100uL/5.0mL/MSVOA_X/MEOH
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 40955

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_X\Method\82X051921W.M
 Title : SW846 8260

37	11.756	1746	1750	1759	rVB	529613	665370	52.17%	5.654%
38	11.841	1759	1764	1767	rBV	25675	40099	3.14%	0.341%
39	11.890	1767	1772	1776	rVB2	326915	399456	31.32%	3.395%
40	11.939	1776	1780	1783	rBV2	53699	67781	5.31%	0.576%

41	11.975	1783	1786	1789	rVB	74202	80034	6.27%	0.680%
42	12.024	1789	1794	1800	rBV2	318291	446353	34.99%	3.793%
43	12.103	1800	1807	1813	rBV3	261094	507788	39.81%	4.315%
44	12.152	1813	1815	1818	rVB	34965	31971	2.51%	0.272%
45	12.250	1826	1831	1833	rBV	116996	165920	13.01%	1.410%

46	12.317	1837	1842	1849	rVB4	141570	251038	19.68%	2.133%
47	12.427	1849	1860	1864	rBV	151082	199894	15.67%	1.699%
48	12.469	1864	1867	1873	rVB2	41196	63493	4.98%	0.540%
49	12.536	1873	1878	1880	rBV	59596	88570	6.94%	0.753%
50	12.555	1880	1881	1884	rVV	43080	38010	2.98%	0.323%

51	12.616	1887	1891	1895	rVB	72914	81282	6.37%	0.691%
52	12.701	1901	1905	1914	rVB6	22551	61789	4.84%	0.525%
53	12.823	1918	1925	1927	rVV6	14201	24825	1.95%	0.211%
54	12.853	1927	1930	1933	rVB	14675	14383	1.13%	0.122%
55	12.920	1933	1941	1944	rBV5	41802	94359	7.40%	0.802%

56	12.994	1945	1953	1958	rVV3	106583	237024	18.58%	2.014%
57	13.048	1958	1962	1968	rVB2	46773	69784	5.47%	0.593%
58	13.170	1977	1982	1986	rBV4	34105	52093	4.08%	0.443%
59	13.219	1986	1990	1992	rVV4	24984	32584	2.55%	0.277%
60	13.268	1993	1998	2001	rVV	229209	281430	22.06%	2.392%

61	13.298	2001	2003	2008	rVV2	83870	97983	7.68%	0.833%
62	13.347	2008	2011	2013	rVV4	13778	18199	1.43%	0.155%
63	13.390	2015	2018	2020	rVV	47658	59931	4.70%	0.509%
64	13.420	2020	2023	2033	rVB8	38077	88554	6.94%	0.753%
65	13.506	2033	2037	2042	rBV7	14922	27915	2.19%	0.237%

66	13.585	2047	2050	2056	rVV4	19144	33292	2.61%	0.283%
67	13.682	2061	2066	2067	rBV2	25836	35259	2.76%	0.300%
68	13.737	2072	2075	2078	rVV3	79492	111366	8.73%	0.946%
69	13.774	2078	2081	2084	rVV2	68821	93963	7.37%	0.798%
70	13.804	2084	2086	2088	rVV2	30448	35293	2.77%	0.300%

71	13.847	2088	2093	2100	rVV3	63617	121221	9.50%	1.030%
72	13.914	2100	2104	2111	rVV8	16537	37112	2.91%	0.315%
73	14.042	2121	2125	2128	rBV	139068	143480	11.25%	1.219%
74	14.073	2128	2130	2133	rBV3	12583	14512	1.14%	0.123%
75	14.219	2151	2154	2161	rVB5	22214	40733	3.19%	0.346%

76	14.438	2183	2190	2193	rBV3	42111	68811	5.39%	0.585%
----	--------	------	------	------	------	-------	-------	-------	--------

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060421\
 Data File : VX022419.D
 Acq On : 04 Jun 2021 17:48
 Operator : JC/MD
 Sample : M2583-04 10X
 Misc : 4.99g/10mL/100uL/5.0mL/MSVOA_X/MEOH
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 40955

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_X\Method\82X051921W.M

Title : SW846 8260

77	14.475	2193	2196	2198	rVV4	14031	18962	1.49%	0.161%
78	14.505	2198	2201	2207	rVB3	28154	51121	4.01%	0.434%
79	14.566	2207	2211	2216	rBV4	15630	24188	1.90%	0.206%
80	14.615	2216	2219	2220	rBV	29926	28380	2.23%	0.241%
81	14.640	2220	2223	2227	rVB	38896	50656	3.97%	0.430%
82	14.762	2239	2243	2245	rBV	52930	56047	4.39%	0.476%
83	14.786	2245	2247	2252	rVB3	19259	27732	2.17%	0.236%
84	14.920	2265	2269	2274	rBV5	17520	33653	2.64%	0.286%
85	15.213	2315	2317	2321	rVB	14248	15949	1.25%	0.136%
86	15.420	2348	2351	2357	rVB8	9974	15266	1.20%	0.130%
87	15.493	2357	2363	2369	rBV4	18034	34015	2.67%	0.289%
88	16.383	2505	2509	2514	rVB3	7960	12880	1.01%	0.109%
89	16.603	2541	2545	2553	rVB7	5936	13855	1.09%	0.118%

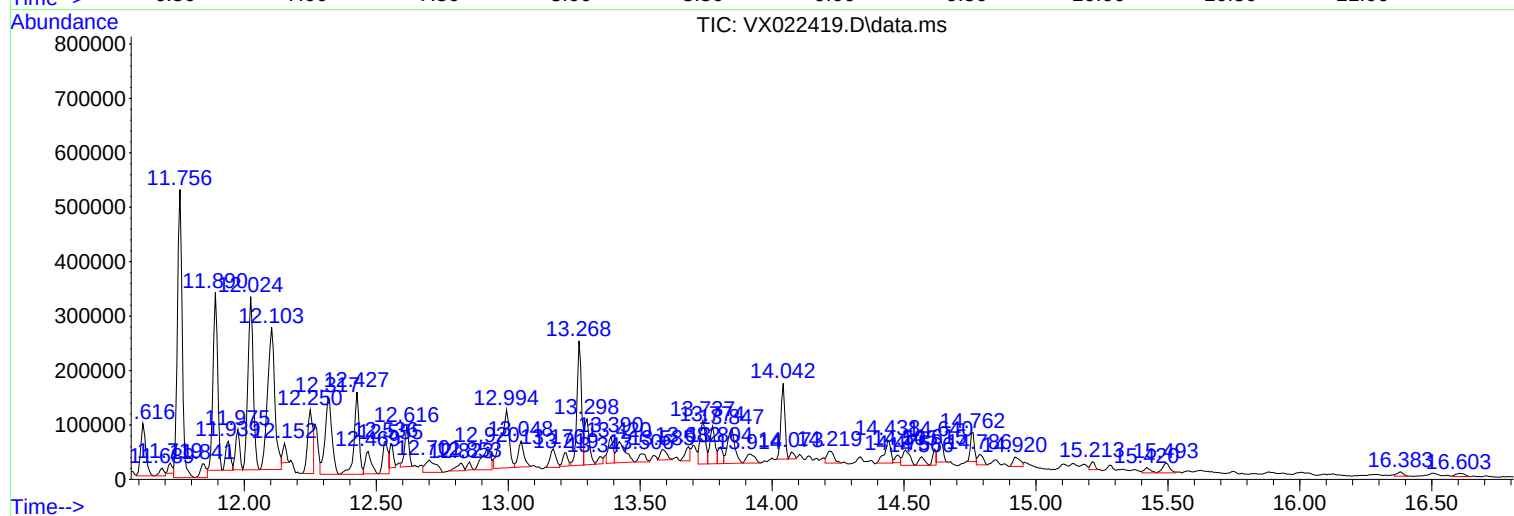
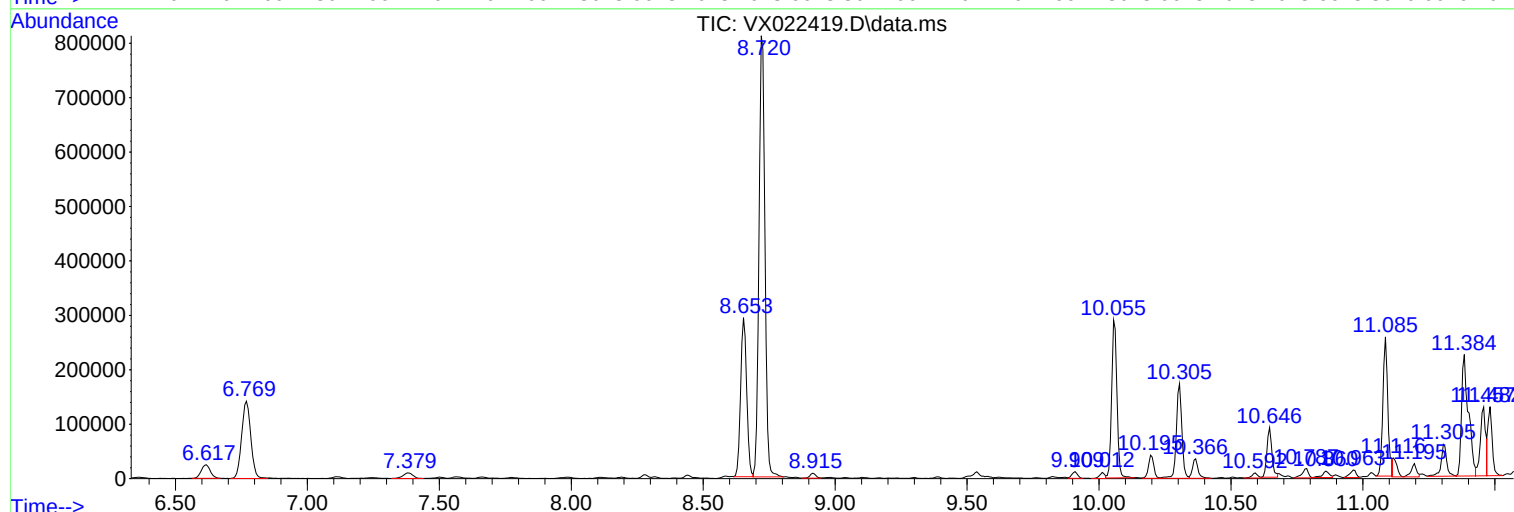
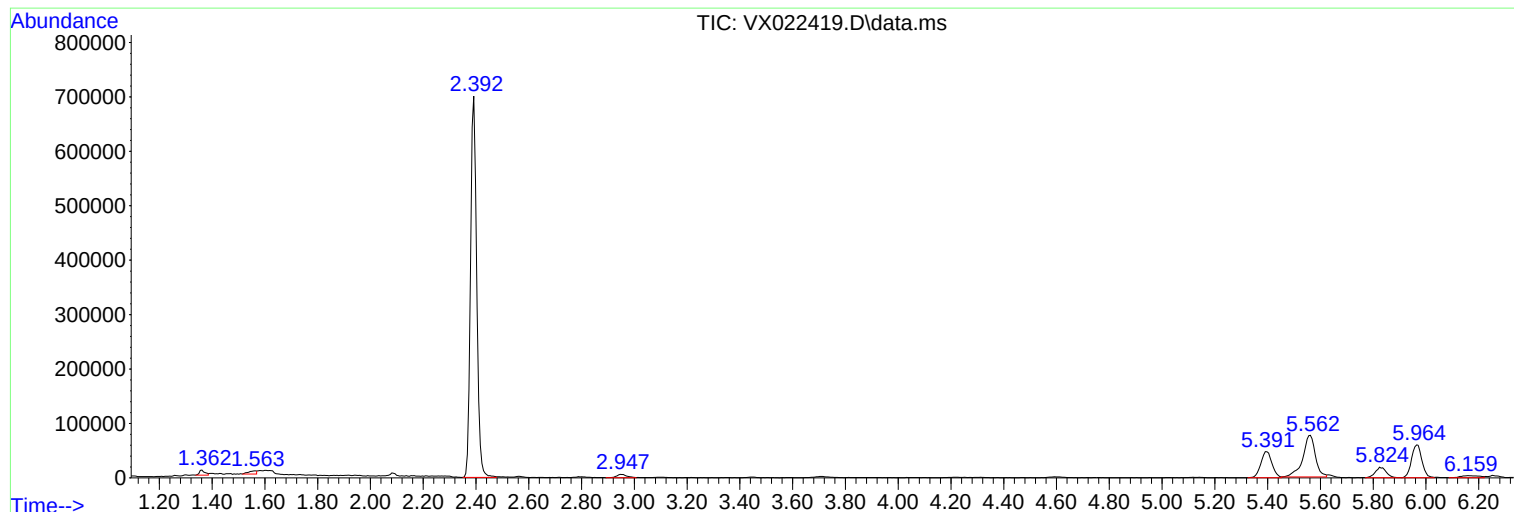
Sum of corrected areas: 11767489

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060421\
Data File : VX022419.D
Acq On : 04 Jun 2021 17:48
Operator : JC/MD
Sample : M2583-04 10X
Misc : 4.99g/10mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
40955

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051921W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060421\
Data File : VX022419.D
Acq On : 04 Jun 2021 17:48
Operator : JC/MD
Sample : M2583-04 10X
Misc : 4.99g/10mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
40955

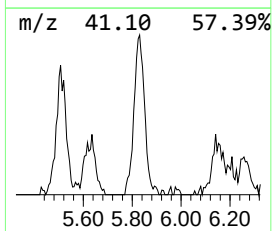
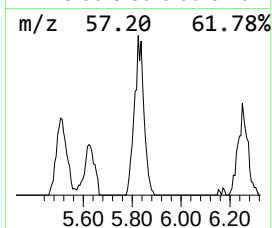
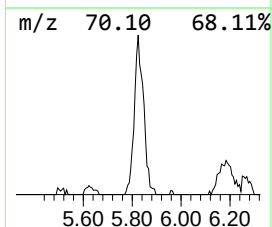
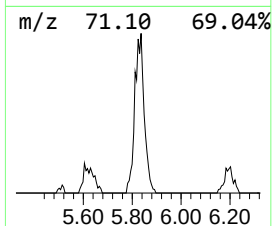
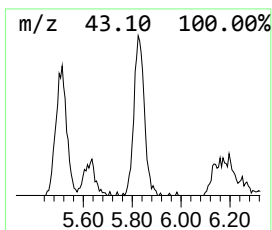
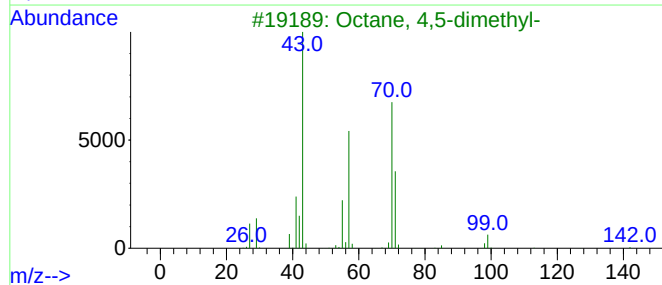
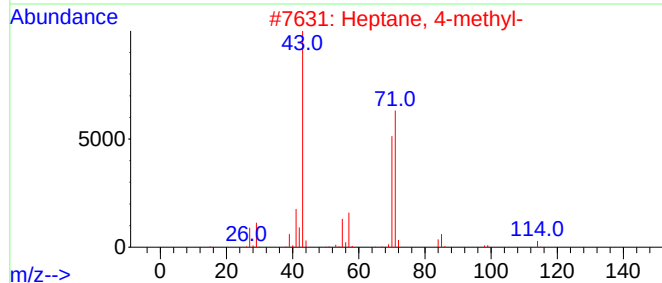
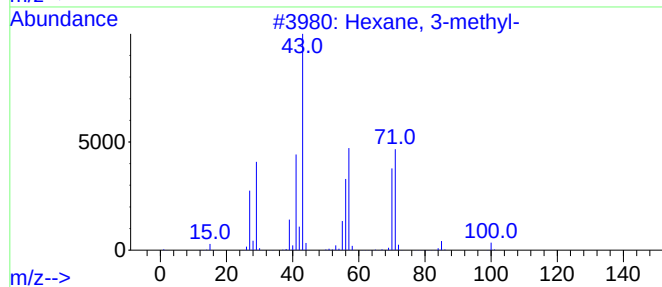
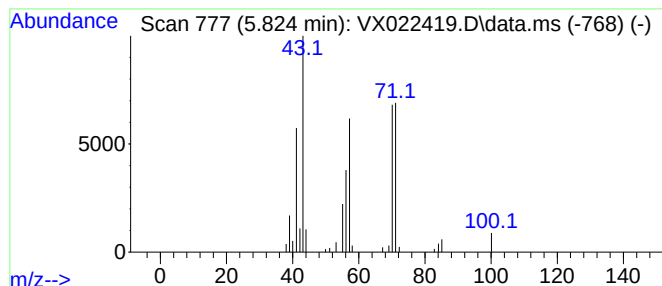
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051921W.M
Quant Title : SW846 8260

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

Peak Number 1 Hexane, 3-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.824	11.00 ug/l	55597	Pentafluorobenzene	5.562

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 3-methyl-	100	C7H16	000589-34-4	59
2			Heptane, 4-methyl-	114	C8H18	000589-53-7	50
3			Octane, 4,5-dimethyl-	142	C10H22	015869-96-2	47
4			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	45
5			Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	45



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060421\
Data File : VX022419.D
Acq On : 04 Jun 2021 17:48
Operator : JC/MD
Sample : M2583-04 10X
Misc : 4.99g/10mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
40955

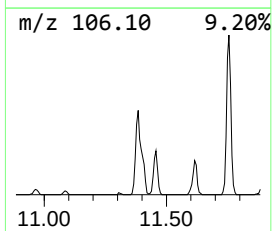
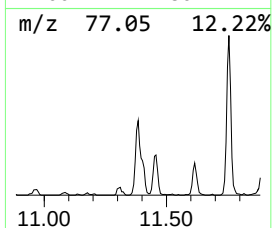
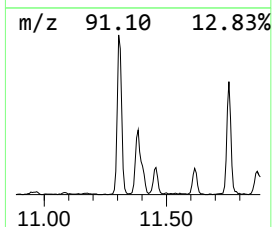
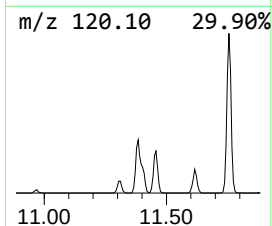
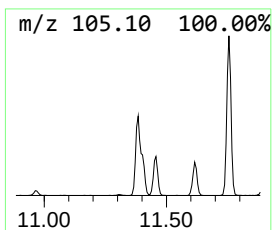
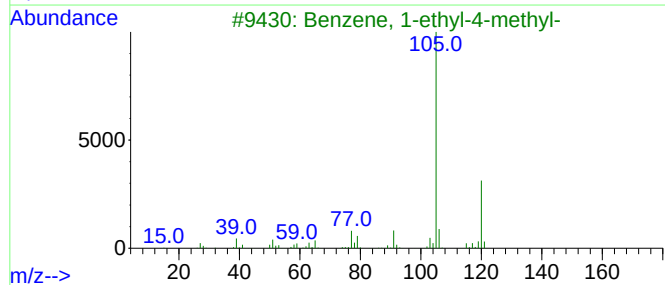
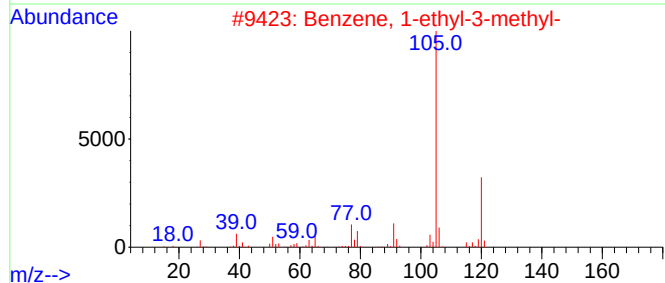
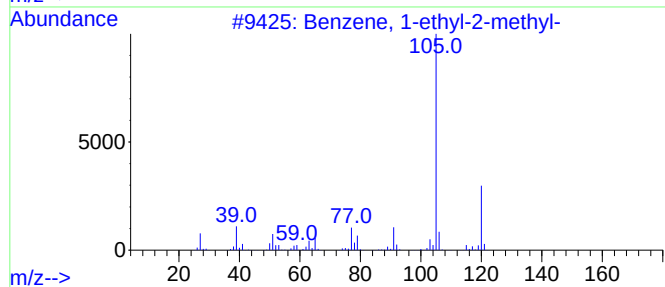
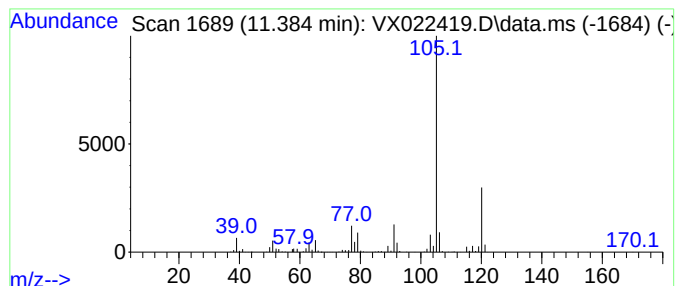
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051921W.M
Quant Title : SW846 8260

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

Peak Number 2 Benzene, 1-ethyl-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.384	45.67 ug/l	407727	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
3			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5			Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060421\
Data File : VX022419.D
Acq On : 04 Jun 2021 17:48
Operator : JC/MD
Sample : M2583-04 10X
Misc : 4.99g/10mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
40955

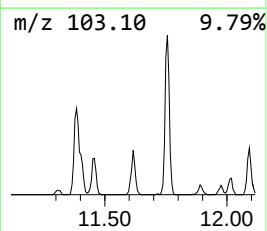
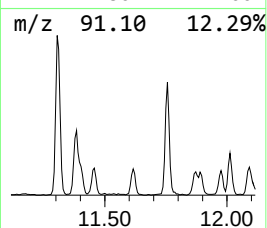
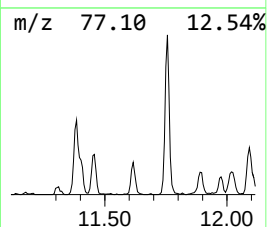
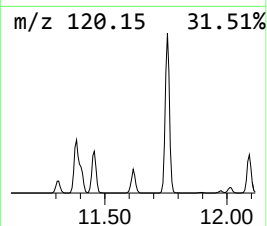
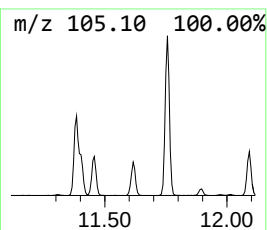
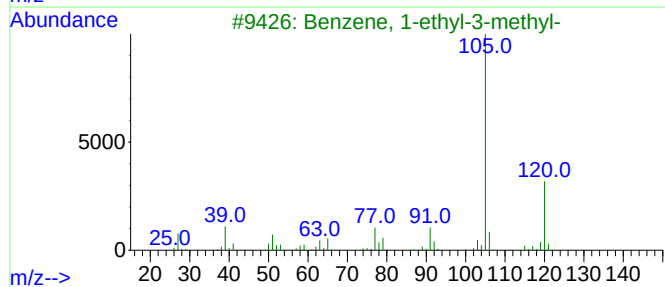
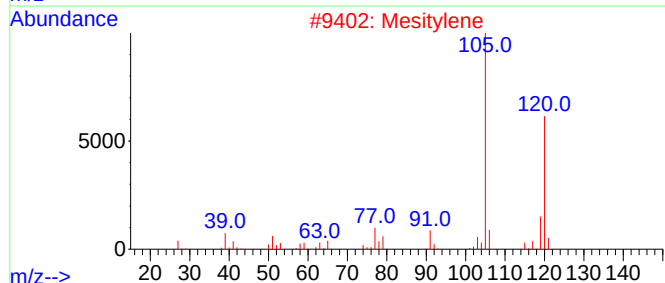
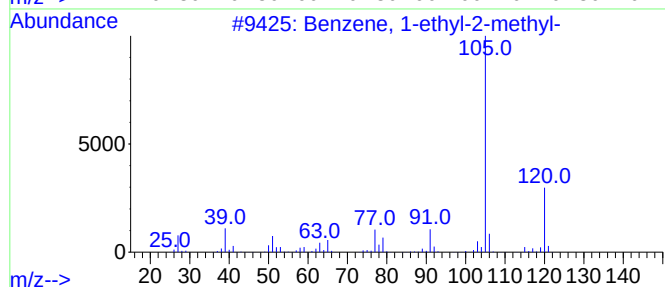
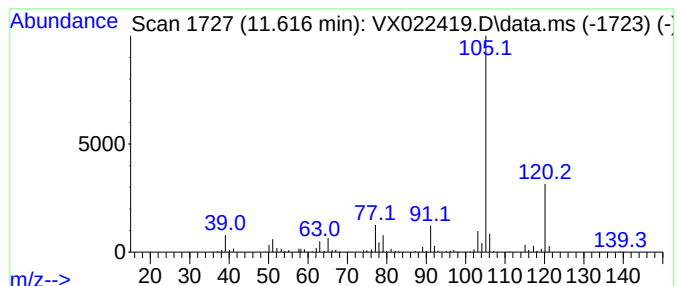
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051921W.M
Quant Title : SW846 8260

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

Peak Number 3 Benzene, 1-ethyl-3-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.616	13.98 ug/l	124789	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2			Mesitylene	120	C9H12	000108-67-8	91
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060421\
Data File : VX022419.D
Acq On : 04 Jun 2021 17:48
Operator : JC/MD
Sample : M2583-04 10X
Misc : 4.99g/10mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
40955

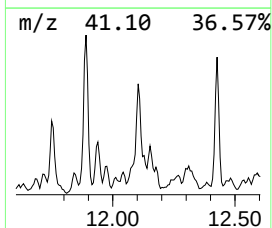
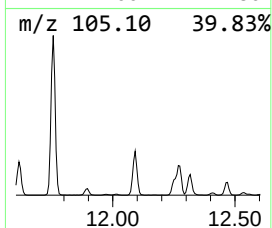
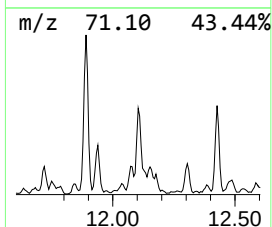
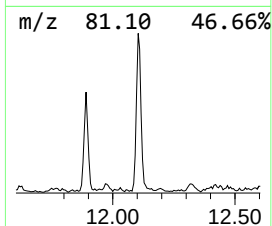
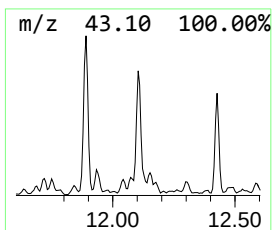
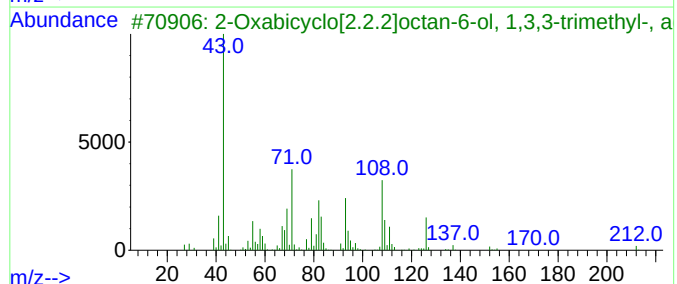
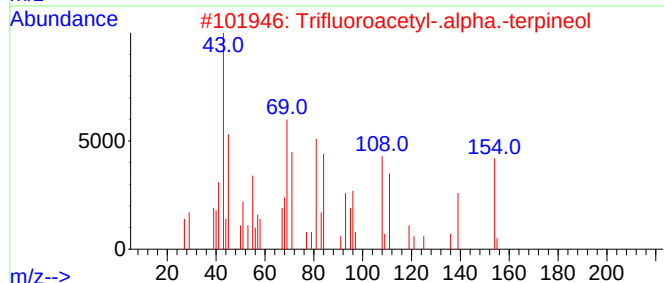
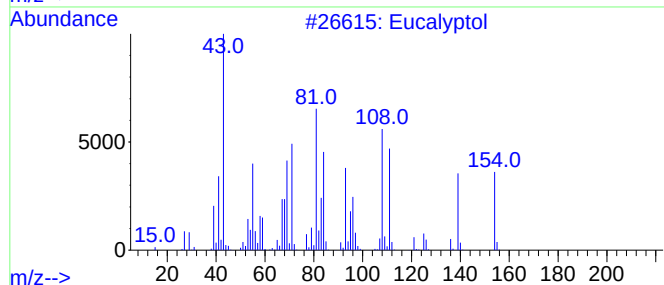
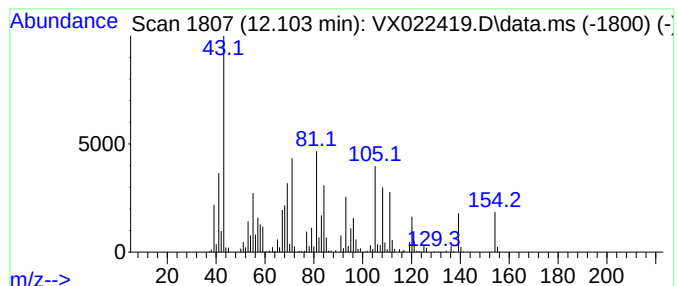
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051921W.M
Quant Title : SW846 8260

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

Peak Number 4 Eucalyptol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.104	56.88 ug/l	507788	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eucalyptol	154	C10H18O	000470-82-6	96
2			Trifluoroacetyl-.alpha.-terpineol	250	C12H17F3O2	1000058-17-6	53
3			2-Oxabicyclo[2.2.2]octan-6-ol, 1...	212	C12H20O3	057709-95-2	43
4			Bicyclo[2.2.1]heptan-2-ol, 2,3,3...	154	C10H18O	000465-31-6	38
5			Ketone, methyl 2,2,3-trimethylcy...	154	C10H18O	017983-22-1	35



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060421\
Data File : VX022419.D
Acq On : 04 Jun 2021 17:48
Operator : JC/MD
Sample : M2583-04 10X
Misc : 4.99g/10mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
40955

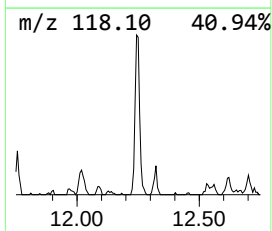
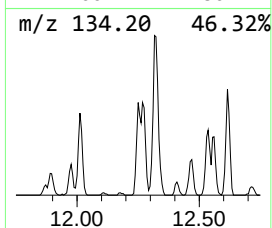
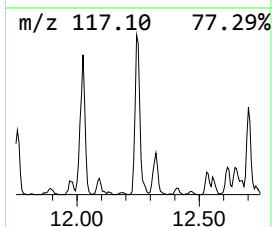
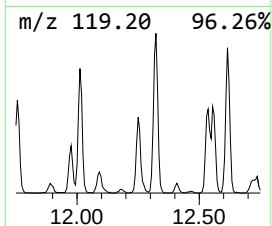
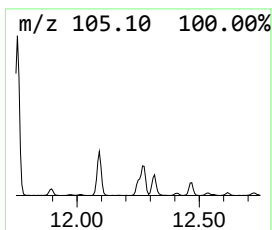
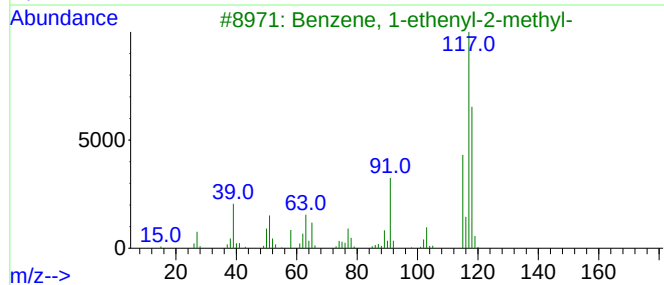
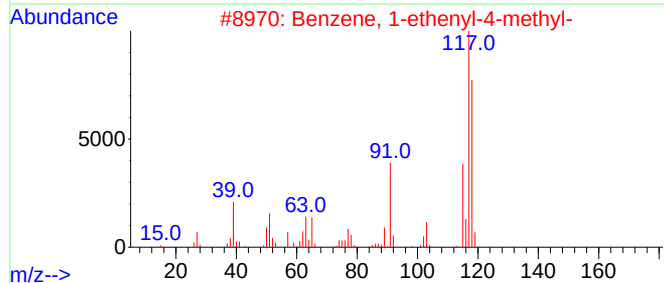
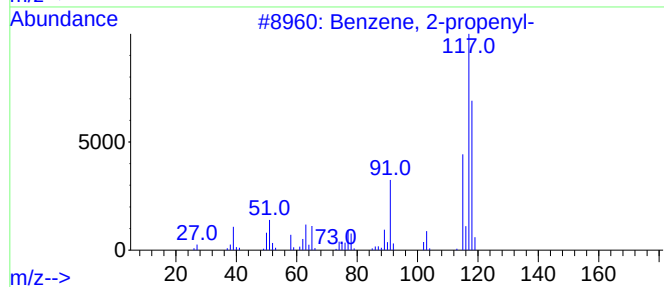
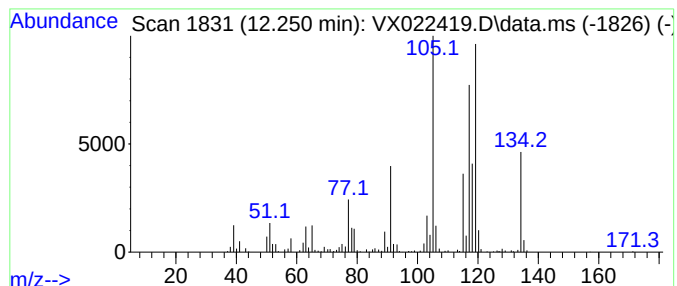
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051921W.M
Quant Title : SW846 8260

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

Peak Number 5 Benzene, 2-propenyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.250	18.59 ug/l	165920	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-propenyl-	118	C9H10	000300-57-2	80
2			Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	80
3			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	80
4			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	78
5			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060421\
Data File : VX022419.D
Acq On : 04 Jun 2021 17:48
Operator : JC/MD
Sample : M2583-04 10X
Misc : 4.99g/10mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
40955

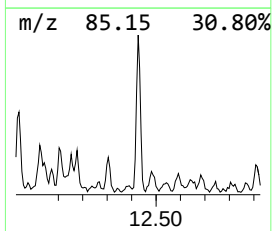
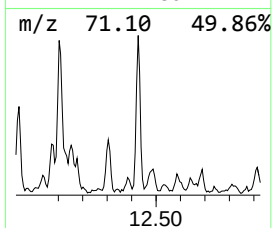
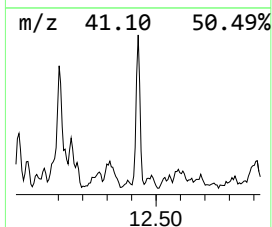
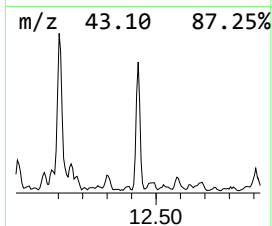
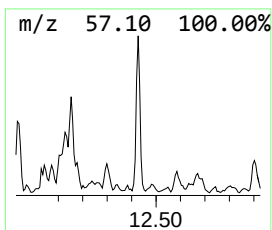
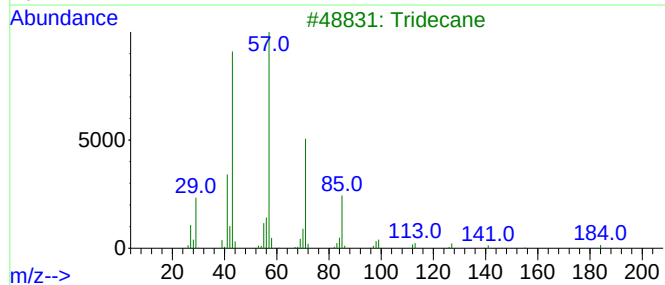
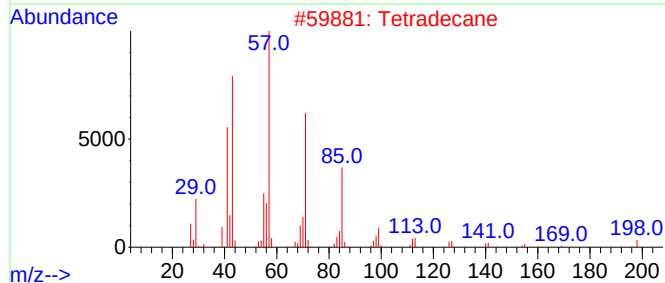
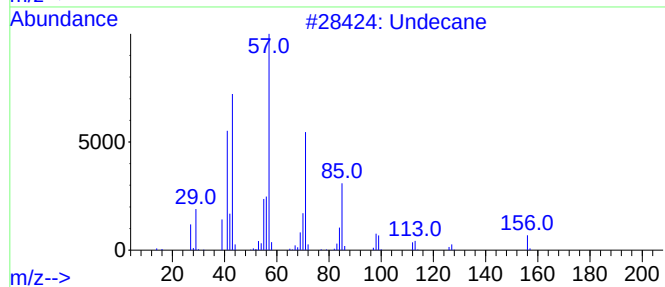
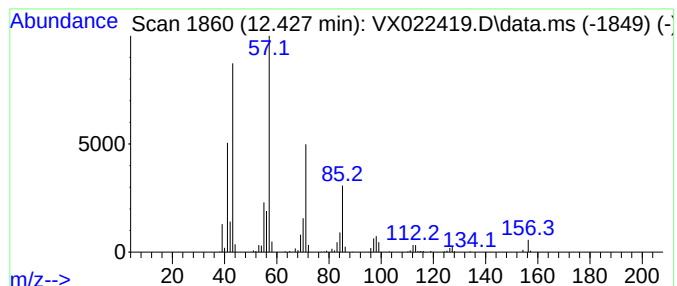
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051921W.M
Quant Title : SW846 8260

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

Peak Number 6 Undecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.427	22.39 ug/l	199894	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane			156	C11H24	001120-21-4	96
2	Tetradecane			198	C14H30	000629-59-4	80
3	Tridecane			184	C13H28	000629-50-5	80
4	Decane, 3,6-dimethyl-			170	C12H26	017312-53-7	64
5	Hexadecane			226	C16H34	000544-76-3	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060421\
Data File : VX022419.D
Acq On : 04 Jun 2021 17:48
Operator : JC/MD
Sample : M2583-04 10X
Misc : 4.99g/10mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
40955

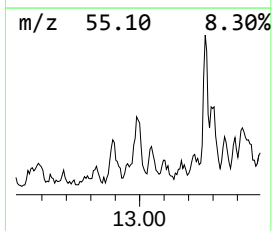
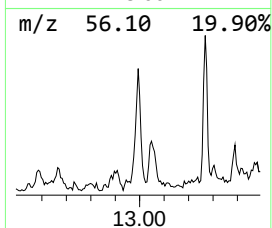
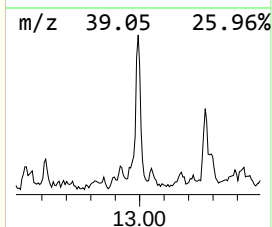
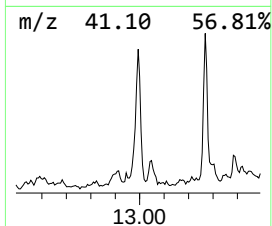
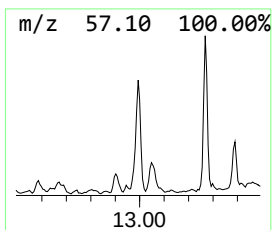
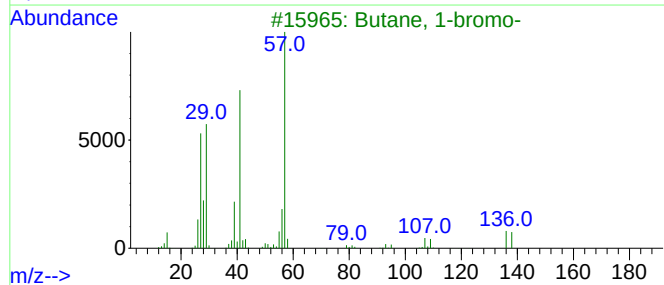
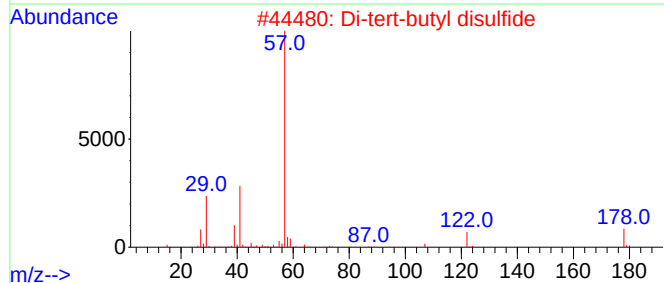
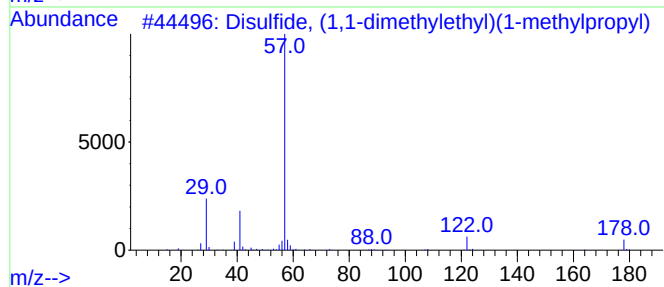
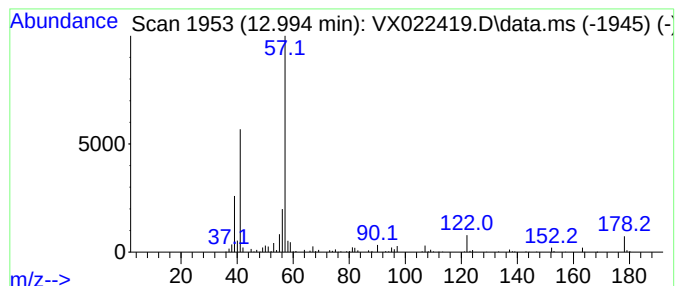
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051921W.M
Quant Title : SW846 8260

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

Peak Number 7 unknown12.994 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.994	26.55 ug/l	237024	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Disulfide, (1,1-dimethylethyl)(1...	178	C8H18S2	072437-46-8	49
2			Di-tert-butyl disulfide	178	C8H18S2	000110-06-5	47
3			Butane, 1-bromo-	136	C4H9Br	000109-65-9	47
4			1-Bromo-3,3-dimethylbutane	164	C6H13Br	001647-23-0	38
5			Oxalic acid, butyl isobutyl ester	202	C10H18O4	1000309-36-9	28



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060421\
Data File : VX022419.D
Acq On : 04 Jun 2021 17:48
Operator : JC/MD
Sample : M2583-04 10X
Misc : 4.99g/10mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
40955

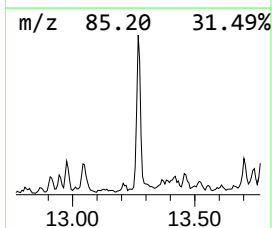
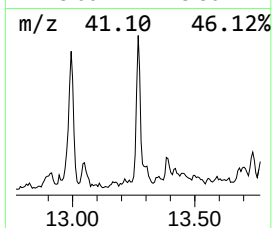
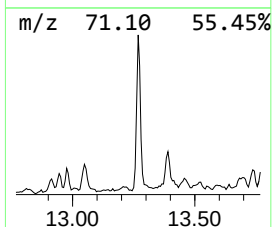
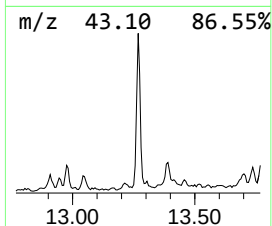
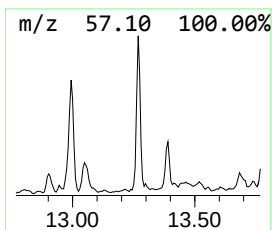
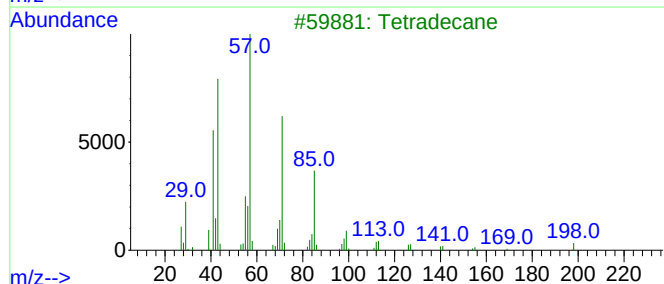
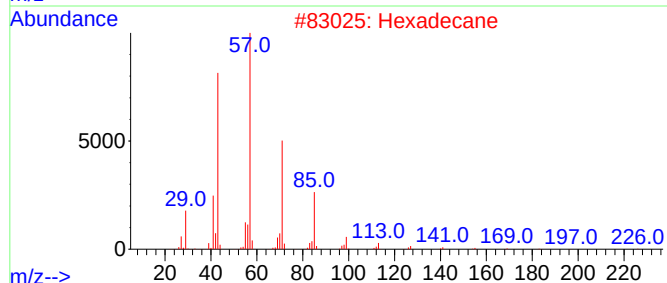
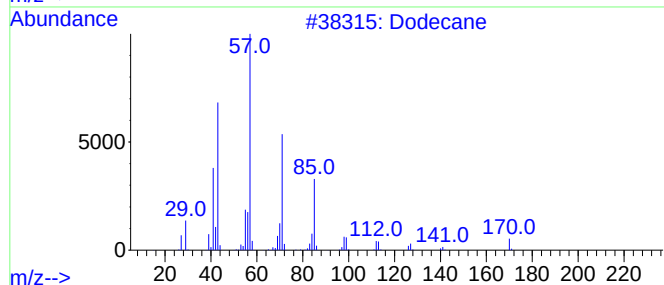
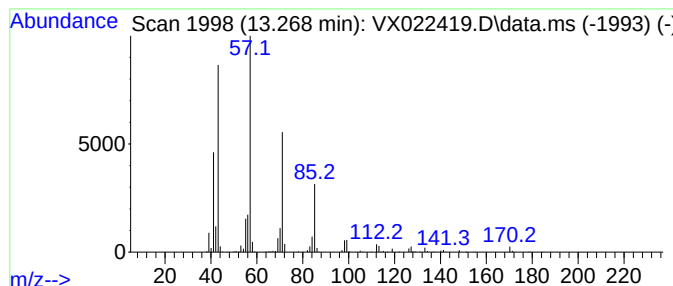
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051921W.M
Quant Title : SW846 8260

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

Peak Number 8 Dodecane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.268	31.53 ug/l	281430	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane	170	C12H26	000112-40-3	96
2			Hexadecane	226	C16H34	000544-76-3	86
3			Tetradecane	198	C14H30	000629-59-4	86
4			Undecane, 2,3-dimethyl-	184	C13H28	017312-77-5	86
5			Tridecane	184	C13H28	000629-50-5	80



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060421\
Data File : VX022419.D
Acq On : 04 Jun 2021 17:48
Operator : JC/MD
Sample : M2583-04 10X
Misc : 4.99g/10mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
40955

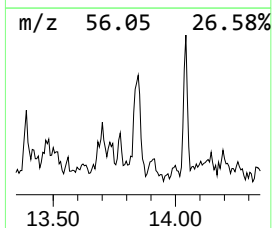
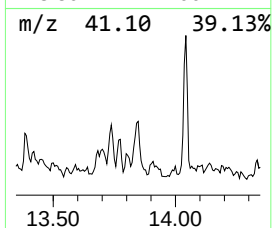
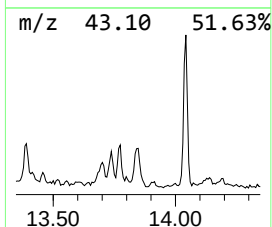
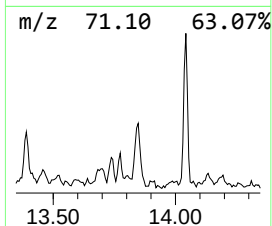
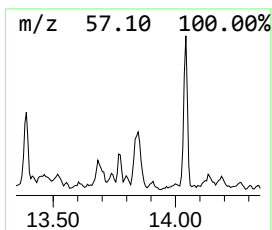
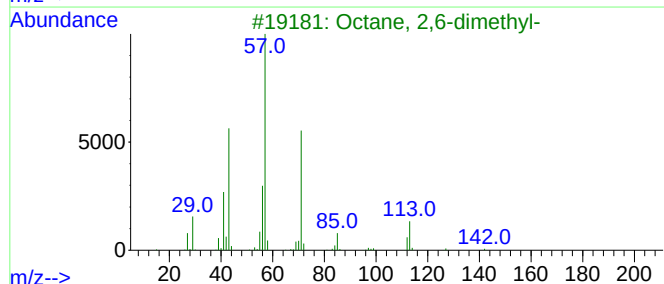
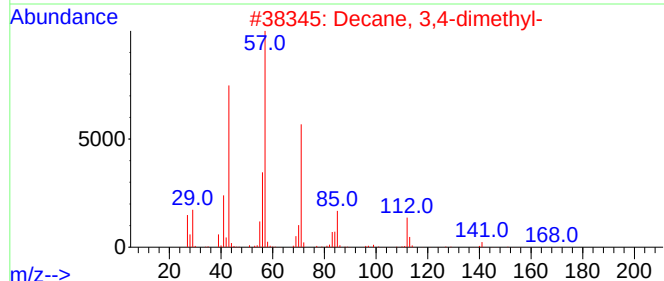
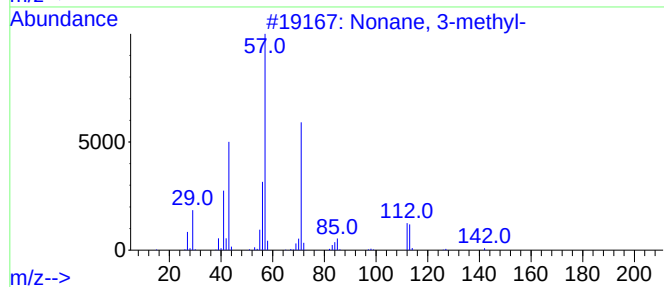
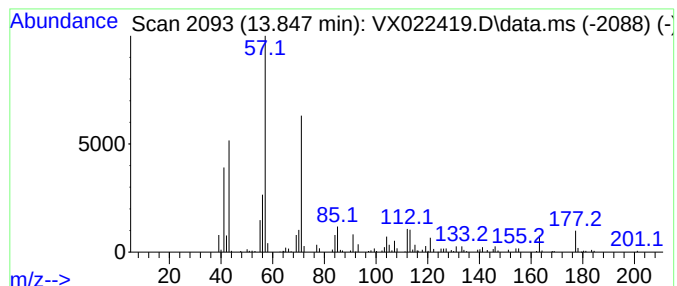
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051921W.M
Quant Title : SW846 8260

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

Peak Number 9 Nonane, 3-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.847	13.58 ug/l	121221	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 3-methyl-	142	C10H22	005911-04-6	72
2			Decane, 3,4-dimethyl-	170	C12H26	017312-45-7	64
3			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	64
4			Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	58
5			Sulfurous acid, decyl 2-ethylhex...	334	C18H38O3S	1000309-19-3	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060421\
Data File : VX022419.D
Acq On : 04 Jun 2021 17:48
Operator : JC/MD
Sample : M2583-04 10X
Misc : 4.99g/10mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
40955

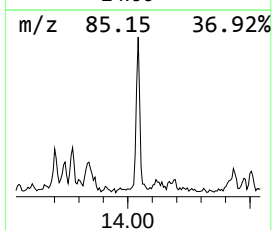
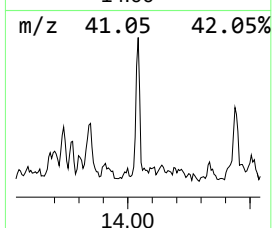
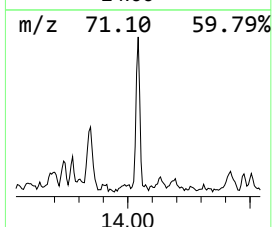
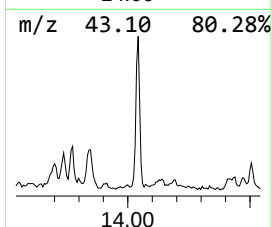
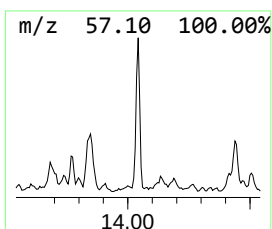
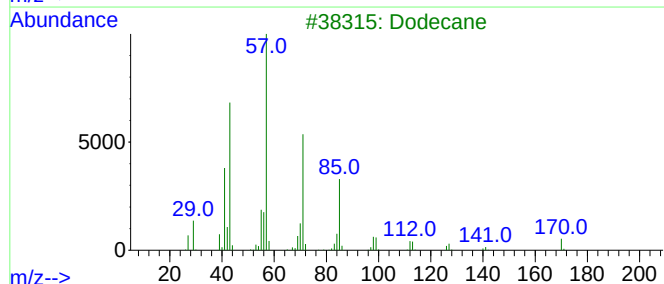
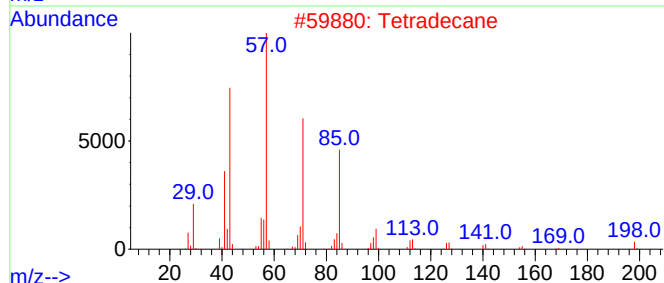
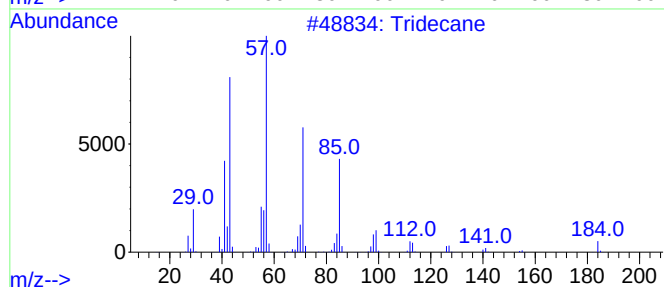
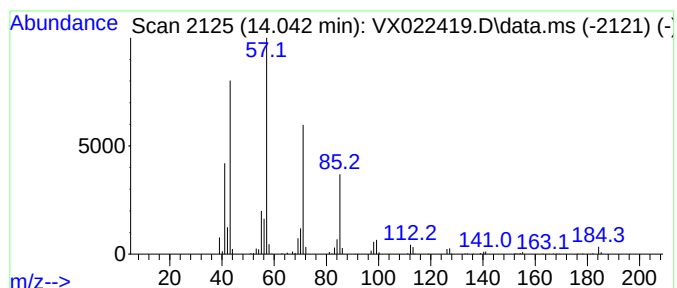
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051921W.M
Quant Title : SW846 8260

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

Peak Number 10 Tridecane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.042	16.07 ug/l	143480	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane	184	C13H28	000629-50-5	94
2			Tetradecane	198	C14H30	000629-59-4	90
3			Dodecane	170	C12H26	000112-40-3	90
4			Hexadecane	226	C16H34	000544-76-3	87
5			Undecane	156	C11H24	001120-21-4	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060421\
 Data File : VX022419.D
 Acq On : 04 Jun 2021 17:48
 Operator : JC/MD
 Sample : M2583-04 10X
 Misc : 4.99g/10mL/100uL/5.0mL/MSVOA_X/MEOH
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 40955

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051921W.M
 Quant Title : SW846 8260

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Hexane, 3-methyl-	5.824	11.0	ug/l	55597	1	5.562	252654	50.0
Benzene, 1-ethy...	11.384	45.7	ug/l	407727	4	12.024	446353	50.0
Benzene, 1-ethy...	11.616	14.0	ug/l	124789	4	12.024	446353	50.0
Eucalyptol	12.104	56.9	ug/l	507788	4	12.024	446353	50.0
Benzene, 2-prop...	12.250	18.6	ug/l	165920	4	12.024	446353	50.0
Undecane	12.427	22.4	ug/l	199894	4	12.024	446353	50.0
unknown12.994	12.994	26.6	ug/l	237024	4	12.024	446353	50.0
Dodecane	13.268	31.5	ug/l	281430	4	12.024	446353	50.0
Nonane, 3-methyl-	13.847	13.6	ug/l	121221	4	12.024	446353	50.0
Tridecane	14.042	16.1	ug/l	143480	4	12.024	446353	50.0