

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX042523\
 Data File : VX035285.D
 Acq On : 25 Apr 2023 15:50
 Operator : JC/MD
 Sample : 02431-01
 Misc : 5.00g/10mL/100uL/5.00mL/MSVOA_X/MEOH
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 AUD-1943

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\82X042523W.M
 Title : SW846 8260

Signal : TIC: VX035285.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.410	208	217	248	rBV	264612	711713	31.69%	1.861%
2	4.196	498	510	519	rBV3	20208	61083	2.72%	0.160%
3	5.111	641	660	674	rBV	35406	134121	5.97%	0.351%
4	5.385	694	705	713	rBV	132425	377654	16.82%	0.987%
5	5.495	713	723	729	rBV	540938	1696486	75.54%	4.436%
6	5.611	737	742	757	rVB2	181809	534614	23.81%	1.398%
7	5.812	763	775	789	rBV	777548	2245714	100.00%	5.872%
8	5.952	789	798	818	rVV	158986	423931	18.88%	1.108%
9	6.178	818	835	842	rVV4	73945	270009	12.02%	0.706%
10	6.245	842	846	854	rVV4	24959	67817	3.02%	0.177%
11	6.348	854	863	876	rVB4	34046	96756	4.31%	0.253%
12	6.604	894	905	920	rBV	192753	468830	20.88%	1.226%
13	6.757	920	930	947	rBV	400662	961475	42.81%	2.514%
14	7.372	1020	1031	1045	rBV2	49002	115913	5.16%	0.303%
15	8.269	1173	1178	1182	rVV	37545	64348	2.87%	0.168%
16	8.598	1225	1232	1235	rBV2	60070	113961	5.07%	0.298%
17	8.647	1235	1240	1246	rVV	872229	1345608	59.92%	3.518%
18	8.720	1246	1252	1258	rVV	602758	952676	42.42%	2.491%
19	8.909	1277	1283	1289	rBV	109304	157279	7.00%	0.411%
20	8.976	1289	1294	1303	rVB2	43989	76688	3.41%	0.201%
21	9.098	1305	1314	1320	rBV	41347	69644	3.10%	0.182%
22	9.384	1351	1361	1369	rBV2	86160	149917	6.68%	0.392%
23	9.506	1375	1381	1386	rBV4	94451	210595	9.38%	0.551%
24	9.561	1386	1390	1394	rVB	96061	135693	6.04%	0.355%
25	9.616	1394	1399	1403	rBV	104501	167612	7.46%	0.438%
26	9.756	1416	1422	1427	rBV	40608	62113	2.77%	0.162%
27	9.829	1427	1434	1438	rBV3	136810	286311	12.75%	0.749%
28	9.903	1442	1446	1451	rVB	244930	353894	15.76%	0.925%
29	10.006	1456	1463	1467	rBV	251315	370127	16.48%	0.968%
30	10.055	1467	1471	1477	rVV	785668	1073981	47.82%	2.808%
31	10.195	1486	1494	1498	rBV2	177672	306011	13.63%	0.800%
32	10.299	1498	1511	1517	rBV2	494019	1135446	50.56%	2.969%
33	10.360	1517	1521	1531	rVB	837240	1120750	49.91%	2.930%
34	10.451	1531	1536	1542	rBV	63351	96815	4.31%	0.253%
35	10.585	1551	1558	1562	rBV2	208479	335204	14.93%	0.876%
36	10.640	1562	1567	1570	rBV	193675	269184	11.99%	0.704%

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37	10.671	1570	1572	1575	rVB	85190	83493	3.72%	0.218%
38	10.780	1582	1590	1594	rBV2	358785	546452	24.33%	1.429%
39	10.854	1598	1602	1606	rVV3	282617	492139	21.91%	1.287%
40	10.896	1606	1609	1613	rVV	153408	223751	9.96%	0.585%
41	10.951	1613	1618	1623	rVV3	106665	231268	10.30%	0.605%
42	11.000	1623	1626	1628	rVV2	71405	90478	4.03%	0.237%
43	11.030	1628	1631	1634	rVV2	104378	135392	6.03%	0.354%
44	11.085	1634	1640	1650	rVB2	837562	1494735	66.56%	3.908%
45	11.189	1650	1657	1660	rBV2	227022	353984	15.76%	0.926%
46	11.213	1660	1661	1669	rVB5	116216	132657	5.91%	0.347%
47	11.305	1672	1676	1679	rVV	146013	191205	8.51%	0.500%
48	11.335	1679	1681	1684	rVV3	66772	94045	4.19%	0.246%
49	11.378	1684	1688	1694	rVV	442080	836246	37.24%	2.186%
50	11.433	1694	1697	1698	rVV	247322	299848	13.35%	0.784%
51	11.451	1698	1700	1701	rVV	251260	255153	11.36%	0.667%
52	11.475	1701	1704	1709	rVB	939500	1143893	50.94%	2.991%
53	11.616	1722	1727	1734	rVB3	268742	474303	21.12%	1.240%
54	11.683	1734	1738	1740	rBV2	70050	85831	3.82%	0.224%
55	11.713	1740	1743	1746	rBV	183160	209522	9.33%	0.548%
56	11.756	1746	1750	1759	rVB	688456	1035334	46.10%	2.707%
57	11.835	1759	1763	1765	rBV3	59444	80574	3.59%	0.211%
58	11.890	1765	1772	1777	rVB2	168177	275172	12.25%	0.719%
59	11.945	1777	1781	1783	rBV	177706	220477	9.82%	0.576%
60	11.969	1783	1785	1789	rVB2	130760	152885	6.81%	0.400%
61	12.024	1789	1794	1800	rBV	824996	1196743	53.29%	3.129%
62	12.091	1800	1805	1810	rVB3	230130	444134	19.78%	1.161%
63	12.170	1815	1818	1824	rVB3	110361	168988	7.52%	0.442%
64	12.250	1824	1831	1832	rBV4	273027	389637	17.35%	1.019%
65	12.268	1832	1834	1838	rVV	306325	366603	16.32%	0.959%
66	12.317	1838	1842	1849	rVB3	512243	888502	39.56%	2.323%
67	12.420	1849	1859	1863	rBV	653408	908361	40.45%	2.375%
68	12.463	1863	1866	1872	rVB	229433	327425	14.58%	0.856%
69	12.536	1873	1878	1879	rBV	198465	257908	11.48%	0.674%
70	12.554	1879	1881	1884	rVV	224725	283873	12.64%	0.742%
71	12.591	1884	1887	1888	rVV3	129608	172414	7.68%	0.451%
72	12.615	1888	1891	1894	rVV	304478	377408	16.81%	0.987%
73	12.646	1894	1896	1899	rVB2	70271	72541	3.23%	0.190%
74	12.725	1899	1909	1913	rBV4	178581	507223	22.59%	1.326%
75	12.817	1917	1924	1927	rVV4	147044	253344	11.28%	0.662%
76	12.847	1927	1929	1932	rVB2	70082	74640	3.32%	0.195%

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77	12.890	1932	1936	1938	rBV3	100527	146432	6.52%	0.383%
78	12.920	1938	1941	1945	rVV	162253	236768	10.54%	0.619%
79	12.963	1945	1948	1955	rVB	282449	472233	21.03%	1.235%
80	13.042	1955	1961	1964	rBV4	153656	296855	13.22%	0.776%
81	13.164	1975	1981	1986	rBV3	204958	346708	15.44%	0.906%
82	13.213	1986	1989	1992	rBV2	134041	151120	6.73%	0.395%
83	13.268	1992	1998	1999	rVV3	254806	360647	16.06%	0.943%
84	13.292	1999	2002	2007	rVV2	451316	625623	27.86%	1.636%
85	13.371	2012	2015	2019	rBV2	57505	73532	3.27%	0.192%
86	13.420	2019	2023	2032	rVB3	270157	490335	21.83%	1.282%
87	13.499	2032	2036	2040	rBV2	86470	133049	5.92%	0.348%
88	13.542	2040	2043	2046	rBV	76566	94191	4.19%	0.246%
89	13.585	2046	2050	2055	rBV3	135648	240807	10.72%	0.630%
90	13.670	2061	2064	2068	rVB3	71157	108292	4.82%	0.283%
91	13.853	2088	2094	2100	rVB5	129487	252686	11.25%	0.661%
92	13.926	2100	2106	2111	rBV5	115997	247413	11.02%	0.647%
93	14.036	2121	2124	2127	rVB3	87687	94183	4.19%	0.246%
94	14.073	2127	2130	2134	rBV2	72780	96338	4.29%	0.252%
95	14.225	2150	2155	2160	rVV3	196118	353647	15.75%	0.925%
96	14.371	2176	2179	2183	rVB	57125	71050	3.16%	0.186%
97	14.481	2193	2197	2205	rVB5	106736	212965	9.48%	0.557%
98	14.615	2216	2219	2221	rBV	121579	148795	6.63%	0.389%
99	14.670	2226	2228	2232	rVB3	55858	63265	2.82%	0.165%
100	15.487	2358	2362	2367	rBV2	52207	81644	3.64%	0.213%

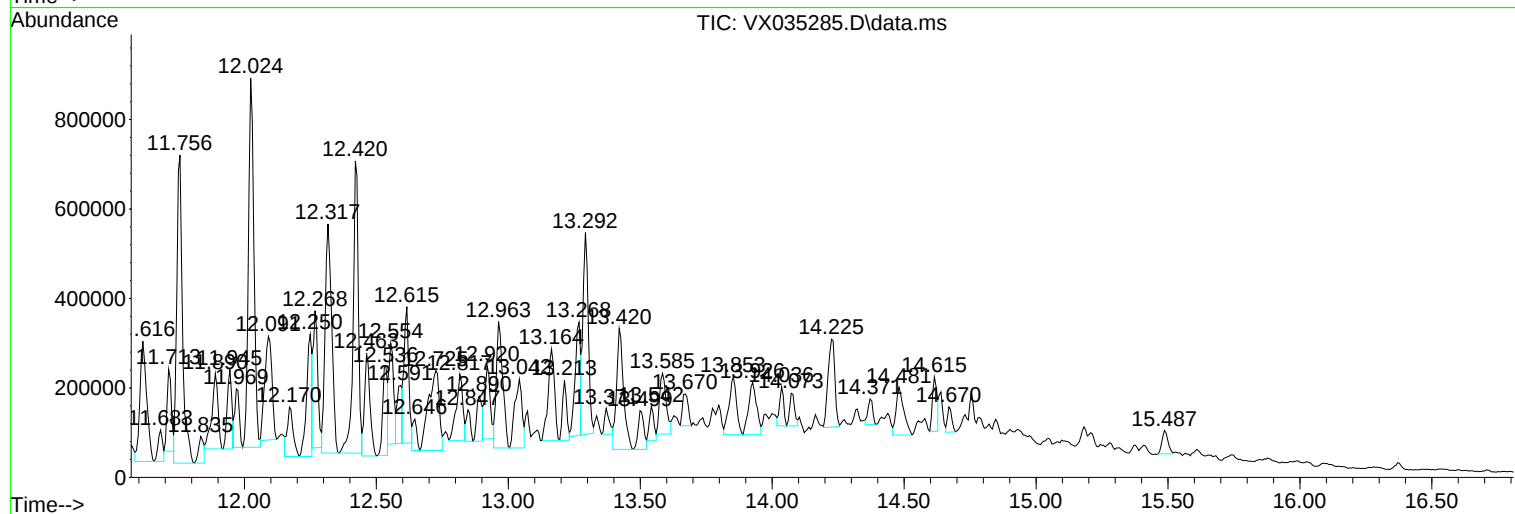
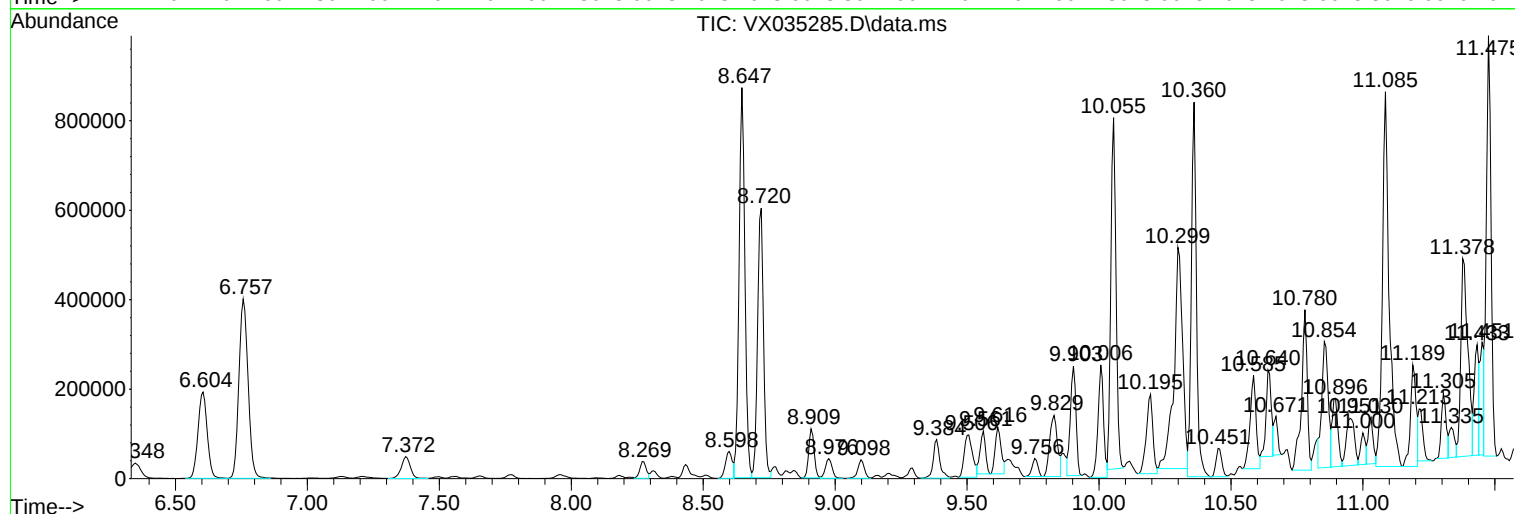
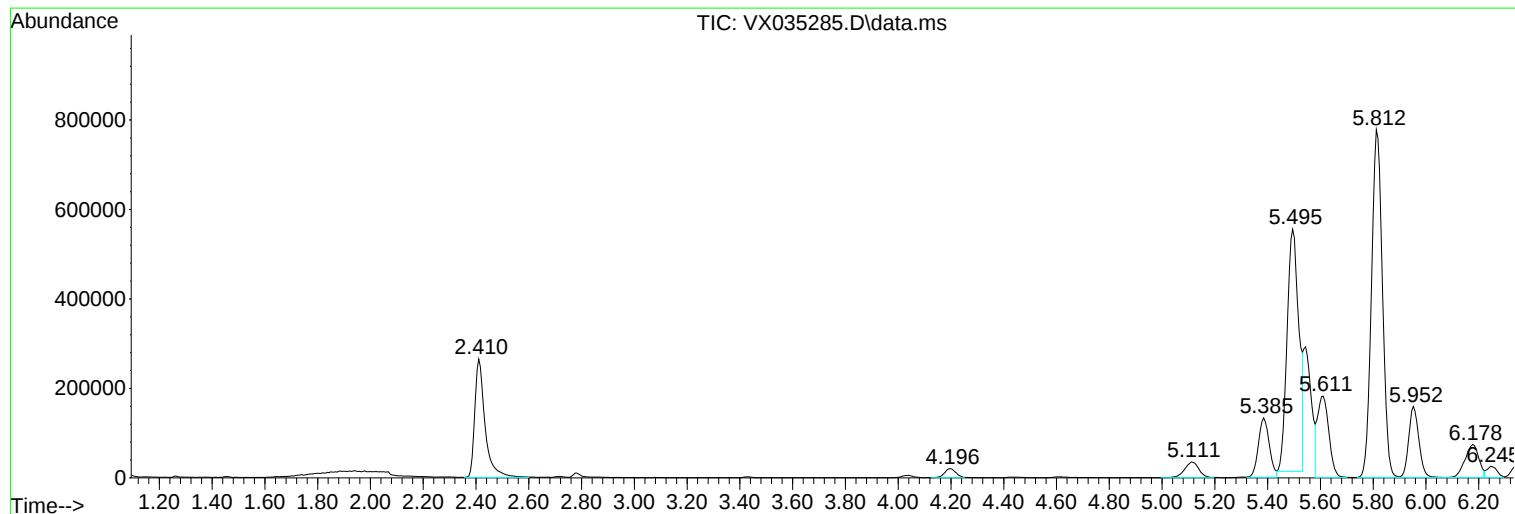
Sum of corrected areas: 38247132

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Instrument :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X042523W.M
Quant Title : SW846 8260

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



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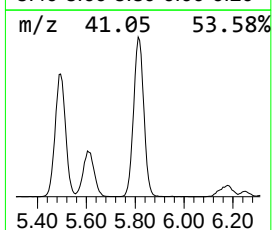
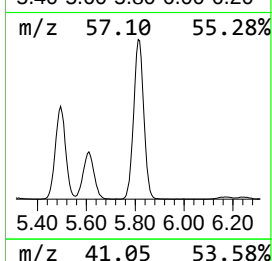
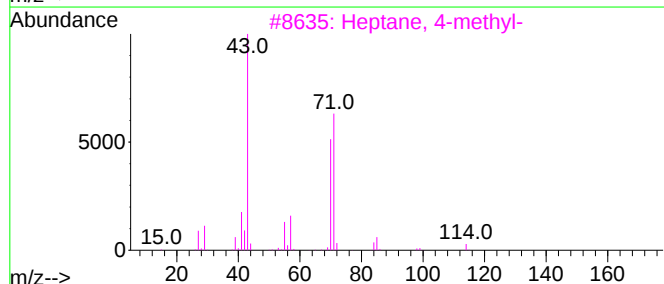
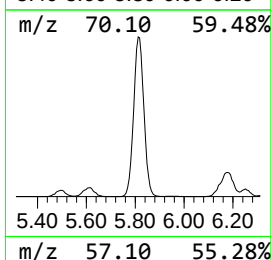
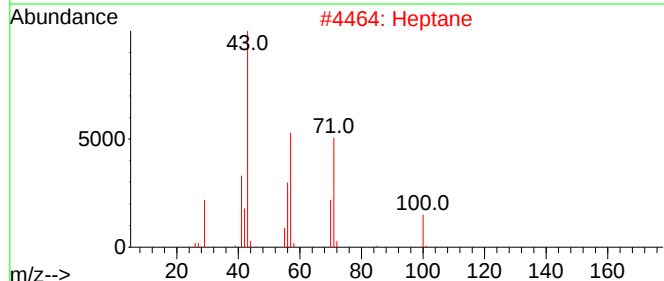
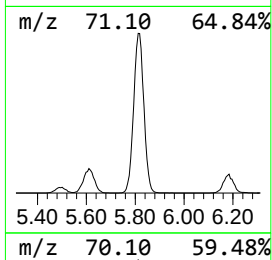
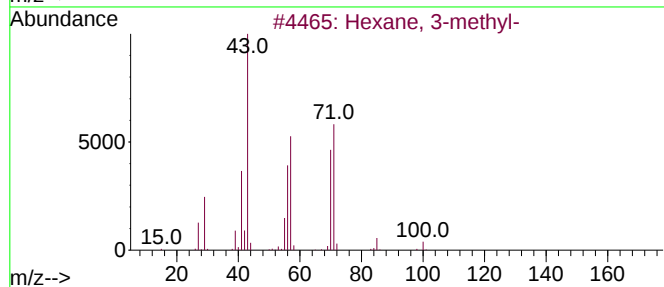
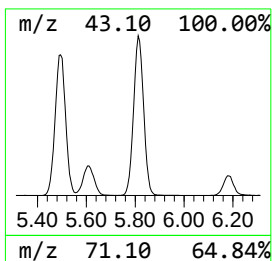
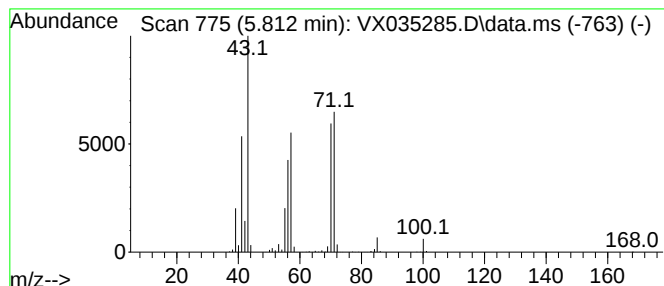
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X042523W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.812	66.19 ug/l	2245710	Pentafluorobenzene	5.544

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 3-methyl-	100	C7H16	000589-34-4	94
2			Heptane	100	C7H16	000142-82-5	80
3			Heptane, 4-methyl-	114	C8H18	000589-53-7	59
4			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	53
5			Heptane, 3,3,4-trimethyl-	142	C10H22	020278-87-9	53



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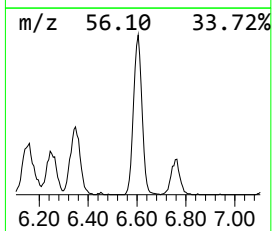
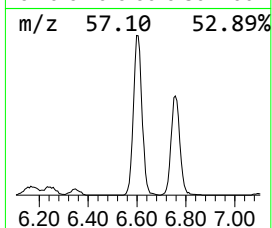
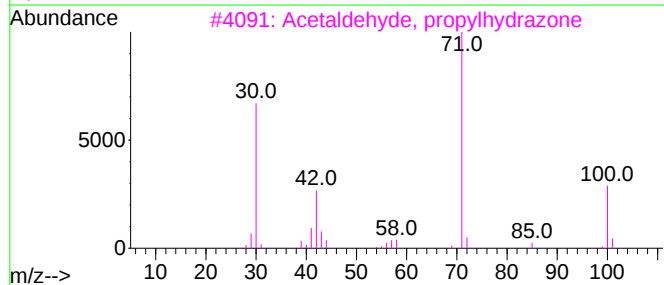
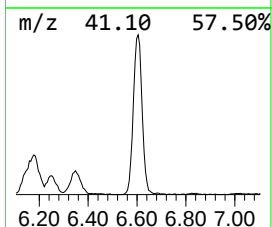
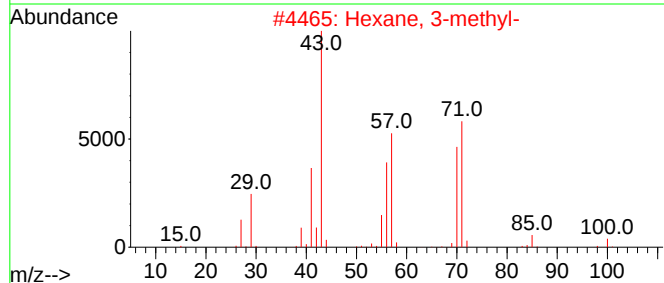
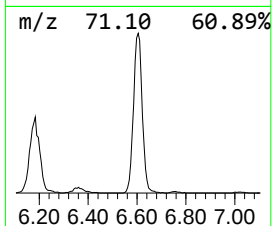
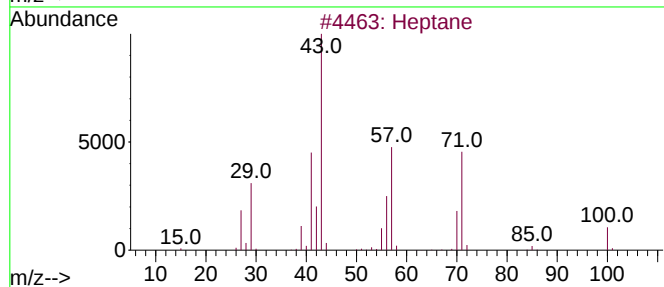
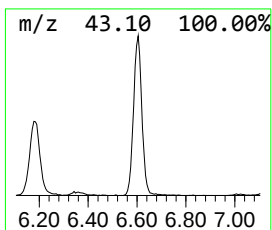
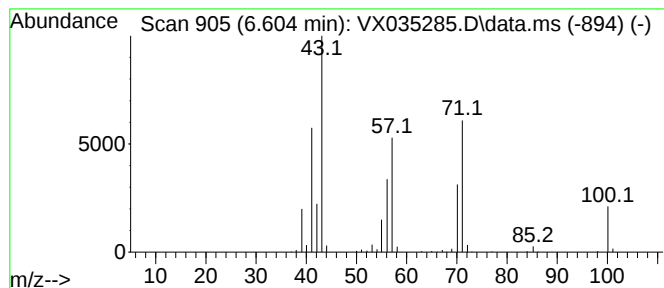
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Quant Title : SW846 8260

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

Peak Number 2 Heptane Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.604	24.38 ug/l	468830	1,4-Difluorobenzene	6.757

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane	100	C7H16	000142-82-5	91
2			Hexane, 3-methyl-	100	C7H16	000589-34-4	59
3			Acetaldehyde, propylhydrazone	100	C5H12N2	007422-88-0	49
4			3-Hexanone	100	C6H12O	000589-38-8	46
5			Pentane, 2-methyl-	86	C6H14	000107-83-5	43



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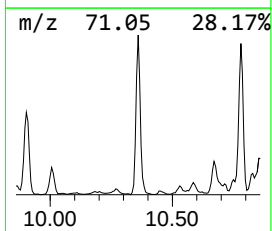
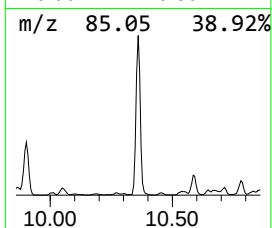
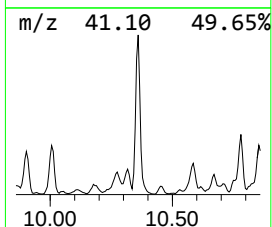
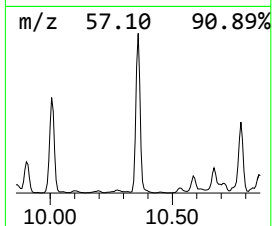
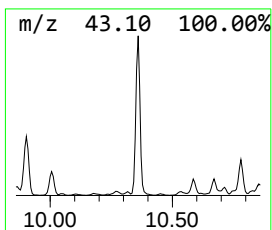
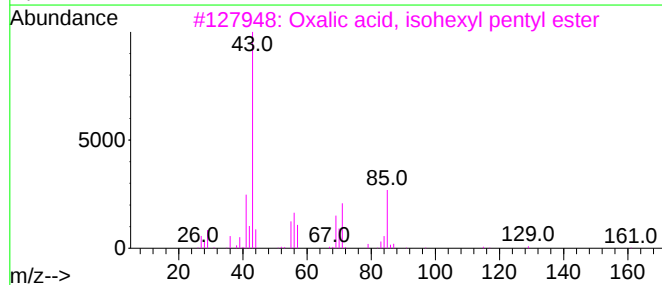
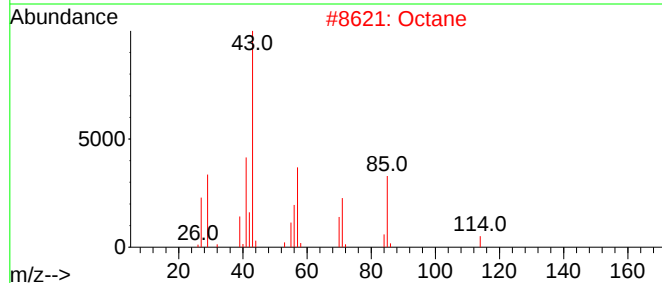
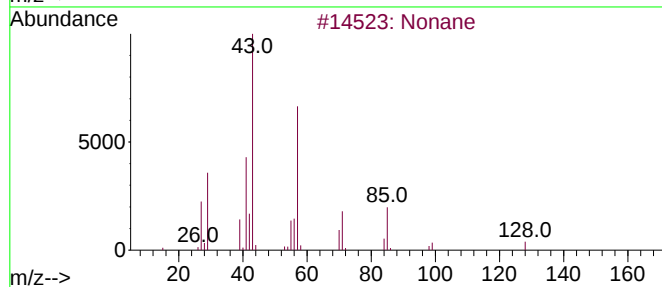
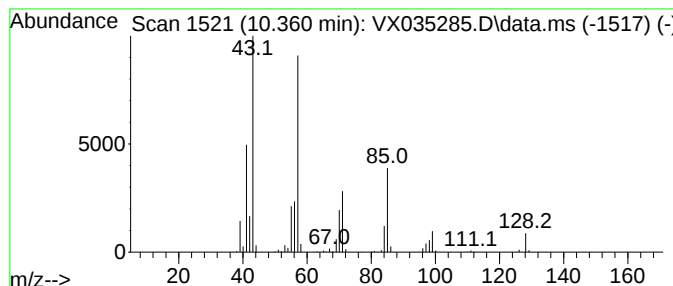
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Quant Title : SW846 8260

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

Peak Number 3 Nonane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.360	52.18 ug/l	1120750	Chlorobenzene-d5	10.055

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane	128	C9H20	000111-84-2	95
2			Octane	114	C8H18	000111-65-9	59
3			Oxalic acid, isohexyl pentyl ester	244	C13H24O4	1000309-32-8	53
4			Sulfurous acid, hexyl pentadecyl...	376	C21H44O3S	1000309-13-7	53
5			4,4-Dimethyl octane	142	C10H22	015869-95-1	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX042523\
Data File : VX035285.D
Acq On : 25 Apr 2023 15:50
Operator : JC/MD
Sample : 02431-01
Misc : 5.00g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
AUD-1943

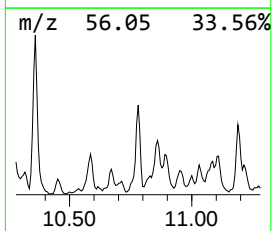
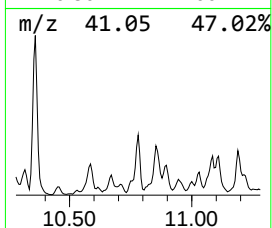
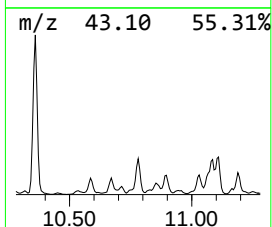
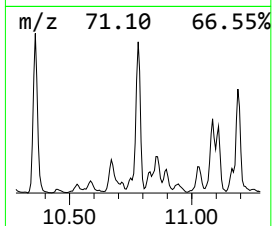
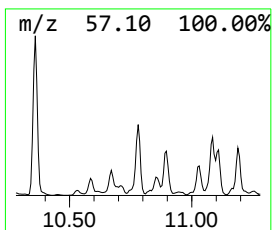
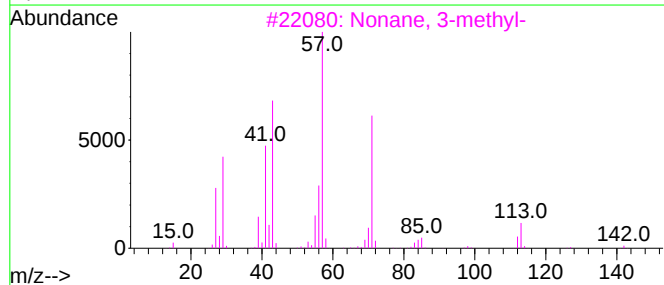
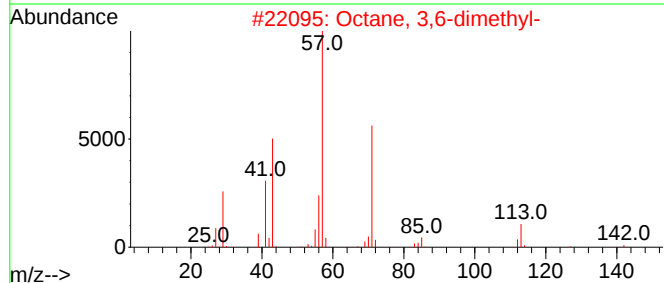
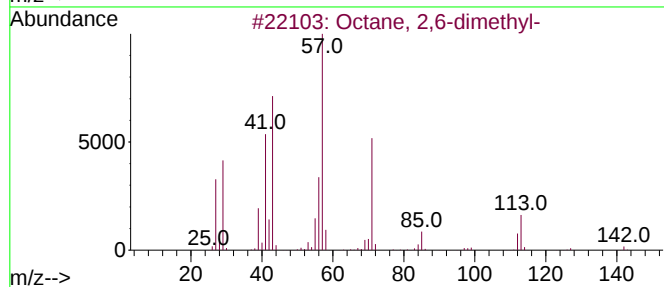
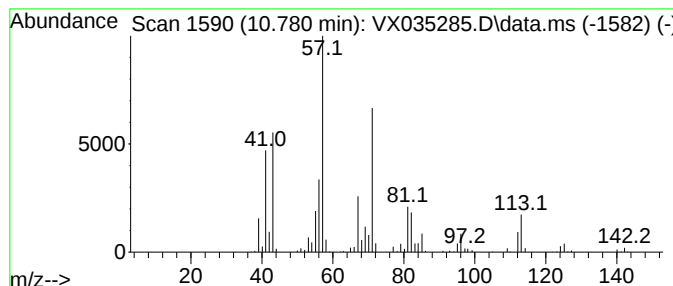
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Quant Title : SW846 8260

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

Peak Number 4 Octane, 2,6-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.780	25.44 ug/l	546452	Chlorobenzene-d5	10.055

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	76
2			Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	70
3			Nonane, 3-methyl-	142	C10H22	005911-04-6	70
4			Undecane, 6,6-dimethyl-	184	C13H28	017312-76-4	50
5			Sulfurous acid, 2-ethylhexyl hex...	278	C14H30O3S	1000309-20-2	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX042523\
Data File : VX035285.D
Acq On : 25 Apr 2023 15:50
Operator : JC/MD
Sample : 02431-01
Misc : 5.00g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
AUD-1943

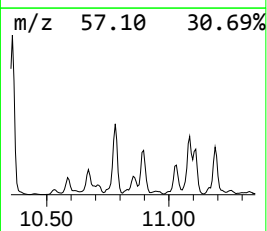
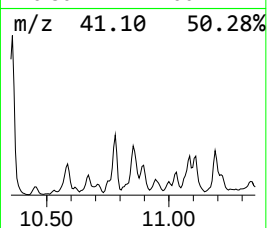
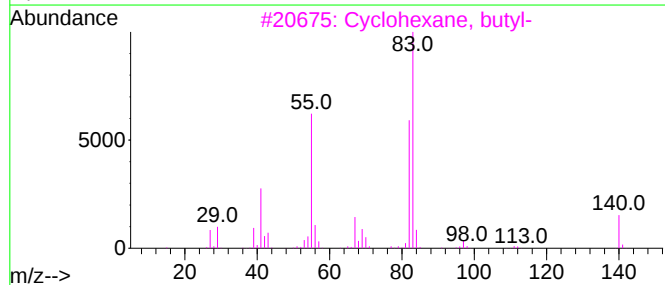
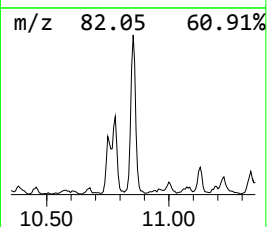
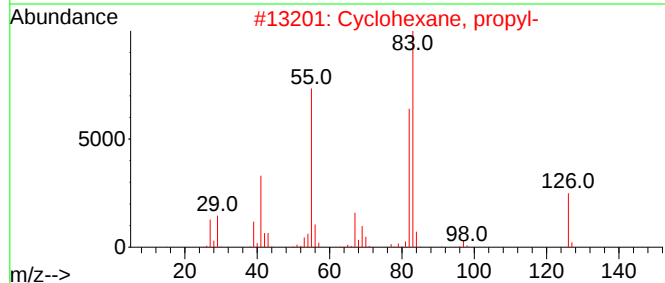
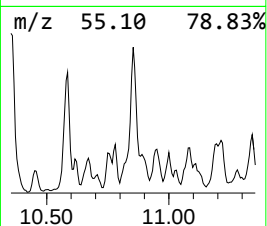
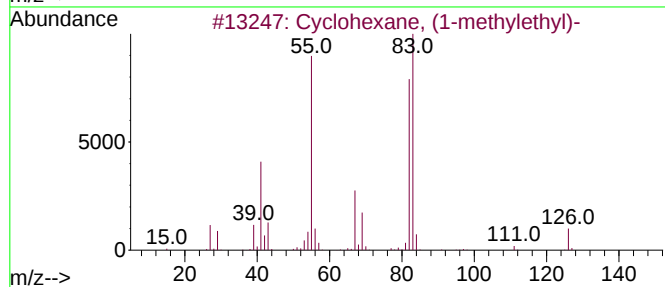
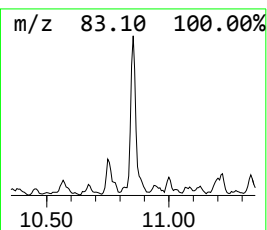
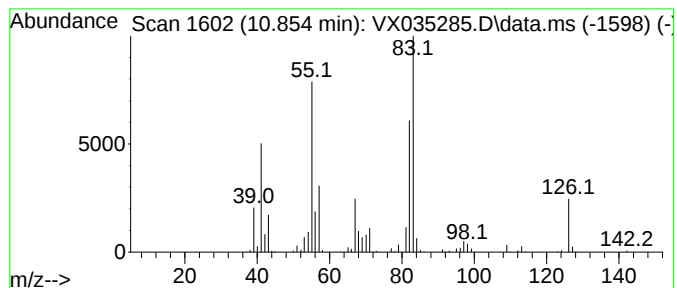
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Quant Title : SW846 8260

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

Peak Number 5 Cyclohexane, (1-methylethyl)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.854	22.91 ug/l	492139	Chlorobenzene-d5	10.055

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	87
2			Cyclohexane, propyl-	126	C9H18	001678-92-8	81
3			Cyclohexane, butyl-	140	C10H20	001678-93-9	53
4			Trichloroacetic acid, 6-chlorohe...	280	C8H12Cl4O2	1000330-89-2	53
5			trans-7-Oxabicyclo[4.3.0]nonane	126	C8H14O	027345-70-6	52



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX042523\
Data File : VX035285.D
Acq On : 25 Apr 2023 15:50
Operator : JC/MD
Sample : 02431-01
Misc : 5.00g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
AUD-1943

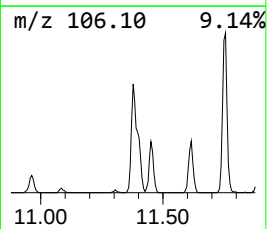
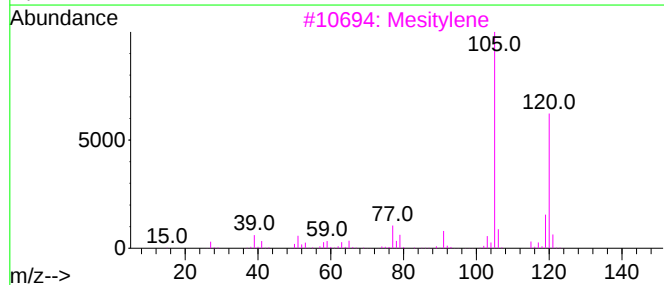
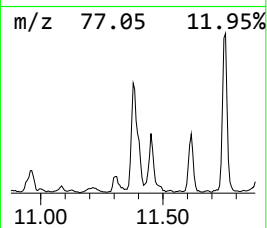
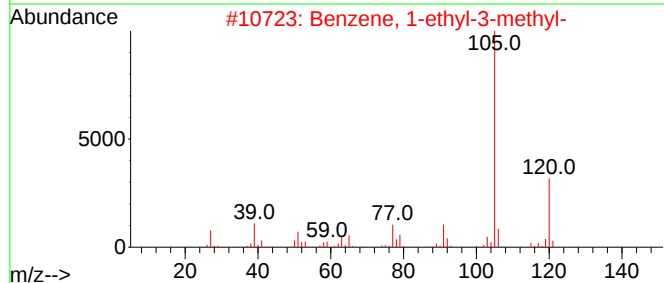
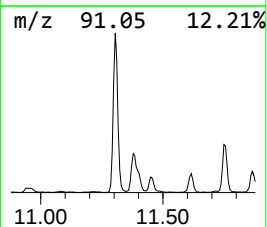
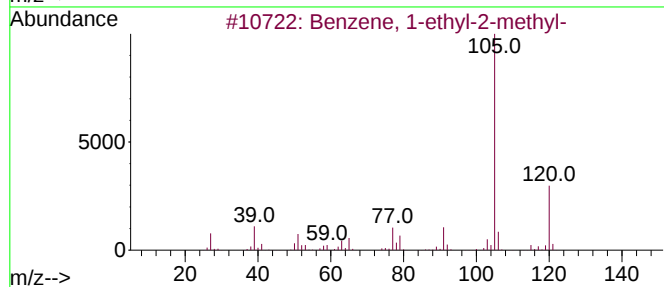
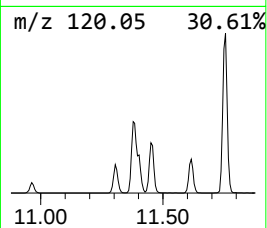
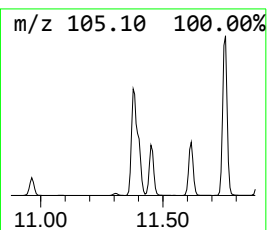
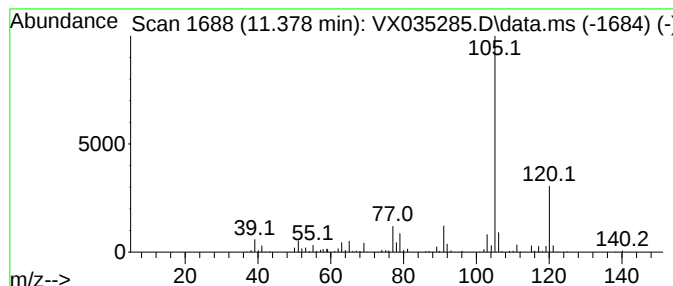
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1-ethyl-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.378	34.94 ug/l	836246	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			Mesitylene	120	C9H12	000108-67-8	91
4			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX042523\
Data File : VX035285.D
Acq On : 25 Apr 2023 15:50
Operator : JC/MD
Sample : 02431-01
Misc : 5.00g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
AUD-1943

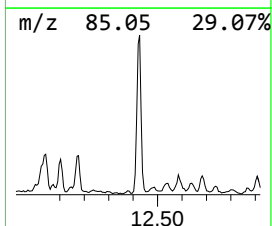
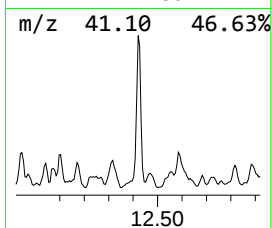
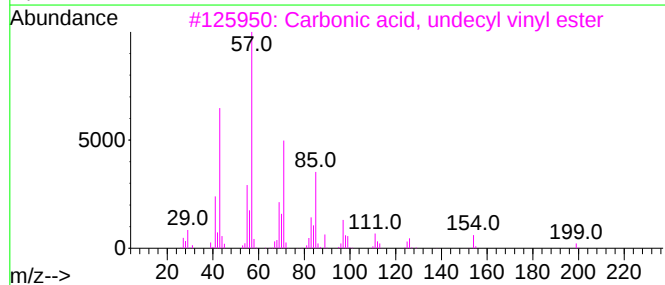
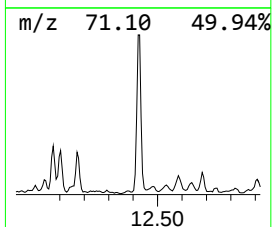
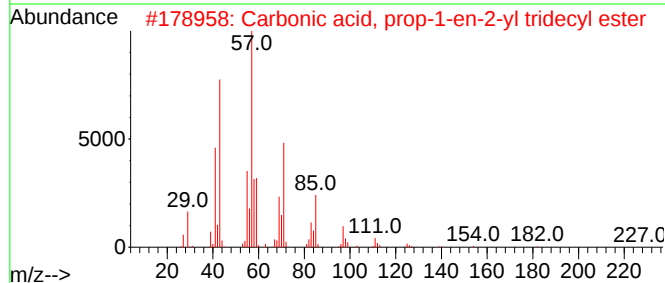
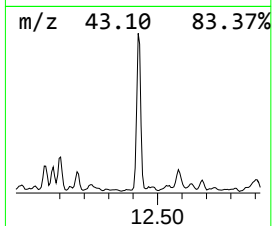
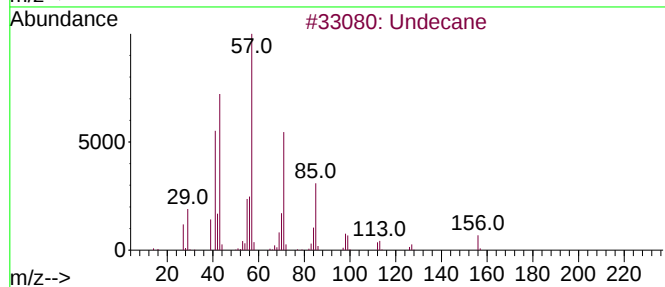
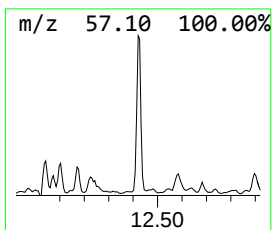
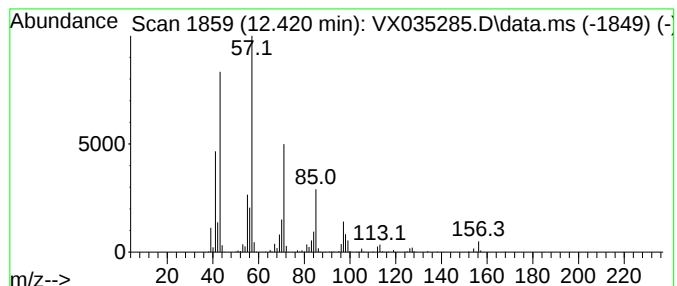
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Quant Title : SW846 8260

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

Peak Number 7 Undecane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.420	37.95 ug/l	908361	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane			156	C11H24	001120-21-4	96
2	Carbonic acid, prop-1-en-2-yl tr...			284	C17H32O3	1000382-90-8	86
3	Carbonic acid, undecyl vinyl ester			242	C14H26O3	1000382-54-9	83
4	Tetradecane			198	C14H30	000629-59-4	80
5	Hexadecane			226	C16H34	000544-76-3	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX042523\
Data File : VX035285.D
Acq On : 25 Apr 2023 15:50
Operator : JC/MD
Sample : 02431-01
Misc : 5.00g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 18 Sample Multiplier: 1

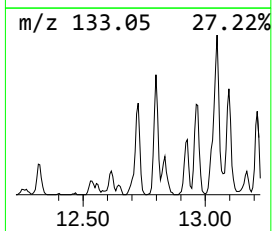
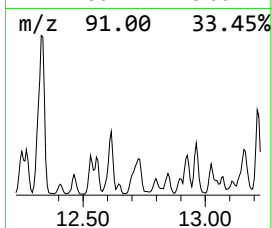
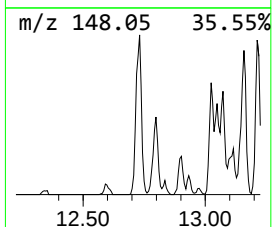
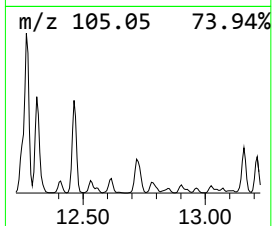
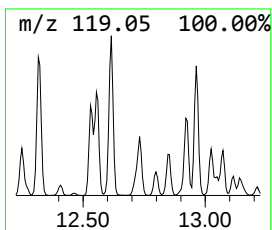
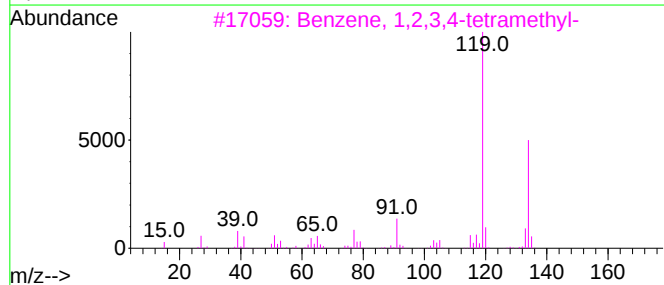
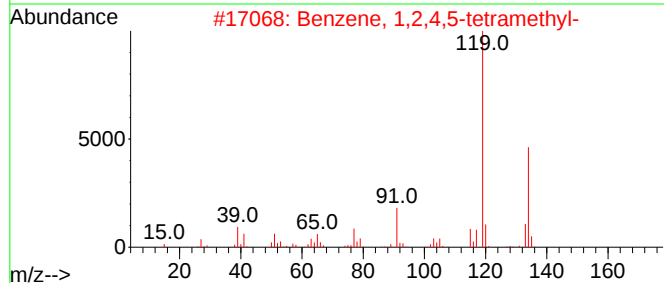
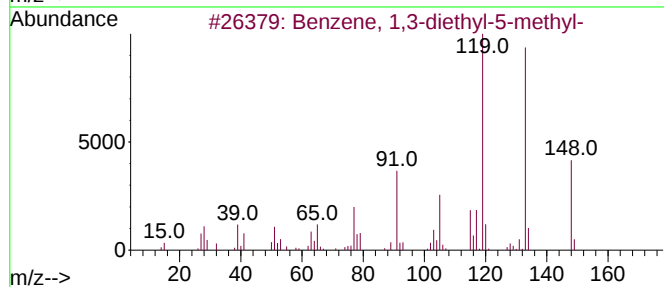
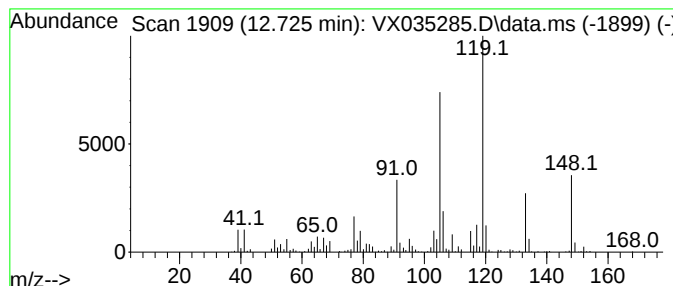
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ClientSampleId :
AUD-1943

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X042523W.M
Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,3-diethyl-5-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD			R.T.
12.725	21.19 ug/l	507223	1,4-Dichlorobenzene-d4			12.024
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Benzene, 1,3-diethyl-5-methyl-		148	C11H16	002050-24-0	64
2	Benzene, 1,2,4,5-tetramethyl-		134	C10H14	000095-93-2	64
3	Benzene, 1,2,3,4-tetramethyl-		134	C10H14	000488-23-3	64
4	p-Cymene		134	C10H14	000099-87-6	60
5	1,3,5-Cycloheptatriene, 3,7,7-tr...		134	C10H14	003479-89-8	60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX042523\
Data File : VX035285.D
Acq On : 25 Apr 2023 15:50
Operator : JC/MD
Sample : 02431-01
Misc : 5.00g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
AUD-1943

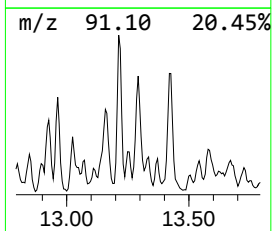
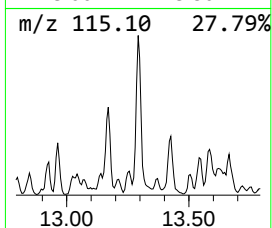
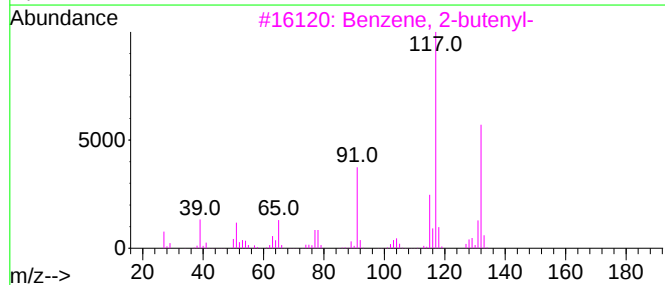
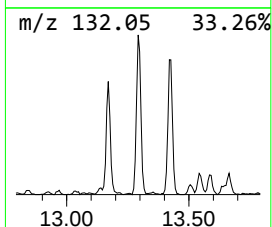
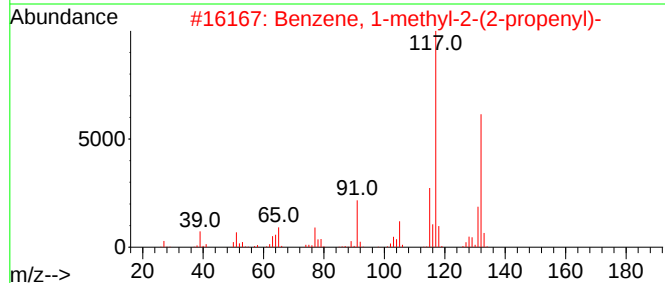
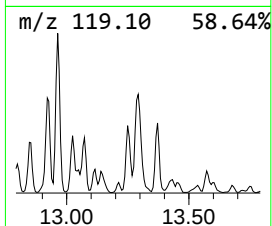
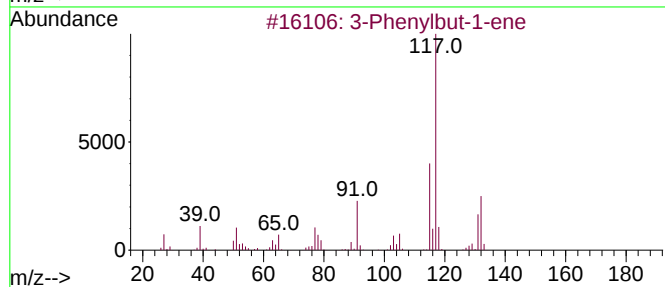
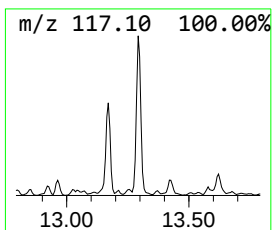
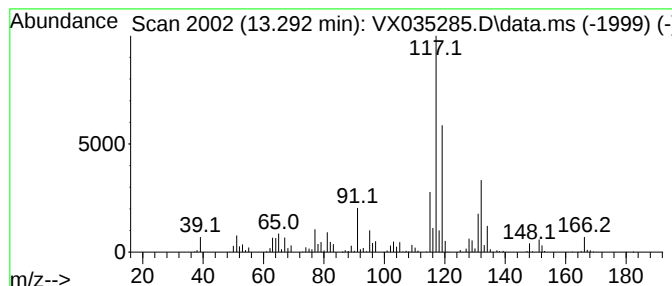
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Quant Title : SW846 8260

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

Peak Number 9 3-Phenylbut-1-ene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	26.14 ug/l	625623	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Phenylbut-1-ene	132	C10H12	000934-10-1	93
2			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	70
3			Benzene, 2-butenyl-	132	C10H12	001560-06-1	64
4			Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	64
5			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX042523\
Data File : VX035285.D
Acq On : 25 Apr 2023 15:50
Operator : JC/MD
Sample : 02431-01
Misc : 5.00g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
AUD-1943

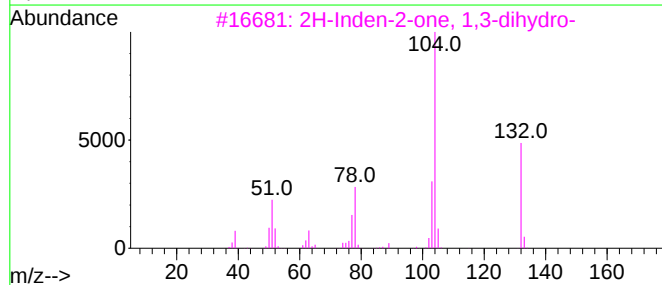
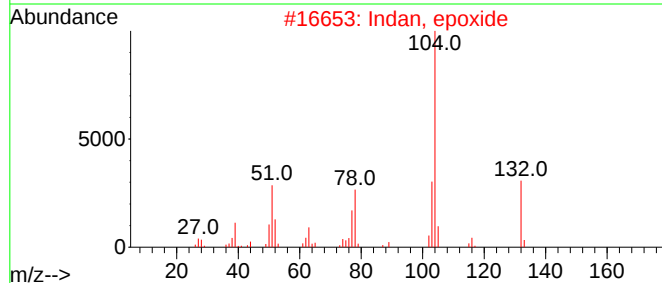
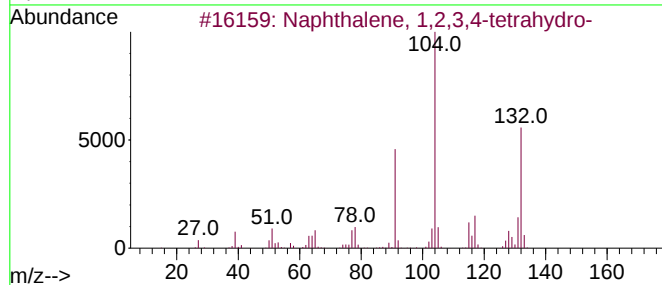
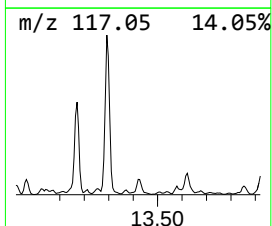
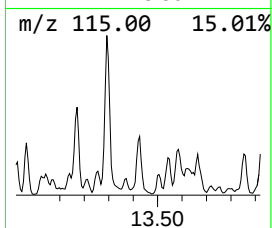
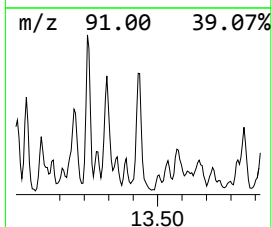
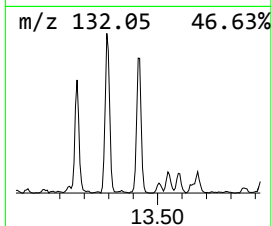
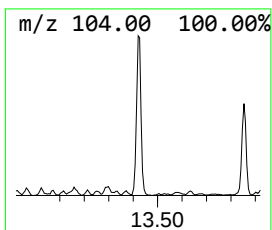
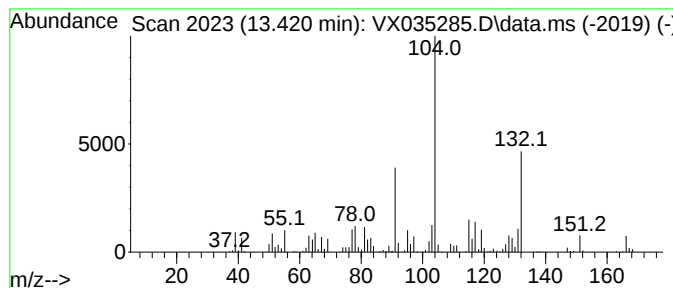
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X042523W.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.420	20.49 ug/l	490335	1,4-Dichlorobenzene-d4	12.024

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	90
2		Indan, epoxide	132	C9H8O	000768-22-9	52
3		2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	50
4		Benzene, cyclobutyl-	132	C10H12	004392-30-7	42
5		Benzene, 3-cyclohexen-1-yl-	158	C12H14	004994-16-5	38



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX042523\
Data File : VX035285.D
Acq On : 25 Apr 2023 15:50
Operator : JC/MD
Sample : 02431-01
Misc : 5.00g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
AUD-1943

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X042523W.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
Hexane, 3-methyl-	5.812	66.2	ug/l	2245710	1	5.544	1696490	50.0
Heptane	6.604	24.4	ug/l	468830	2	6.757	961475	50.0
Nonane	10.360	52.2	ug/l	1120750	3	10.055	1073980	50.0
Octane, 2,6-dim...	10.780	25.4	ug/l	546452	3	10.055	1073980	50.0
Cyclohexane, (1...	10.854	22.9	ug/l	492139	3	10.055	1073980	50.0
Benzene, 1-ethy...	11.378	34.9	ug/l	836246	4	12.024	1196740	50.0
Undecane	12.420	38.0	ug/l	908361	4	12.024	1196740	50.0
Benzene, 1,3-di...	12.725	21.2	ug/l	507223	4	12.024	1196740	50.0
3-Phenylbut-1-ene	13.292	26.1	ug/l	625623	4	12.024	1196740	50.0
Naphthalene, 1,...	13.420	20.5	ug/l	490335	4	12.024	1196740	50.0