

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092223\
 Data File : VX037877.D
 Acq On : 22 Sep 2023 14:57
 Operator : JC/MD
 Sample : 04436-09ME 20X
 Misc : 6.45g/5mL/100uL/5.0mL/MSVOA_X/MEOH
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 B102-08(5-5.5)ME

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\82X091523W.M
 Title : SW846 8260

Signal : TIC: VX037877.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.300	518	527	538	rVB3	35964	105255	7.43%	0.355%
2	5.385	693	705	713	rBV	130815	387357	27.33%	1.305%
3	5.550	713	732	741	rVV2	211082	757527	53.46%	2.553%
4	5.824	765	777	789	rBV3	80349	250917	17.71%	0.846%
5	5.952	789	798	810	rBV	137800	372335	26.27%	1.255%
6	6.159	818	832	840	rBV4	37081	127685	9.01%	0.430%
7	6.251	840	847	856	rVV4	61183	184070	12.99%	0.620%
8	6.348	856	863	877	rVB	57486	164885	11.64%	0.556%
9	6.610	893	906	920	rVB2	39371	103120	7.28%	0.347%
10	6.757	920	930	940	rBV	346730	889268	62.75%	2.997%
11	7.379	1019	1032	1044	rBV	281068	698922	49.32%	2.355%
12	7.653	1072	1077	1088	rVB2	41896	82616	5.83%	0.278%
13	7.775	1089	1097	1104	rVB2	64482	134462	9.49%	0.453%
14	7.952	1120	1126	1142	rVB3	74386	231096	16.31%	0.779%
15	8.104	1142	1151	1153	rBV3	49561	93134	6.57%	0.314%
16	8.141	1153	1157	1161	rVV	67209	113102	7.98%	0.381%
17	8.189	1161	1165	1174	rVB3	53326	107052	7.55%	0.361%
18	8.275	1174	1179	1182	rBV	129756	219473	15.49%	0.740%
19	8.311	1182	1185	1192	rVB4	107054	177886	12.55%	0.599%
20	8.433	1200	1205	1213	rBV	135591	255917	18.06%	0.862%
21	8.506	1213	1217	1226	rVB3	68894	132198	9.33%	0.445%
22	8.598	1226	1232	1235	rBV2	267895	480589	33.91%	1.619%
23	8.647	1235	1240	1250	rVB	800566	1355280	95.64%	4.567%
24	8.775	1250	1261	1265	rBV	101685	189925	13.40%	0.640%
25	8.848	1270	1273	1279	rVB	85398	127246	8.98%	0.429%
26	8.915	1279	1284	1289	rBV3	78990	149618	10.56%	0.504%
27	8.976	1289	1294	1300	rVV	196999	301525	21.28%	1.016%
28	9.098	1306	1314	1320	rBV2	120554	233262	16.46%	0.786%
29	9.159	1320	1324	1329	rVV2	118404	189297	13.36%	0.638%
30	9.384	1356	1361	1365	rBV	145432	210053	14.82%	0.708%
31	9.500	1376	1380	1386	rVV4	170015	365806	25.81%	1.233%
32	9.561	1386	1390	1394	rVV	271220	375515	26.50%	1.265%
33	9.616	1394	1399	1403	rVV2	318561	526248	37.13%	1.773%
34	9.659	1403	1406	1416	rVB5	78232	183666	12.96%	0.619%
35	9.762	1416	1423	1428	rBV3	135587	242571	17.12%	0.817%
36	9.823	1428	1433	1438	rBV3	157280	335436	23.67%	1.130%

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Integration Parameters: RTEINT.P

Integrator: RTE

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37	9.903	1442	1446	1451	rVV	210864	298644	21.07%	1.006%
38	10.006	1455	1463	1467	rBV	260536	382874	27.02%	1.290%
39	10.055	1467	1471	1476	rVV	777340	1038657	73.29%	3.500%
40	10.128	1476	1483	1486	rVV3	68781	148936	10.51%	0.502%
41	10.177	1487	1491	1499	rVB5	56514	117385	8.28%	0.396%
42	10.274	1499	1507	1510	rBV2	145861	279600	19.73%	0.942%
43	10.317	1510	1514	1517	rVV2	172621	261908	18.48%	0.883%
44	10.354	1517	1520	1531	rVB5	120596	281384	19.86%	0.948%
45	10.457	1531	1537	1545	rBV3	92896	170912	12.06%	0.576%
46	10.585	1551	1558	1566	rVB4	244760	470049	33.17%	1.584%
47	10.671	1566	1572	1577	rBV3	92635	192701	13.60%	0.649%
48	10.780	1582	1590	1594	rBV2	440589	704152	49.69%	2.373%
49	10.854	1594	1602	1606	rVV3	322147	550530	38.85%	1.855%
50	10.896	1606	1609	1614	rVB	122077	167822	11.84%	0.566%
51	10.957	1614	1619	1623	rBV2	113137	194165	13.70%	0.654%
52	11.000	1623	1626	1628	rVV3	68665	93923	6.63%	0.316%
53	11.030	1628	1631	1634	rVV	95739	121043	8.54%	0.408%
54	11.079	1634	1639	1650	rVB	831540	1417129	100.00%	4.775%
55	11.189	1650	1657	1659	rBV2	168801	240184	16.95%	0.809%
56	11.219	1659	1662	1666	rVB3	190127	243755	17.20%	0.821%
57	11.305	1671	1676	1679	rBV	220870	286824	20.24%	0.967%
58	11.390	1686	1690	1693	rBV3	78106	108000	7.62%	0.364%
59	11.433	1693	1697	1701	rBV	148740	193891	13.68%	0.653%
60	11.482	1701	1705	1709	rVB3	91503	143113	10.10%	0.482%
61	11.616	1722	1727	1734	rVB5	130489	269187	19.00%	0.907%
62	11.683	1734	1738	1740	rBV3	70959	87868	6.20%	0.296%
63	11.713	1740	1743	1746	rBV	201176	217110	15.32%	0.732%
64	11.866	1765	1768	1769	rVV	80343	102434	7.23%	0.345%
65	11.890	1769	1772	1776	rVV2	173205	248177	17.51%	0.836%
66	11.945	1776	1781	1783	rVV	188550	271759	19.18%	0.916%
67	11.969	1783	1785	1789	rVV3	120970	157527	11.12%	0.531%
68	12.024	1789	1794	1799	rVV	954650	1365019	96.32%	4.600%
69	12.073	1799	1802	1804	rVV	93468	109371	7.72%	0.369%
70	12.103	1804	1807	1812	rVV4	76455	125996	8.89%	0.425%
71	12.170	1812	1818	1825	rVB4	108441	256108	18.07%	0.863%
72	12.244	1825	1830	1836	rBV2	396671	541812	38.23%	1.826%
73	12.311	1837	1841	1848	rBV2	252162	478168	33.74%	1.611%
74	12.463	1863	1866	1872	rVB3	106293	182190	12.86%	0.614%
75	12.609	1887	1890	1894	rVB	405884	432753	30.54%	1.458%
76	12.695	1899	1904	1913	rVB3	304842	628927	44.38%	2.119%

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77	12.817	1920	1924	1930	rVB5	140376	260036	18.35%	0.876%
78	12.890	1931	1936	1938	rBV2	153504	244520	17.25%	0.824%
79	12.920	1938	1941	1947	rVB	293106	380955	26.88%	1.284%
80	12.981	1947	1951	1954	rBV2	91008	110695	7.81%	0.373%
81	13.024	1954	1958	1964	rVB4	137277	258681	18.25%	0.872%
82	13.091	1966	1969	1974	rBV2	64887	118987	8.40%	0.401%
83	13.164	1978	1981	1985	rVB	173916	198440	14.00%	0.669%
84	13.256	1992	1996	1998	rBV2	63769	88681	6.26%	0.299%
85	13.292	1998	2002	2008	rVV2	552085	836724	59.04%	2.819%
86	13.371	2012	2015	2020	rVV2	111035	218518	15.42%	0.736%
87	13.420	2020	2023	2032	rVB7	127304	291904	20.60%	0.984%
88	13.506	2032	2037	2039	rBV2	97416	136538	9.63%	0.460%
89	13.542	2039	2043	2046	rVV	165386	208614	14.72%	0.703%
90	13.579	2046	2049	2053	rVV3	180861	250262	17.66%	0.843%
91	13.634	2054	2058	2060	rVV4	57624	97802	6.90%	0.330%
92	13.664	2060	2063	2068	rVB2	174665	248613	17.54%	0.838%
93	13.841	2088	2092	2099	rVB7	69675	131802	9.30%	0.444%
94	13.926	2099	2106	2110	rBV5	100391	192175	13.56%	0.648%
95	14.073	2125	2130	2133	rBV2	86325	115906	8.18%	0.391%
96	14.213	2149	2153	2162	rVB2	119413	195772	13.81%	0.660%
97	14.371	2176	2179	2183	rVB	74372	83122	5.87%	0.280%
98	14.481	2193	2197	2201	rBV2	50920	85543	6.04%	0.288%
99	14.633	2215	2222	2226	rBV	187565	271996	19.19%	0.917%
100	14.780	2242	2246	2254	rVV	148418	213768	15.08%	0.720%

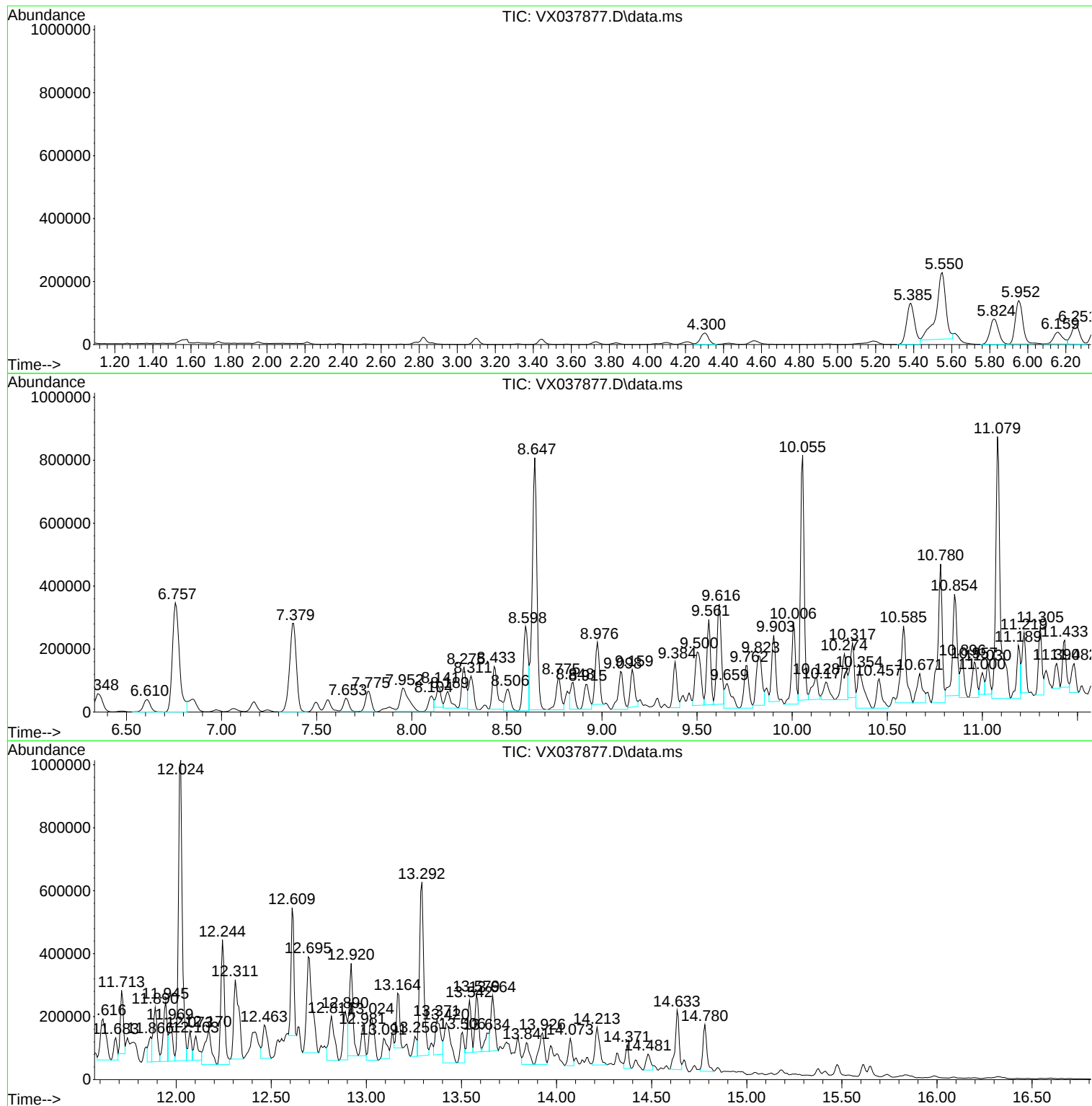
Sum of corrected areas: 29676355

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091523W.M
Quant Title : SW846 8260

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



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B102-08(5-5.5)ME

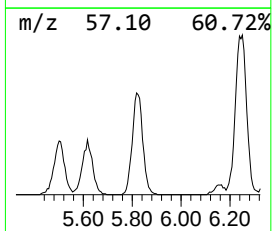
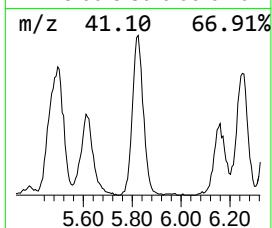
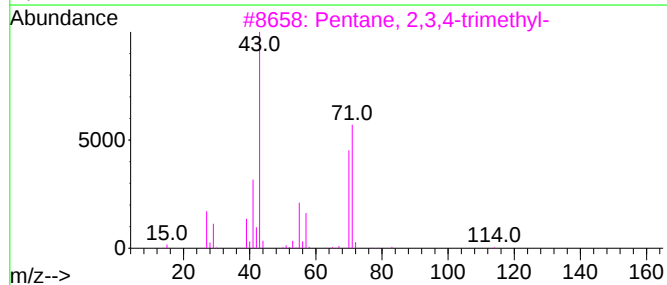
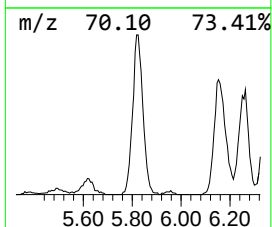
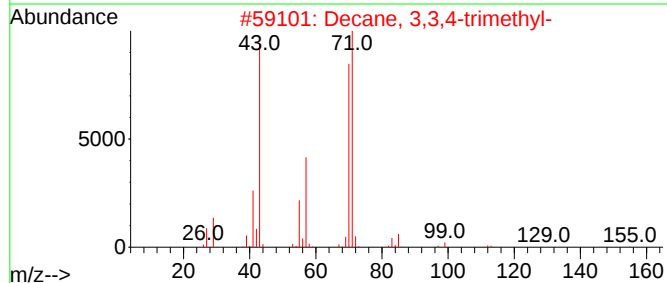
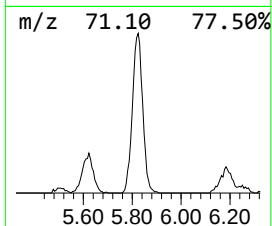
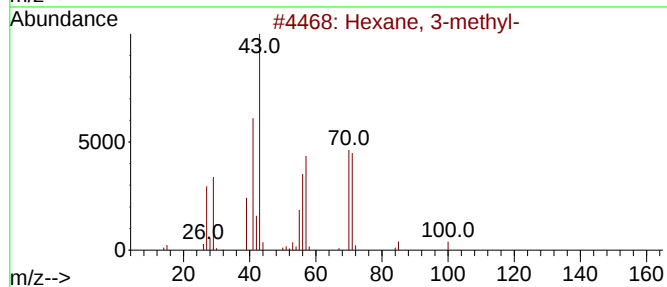
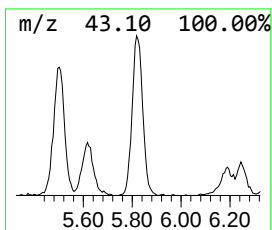
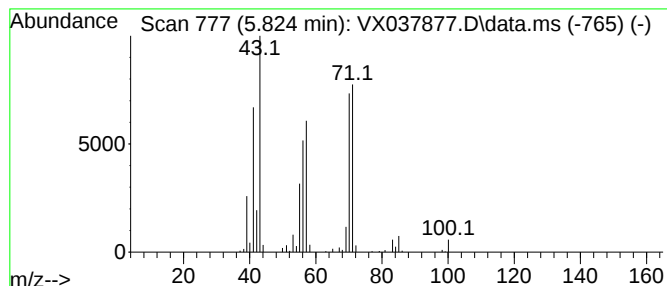
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091523W.M
Quant Title : SW846 8260

TIC Library : C:\Database\NIST20.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	16.56 ug/l	250917	Pentafluorobenzene	5.550

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexane, 3-methyl-	100	C7H16	000589-34-4	96
2			Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	59
3			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	53
4			Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	50
5			Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	50



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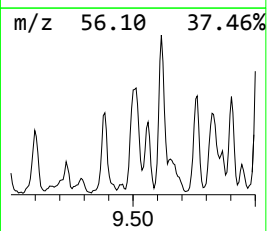
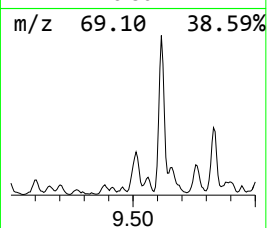
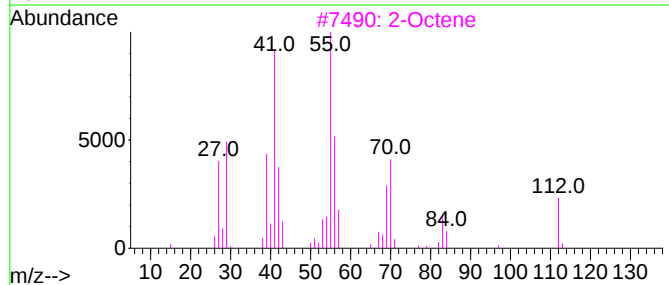
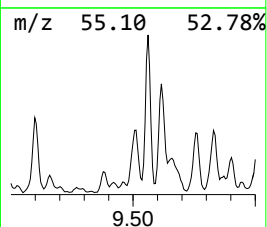
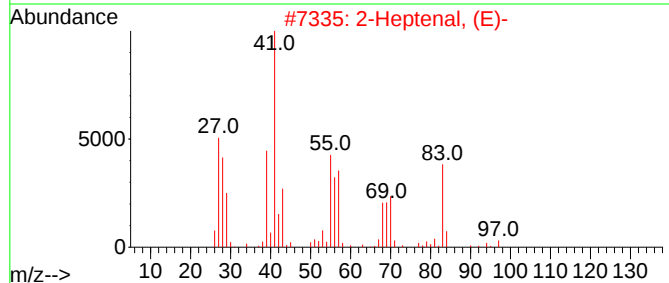
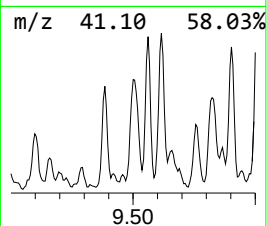
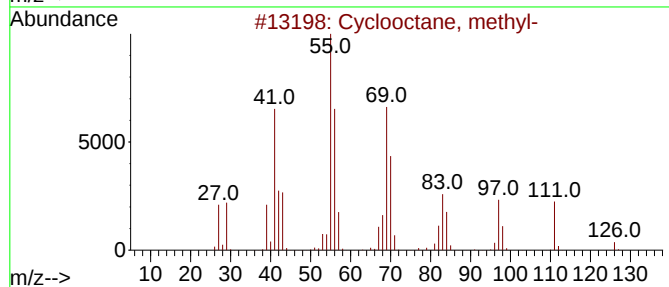
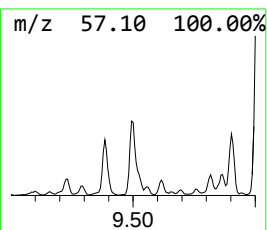
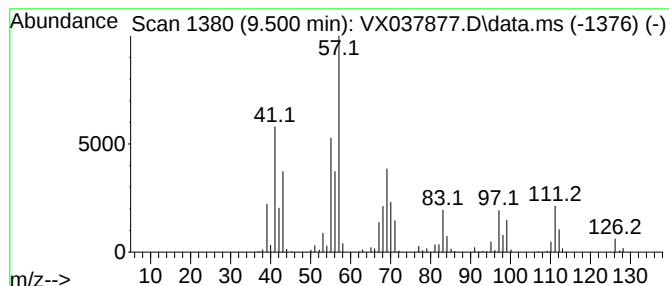
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091523W.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown9.500 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.500	17.61 ug/l	365806	Chlorobenzene-d5	10.055

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclooctane, methyl-	126	C9H18	001502-38-1	49
2			2-Heptenal, (E)-	112	C7H12O	018829-55-5	43
3			2-Octene	112	C8H16	000111-67-1	38
4			3-Octene, (E)-	112	C8H16	014919-01-8	38
5			3-Nonene, (E)-	126	C9H18	020063-92-7	30



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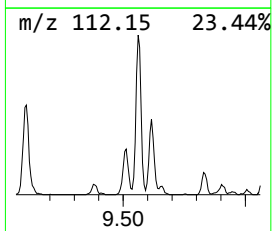
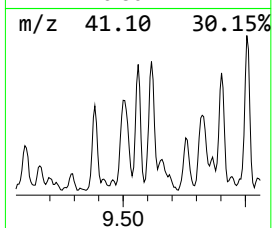
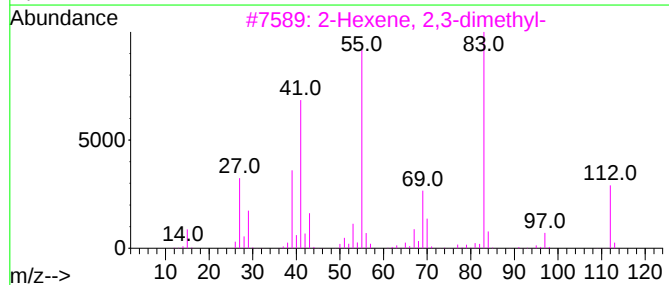
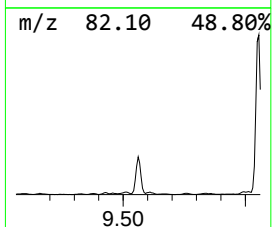
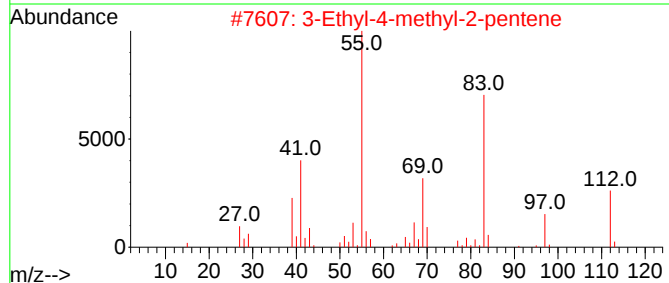
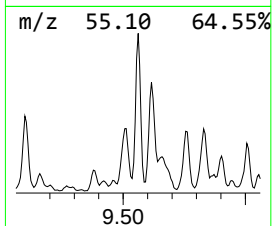
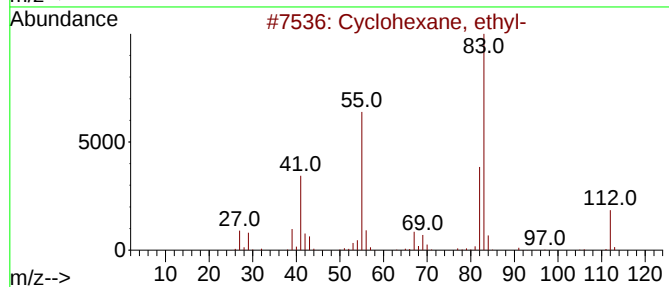
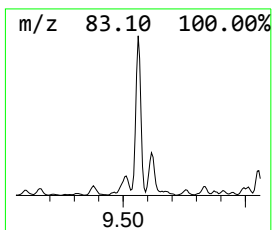
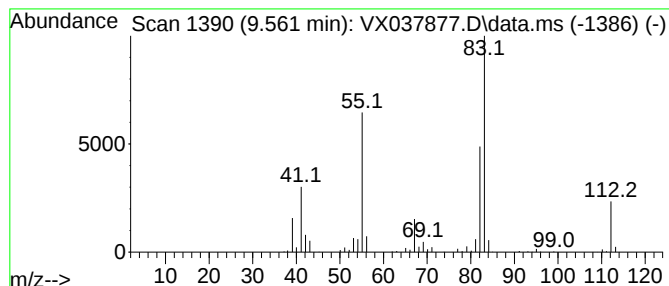
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Quant Title : SW846 8260

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

Peak Number 3 Cyclohexane, ethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.561	18.08 ug/l	375515	Chlorobenzene-d5	10.055

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, ethyl-	112	C8H16	001678-91-7	86
2			3-Ethyl-4-methyl-2-pentene	112	C8H16	019780-68-8	59
3			2-Hexene, 2,3-dimethyl-	112	C8H16	007145-20-2	58
4			3-Hexene, 2,4-dimethyl-	112	C8H16	082085-14-1	53
5			3-Hexene, 3,4-dimethyl-	112	C8H16	030951-95-2	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092223\
Data File : VX037877.D
Acq On : 22 Sep 2023 14:57
Operator : JC/MD
Sample : 04436-09ME 20X
Misc : 6.45g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B102-08(5-5.5)ME

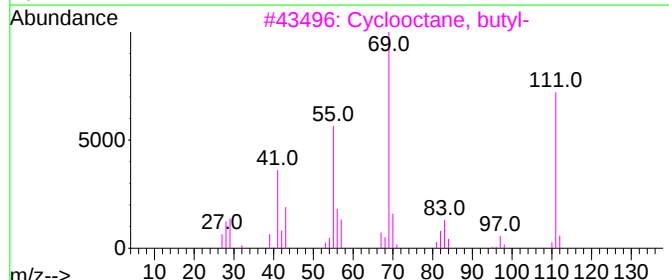
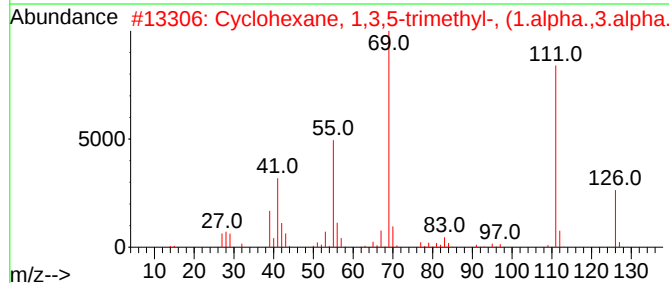
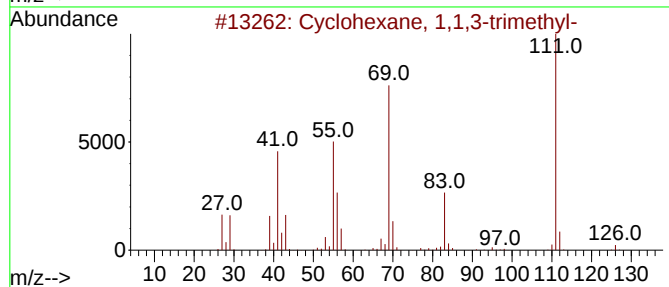
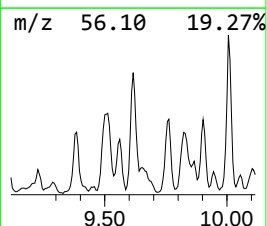
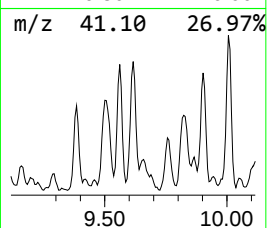
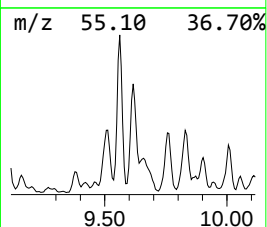
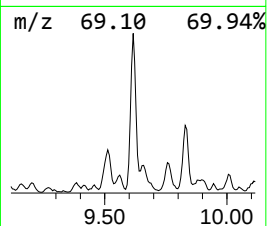
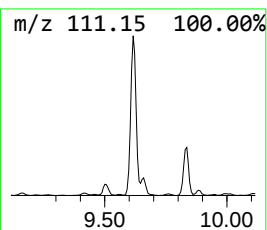
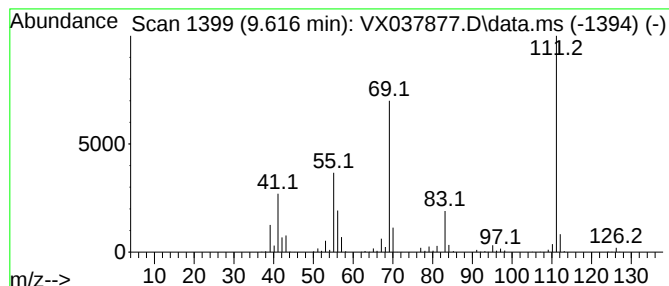
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Quant Title : SW846 8260

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

Peak Number 4 Cyclohexane, 1,1,3-trimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.616	25.33 ug/l	526248	Chlorobenzene-d5	10.055

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	95
2			Cyclohexane, 1,3,5-trimethyl-, (...)	126	C9H18	001795-27-3	64
3			Cyclooctane, butyl-	168	C12H24	016538-93-5	64
4			Cyclooctane, (1-methylpropyl)-	168	C12H24	016538-89-9	59
5			4-Amino-6-hydroxypyrimidine	111	C4H5N3O	001193-22-2	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092223\
Data File : VX037877.D
Acq On : 22 Sep 2023 14:57
Operator : JC/MD
Sample : 04436-09ME 20X
Misc : 6.45g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B102-08(5-5.5)ME

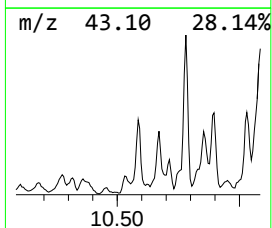
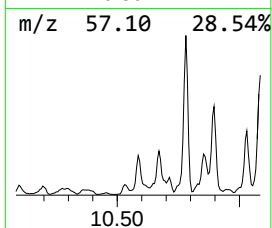
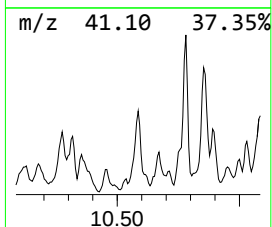
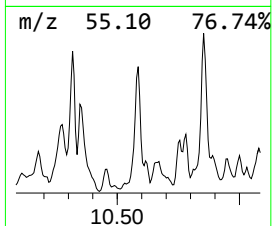
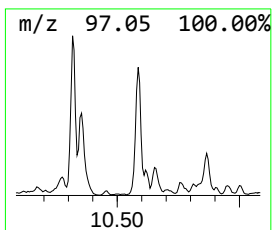
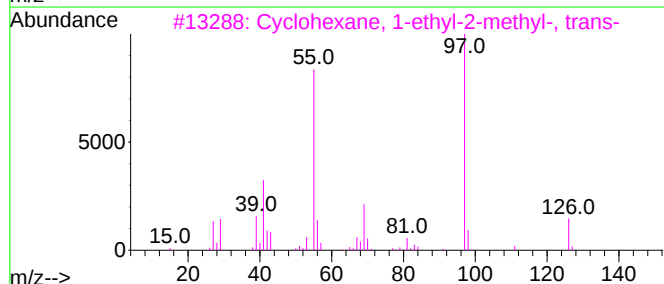
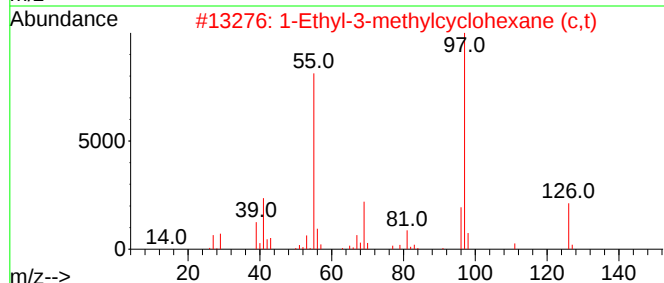
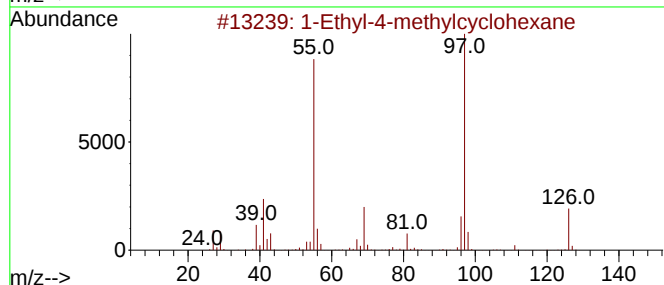
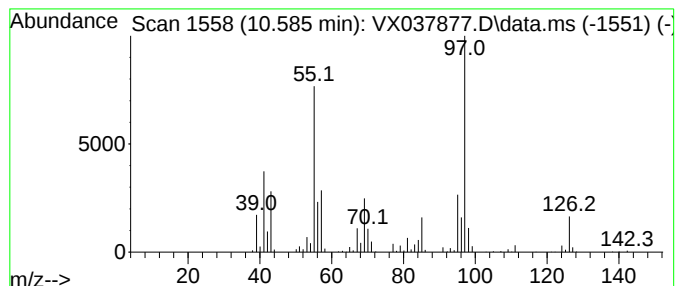
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Quant Title : SW846 8260

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

Peak Number 5 1-Ethyl-4-methylcyclohexane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.585	22.63 ug/l	470049	Chlorobenzene-d5	10.055

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Ethyl-4-methylcyclohexane	126	C9H18	003728-56-1	70
2			1-Ethyl-3-methylcyclohexane (c,t)	126	C9H18	003728-55-0	70
3			Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-78-8	70
4			cis-1-Ethyl-3-methyl-cyclohexane	126	C9H18	019489-10-2	70
5			Cyclohexane, 1-ethyl-2-methyl-	126	C9H18	003728-54-9	62



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092223\
Data File : VX037877.D
Acq On : 22 Sep 2023 14:57
Operator : JC/MD
Sample : 04436-09ME 20X
Misc : 6.45g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B102-08(5-5.5)ME

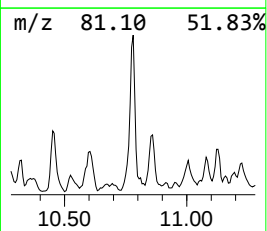
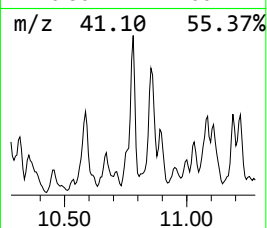
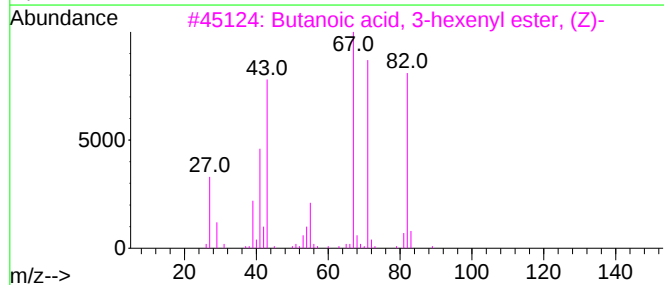
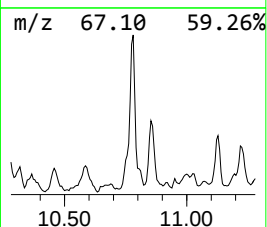
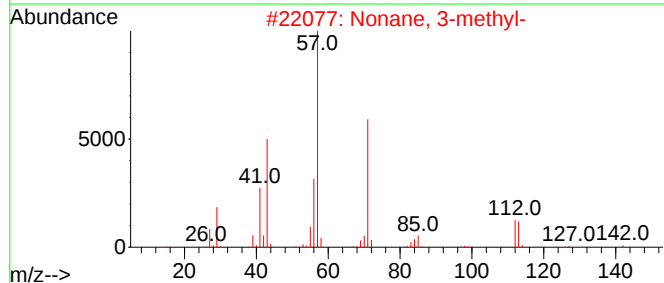
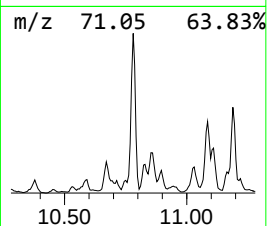
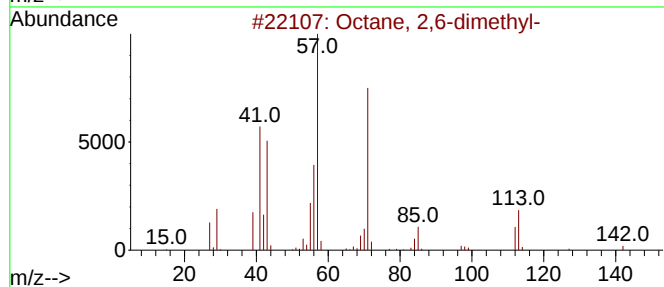
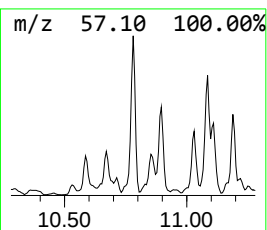
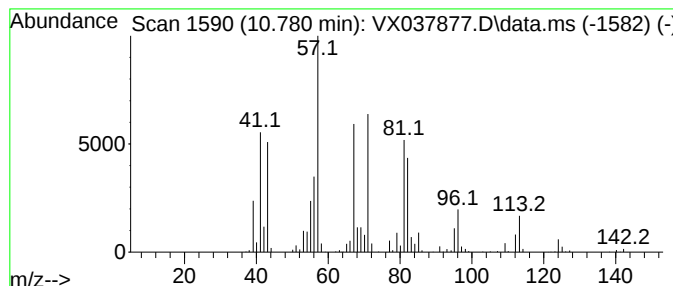
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.780	33.90 ug/l	704152	Chlorobenzene-d5	10.055

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	86
2			Nonane, 3-methyl-	142	C10H22	005911-04-6	46
3			Butanoic acid, 3-hexenyl ester, ...	170	C10H18O2	016491-36-4	43
4			Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	38
5			3,4-Nonadiene	124	C9H16	037050-03-6	27



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092223\
Data File : VX037877.D
Acq On : 22 Sep 2023 14:57
Operator : JC/MD
Sample : 04436-09ME 20X
Misc : 6.45g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 16 Sample Multiplier: 1

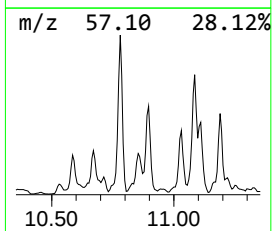
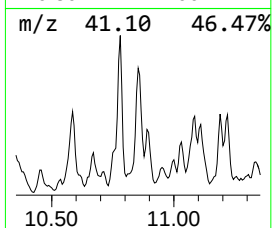
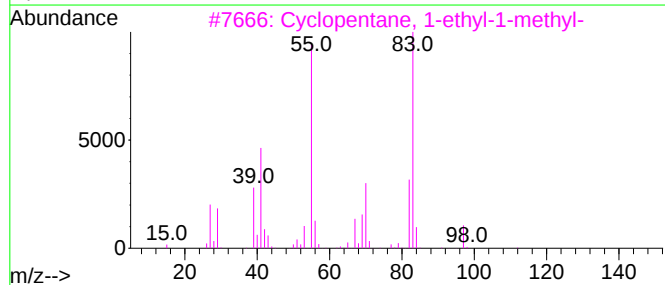
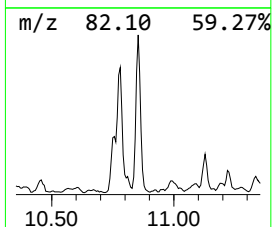
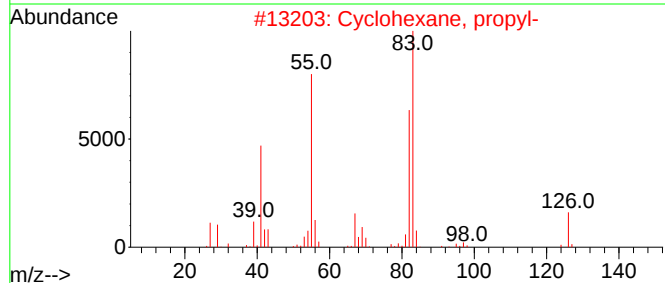
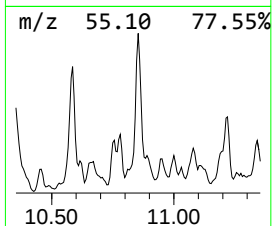
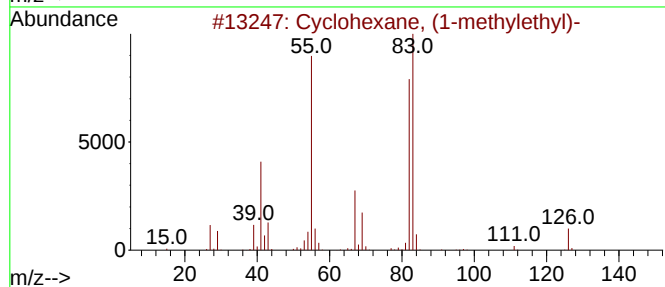
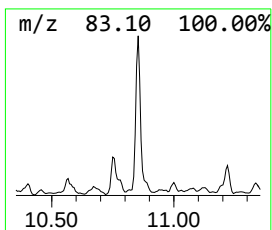
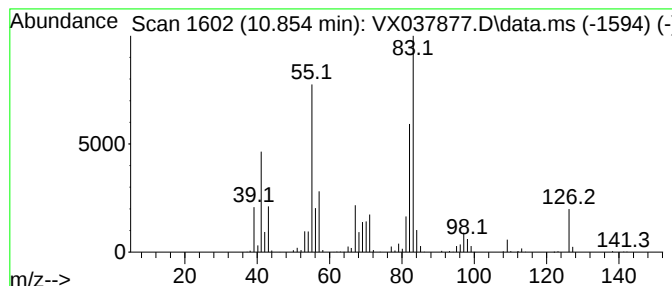
Instrument :
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ClientSampleId :
B102-08(5-5.5)ME

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091523W.M
Quant Title : SW846 8260

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

Peak Number 7 Cyclohexane, (1-methylethyl)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.854	26.50 ug/l	550530	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	83
3	Cyclopentane, 1-ethyl-1-methyl-	112	C8H16	016747-50-5	58
4	Cyclohexane, ethyl-	112	C8H16	001678-91-7	53
5	Oxalic acid, cyclohexyl dodecyl ...	340	C20H36O4	1000309-31-4	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092223\
Data File : VX037877.D
Acq On : 22 Sep 2023 14:57
Operator : JC/MD
Sample : 04436-09ME 20X
Misc : 6.45g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B102-08(5-5.5)ME

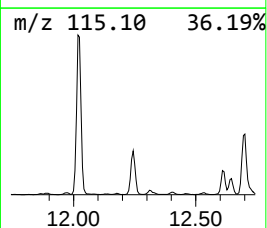
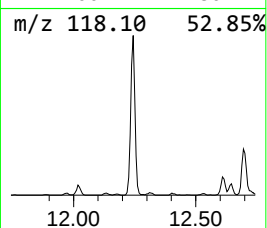
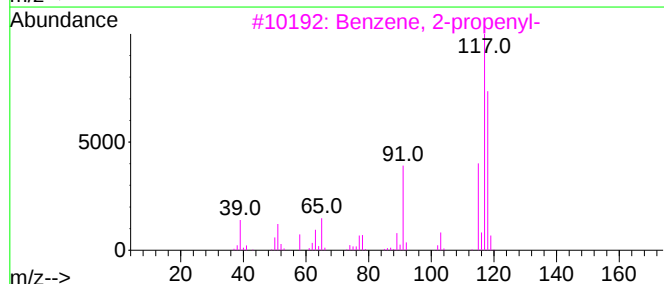
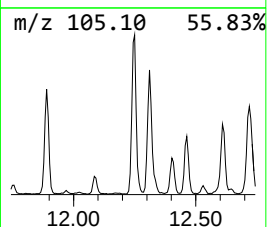
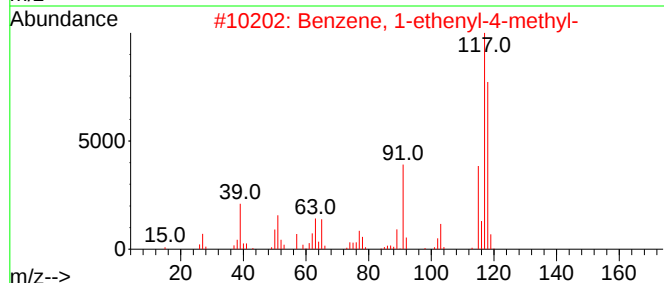
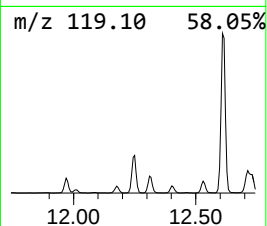
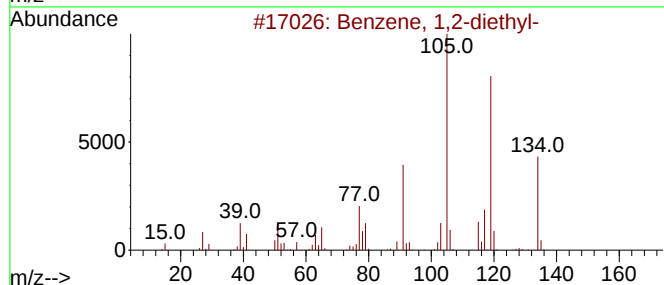
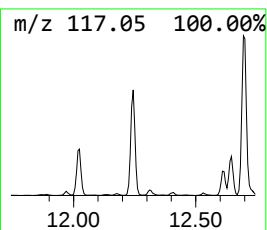
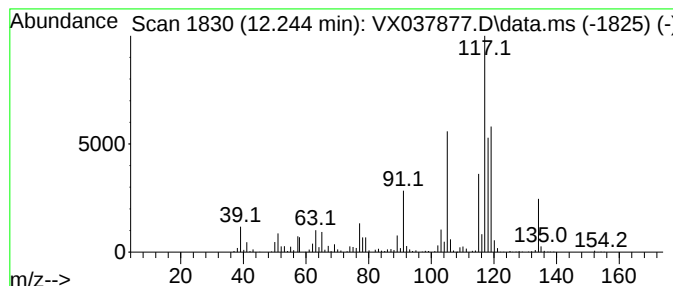
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2-diethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	19.85 ug/l	541812	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	83
2			Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	83
3			Benzene, 2-propenyl-	118	C9H10	000300-57-2	80
4			Benzene, 1-propenyl-	118	C9H10	000637-50-3	60
5			Benzene, cyclopropyl-	118	C9H10	000873-49-4	60



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092223\
Data File : VX037877.D
Acq On : 22 Sep 2023 14:57
Operator : JC/MD
Sample : 04436-09ME 20X
Misc : 6.45g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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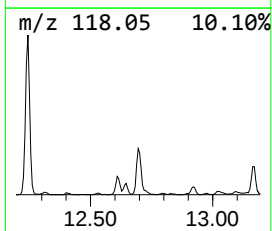
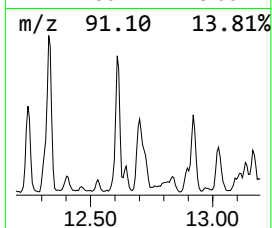
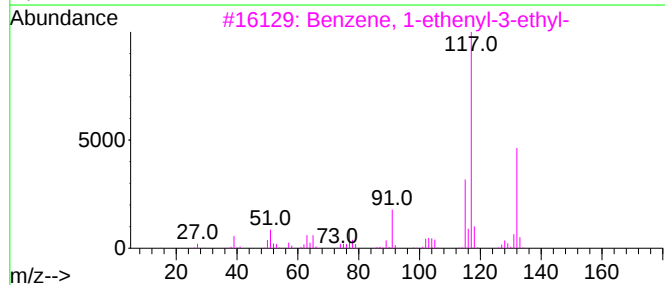
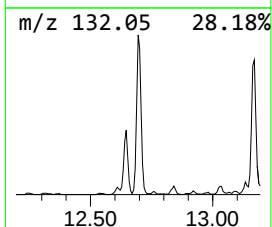
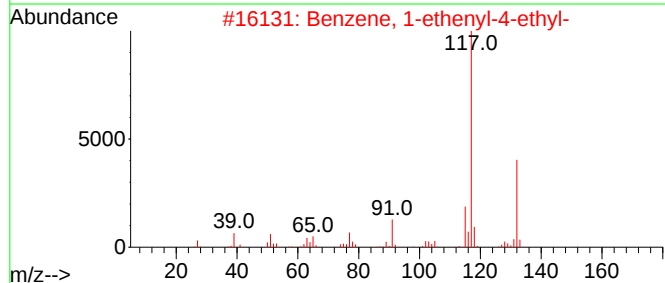
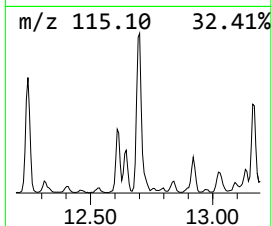
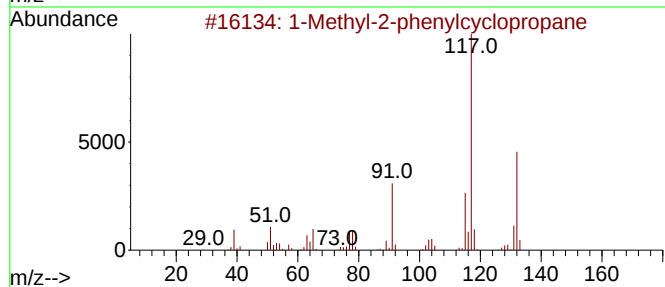
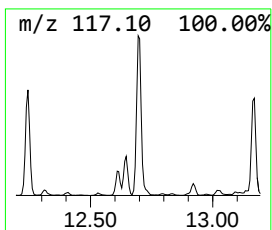
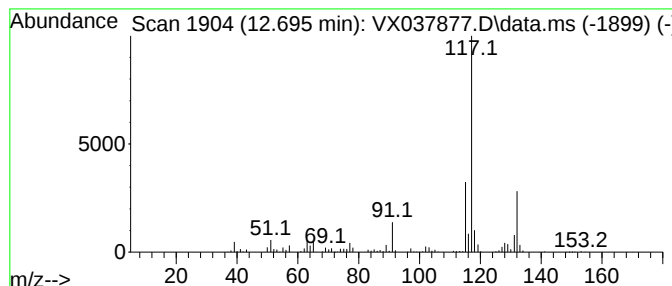
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Quant Title : SW846 8260

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

Peak Number 9 1-Methyl-2-phenylcyclopropane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.695	23.04 ug/l	628927	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	87
2			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
3			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
4			Indan, 1-methyl-	132	C10H12	000767-58-8	87
5			Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	86



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092223\
Data File : VX037877.D
Acq On : 22 Sep 2023 14:57
Operator : JC/MD
Sample : 04436-09ME 20X
Misc : 6.45g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B102-08(5-5.5)ME

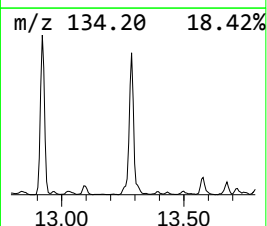
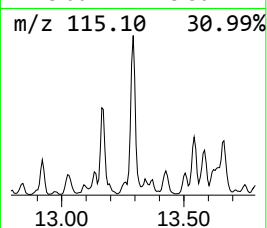
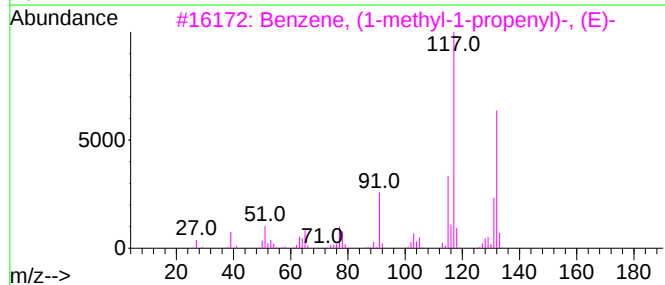
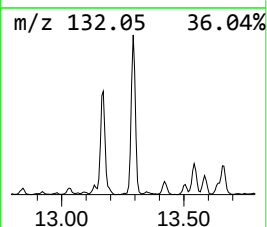
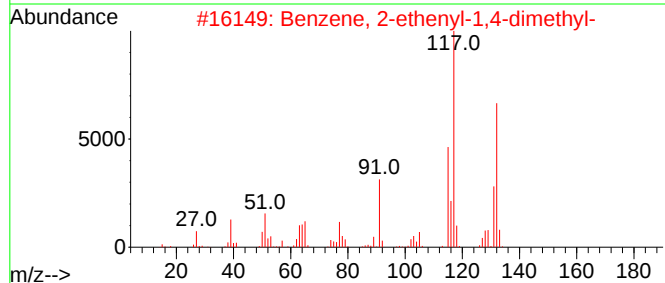
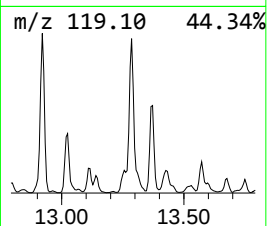
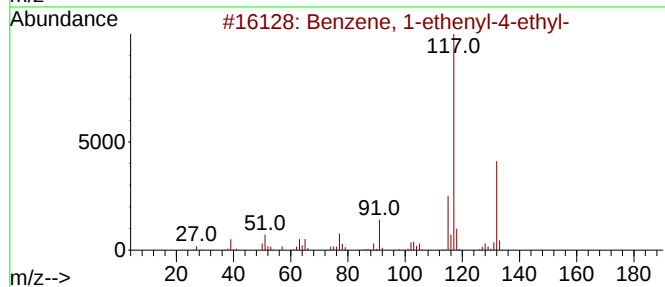
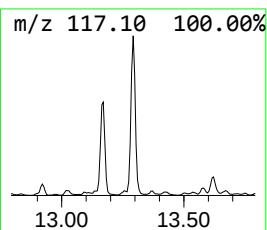
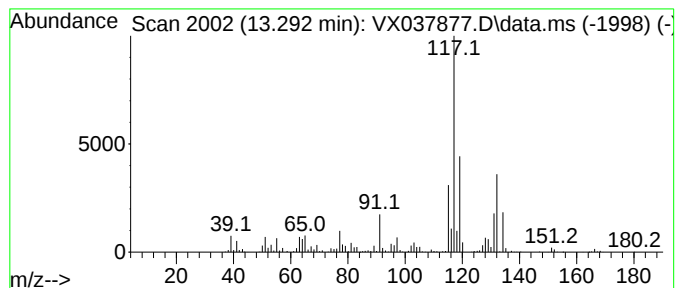
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Quant Title : SW846 8260

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

Peak Number 10 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	30.65 ug/l	836724	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	92
2			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	91
3			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	86
4			Indan, 1-methyl-	132	C10H12	000767-58-8	70
5			Benzene, 2-butenyl-	132	C10H12	001560-06-1	70



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092223\
Data File : VX037877.D
Acq On : 22 Sep 2023 14:57
Operator : JC/MD
Sample : 04436-09ME 20X
Misc : 6.45g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B102-08(5-5.5)ME

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091523W.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-	5.824	16.6	ug/l	250917	1	5.550	757527	50.0
unknown9.500	9.500	17.6	ug/l	365806	3	10.055	1038660	50.0
Cyclohexane, et...	9.561	18.1	ug/l	375515	3	10.055	1038660	50.0
Cyclohexane, 1,...	9.616	25.3	ug/l	526248	3	10.055	1038660	50.0
1-Ethyl-4-methy...	10.585	22.6	ug/l	470049	3	10.055	1038660	50.0
Octane, 2,6-dim...	10.780	33.9	ug/l	704152	3	10.055	1038660	50.0
Cyclohexane, (1...	10.854	26.5	ug/l	550530	3	10.055	1038660	50.0
Benzene, 1,2-di...	12.244	19.9	ug/l	541812	4	12.024	1365020	50.0
1-Methyl-2-phen...	12.695	23.0	ug/l	628927	4	12.024	1365020	50.0
Benzene, 1-ethe...	13.292	30.6	ug/l	836724	4	12.024	1365020	50.0