

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX080723\
 Data File : VX036849.D
 Acq On : 07 Aug 2023 20:05
 Operator : JC/MD
 Sample : 03903-12
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-01W

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: VX036849.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.782	270	278	283	rBV	32887	76819	5.01%	0.251%
2	3.099	320	330	344	rBV	74877	172835	11.28%	0.564%
3	4.214	498	513	520	rVV2	18645	58544	3.82%	0.191%
4	4.306	520	528	539	rVV2	25873	73982	4.83%	0.241%
5	5.391	693	706	715	rBV	89558	261254	17.05%	0.852%
6	5.556	716	733	739	rVV	142014	480903	31.39%	1.568%
7	5.623	740	744	763	rVV	61958	173163	11.30%	0.565%
8	5.830	765	778	790	rVV4	41595	136990	8.94%	0.447%
9	5.958	790	799	811	rVV	93244	244079	15.93%	0.796%
10	6.159	814	832	840	rVV3	65749	219928	14.36%	0.717%
11	6.257	840	848	857	rVV3	108525	324508	21.18%	1.058%
12	6.360	857	865	876	rVB3	45672	131207	8.56%	0.428%
13	6.763	921	931	941	rBV	232794	572673	37.38%	1.868%
14	7.177	991	999	1005	rVB	25864	55962	3.65%	0.183%
15	7.372	1019	1031	1042	rVB3	130850	364030	23.76%	1.187%
16	7.500	1042	1052	1057	rBV3	28700	63410	4.14%	0.207%
17	7.561	1057	1062	1072	rVB2	32346	67982	4.44%	0.222%
18	7.775	1091	1097	1105	rVB	73419	146958	9.59%	0.479%
19	7.958	1122	1127	1129	rBV2	55034	98270	6.41%	0.321%
20	8.104	1143	1151	1159	rBV	99869	208289	13.60%	0.679%
21	8.189	1159	1165	1175	rVV4	47324	126533	8.26%	0.413%
22	8.311	1175	1185	1193	rVV4	98908	245776	16.04%	0.802%
23	8.439	1201	1206	1209	rBV	49070	78266	5.11%	0.255%
24	8.512	1213	1218	1225	rVB2	52999	115826	7.56%	0.378%
25	8.604	1225	1233	1236	rBV2	289603	550318	35.92%	1.795%
26	8.646	1236	1240	1251	rVB	555588	952111	62.15%	3.105%
27	8.775	1251	1261	1265	rBV	172673	301004	19.65%	0.982%
28	8.817	1265	1268	1271	rVV	59619	95017	6.20%	0.310%
29	8.848	1271	1273	1278	rVV3	47498	61298	4.00%	0.200%
30	8.921	1280	1285	1289	rVV	39634	64383	4.20%	0.210%
31	8.976	1289	1294	1300	rVV	293165	446015	29.11%	1.455%
32	9.104	1307	1315	1320	rBV	269909	455157	29.71%	1.485%
33	9.159	1320	1324	1329	rBV	191910	296093	19.33%	0.966%
34	9.384	1356	1361	1364	rBV2	42537	65760	4.29%	0.214%
35	9.427	1364	1368	1371	rBV	62694	87438	5.71%	0.285%
36	9.506	1376	1381	1386	rVB2	167921	285647	18.65%	0.932%

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37	9.561	1386	1390	1395	rVB	157388	215550	14.07%	0.703%
38	9.616	1395	1399	1403	rBV	118143	183613	11.99%	0.599%
39	9.659	1403	1406	1417	rVB2	63790	130166	8.50%	0.425%
40	9.762	1417	1423	1429	rBV3	76267	136048	8.88%	0.444%
41	9.835	1429	1435	1439	rBV3	70227	135936	8.87%	0.443%
42	10.055	1466	1471	1476	rVB	491952	639050	41.72%	2.084%
43	10.128	1476	1483	1486	rBV	187513	283325	18.49%	0.924%
44	10.274	1502	1507	1511	rBV2	131180	238491	15.57%	0.778%
45	10.317	1512	1514	1517	rVV	95681	110988	7.25%	0.362%
46	10.354	1517	1520	1531	rVB4	55121	123540	8.06%	0.403%
47	10.457	1531	1537	1545	rVB2	95612	159551	10.42%	0.520%
48	10.591	1550	1559	1566	rBV3	141335	378939	24.74%	1.236%
49	10.658	1566	1570	1579	rVB3	32914	69300	4.52%	0.226%
50	10.780	1579	1590	1593	rBV2	197325	359637	23.48%	1.173%
51	10.853	1598	1602	1609	rBV3	149715	238769	15.59%	0.779%
52	10.963	1615	1620	1623	rBV	143993	196153	12.80%	0.640%
53	11.079	1634	1639	1643	rBV	404077	527284	34.42%	1.720%
54	11.128	1643	1647	1651	rVB	132558	169985	11.10%	0.554%
55	11.219	1659	1662	1667	rVB2	71648	94818	6.19%	0.309%
56	11.305	1667	1676	1679	rBV	122228	151020	9.86%	0.493%
57	11.396	1686	1691	1695	rBV3	42093	69792	4.56%	0.228%
58	11.615	1722	1727	1734	rVB	136544	190627	12.44%	0.622%
59	11.750	1746	1749	1754	rVV	79449	107583	7.02%	0.351%
60	11.865	1761	1768	1770	rBV	276382	399645	26.09%	1.303%
61	11.890	1770	1772	1777	rVB	360337	375527	24.51%	1.225%
62	12.018	1789	1793	1797	rBV	513553	661276	43.17%	2.157%
63	12.054	1797	1799	1803	rVV	85240	113216	7.39%	0.369%
64	12.176	1815	1819	1824	rVB	100942	122441	7.99%	0.399%
65	12.243	1825	1830	1836	rBV	1201932	1488793	97.19%	4.856%
66	12.310	1836	1841	1850	rVB2	606044	916829	59.85%	2.990%
67	12.402	1851	1856	1862	rVB	259681	326157	21.29%	1.064%
68	12.469	1862	1867	1872	rVB2	39796	61104	3.99%	0.199%
69	12.609	1883	1890	1894	rBV	1235129	1531908	100.00%	4.996%
70	12.646	1894	1896	1900	rVB2	146885	175191	11.44%	0.571%
71	12.695	1900	1904	1912	rVB	982267	1397231	91.21%	4.557%
72	12.762	1912	1915	1917	rBV	49088	63162	4.12%	0.206%
73	12.841	1924	1928	1933	rVB	203567	287862	18.79%	0.939%
74	12.920	1933	1941	1945	rBV	976050	1239992	80.94%	4.044%
75	12.963	1945	1948	1954	rVV2	59722	96645	6.31%	0.315%
76	13.024	1954	1958	1965	rVB3	235398	362622	23.67%	1.183%

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77	13.097	1965	1970	1971	rBV	83712	104723	6.84%	0.342%
78	13.134	1973	1976	1979	rVV	236631	324841	21.20%	1.059%
79	13.170	1979	1982	1985	rVV	396633	534629	34.90%	1.744%
80	13.194	1985	1986	1991	rVV2	110968	92046	6.01%	0.300%
81	13.255	1991	1996	1998	rVV2	122305	165583	10.81%	0.540%
82	13.292	1998	2002	2008	rVV2	630617	920103	60.06%	3.001%
83	13.365	2008	2014	2018	rVV2	160428	335035	21.87%	1.093%
84	13.426	2020	2024	2032	rVB2	73773	118799	7.75%	0.387%
85	13.505	2032	2037	2039	rBV	151662	192982	12.60%	0.629%
86	13.542	2039	2043	2046	rVV	483611	677959	44.26%	2.211%
87	13.579	2046	2049	2053	rVV3	534451	777000	50.72%	2.534%
88	13.664	2053	2063	2070	rVV3	660375	1484396	96.90%	4.841%
89	13.749	2073	2077	2079	rVV	49010	59069	3.86%	0.193%
90	13.792	2079	2084	2090	rVV3	44830	105673	6.90%	0.345%
91	13.853	2090	2094	2098	rVB	77154	104057	6.79%	0.339%
92	13.920	2099	2105	2110	rBV2	203652	307724	20.09%	1.004%
93	13.969	2110	2113	2117	rVV	110717	132290	8.64%	0.431%
94	14.072	2126	2130	2134	rBV	237731	284413	18.57%	0.928%
95	14.213	2148	2153	2158	rBV	332018	471827	30.80%	1.539%
96	14.316	2167	2170	2176	rBV	130582	179115	11.69%	0.584%
97	14.371	2176	2179	2183	rVB	181717	214754	14.02%	0.700%
98	14.481	2192	2197	2201	rBV2	69129	111382	7.27%	0.363%
99	14.633	2219	2222	2226	rVV	101577	140111	9.15%	0.457%
100	14.780	2240	2246	2250	rBV2	93481	133835	8.74%	0.437%

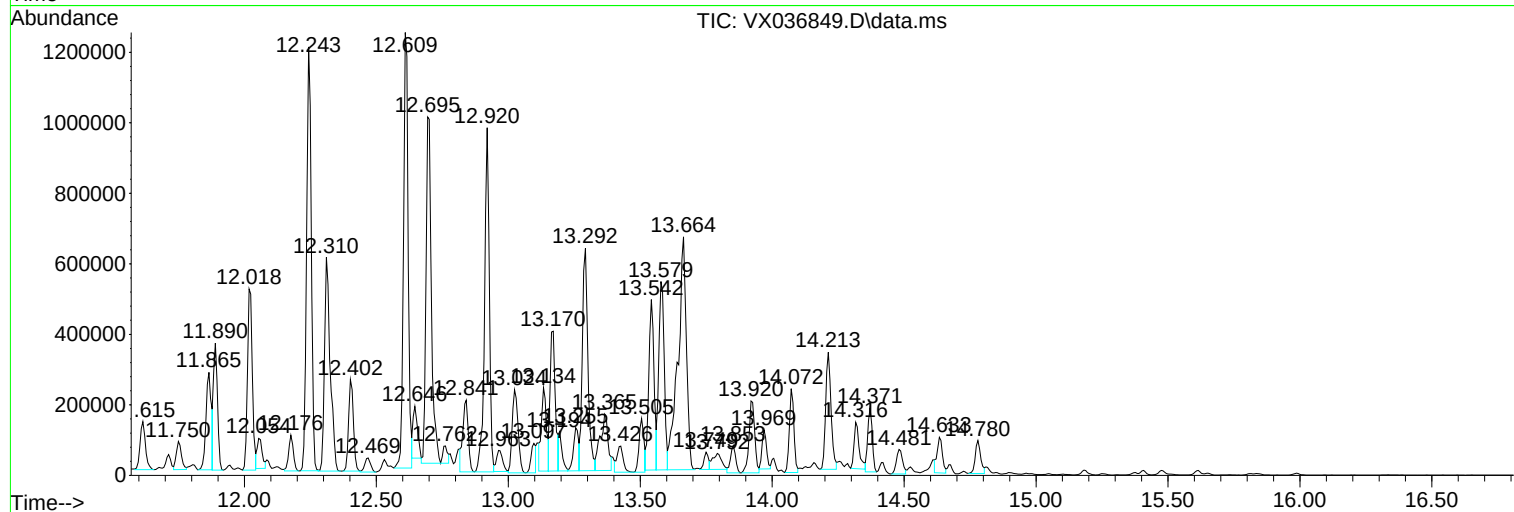
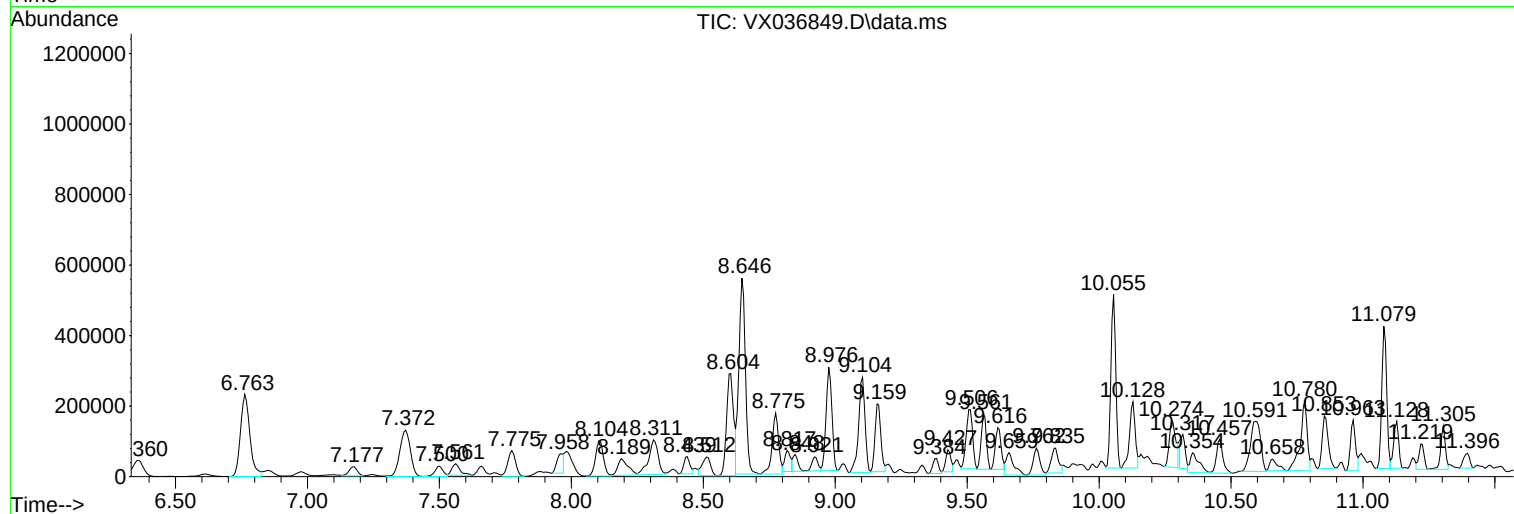
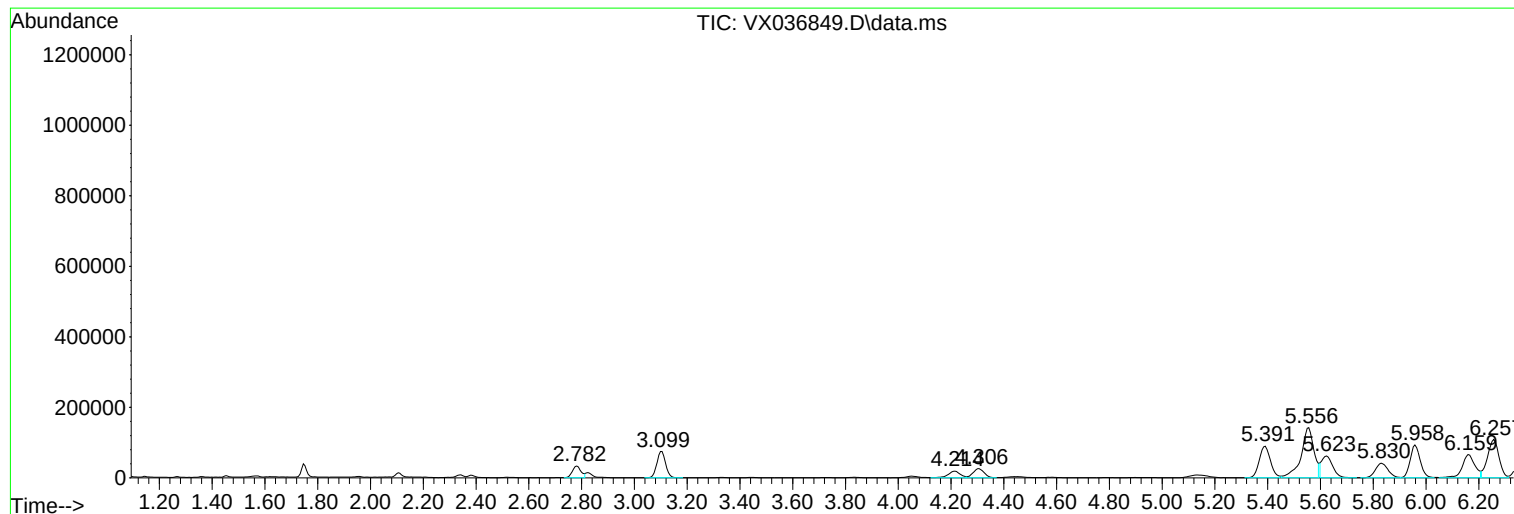
Sum of corrected areas: 30660540

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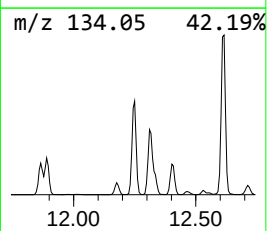
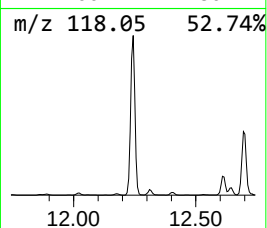
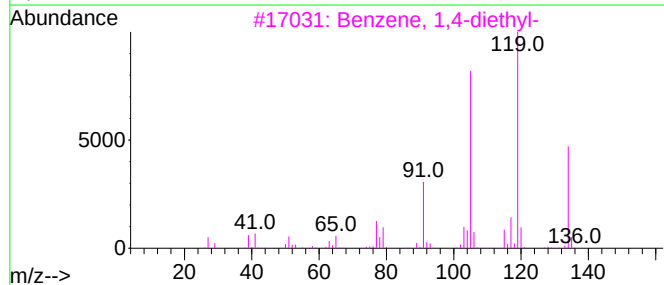
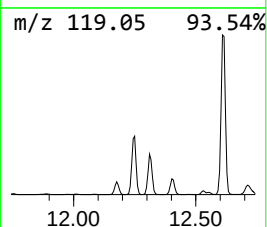
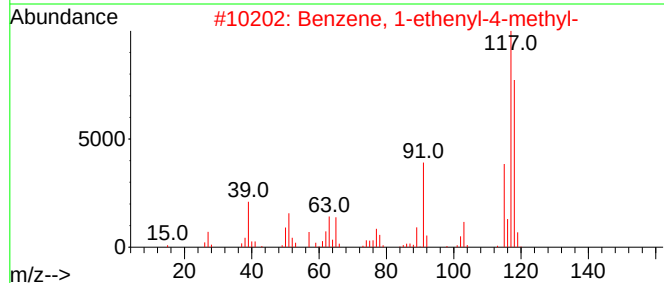
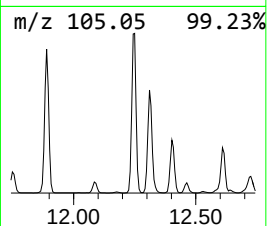
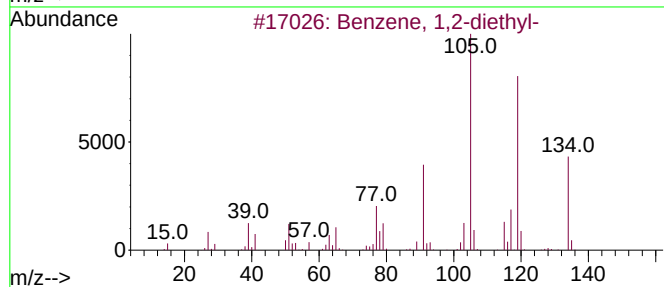
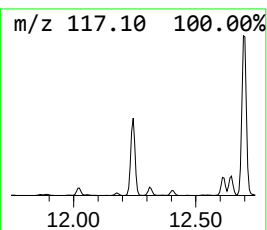
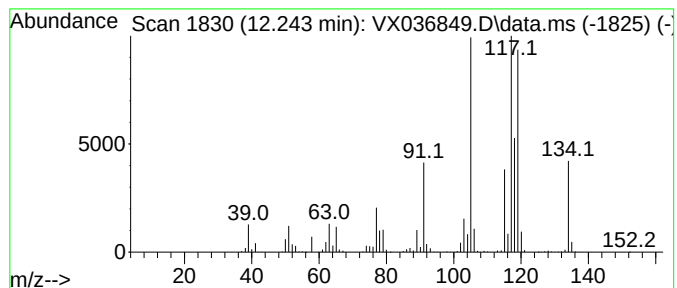
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.243	112.57 ug/l	1488790	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	87
2			Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	60
3			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	55
4			(Z)-1-Phenylpropene	118	C9H10	000766-90-5	55
5			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	50



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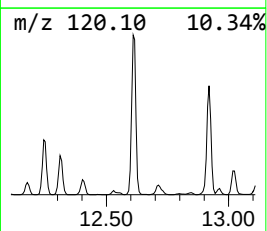
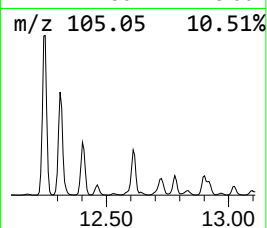
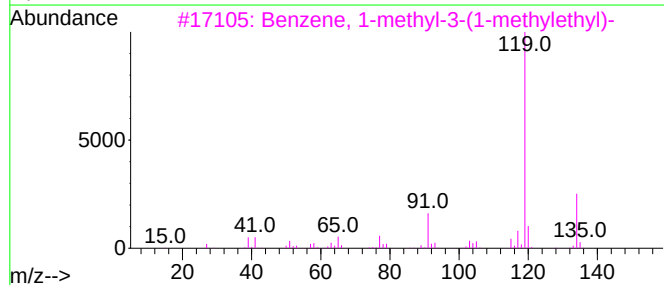
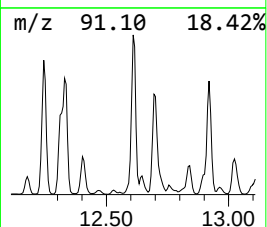
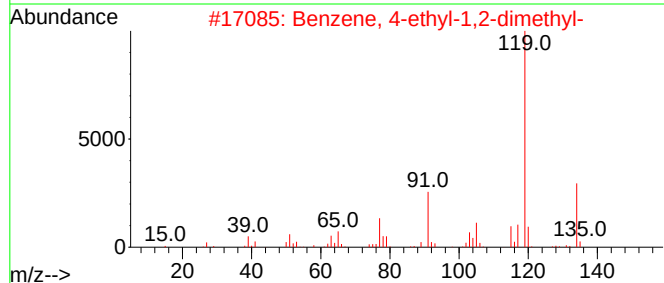
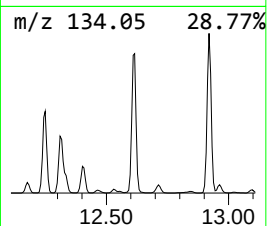
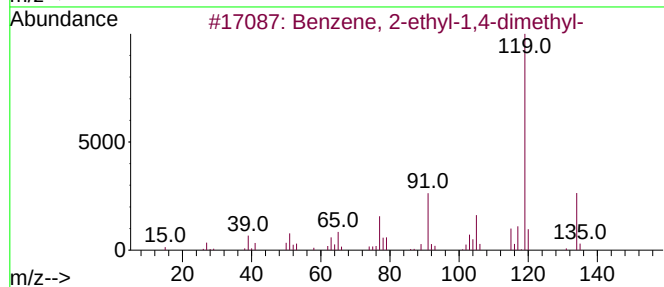
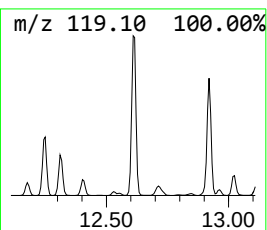
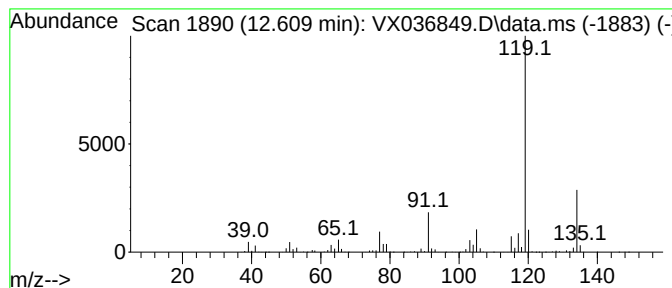
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.609	115.83 ug/l	1531910	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
3			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
4			p-Cymene	134	C10H14	000099-87-6	95
5			o-Cymene	134	C10H14	000527-84-4	95



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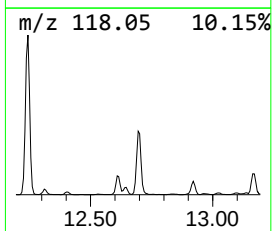
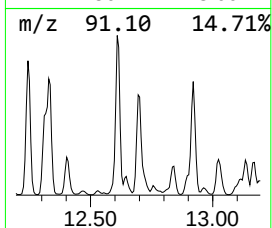
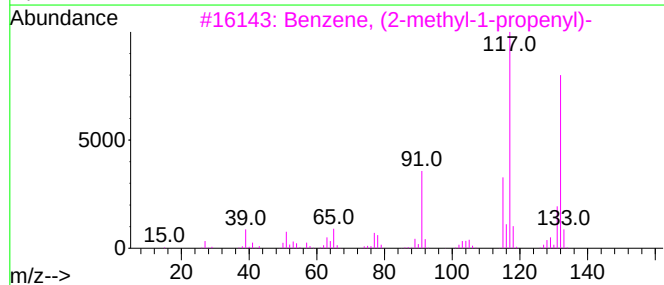
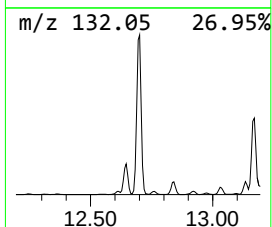
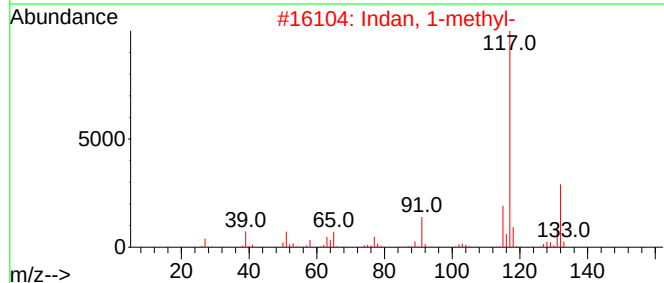
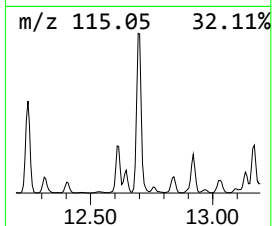
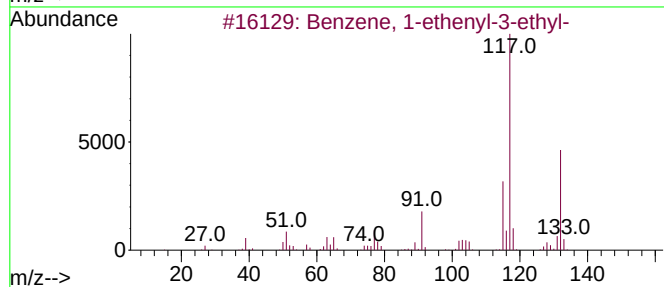
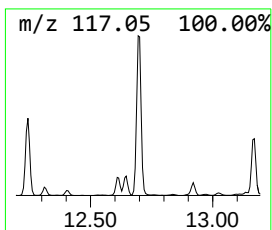
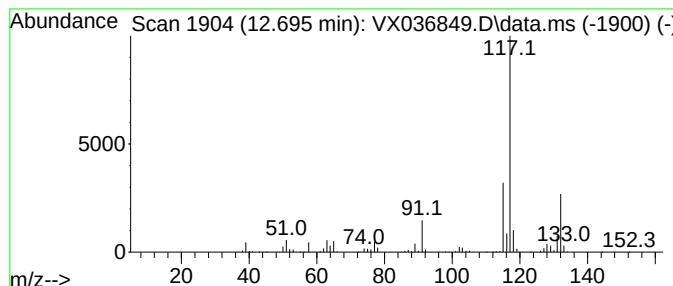
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.695	105.65 ug/l	1397230	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	91
2		Indan, 1-methyl-	132	C10H12	000767-58-8	90
3		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	90
4		1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	87
5		1-Phenyl-1-butene	132	C10H12	000824-90-8	87



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX080723\
Data File : VX036849.D
Acq On : 07 Aug 2023 20:05
Operator : JC/MD
Sample : 03903-12
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-01W

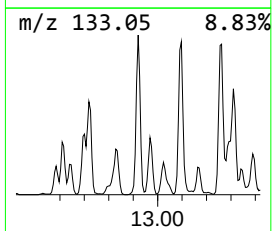
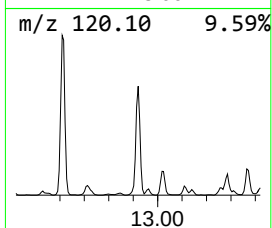
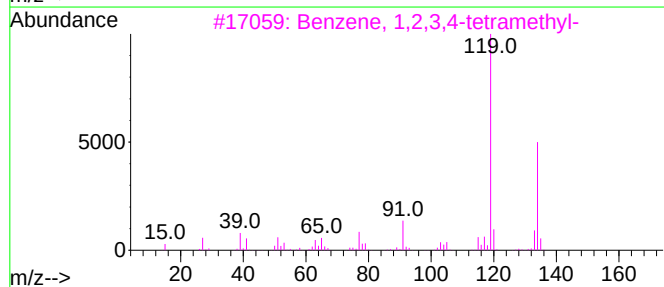
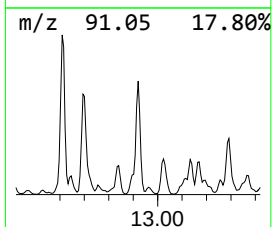
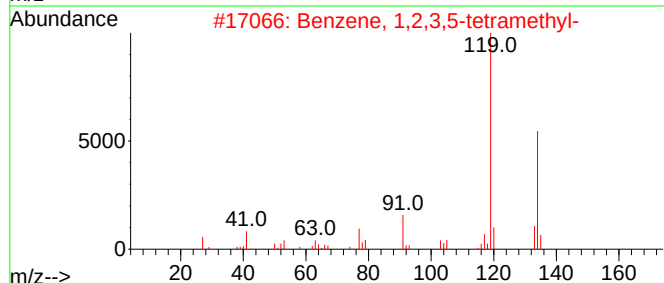
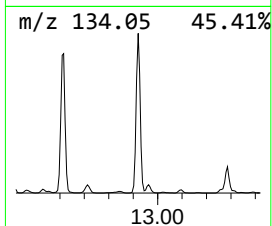
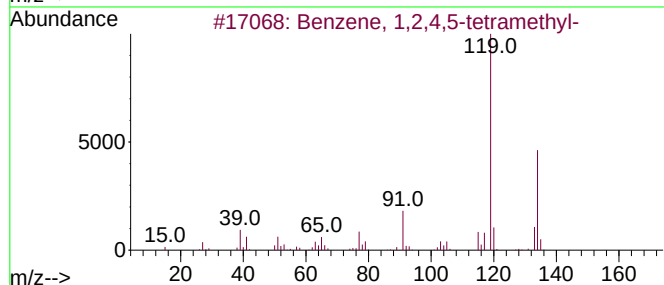
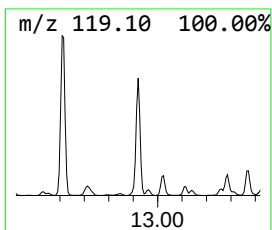
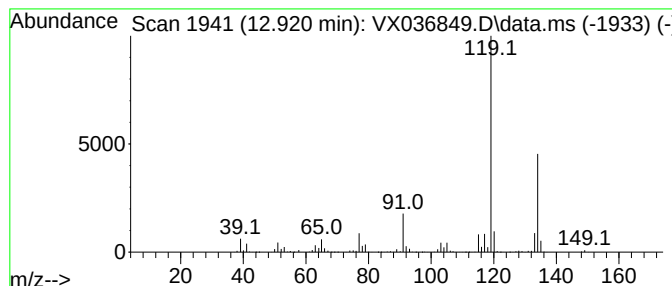
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.920	93.76 ug/l	1239990	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, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
3			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
4			o-Cymene	134	C10H14	000527-84-4	95
5			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX080723\
Data File : VX036849.D
Acq On : 07 Aug 2023 20:05
Operator : JC/MD
Sample : 03903-12
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-01W

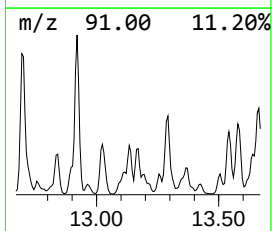
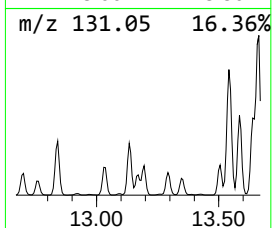
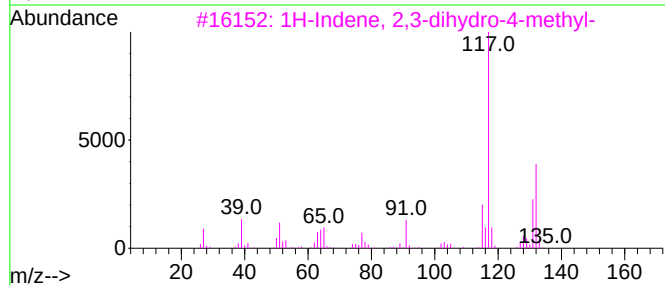
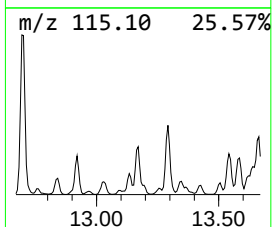
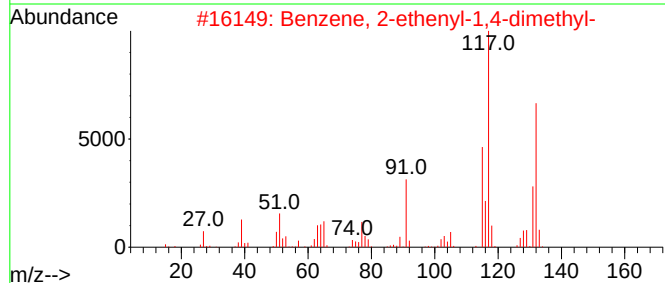
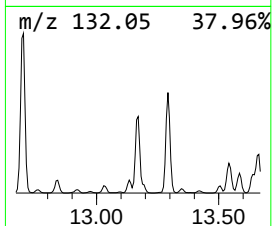
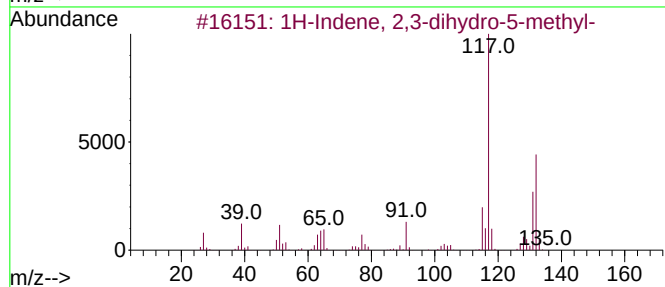
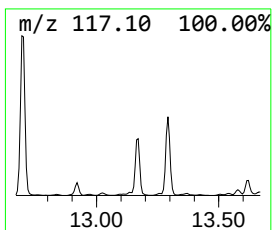
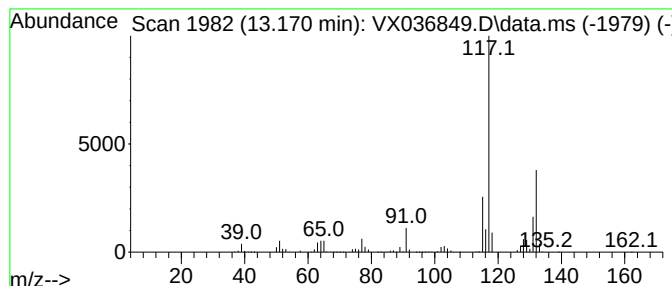
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Quant Title : SW846 8260

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

Peak Number 5 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.170	40.42 ug/l	534629	1,4-Dichlorobenzene-d4	12.024

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93	
2	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	91	
3	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90	
4	Indan, 1-methyl-	132	C10H12	000767-58-8	90	
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	90	



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX080723\
Data File : VX036849.D
Acq On : 07 Aug 2023 20:05
Operator : JC/MD
Sample : 03903-12
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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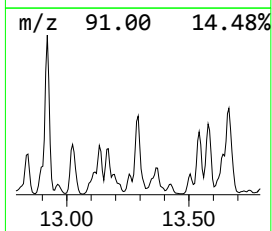
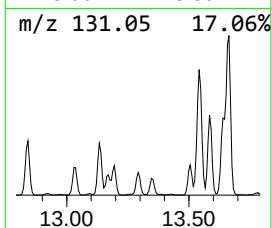
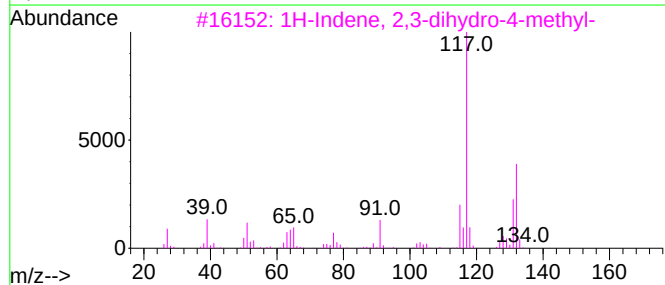
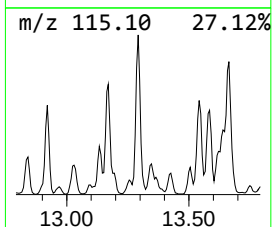
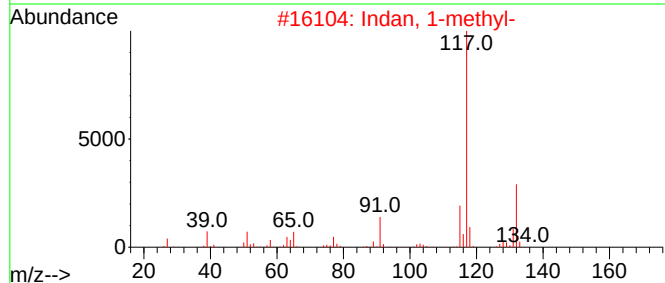
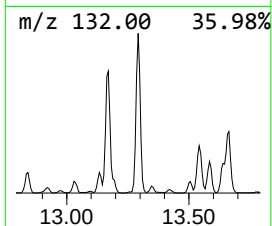
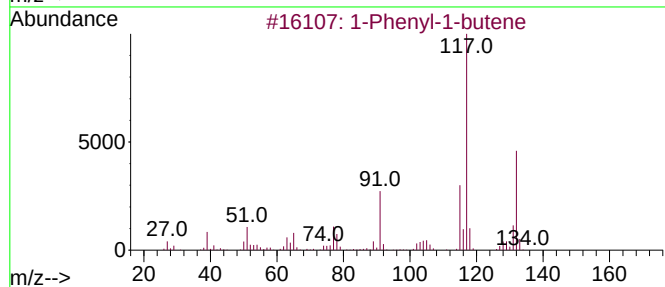
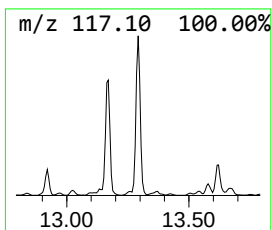
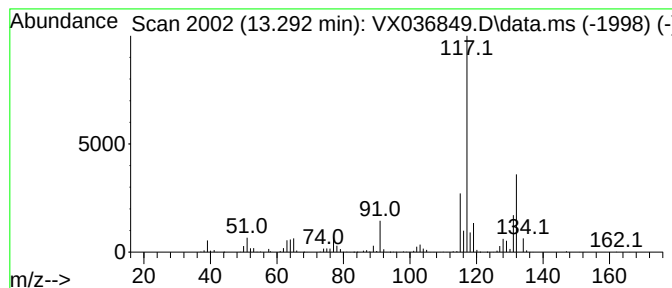
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Quant Title : SW846 8260

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

Peak Number 6 1-Phenyl-1-butene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	69.57 ug/l	920103	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenyl-1-butene	132	C10H12	000824-90-8	87
2			Indan, 1-methyl-	132	C10H12	000767-58-8	87
3			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	87
4			3-Phenylbut-1-ene	132	C10H12	000934-10-1	87
5			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	86



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX080723\
Data File : VX036849.D
Acq On : 07 Aug 2023 20:05
Operator : JC/MD
Sample : 03903-12
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-01W

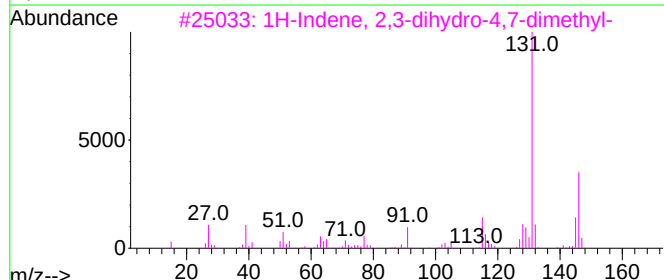
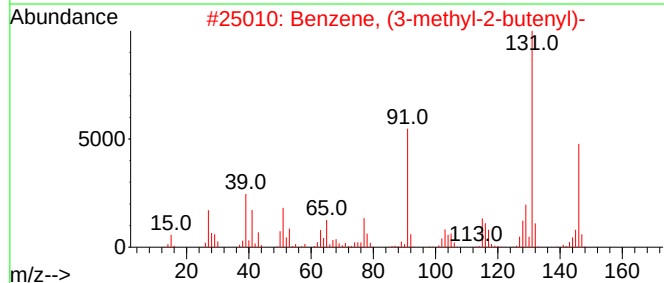
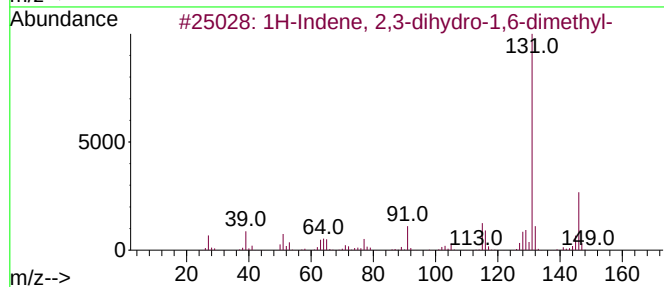
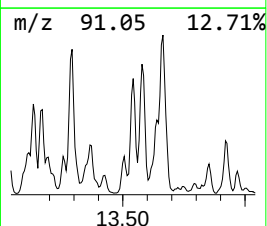
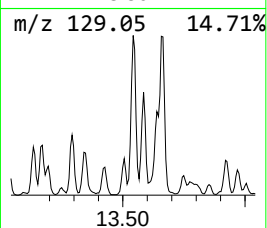
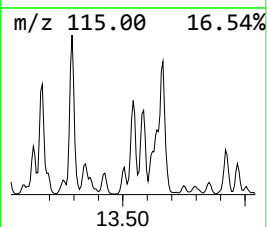
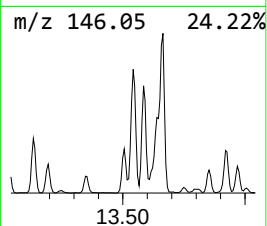
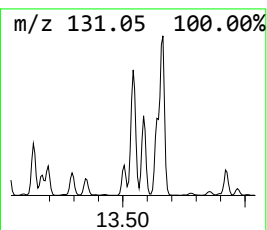
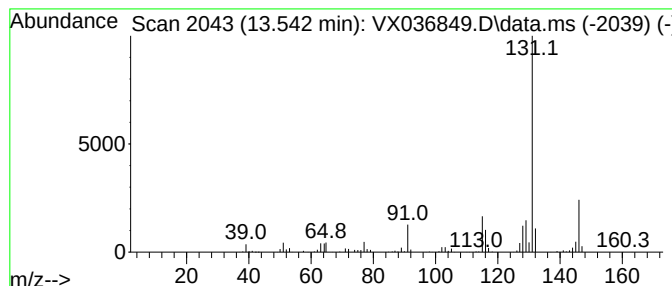
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Quant Title : SW846 8260

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

Peak Number 7 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.542	51.26 ug/l	677959	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	95
2			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	94
3			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	91
4			2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	91
5			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX080723\
Data File : VX036849.D
Acq On : 07 Aug 2023 20:05
Operator : JC/MD
Sample : 03903-12
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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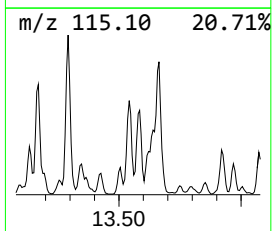
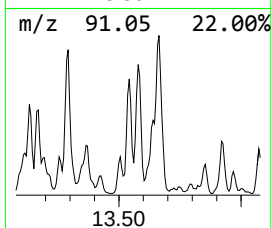
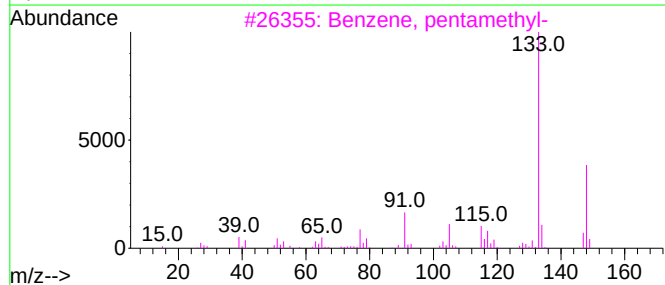
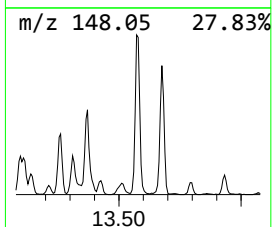
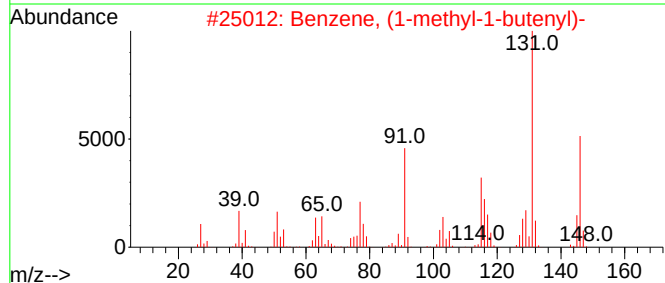
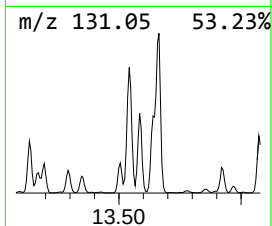
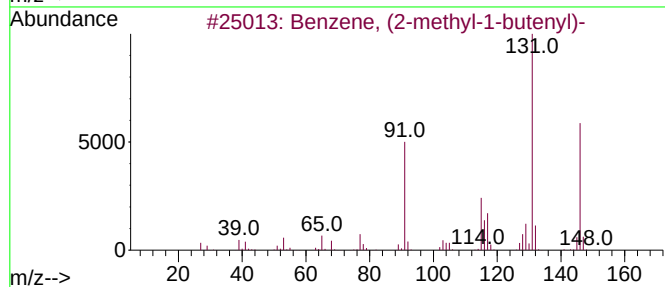
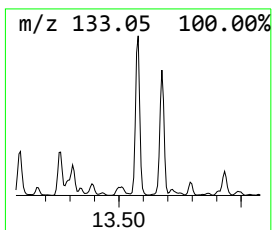
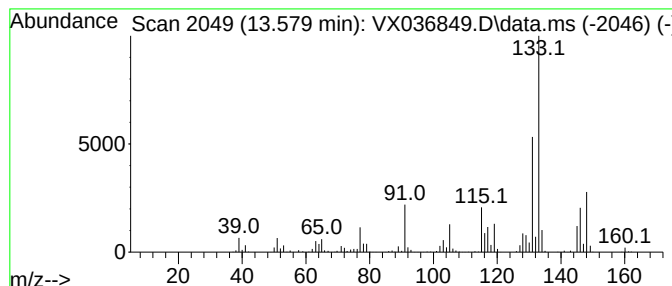
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, (2-methyl-1-butenyl)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.579	58.75 ug/l	777000	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	83
2			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	80
3			Benzene, pentamethyl-	148	C11H16	000700-12-9	53
4			Benzene, 1,3-dimethyl-5-(1-methy...	148	C11H16	004706-90-5	49
5			Benzene, 1,4-dimethyl-2-(1-methy...	148	C11H16	004132-72-3	46



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX080723\
Data File : VX036849.D
Acq On : 07 Aug 2023 20:05
Operator : JC/MD
Sample : 03903-12
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 27 Sample Multiplier: 1

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ClientSampleId :
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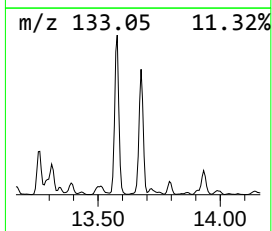
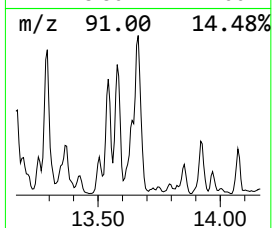
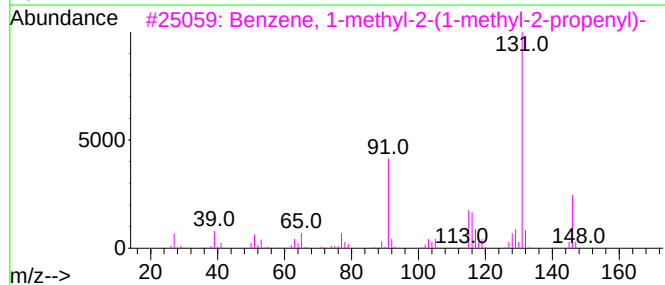
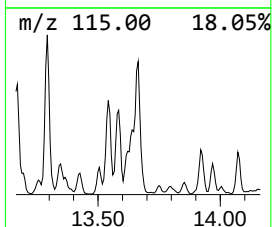
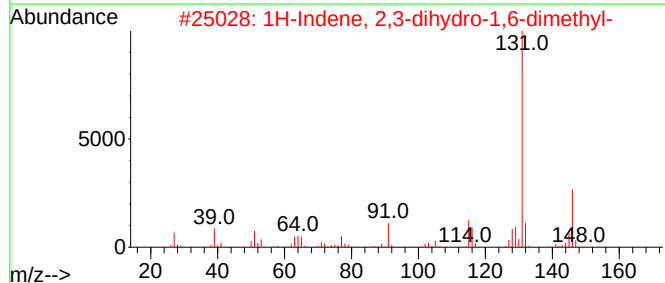
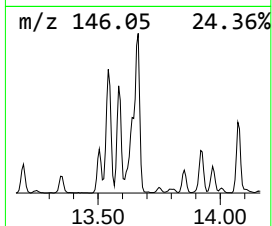
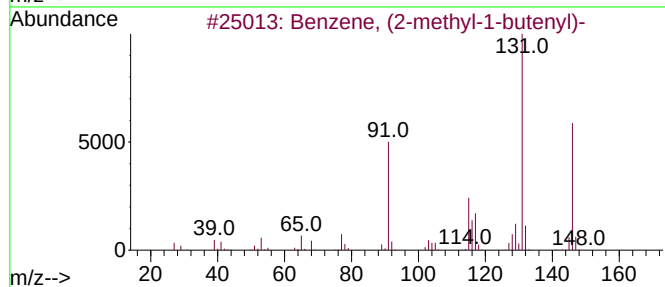
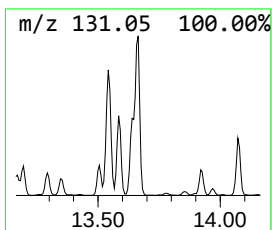
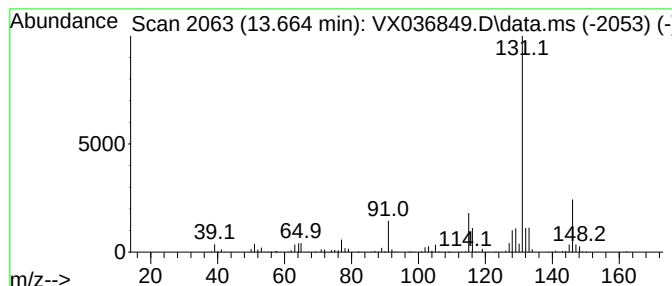
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Quant Title : SW846 8260

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

Peak Number 9 Benzene, 1-methyl-2-(1-meth... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.664	112.24 ug/l	1484400	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	93
2			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	90
3			Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	90
4			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	90
5			Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX080723\
Data File : VX036849.D
Acq On : 07 Aug 2023 20:05
Operator : JC/MD
Sample : 03903-12
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 27 Sample Multiplier: 1

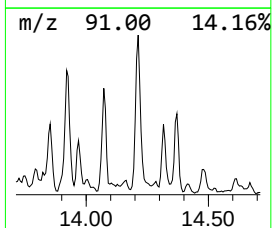
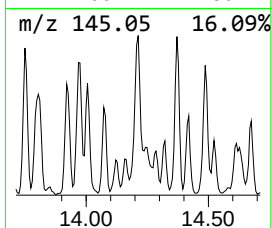
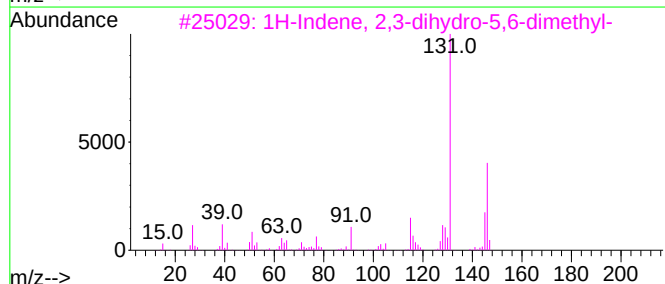
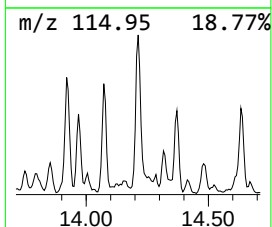
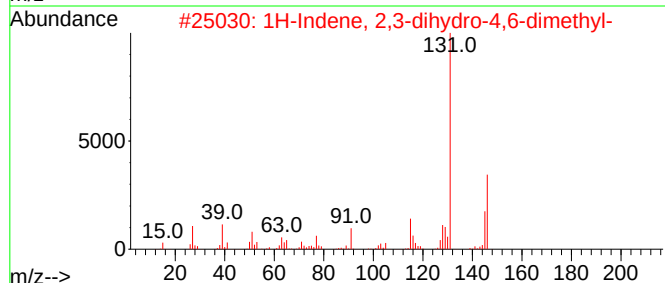
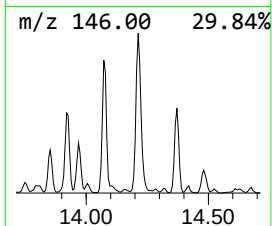
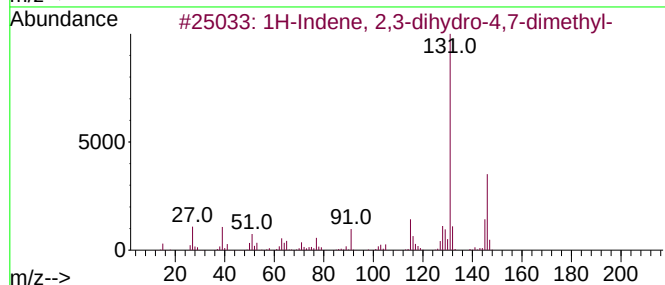
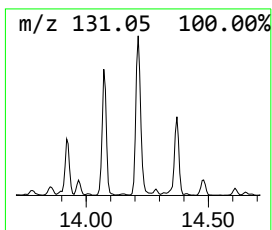
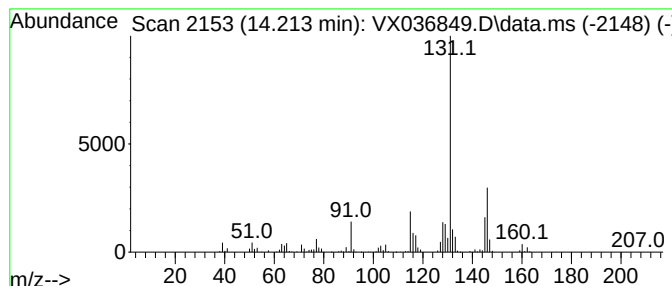
Instrument :
MSVOA_X
ClientSampleId :
MW-01W

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 10

R.T.	EstConc	Area	Relative to ISTD			R.T.
14.213	35.68 ug/l	471827	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	95
2	1H-Indene, 2,3-dihydro-4,6-dimet...		146	C11H14	001685-82-1	93
3	1H-Indene, 2,3-dihydro-5,6-dimet...		146	C11H14	001075-22-5	91
4	Benzene, (1,2-dimethyl-1-propenyl)-		146	C11H14	000769-57-3	90
5	1H-Indene, 2,3-dihydro-1,3-dimet...		146	C11H14	004175-53-5	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX080723\
Data File : VX036849.D
Acq On : 07 Aug 2023 20:05
Operator : JC/MD
Sample : 03903-12
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW-01W

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
Benzene, 1,2-di...	12.243	112.6	ug/l	1488790	4	12.024	661276	50.0
Benzene, 2-ethy...	12.609	115.8	ug/l	1531910	4	12.024	661276	50.0
Benzene, 1-ethe...	12.695	105.7	ug/l	1397230	4	12.024	661276	50.0
Benzene, 1,2,4,...	12.920	93.8	ug/l	1239990	4	12.024	661276	50.0
1H-Indene, 2,3-...	13.170	40.4	ug/l	534629	4	12.024	661276	50.0
1-Phenyl-1-butene	13.292	69.6	ug/l	920103	4	12.024	661276	50.0
1H-Indene, 2,3-...	13.542	51.3	ug/l	677959	4	12.024	661276	50.0
Benzene, (2-met...	13.579	58.8	ug/l	777000	4	12.024	661276	50.0
Benzene, 1-meth...	13.664	112.2	ug/l	1484400	4	12.024	661276	50.0
1H-Indene, 2,3-...	14.213	35.7	ug/l	471827	4	12.024	661276	50.0