

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX080823\
 Data File : VX036870.D
 Acq On : 08 Aug 2023 17:39
 Operator : JC/MD
 Sample : 03903-20
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-10W

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\82X072123W.M
 Title : SW846 8260

Signal : TIC: VX036870.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.825	281	285	303	rVB	23662	53981	3.37%	0.175%
2	3.105	320	331	342	rVB	36125	83092	5.19%	0.269%
3	5.385	691	705	716	rBV	87935	257438	16.07%	0.834%
4	5.556	716	733	740	rVV2	140350	493981	30.84%	1.601%
5	5.623	740	744	758	rVB2	45976	127240	7.94%	0.412%
6	5.830	764	778	791	rBV3	52158	177280	11.07%	0.575%
7	5.958	791	799	813	rVV	90919	243036	15.17%	0.788%
8	6.159	816	832	840	rVV5	52266	174639	10.90%	0.566%
9	6.257	840	848	857	rVV3	82849	247492	15.45%	0.802%
10	6.355	857	864	876	rVB3	42267	120334	7.51%	0.390%
11	6.763	921	931	940	rVV	230841	551747	34.45%	1.788%
12	7.373	1021	1031	1042	rVB3	181865	482650	30.13%	1.564%
13	7.501	1042	1052	1057	rBV2	24542	55614	3.47%	0.180%
14	7.562	1057	1062	1072	rVB2	32463	69560	4.34%	0.225%
15	7.775	1090	1097	1106	rVB3	86459	180413	11.26%	0.585%
16	7.958	1121	1127	1143	rVB2	80123	241020	15.05%	0.781%
17	8.104	1143	1151	1159	rBV2	64995	144161	9.00%	0.467%
18	8.190	1159	1165	1175	rVV4	41508	101100	6.31%	0.328%
19	8.275	1175	1179	1181	rVV2	31580	48645	3.04%	0.158%
20	8.318	1181	1186	1192	rVV3	85176	172516	10.77%	0.559%
21	8.433	1201	1205	1209	rVV	46795	80857	5.05%	0.262%
22	8.513	1213	1218	1225	rVB2	52350	109862	6.86%	0.356%
23	8.604	1225	1233	1236	rBV2	249678	493992	30.84%	1.601%
24	8.647	1236	1240	1251	rVB	544170	910454	56.84%	2.951%
25	8.775	1251	1261	1265	rBV	176254	301762	18.84%	0.978%
26	8.817	1265	1268	1271	rVV	59915	98095	6.12%	0.318%
27	8.848	1271	1273	1279	rVV3	48456	64809	4.05%	0.210%
28	8.921	1280	1285	1289	rVV2	38273	60856	3.80%	0.197%
29	8.976	1289	1294	1300	rVB	312783	497840	31.08%	1.613%
30	9.104	1307	1315	1320	rVB	240012	403254	25.18%	1.307%
31	9.159	1320	1324	1328	rBV	106062	159806	9.98%	0.518%
32	9.384	1355	1361	1365	rBV2	53027	89948	5.62%	0.292%
33	9.506	1376	1381	1386	rVV2	167253	292496	18.26%	0.948%
34	9.561	1386	1390	1394	rVB	153905	205269	12.82%	0.665%
35	9.616	1394	1399	1403	rBV	164706	251897	15.73%	0.816%
36	9.659	1403	1406	1416	rVB3	67496	133565	8.34%	0.433%

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37	9.762	1417	1423	1428	rBV3	54246	95231	5.95%	0.309%
38	9.829	1428	1434	1438	rBV2	104239	192975	12.05%	0.625%
39	9.903	1443	1446	1455	rVB3	33293	62908	3.93%	0.204%
40	10.055	1466	1471	1476	rVB	485701	636598	39.75%	2.063%
41	10.128	1476	1483	1487	rBV	139503	216927	13.54%	0.703%
42	10.195	1487	1494	1498	rVB2	128690	197207	12.31%	0.639%
43	10.299	1503	1511	1517	rVV3	370746	825925	51.57%	2.677%
44	10.354	1517	1520	1532	rVB5	67211	145935	9.11%	0.473%
45	10.457	1532	1537	1544	rVB	90400	131804	8.23%	0.427%
46	10.585	1546	1558	1566	rBV3	166474	452145	28.23%	1.465%
47	10.653	1566	1569	1577	rVB4	34313	78451	4.90%	0.254%
48	10.781	1586	1590	1594	rVB	196925	252708	15.78%	0.819%
49	10.854	1598	1602	1607	rBV3	160334	243162	15.18%	0.788%
50	10.963	1615	1620	1623	rBV	94494	132033	8.24%	0.428%
51	11.079	1635	1639	1643	rBV	420907	524289	32.73%	1.699%
52	11.128	1643	1647	1650	rVB2	103242	131273	8.20%	0.425%
53	11.219	1659	1662	1666	rVB3	88268	111724	6.98%	0.362%
54	11.305	1672	1676	1679	rBV	207814	243715	15.22%	0.790%
55	11.378	1684	1688	1695	rBV	271693	498575	31.13%	1.616%
56	11.451	1695	1700	1703	rBV	133797	168029	10.49%	0.545%
57	11.616	1722	1727	1734	rVB	156280	239168	14.93%	0.775%
58	11.750	1745	1749	1754	rVB	528496	636196	39.72%	2.062%
59	11.866	1760	1768	1770	rBV	155894	247529	15.45%	0.802%
60	11.890	1770	1772	1776	rVB	196008	208539	13.02%	0.676%
61	12.024	1789	1794	1799	rBV	513557	704260	43.97%	2.282%
62	12.085	1801	1804	1808	rVB	96301	109130	6.81%	0.354%
63	12.177	1815	1819	1825	rVB	79039	114452	7.15%	0.371%
64	12.244	1825	1830	1837	rBV	1100004	1467490	91.62%	4.756%
65	12.311	1837	1841	1850	rVB2	480585	746990	46.64%	2.421%
66	12.402	1851	1856	1862	rBV	151951	206394	12.89%	0.669%
67	12.463	1862	1866	1872	rVB2	56922	92405	5.77%	0.299%
68	12.530	1872	1877	1879	rBV	65166	86145	5.38%	0.279%
69	12.609	1886	1890	1894	rVV	1069967	1312692	81.96%	4.254%
70	12.646	1894	1896	1901	rVB2	215961	255391	15.94%	0.828%
71	12.701	1901	1905	1912	rVB	626351	901130	56.26%	2.920%
72	12.762	1912	1915	1920	rBV2	37047	51122	3.19%	0.166%
73	12.811	1920	1923	1925	rBV	84879	107907	6.74%	0.350%
74	12.841	1925	1928	1933	rVB2	142928	197038	12.30%	0.639%
75	12.920	1933	1941	1945	rBV	1266328	1601702	100.00%	5.191%
76	12.963	1945	1948	1954	rVB2	76787	112778	7.04%	0.365%

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77	13.024	1954	1958	1964	rBV3	216117	313206	19.55%	1.015%
78	13.115	1967	1973	1974	rBV2	59573	114707	7.16%	0.372%
79	13.134	1974	1976	1979	rBV	99761	99581	6.22%	0.323%
80	13.170	1979	1982	1985	rBV	299685	315895	19.72%	1.024%
81	13.256	1991	1996	1998	rBV2	93972	129721	8.10%	0.420%
82	13.292	1998	2002	2008	rVV2	1011557	1335663	83.39%	4.329%
83	13.372	2008	2015	2018	rVV2	154739	272357	17.00%	0.883%
84	13.420	2020	2023	2033	rVB3	95510	153587	9.59%	0.498%
85	13.506	2033	2037	2039	rBV2	102960	134587	8.40%	0.436%
86	13.542	2039	2043	2046	rVV	301942	404585	25.26%	1.311%
87	13.579	2046	2049	2053	rVV3	477035	700352	43.73%	2.270%
88	13.664	2053	2063	2069	rVB3	481087	1063968	66.43%	3.448%
89	13.774	2074	2081	2090	rVV2	184039	376800	23.52%	1.221%
90	13.853	2090	2094	2099	rVB	69814	104395	6.52%	0.338%
91	13.926	2099	2106	2110	rBV2	171542	264312	16.50%	0.857%
92	13.969	2110	2113	2117	rVV	142675	175230	10.94%	0.568%
93	14.006	2117	2119	2126	rVB2	49384	60463	3.77%	0.196%
94	14.073	2126	2130	2135	rBV	172567	217873	13.60%	0.706%
95	14.213	2148	2153	2163	rVV	282719	456366	28.49%	1.479%
96	14.323	2167	2171	2175	rVV	120980	169766	10.60%	0.550%
97	14.371	2175	2179	2183	rVV	160223	197762	12.35%	0.641%
98	14.481	2192	2197	2202	rBV2	78801	122577	7.65%	0.397%
99	14.634	2214	2222	2226	rVV	259676	378248	23.62%	1.226%
100	14.780	2241	2246	2251	rBV	284991	373917	23.34%	1.212%

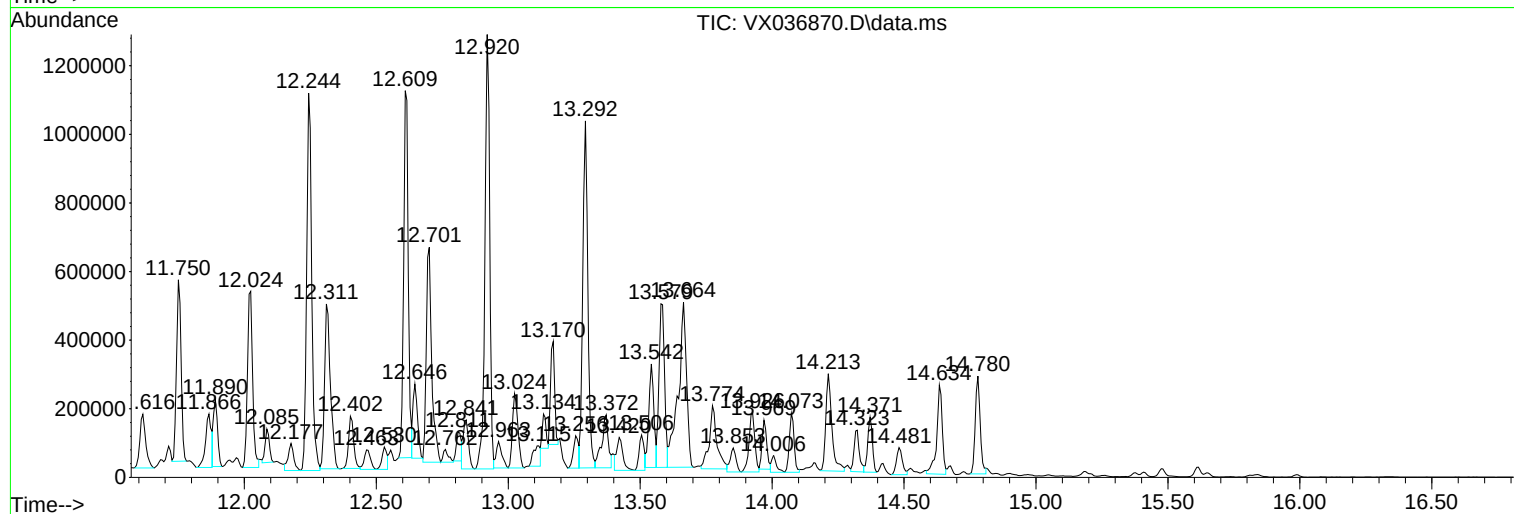
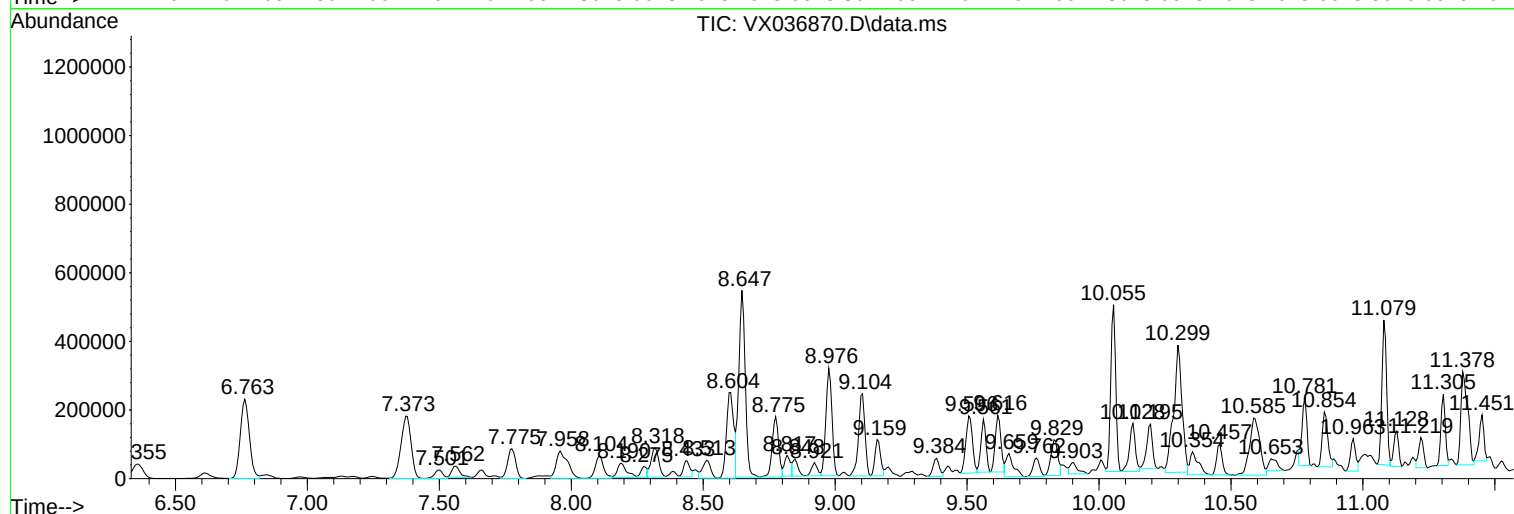
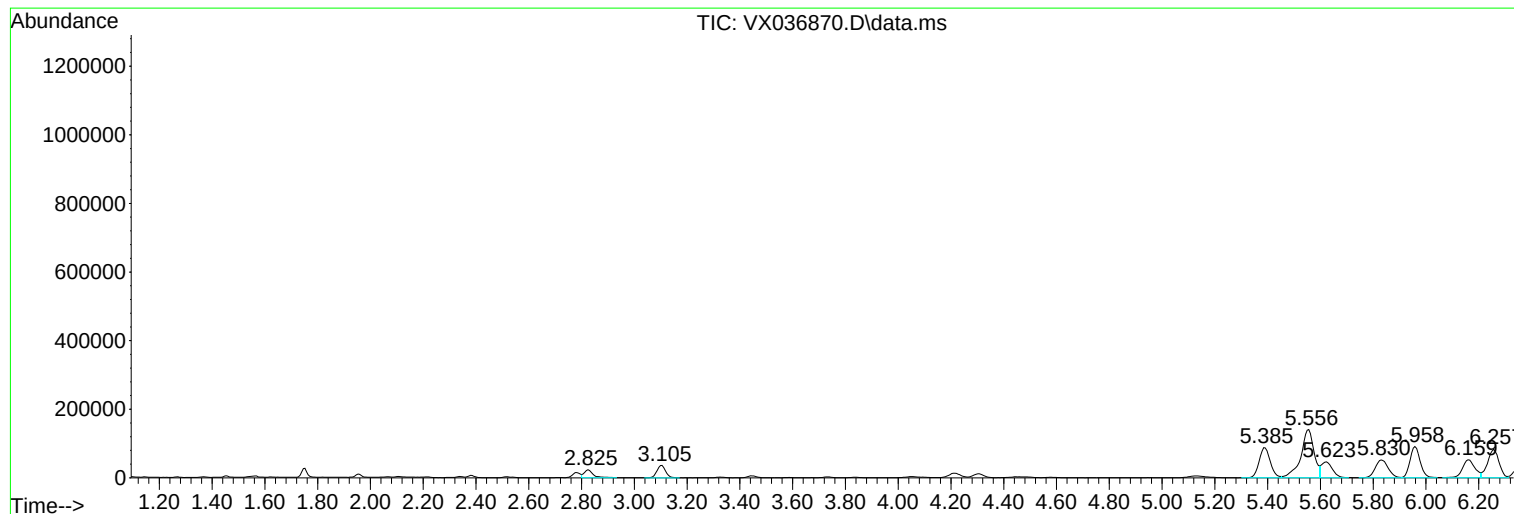
Sum of corrected areas: 30856691

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Instrument :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
Quant Title : SW846 8260

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



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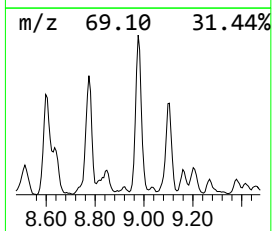
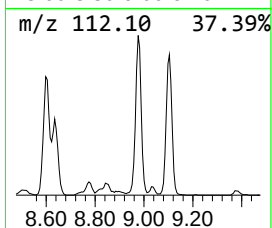
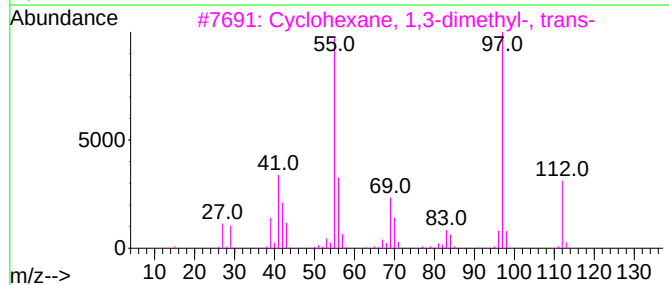
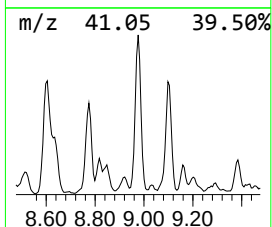
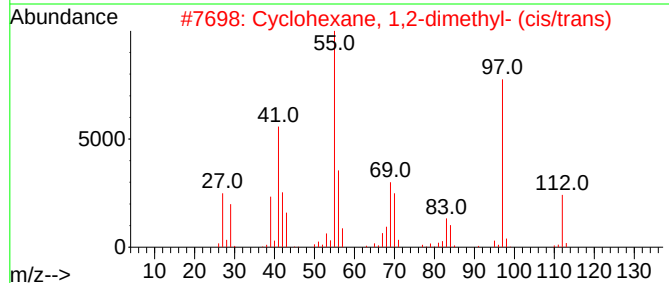
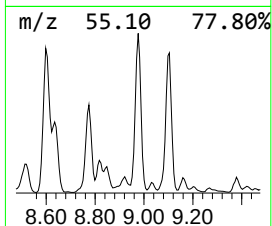
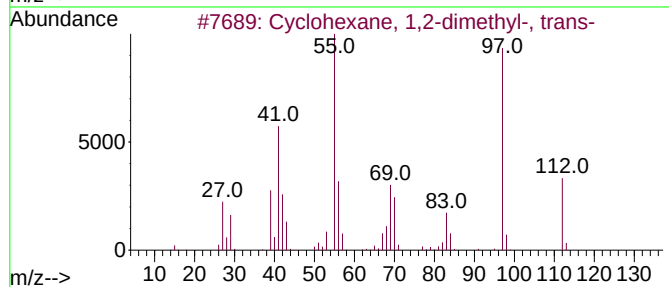
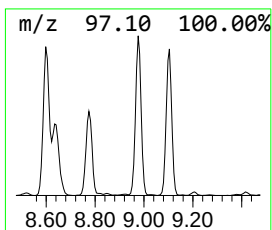
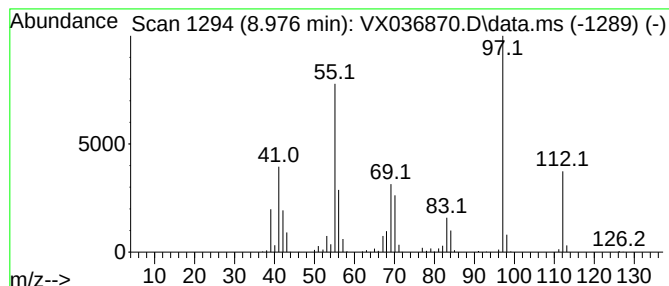
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
Quant Title : SW846 8260

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

Peak Number 1 Cyclohexane, 1,2-dimethyl-,... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.976	39.10 ug/l	497840	Chlorobenzene-d5	10.055

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,2-dimethyl-, trans-	112	C8H16	006876-23-9	96
2			Cyclohexane, 1,2-dimethyl- (cis/...	112	C8H16	000583-57-3	91
3			Cyclohexane, 1,3-dimethyl-, trans-	112	C8H16	002207-03-6	90
4			Cyclohexane, 1,4-dimethyl-, trans-	112	C8H16	002207-04-7	87
5			Cyclohexane, 1,4-dimethyl-	112	C8H16	000589-90-2	81



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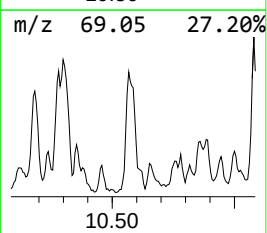
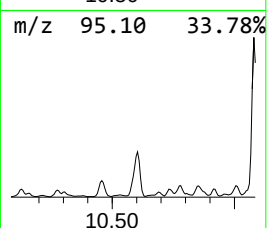
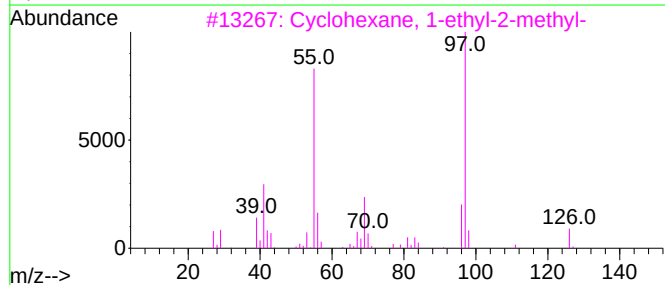
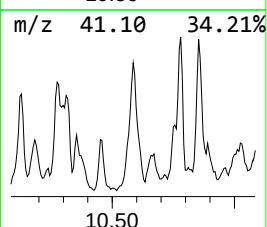
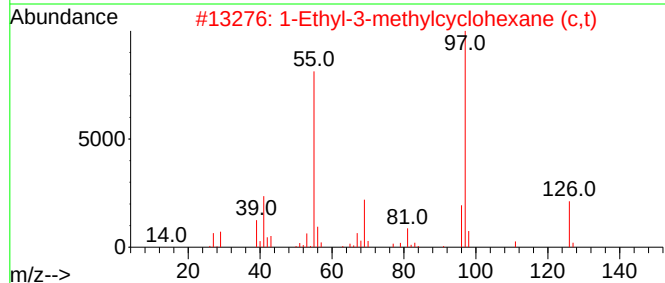
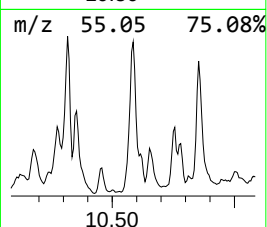
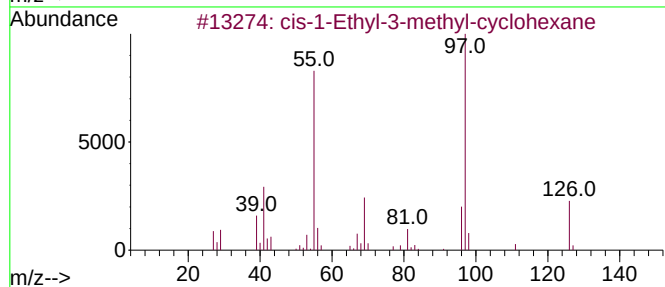
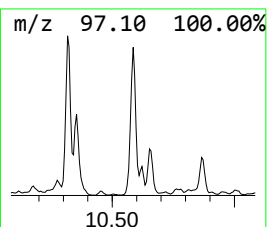
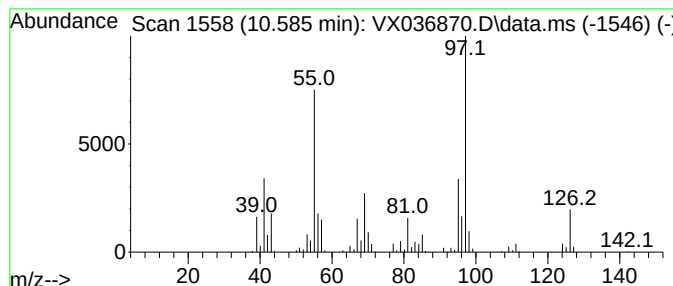
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Quant Title : SW846 8260

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

Peak Number 2 cis-1-Ethyl-3-methyl-cyclo... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.586	35.51 ug/l	452145	Chlorobenzene-d5	10.055

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			cis-1-Ethyl-3-methyl-cyclohexane	126	C9H18	019489-10-2	93
2			1-Ethyl-3-methylcyclohexane (c,t)	126	C9H18	003728-55-0	81
3			Cyclohexane, 1-ethyl-2-methyl-	126	C9H18	003728-54-9	76
4			1-Ethyl-4-methylcyclohexane	126	C9H18	003728-56-1	74
5			Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	006236-88-0	72



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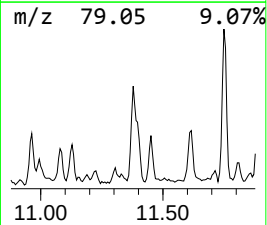
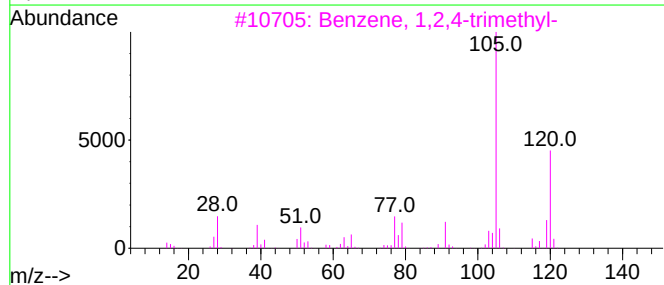
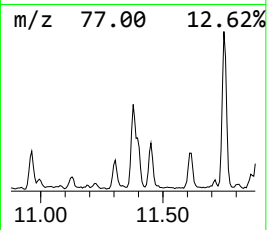
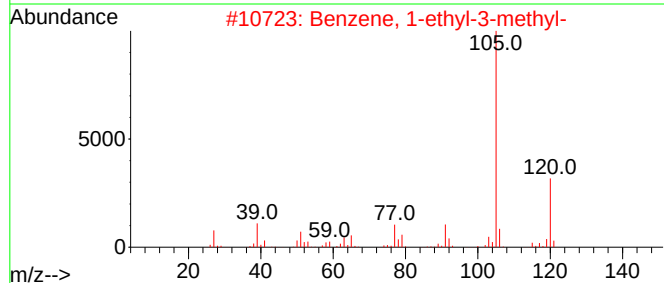
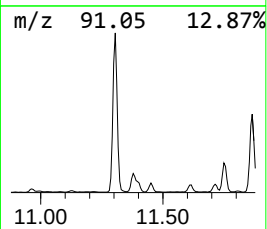
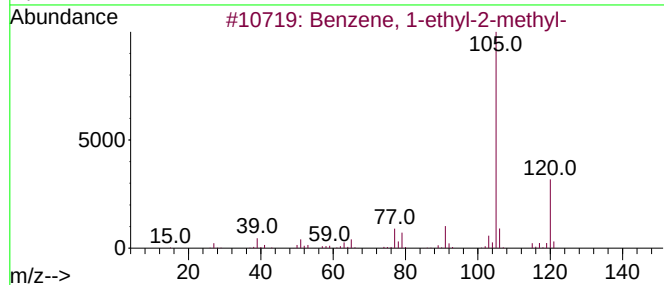
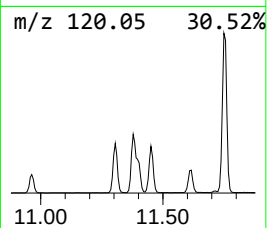
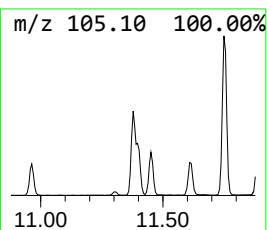
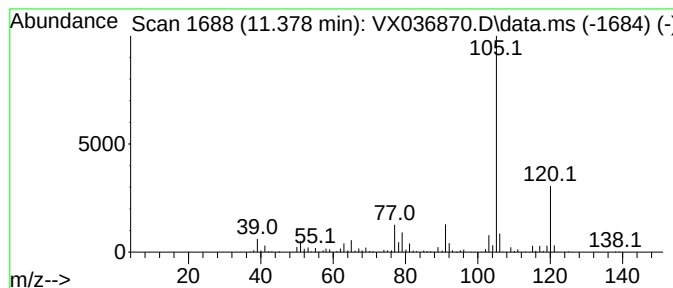
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Quant Title : SW846 8260

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

Peak Number 3 Benzene, 1-ethyl-2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.378	35.40 ug/l	498575	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	95
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4			Mesitylene	120	C9H12	000108-67-8	91
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX080823\
Data File : VX036870.D
Acq On : 08 Aug 2023 17:39
Operator : JC/MD
Sample : 03903-20
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-10W

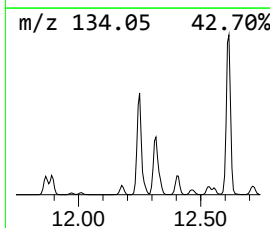
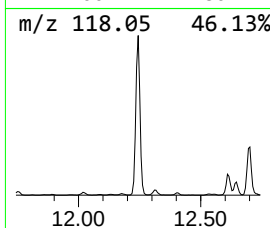
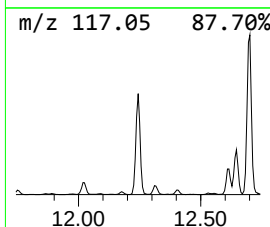
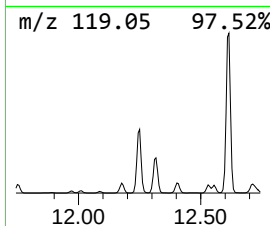
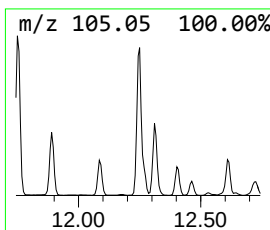
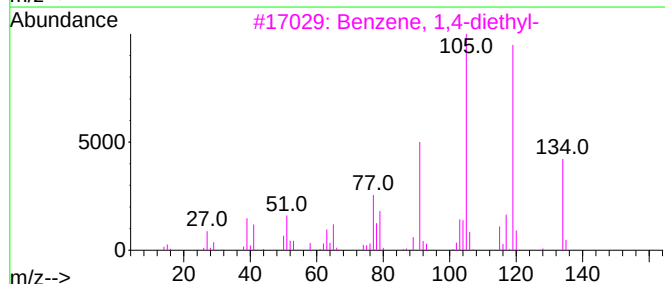
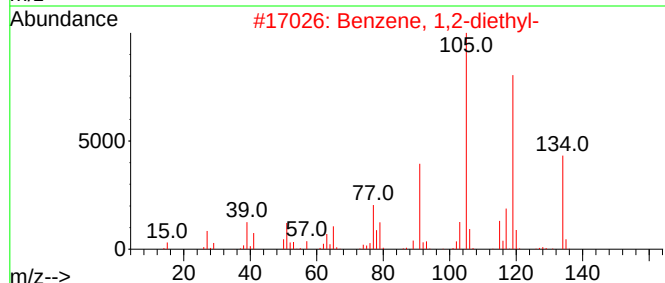
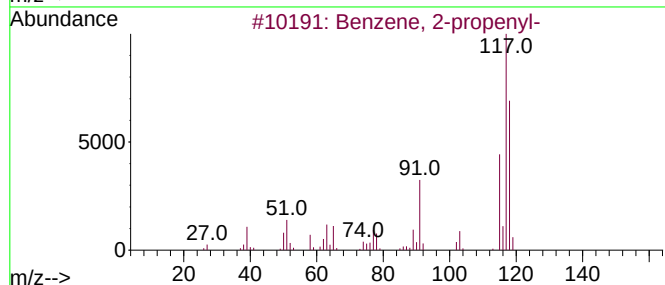
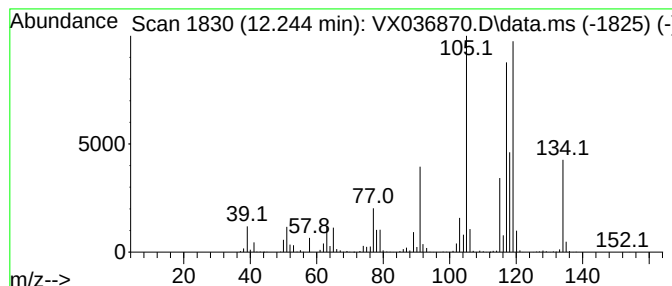
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 2-propenyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	104.19 ug/l	1467490	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, 1,2-diethyl-	134	C10H14	000135-01-3	78
3			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	60
4			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	55
5			Benzene, cyclopropyl-	118	C9H10	000873-49-4	46



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX080823\
Data File : VX036870.D
Acq On : 08 Aug 2023 17:39
Operator : JC/MD
Sample : 03903-20
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-10W

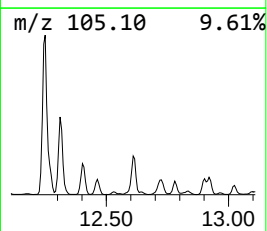
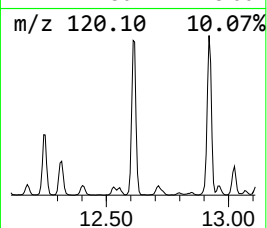
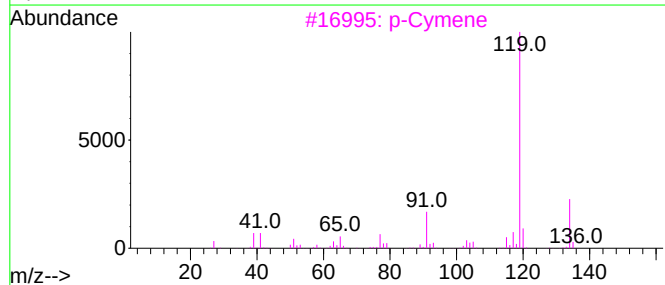
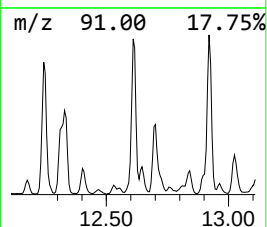
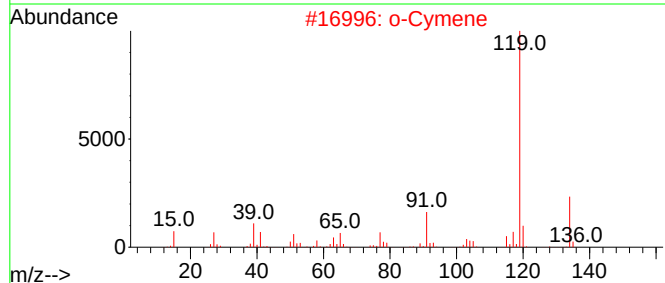
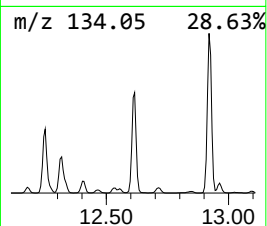
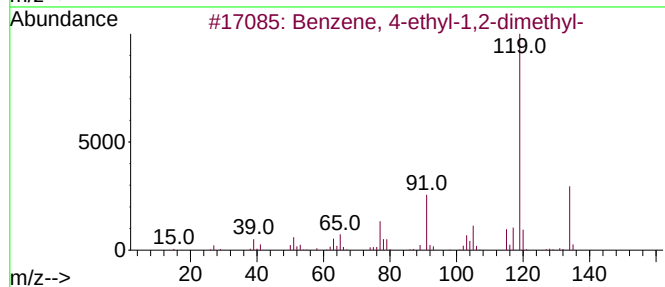
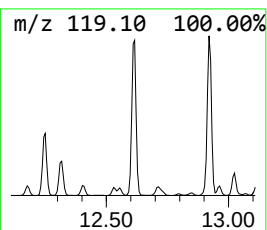
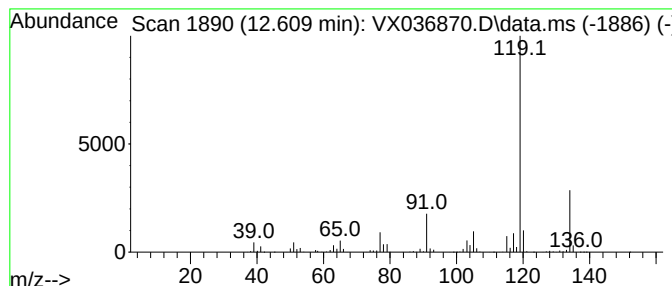
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.610	93.20 ug/l	1312690	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2			o-Cymene	134	C10H14	000527-84-4	95
3			p-Cymene	134	C10H14	000099-87-6	95
4			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX080823\
Data File : VX036870.D
Acq On : 08 Aug 2023 17:39
Operator : JC/MD
Sample : 03903-20
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-10W

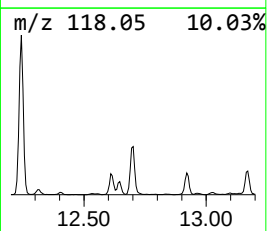
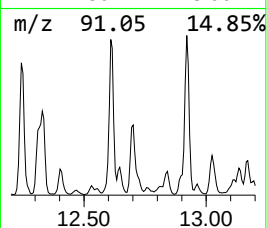
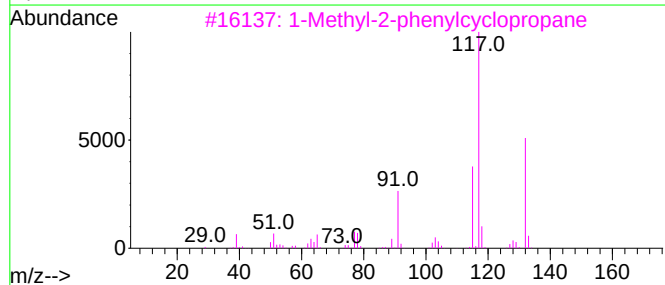
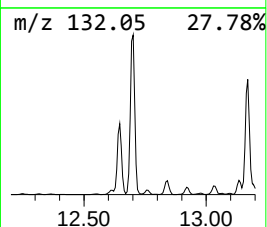
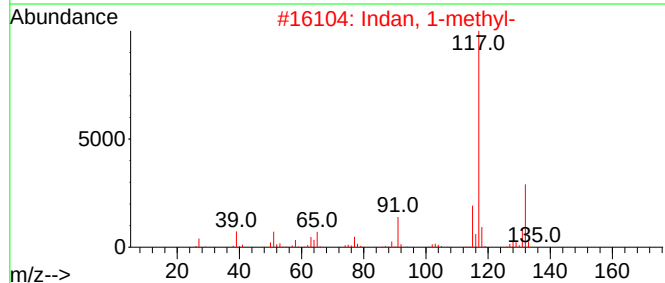
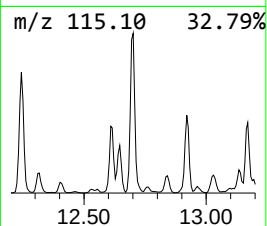
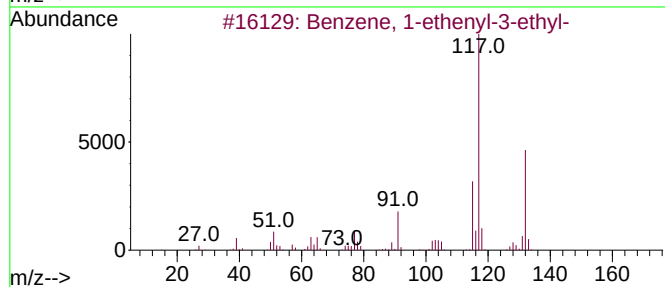
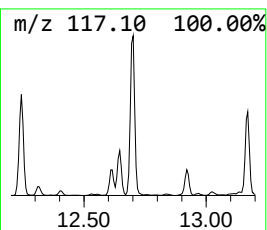
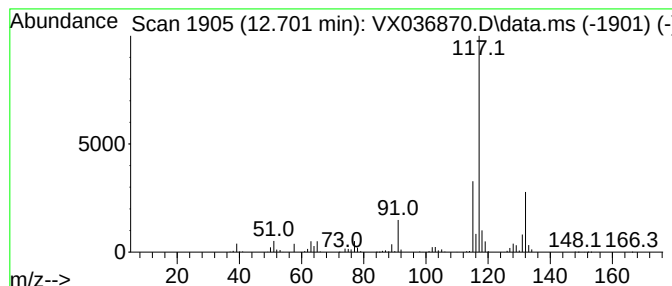
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	63.98 ug/l	901130	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90
2			Indan, 1-methyl-	132	C10H12	000767-58-8	87
3			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	87
4			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	83
5			Benzene, 2-butenyl-	132	C10H12	001560-06-1	83



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX080823\
Data File : VX036870.D
Acq On : 08 Aug 2023 17:39
Operator : JC/MD
Sample : 03903-20
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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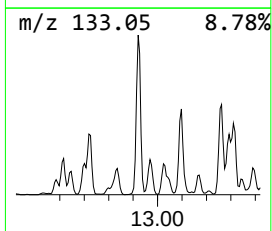
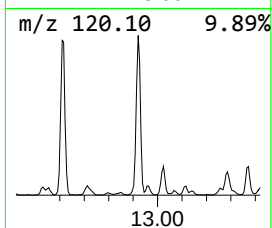
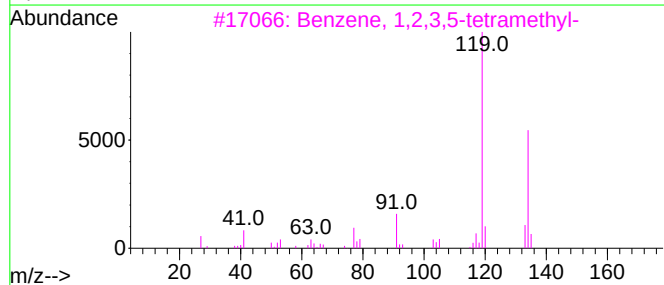
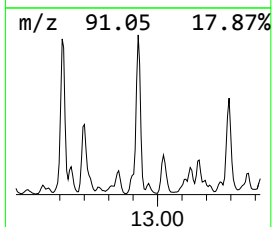
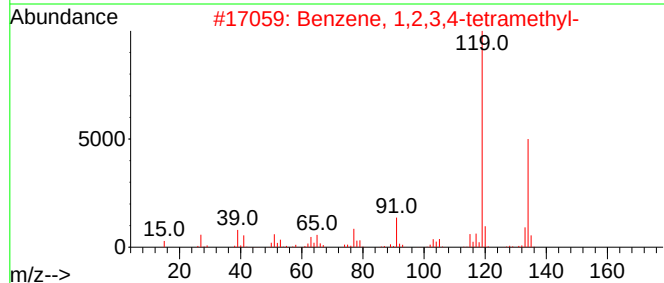
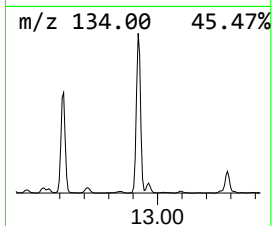
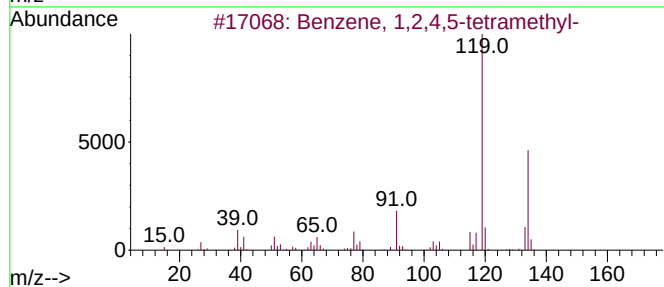
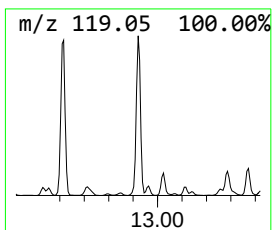
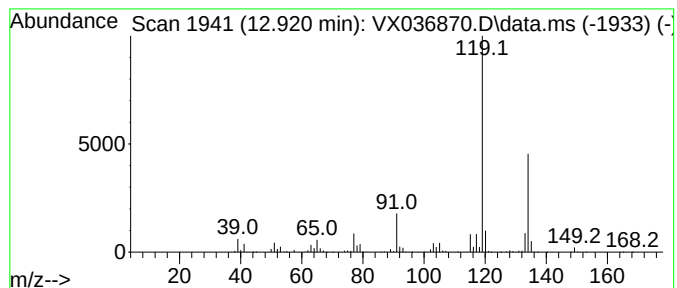
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.920	113.72 ug/l	1601700	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
4			o-Cymene	134	C10H14	000527-84-4	95
5			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX080823\
Data File : VX036870.D
Acq On : 08 Aug 2023 17:39
Operator : JC/MD
Sample : 03903-20
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-10W

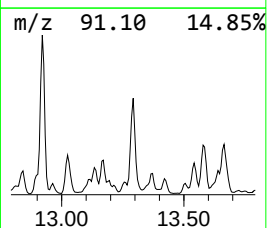
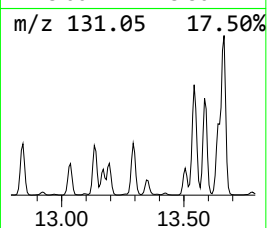
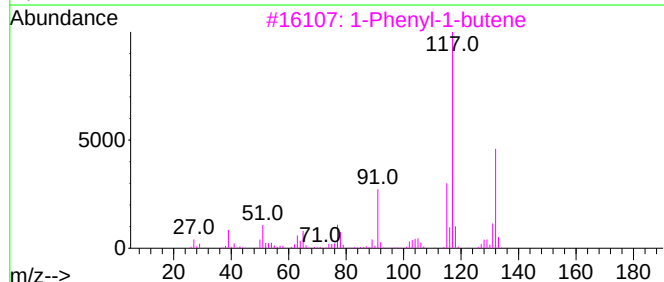
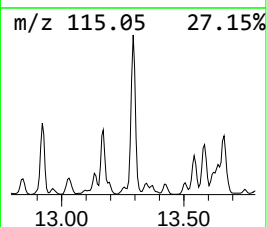
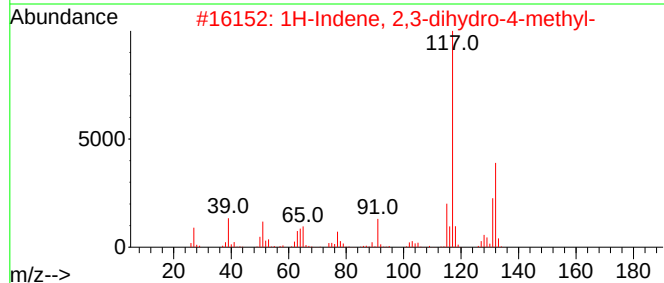
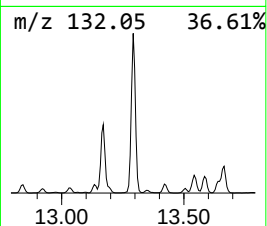
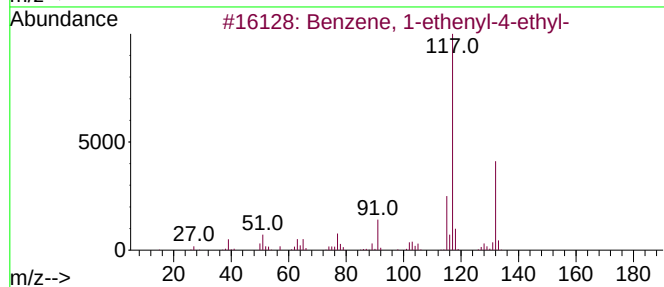
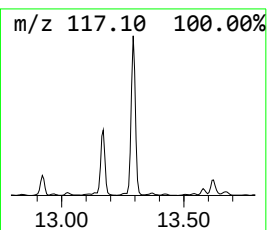
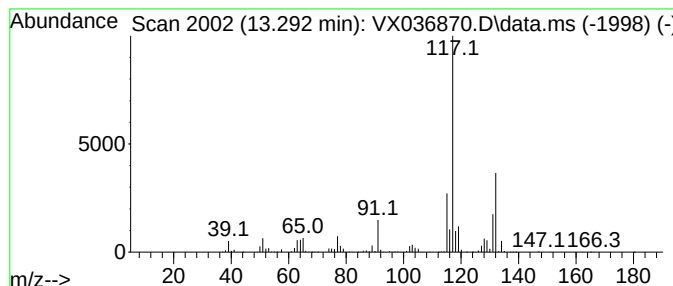
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	94.83 ug/l	1335660	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	90
2			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	87
3			1-Phenyl-1-butene	132	C10H12	000824-90-8	87
4			3-Phenylbut-1-ene	132	C10H12	000934-10-1	87
5			Indan, 1-methyl-	132	C10H12	000767-58-8	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX080823\
Data File : VX036870.D
Acq On : 08 Aug 2023 17:39
Operator : JC/MD
Sample : 03903-20
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-10W

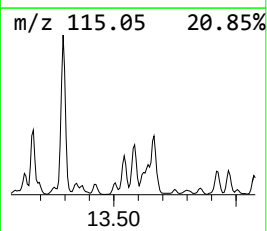
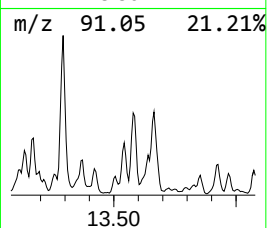
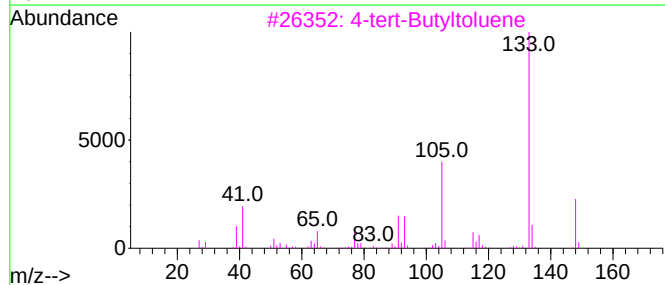
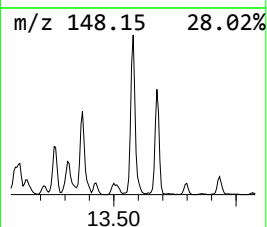
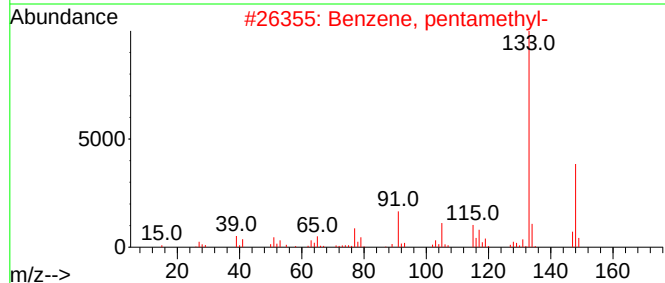
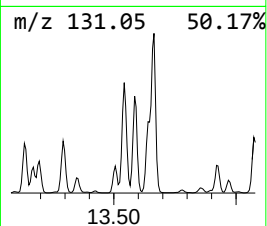
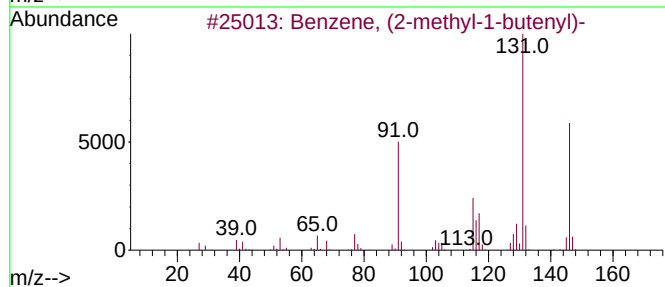
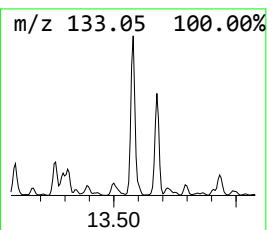
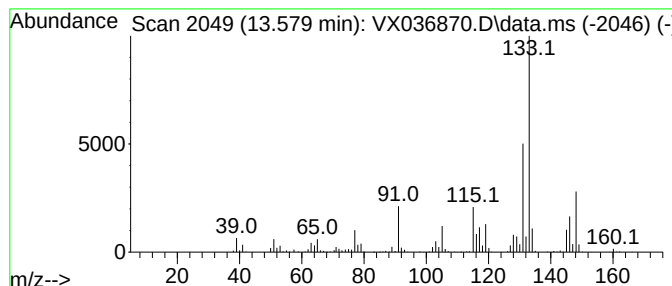
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, (2-methyl-1-butenyl)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.579	49.72 ug/l	700352	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	70
2			Benzene, pentamethyl-	148	C11H16	000700-12-9	64
3			4-tert-Butyltoluene	148	C11H16	000098-51-1	53
4			1,5,6,7-Tetramethylbicyclo[3.2.0]...	148	C11H16	134329-46-7	49
5			Ethanone, 1-(4-ethylphenyl)-	148	C10H12O	000937-30-4	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX080823\
Data File : VX036870.D
Acq On : 08 Aug 2023 17:39
Operator : JC/MD
Sample : 03903-20
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-10W

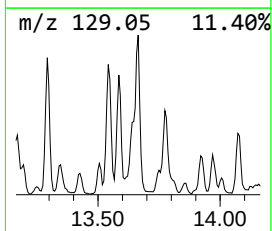
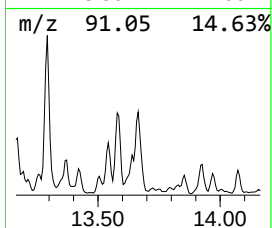
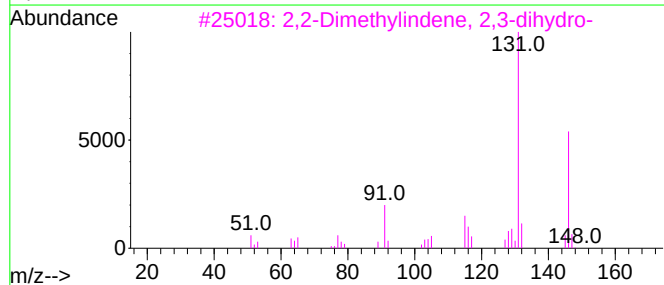
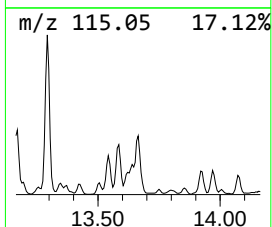
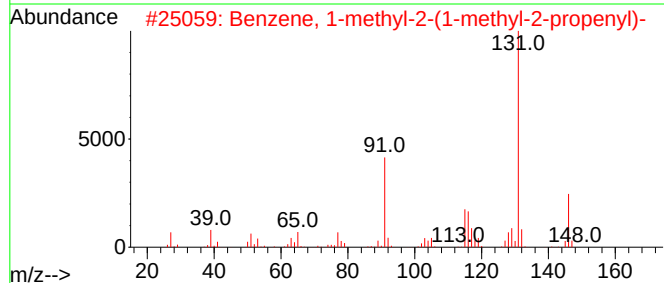
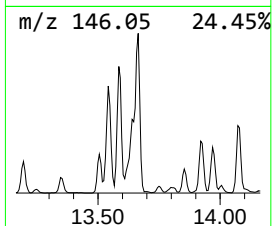
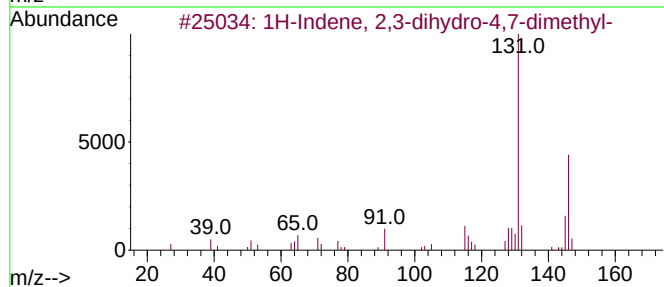
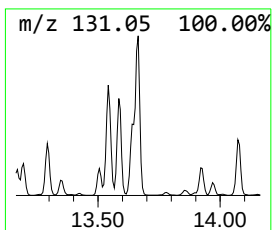
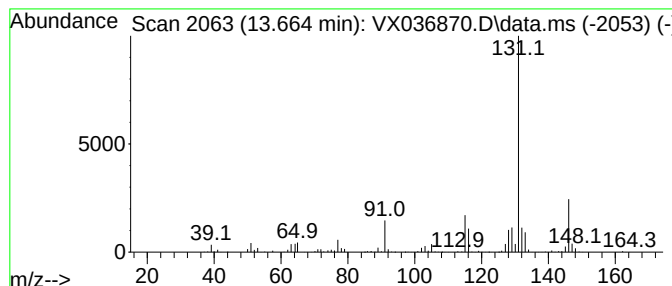
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
Quant Title : SW846 8260

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

Peak Number 10 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.664	75.54 ug/l	1063970	1,4-Dichlorobenzene-d4	12.024

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14		006682-71-9	90
2	Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14		097664-19-2	90
3	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14		020836-11-7	90
4	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14		017057-82-8	90
5	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14		017059-48-2	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX080823\
Data File : VX036870.D
Acq On : 08 Aug 2023 17:39
Operator : JC/MD
Sample : 03903-20
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 20 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-10W

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
Quant Title : SW846 8260

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Cyclohexane, 1,...	8.976	39.1	ug/l	497840	3	10.055	636598	50.0
cis-1-Ethyl-3-m...	10.586	35.5	ug/l	452145	3	10.055	636598	50.0
Benzene, 1-ethy...	11.378	35.4	ug/l	498575	4	12.024	704260	50.0
Benzene, 2-prop...	12.244	104.2	ug/l	1467490	4	12.024	704260	50.0
Benzene, 4-ethy...	12.610	93.2	ug/l	1312690	4	12.024	704260	50.0
Benzene, 1-ethe...	12.701	64.0	ug/l	901130	4	12.024	704260	50.0
Benzene, 1,2,4,...	12.920	113.7	ug/l	1601700	4	12.024	704260	50.0
Benzene, 1-ethe...	13.292	94.8	ug/l	1335660	4	12.024	704260	50.0
Benzene, (2-met...	13.579	49.7	ug/l	700352	4	12.024	704260	50.0
1H-Indene, 2,3-...	13.664	75.5	ug/l	1063970	4	12.024	704260	50.0