

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX081821\
 Data File : VX023827.D
 Acq On : 18 Aug 2021 19:22
 Operator : JC/MD
 Sample : M3414-02 10X
 Misc : 5.04g/10mL/100uL/5.00mL/MSVOA_X/MEOH
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 B-2

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\82X081721W.M
 Title : SW846 8260

Signal : TIC: VX023827.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.569	71	79	80	rBV4	7075	14703	1.82%	0.190%
2	5.397	695	707	719	rBV	88828	255509	31.55%	3.298%
3	5.562	719	734	756	rVB	145745	445911	55.06%	5.756%
4	5.830	769	778	788	rBV4	8412	25004	3.09%	0.323%
5	5.964	788	800	813	rBV	97142	264273	32.63%	3.411%
6	6.257	840	848	857	rVB4	6578	20803	2.57%	0.269%
7	6.610	898	906	914	rBV6	5696	14402	1.78%	0.186%
8	6.769	922	932	945	rBV	237022	569374	70.30%	7.349%
9	7.385	1023	1033	1041	rVB5	13544	36095	4.46%	0.466%
10	7.561	1057	1062	1069	rVV3	5792	12231	1.51%	0.158%
11	7.781	1090	1098	1103	rBV5	4380	9555	1.18%	0.123%
12	7.988	1123	1132	1140	rVB4	7350	19839	2.45%	0.256%
13	8.110	1143	1152	1156	rBV2	5908	12509	1.54%	0.161%
14	8.195	1161	1166	1171	rVV4	6841	13412	1.66%	0.173%
15	8.281	1174	1180	1183	rVV2	17181	30736	3.79%	0.397%
16	8.317	1183	1186	1193	rVB3	10470	17448	2.15%	0.225%
17	8.439	1201	1206	1216	rBV	22866	42478	5.24%	0.548%
18	8.604	1227	1233	1235	rBV2	14614	23817	2.94%	0.307%
19	8.653	1235	1241	1250	rVB	514026	809935	100.00%	10.455%
20	8.775	1253	1261	1265	rBV4	7420	12423	1.53%	0.160%
21	8.915	1278	1284	1290	rBV2	14948	25323	3.13%	0.327%
22	8.982	1290	1295	1303	rVB5	8268	16296	2.01%	0.210%
23	9.104	1306	1315	1321	rBV3	10569	19793	2.44%	0.255%
24	9.293	1340	1346	1352	rVB2	5856	9670	1.19%	0.125%
25	9.390	1354	1362	1371	rBV3	9393	18095	2.23%	0.234%
26	9.506	1375	1381	1387	rBV4	18888	39626	4.89%	0.511%
27	9.567	1387	1391	1395	rVB	13069	18201	2.25%	0.235%
28	9.622	1395	1400	1404	rBV3	7692	14568	1.80%	0.188%
29	9.762	1417	1423	1427	rBV6	4679	8240	1.02%	0.106%
30	9.823	1428	1433	1437	rBV3	8676	17620	2.18%	0.227%
31	9.909	1443	1447	1452	rVB	45647	64572	7.97%	0.833%
32	10.012	1457	1464	1467	rBV	39804	57260	7.07%	0.739%
33	10.055	1467	1471	1478	rVB	488974	665667	82.19%	8.592%
34	10.195	1489	1494	1500	rVB	15316	24464	3.02%	0.316%
35	10.274	1500	1507	1509	rVV5	8713	17689	2.18%	0.228%
36	10.311	1510	1513	1518	rVV2	11677	20682	2.55%	0.267%

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37	10.366	1518	1522	1531	rVB7	8937	23260	2.87%	0.300%
38	10.591	1553	1559	1563	rVV3	16548	25268	3.12%	0.326%
39	10.677	1567	1573	1578	rBV	14191	24137	2.98%	0.312%
40	10.780	1583	1590	1595	rBV	19520	32316	3.99%	0.417%
41	10.860	1595	1603	1608	rVV3	16598	28774	3.55%	0.371%
42	10.963	1614	1620	1625	rBV	16460	24436	3.02%	0.315%
43	11.030	1625	1631	1635	rVV	17365	28158	3.48%	0.363%
44	11.085	1635	1640	1651	rVV	422179	584124	72.12%	7.540%
45	11.195	1654	1658	1666	rVB	38546	52831	6.52%	0.682%
46	11.305	1672	1676	1684	rVB	51609	66316	8.19%	0.856%
47	11.402	1686	1692	1695	rBV	10347	14894	1.84%	0.192%
48	11.481	1703	1705	1711	rVB4	6966	10148	1.25%	0.131%
49	11.634	1719	1730	1735	rBV3	14842	34938	4.31%	0.451%
50	11.683	1735	1738	1741	rVV	13139	17321	2.14%	0.224%
51	11.719	1741	1744	1746	rVV2	15450	18768	2.32%	0.242%
52	11.756	1746	1750	1760	rVB2	37481	69445	8.57%	0.896%
53	11.841	1760	1764	1766	rBV2	16949	21797	2.69%	0.281%
54	11.866	1766	1768	1770	rVV	15820	19373	2.39%	0.250%
55	11.896	1770	1773	1777	rVB	22012	27054	3.34%	0.349%
56	12.024	1788	1794	1800	rBV	407875	544562	67.24%	7.029%
57	12.079	1800	1803	1805	rVV	26388	35889	4.43%	0.463%
58	12.103	1805	1807	1811	rVB	29788	30596	3.78%	0.395%
59	12.176	1816	1819	1826	rVB2	36033	42659	5.27%	0.551%
60	12.250	1826	1831	1837	rBV	91509	118327	14.61%	1.527%
61	12.335	1837	1845	1849	rVB2	44375	89802	11.09%	1.159%
62	12.408	1849	1857	1863	rBV4	13629	29256	3.61%	0.378%
63	12.469	1863	1867	1872	rVB4	9326	14878	1.84%	0.192%
64	12.536	1874	1878	1883	rBV2	10481	19959	2.46%	0.258%
65	12.615	1883	1891	1894	rVV	189243	231712	28.61%	2.991%
66	12.646	1894	1896	1900	rVB2	29036	31872	3.94%	0.411%
67	12.701	1900	1905	1915	rVB3	59330	122618	15.14%	1.583%
68	12.786	1915	1919	1921	rBV2	7005	8573	1.06%	0.111%
69	12.926	1931	1942	1947	rBV	115869	200706	24.78%	2.591%
70	12.975	1947	1950	1954	rVB3	19722	25091	3.10%	0.324%
71	13.030	1954	1959	1966	rVB3	37756	74332	9.18%	0.959%
72	13.115	1966	1973	1976	rBV3	28060	48702	6.01%	0.629%
73	13.170	1979	1982	1987	rVB	59251	67007	8.27%	0.865%
74	13.262	1994	1997	1999	rVV	22442	29216	3.61%	0.377%
75	13.292	1999	2002	2009	rVV4	142982	219142	27.06%	2.829%
76	13.371	2011	2015	2021	rVV2	44776	64003	7.90%	0.826%

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77	13.426	2021	2024	2027	rVB4	12707	15306	1.89%	0.198%
78	13.512	2033	2038	2040	rBV	24629	35071	4.33%	0.453%
79	13.548	2040	2044	2046	rVV	42709	58133	7.18%	0.750%
80	13.585	2046	2050	2054	rVB3	68721	98328	12.14%	1.269%
81	13.676	2060	2065	2072	rVB3	45781	98705	12.19%	1.274%
82	13.737	2072	2075	2078	rVV4	8824	14575	1.80%	0.188%
83	13.780	2078	2082	2088	rVV2	26533	46683	5.76%	0.603%
84	13.835	2088	2091	2099	rVB8	16009	33381	4.12%	0.431%
85	13.926	2099	2106	2111	rBV3	29148	53335	6.59%	0.688%
86	13.975	2111	2114	2117	rVV	17927	21689	2.68%	0.280%
87	14.079	2127	2131	2135	rBV	41743	52112	6.43%	0.673%
88	14.164	2142	2145	2149	rVB3	7052	8711	1.08%	0.112%
89	14.219	2149	2154	2164	rVB2	42528	70449	8.70%	0.909%
90	14.322	2167	2171	2176	rBV	21553	29271	3.61%	0.378%
91	14.377	2176	2180	2183	rVB	30382	36123	4.46%	0.466%
92	14.420	2183	2187	2192	rVB5	7195	11797	1.46%	0.152%
93	14.487	2192	2198	2203	rBV4	8596	19323	2.39%	0.249%
94	14.639	2215	2223	2228	rBV	83299	113076	13.96%	1.460%
95	14.780	2241	2246	2251	rBV	37881	52034	6.42%	0.672%
96	15.481	2354	2361	2368	rBV4	5173	8269	1.02%	0.107%
97	15.615	2379	2383	2387	rBV2	7109	10318	1.27%	0.133%

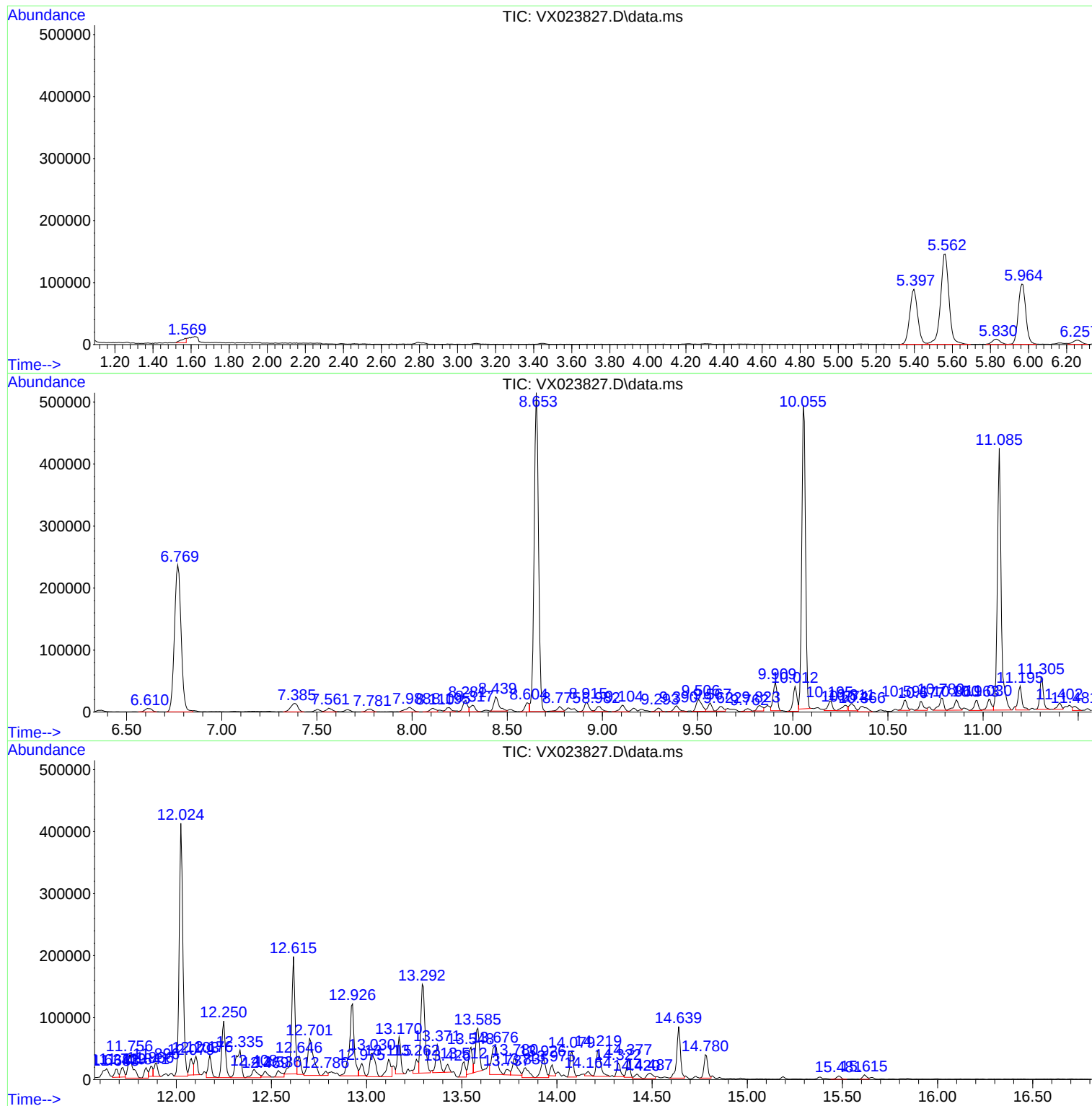
Sum of corrected areas: 7747172

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ALS Vial : 25 Sample Multiplier: 1

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MSVOA_X
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B-2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081721W.M
Quant Title : SW846 8260

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



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Instrument :
MSVOA_X
ClientSampled :
B-2

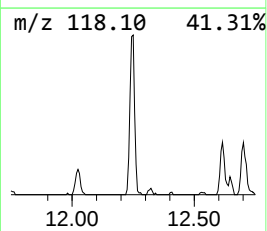
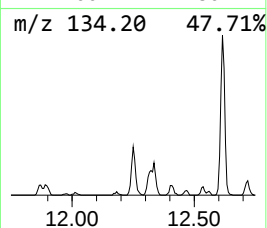
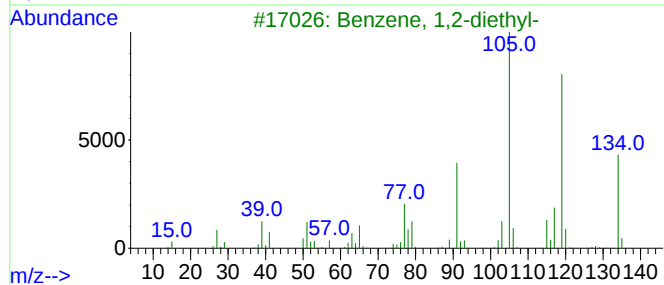
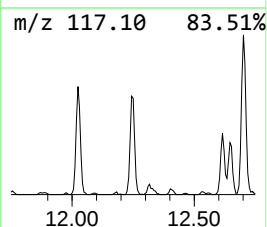
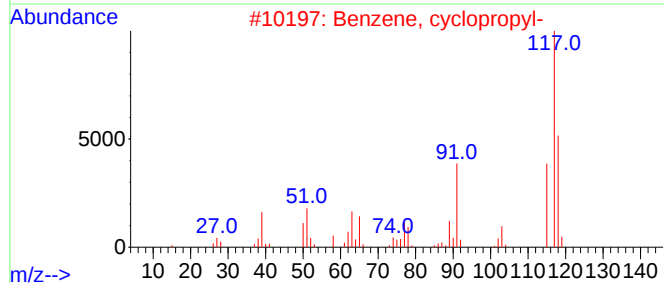
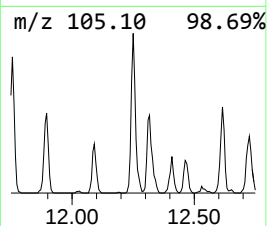
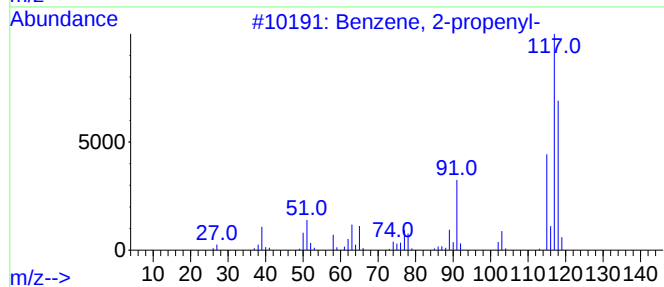
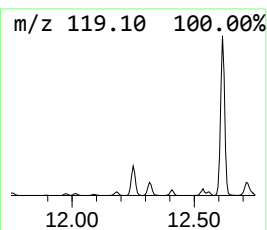
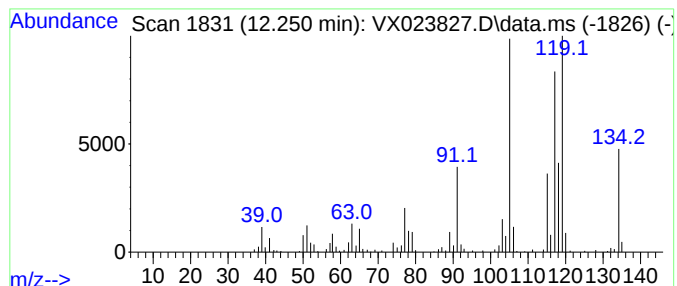
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081721W.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 2-propenyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.250	10.86 ug/l	118327	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, cyclopropyl-	118	C9H10	000873-49-4	80
3			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	60
4			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	55
5			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	55



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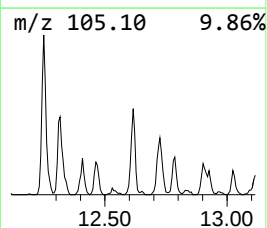
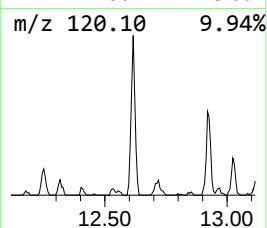
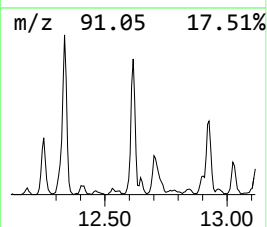
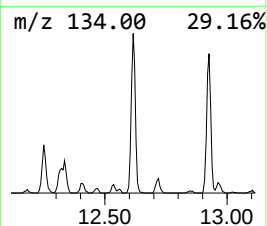
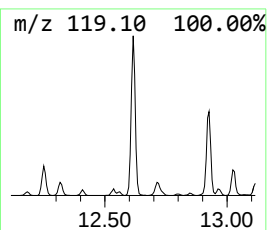
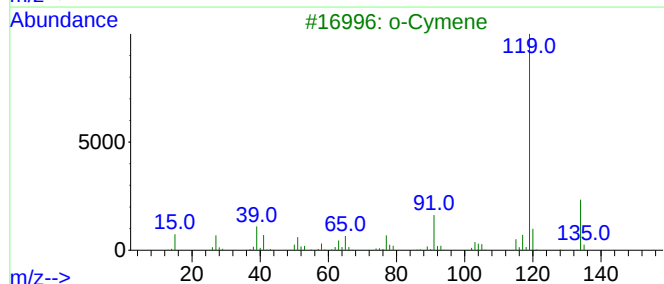
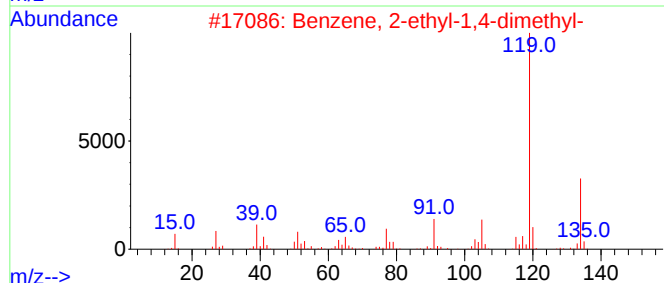
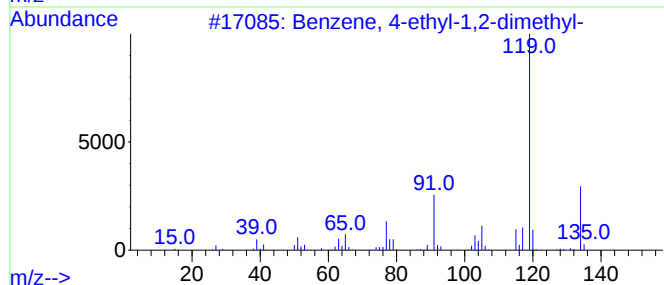
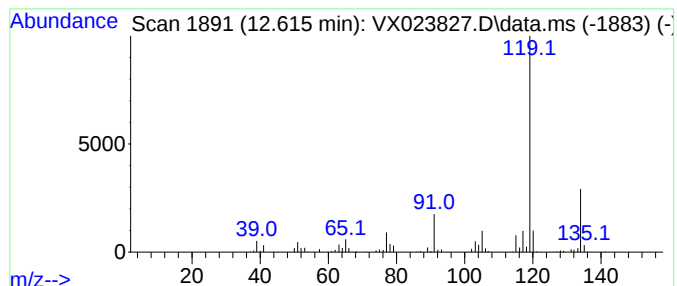
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081721W.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.615	21.28 ug/l	231712	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	97
2			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94
3			o-Cymene	134	C10H14	000527-84-4	94
4			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94
5			p-Cymene	134	C10H14	000099-87-6	94



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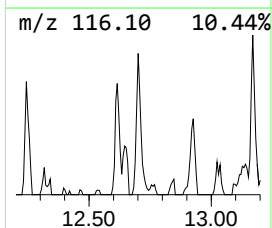
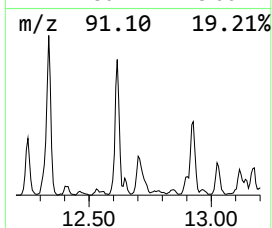
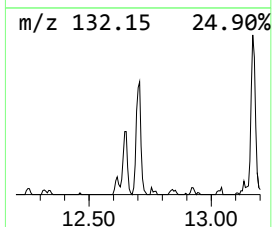
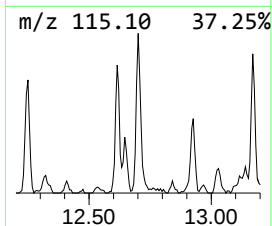
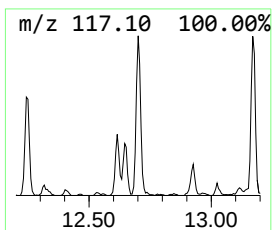
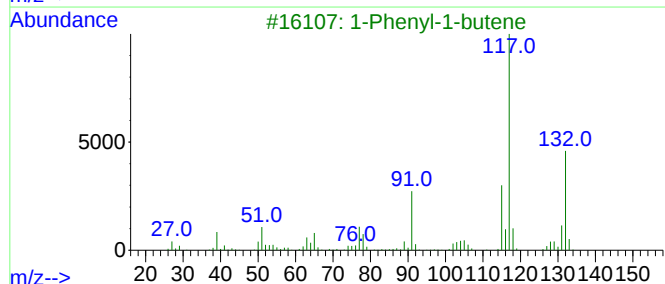
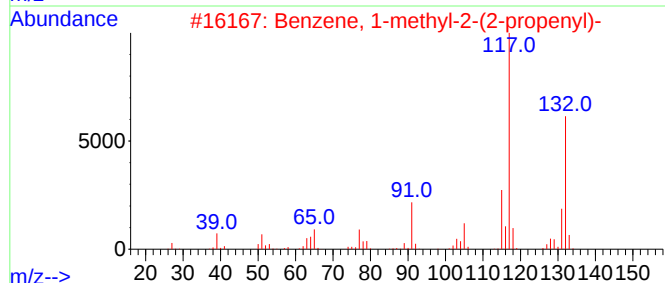
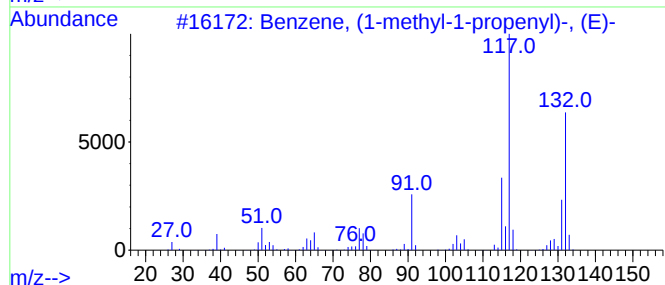
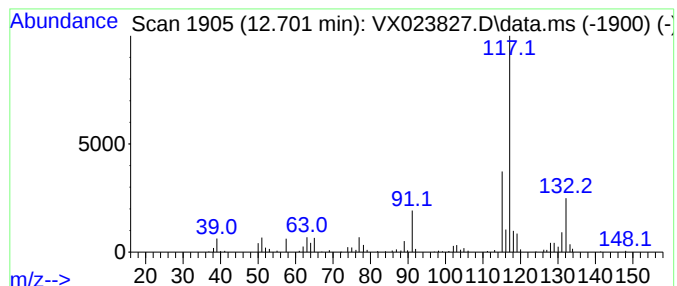
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Quant Title : SW846 8260

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

Peak Number 3 Benzene, (1-methyl-1-propen... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	11.26 ug/l	122618	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	91
2			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	91
3			1-Phenyl-1-butene	132	C10H12	000824-90-8	91
4			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	91
5			Benzene, 2-butenyl-	132	C10H12	001560-06-1	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX081821\
Data File : VX023827.D
Acq On : 18 Aug 2021 19:22
Operator : JC/MD
Sample : M3414-02 10X
Misc : 5.04g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B-2

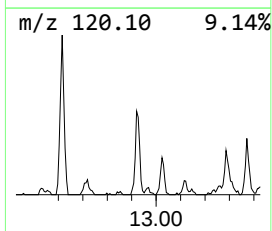
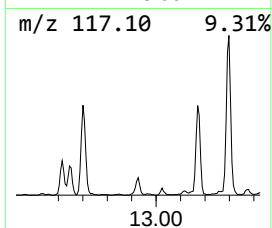
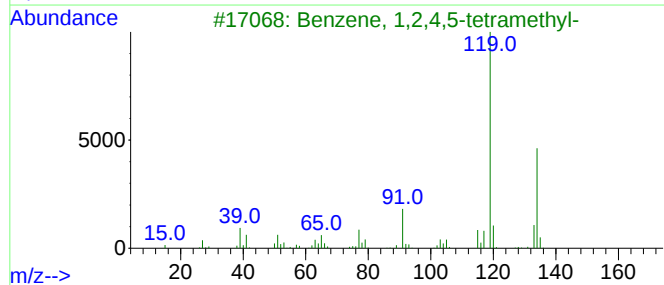
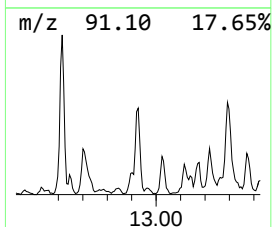
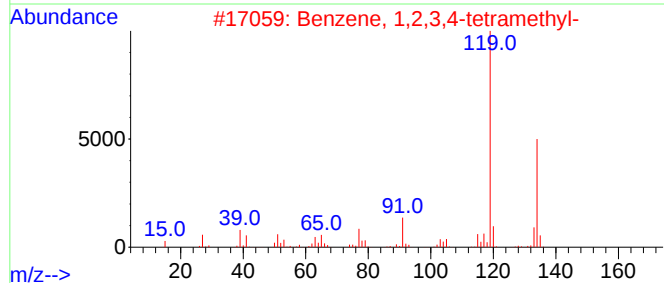
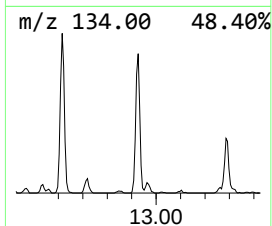
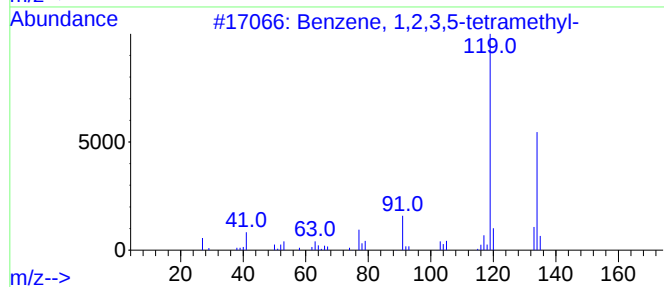
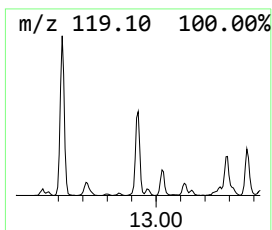
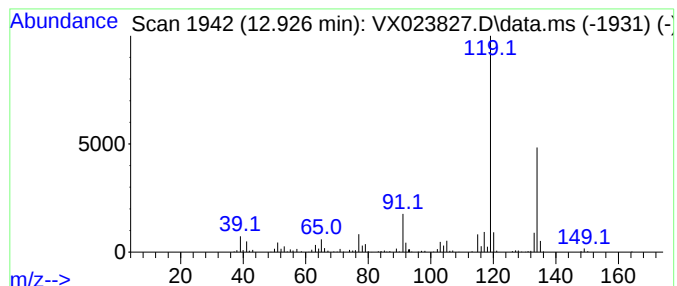
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.926	18.43 ug/l	200706	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
2			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
4			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94
5			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX081821\
Data File : VX023827.D
Acq On : 18 Aug 2021 19:22
Operator : JC/MD
Sample : M3414-02 10X
Misc : 5.04g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B-2

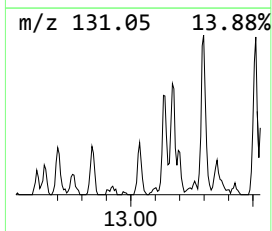
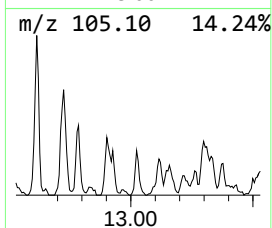
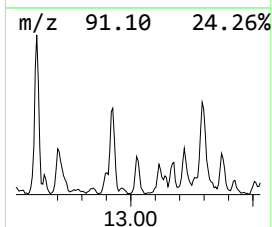
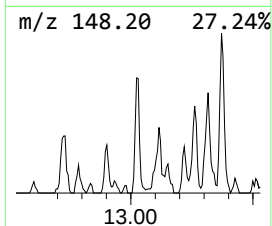
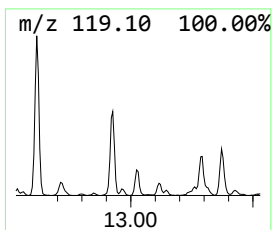
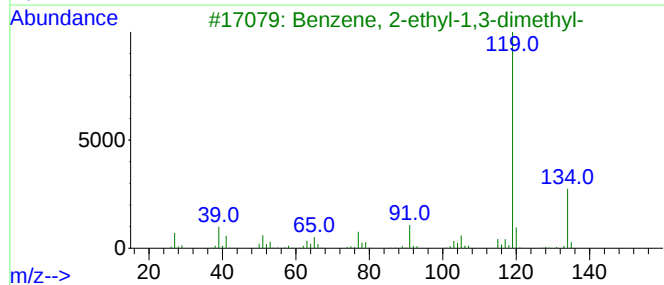
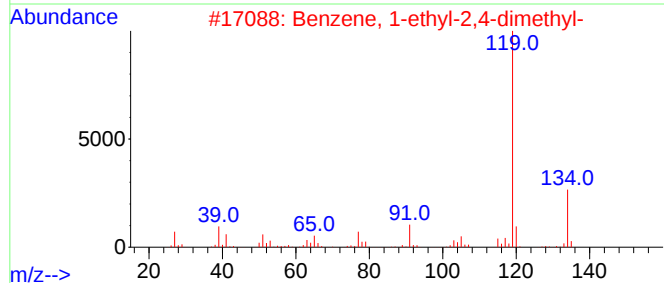
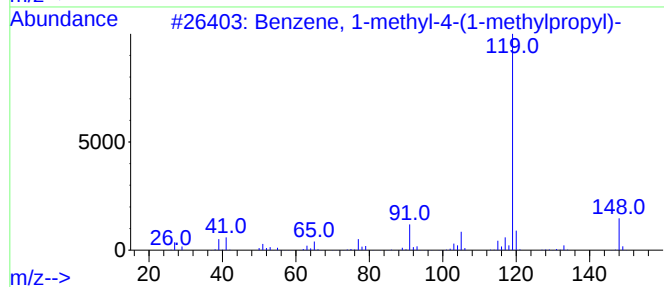
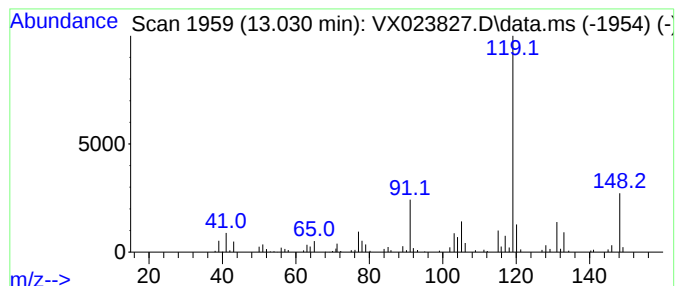
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1-methyl-4-(1-meth... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.030	6.82 ug/l	74332	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	81
2			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	68
3			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	68
4			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	68
5			o-Cymene	134	C10H14	000527-84-4	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX081821\
Data File : VX023827.D
Acq On : 18 Aug 2021 19:22
Operator : JC/MD
Sample : M3414-02 10X
Misc : 5.04g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B-2

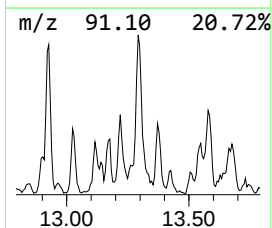
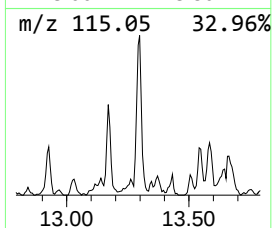
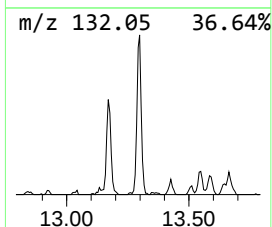
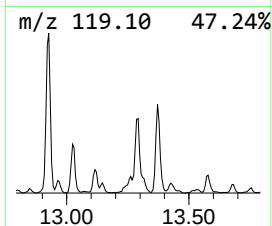
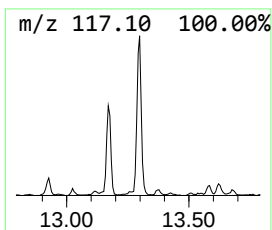
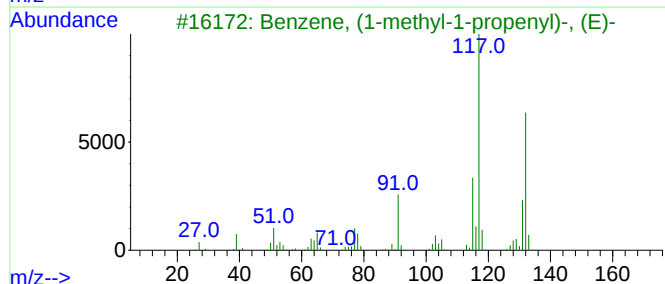
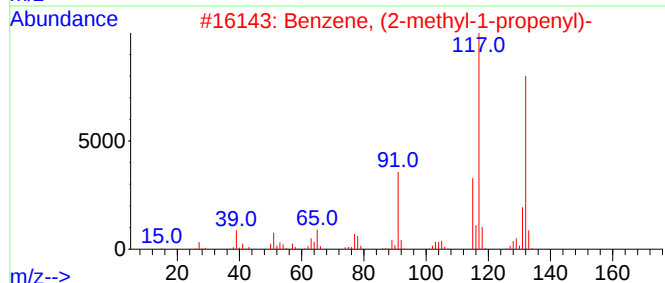
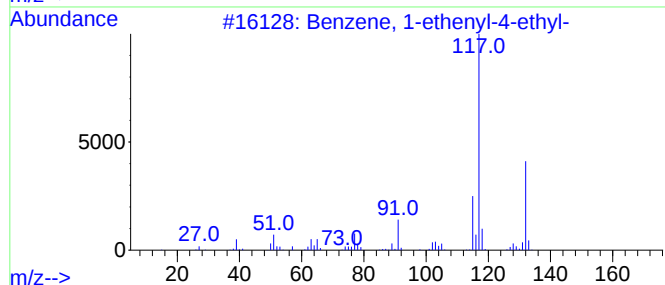
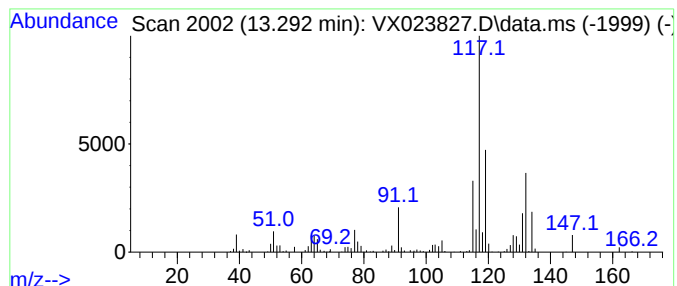
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	20.12 ug/l	219142	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	64
2			Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	62
3			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	62
4			1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	60
5			Indan, 1-methyl-	132	C10H12	000767-58-8	52



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX081821\
Data File : VX023827.D
Acq On : 18 Aug 2021 19:22
Operator : JC/MD
Sample : M3414-02 10X
Misc : 5.04g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
B-2

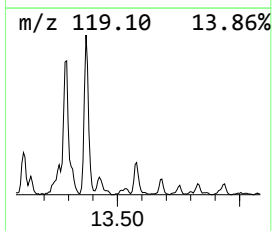
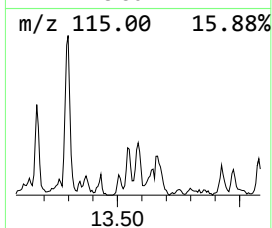
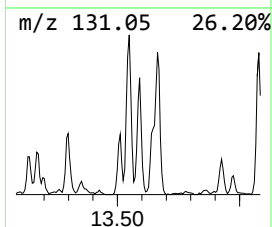
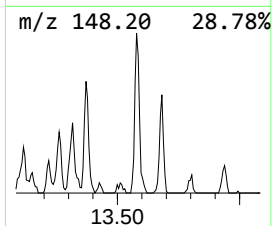
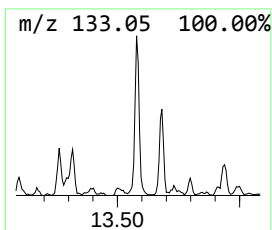
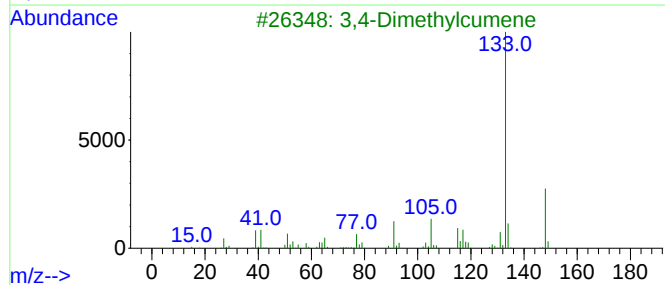
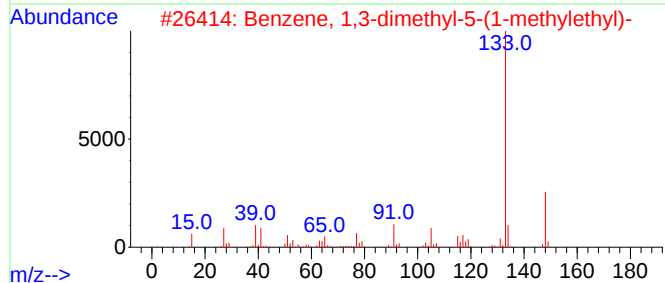
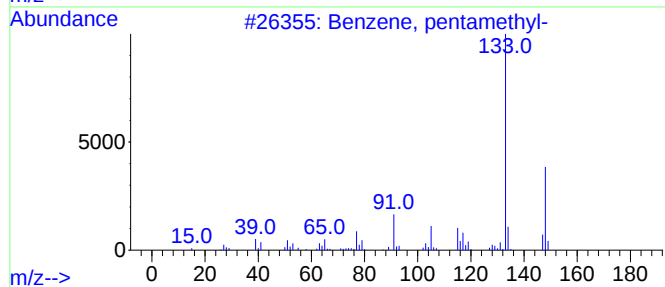
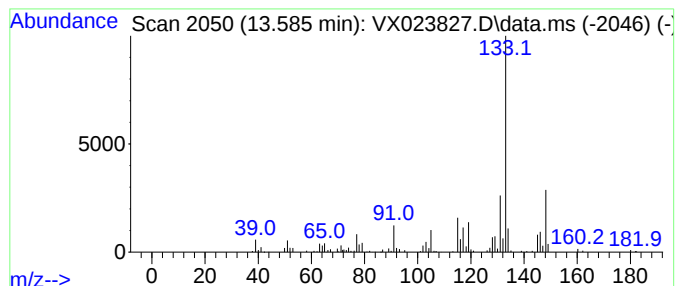
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, pentamethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.585	9.03 ug/l	98328	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, pentamethyl-	148	C11H16	000700-12-9	76
2			Benzene, 1,3-dimethyl-5-(1-methy...	148	C11H16	004706-90-5	68
3			3,4-Dimethylcumene	148	C11H16	1000370-34-1	64
4			Benzene, 1,4-diethyl-2-methyl-	148	C11H16	013632-94-5	64
5			4-tert-Butyltoluene	148	C11H16	000098-51-1	62



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX081821\
Data File : VX023827.D
Acq On : 18 Aug 2021 19:22
Operator : JC/MD
Sample : M3414-02 10X
Misc : 5.04g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B-2

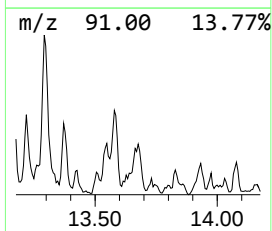
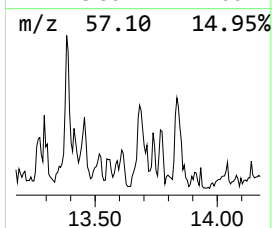
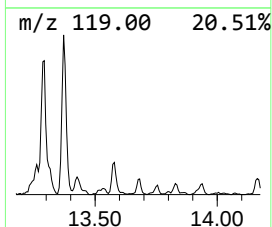
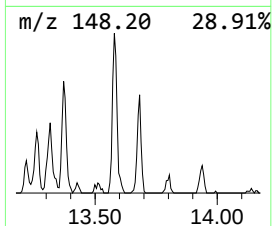
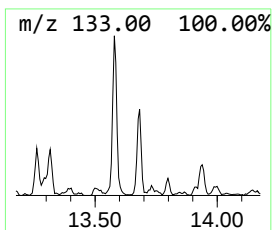
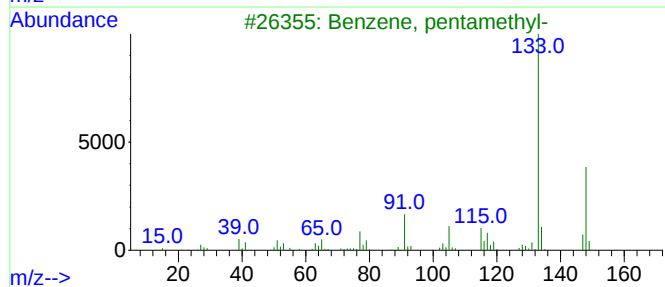
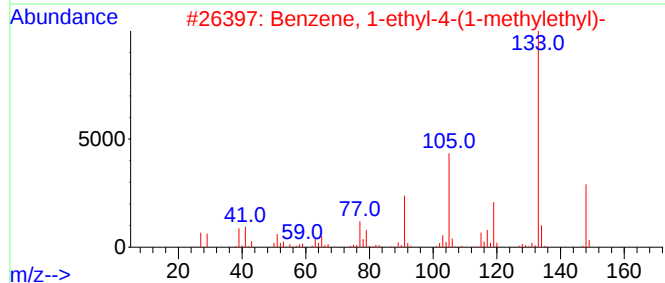
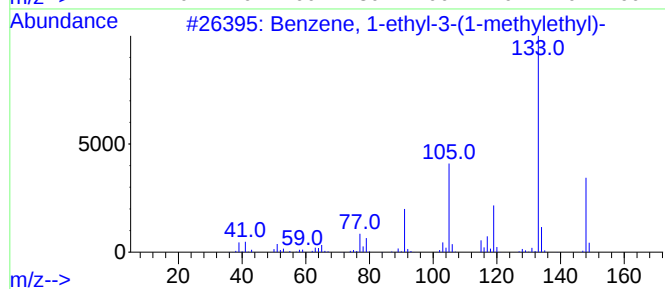
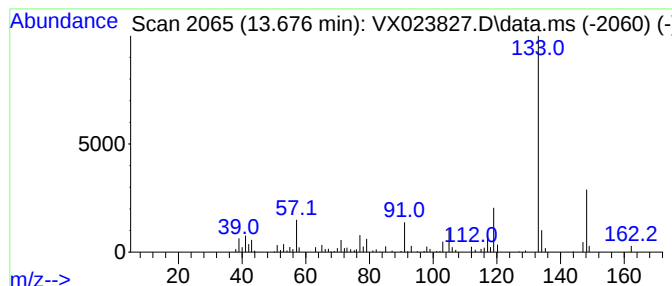
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1-ethyl-3-(1-methy... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.676	9.06 ug/l	98705	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3-(1-methylethyl)-	148	C11H16	004920-99-4	83
2			Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	83
3			Benzene, pentamethyl-	148	C11H16	000700-12-9	80
4			Benzene, 2,4-dimethyl-1-(1-methy...	148	C11H16	004706-89-2	74
5			Benzene, 1,4-dimethyl-2-(1-methy...	148	C11H16	004132-72-3	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX081821\
Data File : VX023827.D
Acq On : 18 Aug 2021 19:22
Operator : JC/MD
Sample : M3414-02 10X
Misc : 5.04g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B-2

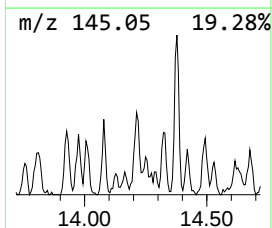
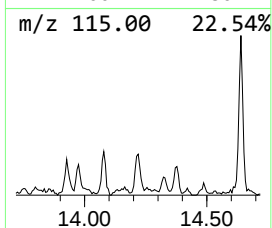
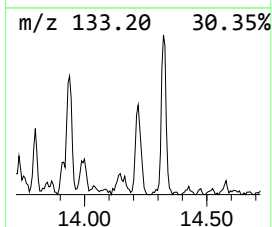
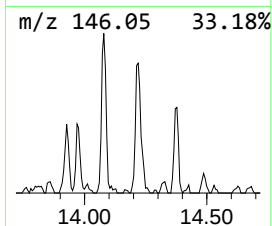
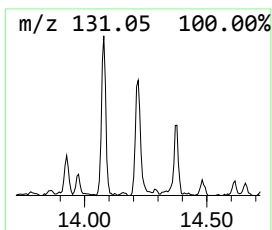
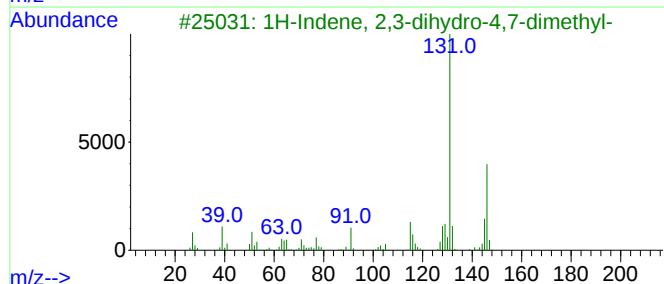
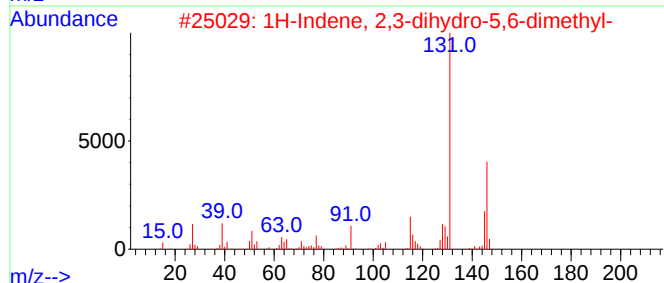
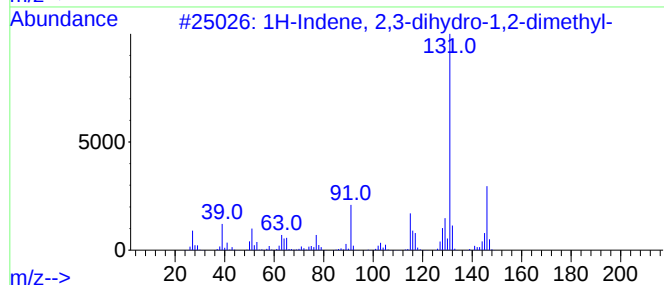
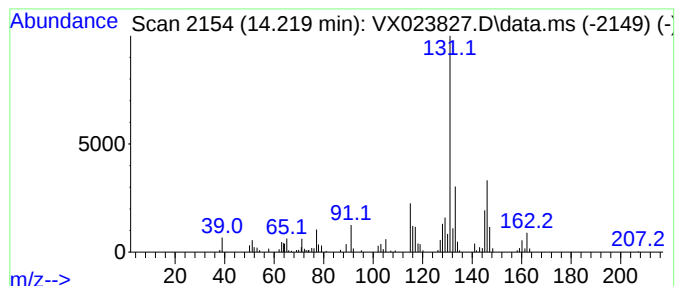
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Quant Title : SW846 8260

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

Peak Number 9 1H-Indene, 2,3-dihydro-1,2-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.219	6.47 ug/l	70449	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	90
2			1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	81
3			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	81
4			Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	81
5			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	76



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX081821\
Data File : VX023827.D
Acq On : 18 Aug 2021 19:22
Operator : JC/MD
Sample : M3414-02 10X
Misc : 5.04g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 25 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B-2

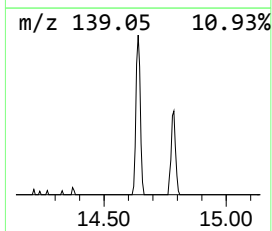
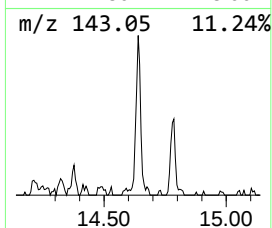
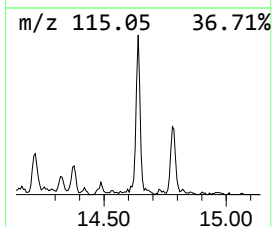
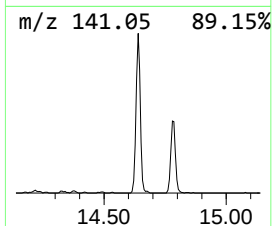
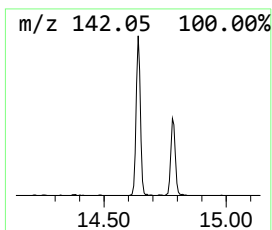
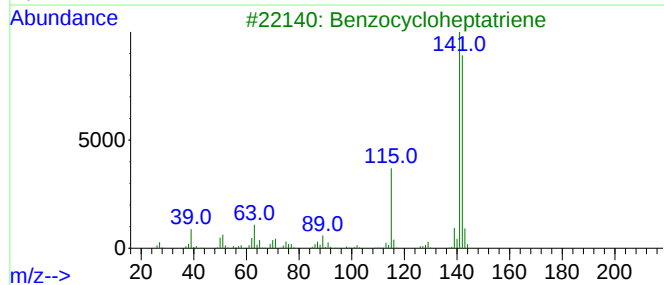
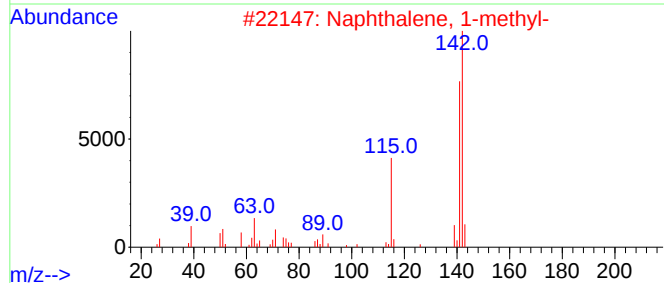
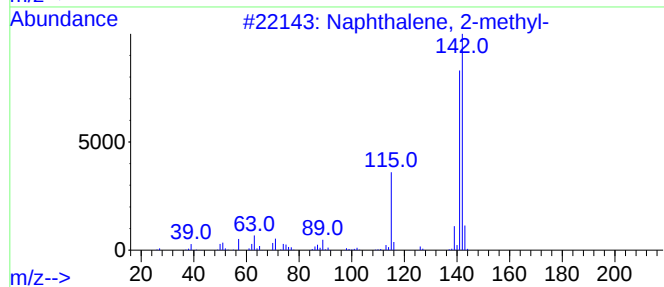
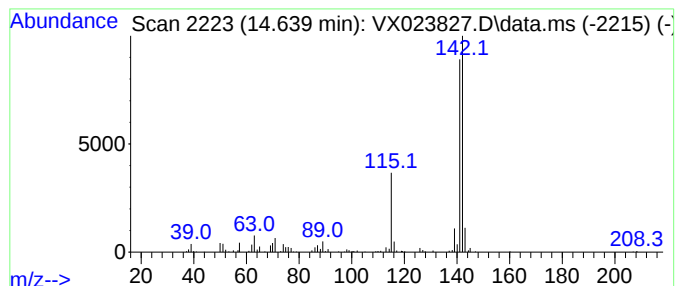
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Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.639	10.38 ug/l	113076	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	97
2			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
3			Benzocycloheptatriene	142	C11H10	000264-09-5	91
4			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
5			Naphthalene, 1-[(2-propenyloxy)m...	198	C14H14O	074685-39-5	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX081821\
 Data File : VX023827.D
 Acq On : 18 Aug 2021 19:22
 Operator : JC/MD
 Sample : M3414-02 10X
 Misc : 5.04g/10mL/100uL/5.00mL/MSVOA_X/MEOH
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 B-2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081721W.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, 2-prop...	12.250	10.9	ug/l	118327	4	12.024	544562	50.0
Benzene, 4-ethy...	12.615	21.3	ug/l	231712	4	12.024	544562	50.0
Benzene, (1-met...	12.701	11.3	ug/l	122618	4	12.024	544562	50.0
Benzene, 1,2,3...	12.926	18.4	ug/l	200706	4	12.024	544562	50.0
Benzene, 1-meth...	13.030	6.8	ug/l	74332	4	12.024	544562	50.0
Benzene, 1-ethe...	13.292	20.1	ug/l	219142	4	12.024	544562	50.0
Benzene, pentam...	13.585	9.0	ug/l	98328	4	12.024	544562	50.0
Benzene, 1-ethy...	13.676	9.1	ug/l	98705	4	12.024	544562	50.0
1H-Indene, 2,3-...	14.219	6.5	ug/l	70449	4	12.024	544562	50.0
Naphthalene, 1-...	14.639	10.4	ug/l	113076	4	12.024	544562	50.0