

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX050123\
 Data File : VX035440.D
 Acq On : 01 May 2023 15:08
 Operator : JC/MD
 Sample : 02558-05
 Misc : 4.49g/10mL/100uL/5.00mL/MSVOA_X/MEOH
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 P001-SS012-0006-01

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

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

Signal : TIC: VX035440.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.385	695	705	720	rBV	122545	352465	12.50%	0.489%
2	5.544	720	731	747	rVB	242147	682692	24.21%	0.948%
3	5.952	787	798	815	rBV2	143029	384817	13.65%	0.534%
4	6.757	920	930	945	rBV	366533	880258	31.22%	1.222%
5	8.647	1234	1240	1248	rVB	793568	1198279	42.50%	1.663%
6	10.006	1456	1463	1466	rBV	37933	52102	1.85%	0.072%
7	10.055	1466	1471	1481	rVB	745477	1035219	36.72%	1.437%
8	10.195	1487	1494	1498	rVV	34668	54884	1.95%	0.076%
9	10.299	1498	1511	1517	rVV2	138750	274735	9.74%	0.381%
10	10.360	1517	1521	1531	rVB2	106951	152863	5.42%	0.212%
11	10.585	1551	1558	1562	rBV3	58686	96658	3.43%	0.134%
12	10.640	1563	1567	1571	rBV	88704	123915	4.39%	0.172%
13	10.781	1582	1590	1594	rBV2	134439	202984	7.20%	0.282%
14	10.854	1594	1602	1606	rVV3	113520	206882	7.34%	0.287%
15	10.890	1606	1608	1613	rVB2	62594	85003	3.01%	0.118%
16	10.957	1613	1619	1623	rBV3	35801	71654	2.54%	0.099%
17	11.030	1628	1631	1634	rVV2	59529	80472	2.85%	0.112%
18	11.085	1634	1640	1650	rVB2	698843	1188477	42.15%	1.650%
19	11.189	1650	1657	1660	rBV	156744	253002	8.97%	0.351%
20	11.219	1660	1662	1666	rVB3	114863	124089	4.40%	0.172%
21	11.305	1672	1676	1679	rBV	74155	98779	3.50%	0.137%
22	11.378	1684	1688	1694	rVV	410209	760929	26.99%	1.056%
23	11.451	1694	1700	1702	rVV2	358117	627839	22.27%	0.872%
24	11.476	1702	1704	1709	rVB	578400	666642	23.64%	0.925%
25	11.616	1722	1727	1734	rVB2	287419	489633	17.37%	0.680%
26	11.683	1734	1738	1740	rBV2	103898	124703	4.42%	0.173%
27	11.713	1740	1743	1746	rBV	361852	390078	13.84%	0.541%
28	11.756	1746	1750	1759	rVB	963920	1492826	52.95%	2.072%
29	11.835	1759	1763	1766	rBV2	105520	170165	6.04%	0.236%
30	11.890	1766	1772	1777	rVB2	215109	317774	11.27%	0.441%
31	11.945	1777	1781	1783	rBV	265996	327145	11.60%	0.454%
32	11.975	1783	1786	1789	rVB2	208687	244977	8.69%	0.340%
33	12.024	1789	1794	1799	rBV3	843168	1305816	46.31%	1.813%
34	12.091	1799	1805	1811	rVB3	579570	1132918	40.18%	1.573%
35	12.171	1811	1818	1826	rVB3	304244	618736	21.95%	0.859%
36	12.250	1826	1831	1832	rBV2	517377	696509	24.70%	0.967%

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX050123\
 Data File : VX035440.D
 Acq On : 01 May 2023 15:08
 Operator : JC/MD
 Sample : 02558-05
 Misc : 4.49g/10mL/100uL/5.00mL/MSVOA_X/MEOH
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 P001-SS012-0006-01

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

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

37	12.268	1832	1834	1838	rVB	571878	621190	22.03%	0.862%
38	12.317	1838	1842	1848	rVB3	1093239	1772443	62.86%	2.460%
39	12.420	1848	1859	1863	rBV	970320	1525086	54.09%	2.117%
40	12.463	1863	1866	1872	rVB2	503417	766066	27.17%	1.063%

41	12.536	1873	1878	1879	rBV	544393	688293	24.41%	0.955%
42	12.555	1879	1881	1884	rVV	549181	616746	21.87%	0.856%
43	12.616	1888	1891	1894	rVB	716640	759909	26.95%	1.055%
44	12.701	1899	1905	1907	rBV2	412103	850183	30.15%	1.180%
45	12.817	1917	1924	1927	rBV3	337856	630113	22.35%	0.875%

46	12.847	1927	1929	1932	rVB2	202967	200744	7.12%	0.279%
47	12.890	1932	1936	1938	rBV2	397843	608486	21.58%	0.845%
48	12.920	1938	1941	1944	rVV	634826	915224	32.46%	1.270%
49	12.963	1944	1948	1955	rVB	995156	1764927	62.60%	2.450%
50	13.042	1955	1961	1964	rBV4	529697	883024	31.32%	1.226%

51	13.073	1964	1966	1968	rVB	170845	133923	4.75%	0.186%
52	13.164	1974	1981	1985	rBV4	621634	1052009	37.31%	1.460%
53	13.213	1985	1989	1992	rBV3	344777	388224	13.77%	0.539%
54	13.268	1992	1998	1999	rBV3	936310	1312041	46.54%	1.821%
55	13.292	1999	2002	2007	rVB3	1707828	2391335	84.82%	3.320%

56	13.384	2013	2017	2020	rBV2	540952	798760	28.33%	1.109%
57	13.420	2020	2023	2032	rVB3	917621	1937870	68.73%	2.690%
58	13.506	2032	2037	2040	rBV2	368491	568243	20.15%	0.789%
59	13.542	2040	2043	2046	rBV	360532	472184	16.75%	0.655%
60	13.585	2046	2050	2056	rBV3	554201	1017415	36.09%	1.412%

61	13.670	2060	2064	2068	rVV3	423025	792722	28.12%	1.100%
62	13.737	2072	2075	2078	rVV2	331420	423271	15.01%	0.588%
63	13.774	2078	2081	2084	rVV	640999	839035	29.76%	1.165%
64	13.847	2088	2093	2099	rVB4	1045194	1850498	65.63%	2.569%
65	13.926	2099	2106	2111	rBV6	715278	1426200	50.58%	1.980%

66	13.969	2111	2113	2116	rBV	198669	278913	9.89%	0.387%
67	14.042	2121	2125	2127	rVB3	573378	630488	22.36%	0.875%
68	14.079	2127	2131	2134	rBV2	559880	737499	26.16%	1.024%
69	14.140	2138	2141	2143	rBV3	172180	168786	5.99%	0.234%
70	14.164	2143	2145	2147	rVV	141359	137641	4.88%	0.191%

71	14.225	2150	2155	2163	rVB4	1025024	2085809	73.98%	2.895%
72	14.323	2167	2171	2176	rBV2	423138	659062	23.38%	0.915%
73	14.371	2176	2179	2183	rVB	592766	671850	23.83%	0.933%
74	14.420	2183	2187	2193	rVB4	333372	685888	24.33%	0.952%
75	14.481	2193	2197	2206	rBV5	766940	1749397	62.05%	2.428%

76	14.566	2206	2211	2215	rBV5	224103	489531	17.36%	0.680%
----	--------	------	------	------	------	--------	--------	--------	--------

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX050123\
 Data File : VX035440.D
 Acq On : 01 May 2023 15:08
 Operator : JC/MD
 Sample : 02558-05
 Misc : 4.49g/10mL/100uL/5.00mL/MSVOA_X/MEOH
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 P001-SS012-0006-01

Integration Parameters: RTEINT.P

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

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

77	14.615	2215	2219	2220	rBV	1425917	1474999	52.31%	2.048%
78	14.640	2220	2223	2226	rVV	2283588	2819458	100.00%	3.914%
79	14.676	2226	2229	2233	rVB2	535317	716911	25.43%	0.995%
80	14.731	2233	2238	2240	rBV2	339234	555040	19.69%	0.770%
81	14.755	2240	2242	2243	rVV	408114	366384	12.99%	0.509%
82	14.780	2243	2246	2250	rVV	1359373	1785884	63.34%	2.479%
83	14.847	2254	2257	2261	rVB2	429590	513228	18.20%	0.712%
84	14.902	2264	2266	2270	rBV2	127588	172866	6.13%	0.240%
85	15.048	2286	2290	2293	rVB2	150157	218778	7.76%	0.304%
86	15.182	2308	2312	2315	rBV	553122	870641	30.88%	1.209%
87	15.377	2339	2344	2347	rBV	633084	903667	32.05%	1.254%
88	15.414	2347	2350	2355	rVV3	302089	473917	16.81%	0.658%
89	15.481	2356	2361	2369	rVB	1340711	2050965	72.74%	2.847%
90	15.615	2378	2383	2386	rBV	1312017	1965955	69.73%	2.729%
91	15.652	2386	2389	2398	rVB	882520	1402292	49.74%	1.947%
92	15.743	2400	2404	2409	rVB3	180151	295438	10.48%	0.410%
93	15.841	2412	2420	2430	rVB3	310792	983122	34.87%	1.365%
94	15.987	2440	2444	2456	rVB	200483	434453	15.41%	0.603%
95	16.091	2457	2461	2469	rBV3	72783	132804	4.71%	0.184%
96	16.261	2486	2489	2492	rBV2	32981	53941	1.91%	0.075%
97	16.353	2500	2504	2512	rVB	113679	211756	7.51%	0.294%
98	16.554	2534	2537	2541	rVV5	40086	85151	3.02%	0.118%
99	16.597	2541	2544	2550	rVV3	67674	129137	4.58%	0.179%
100	16.658	2550	2554	2568	rVB2	56777	153459	5.44%	0.213%

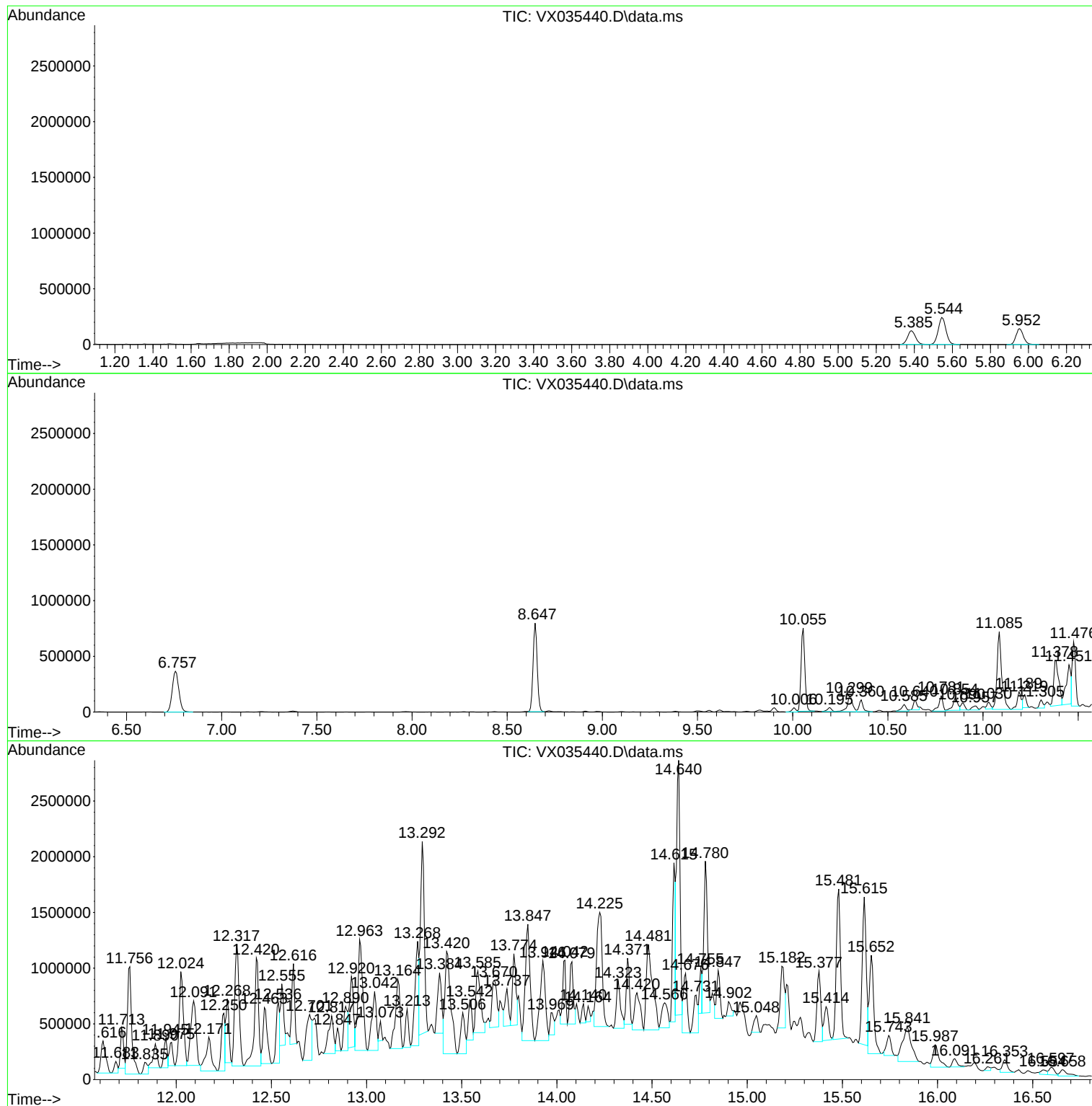
Sum of corrected areas: 72038193

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX050123\
Data File : VX035440.D
Acq On : 01 May 2023 15:08
Operator : JC/MD
Sample : 02558-05
Misc : 4.49g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
P001-SS012-0006-01

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

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX050123\
Data File : VX035440.D
Acq On : 01 May 2023 15:08
Operator : JC/MD
Sample : 02558-05
Misc : 4.49g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
P001-SS012-0006-01

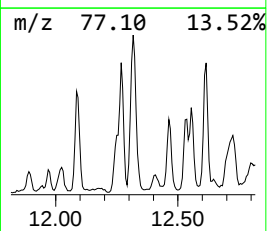
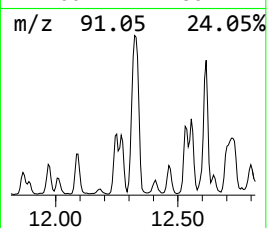
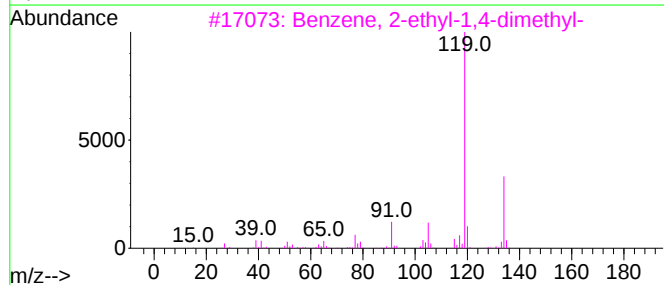
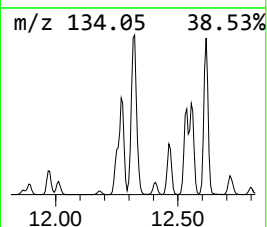
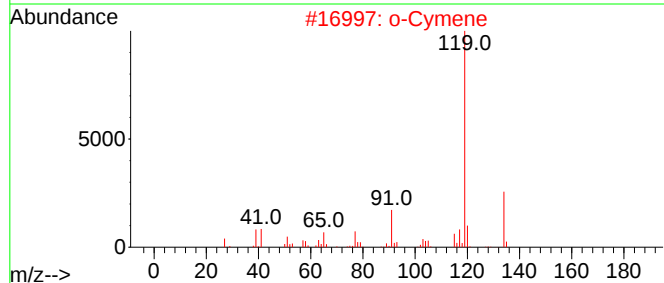
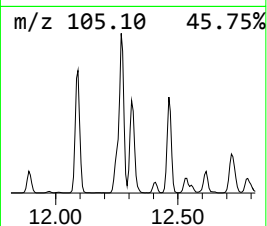
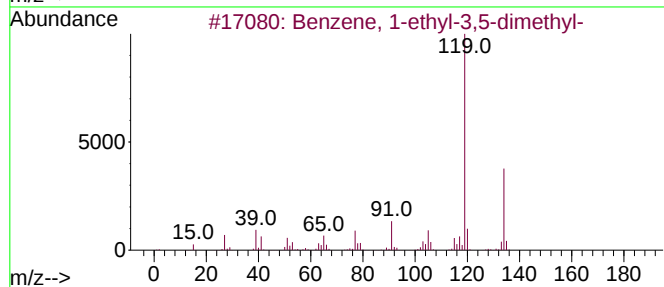
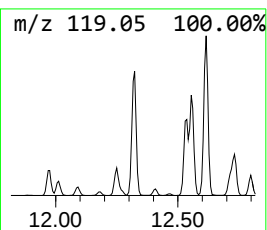
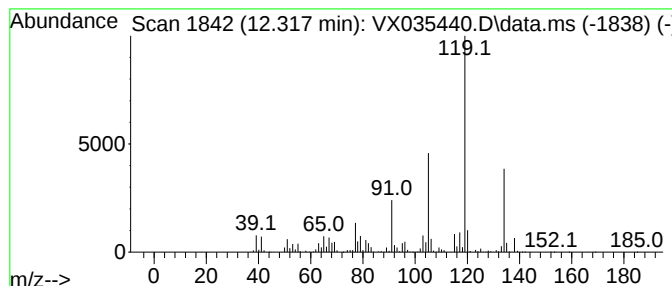
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X042523W.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.317	67.87 ug/l	1772440	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	95
2			o-Cymene	134	C10H14	000527-84-4	95
3			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94
4			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94
5			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	93



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX050123\
Data File : VX035440.D
Acq On : 01 May 2023 15:08
Operator : JC/MD
Sample : 02558-05
Misc : 4.49g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
P001-SS012-0006-01

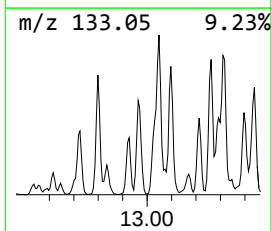
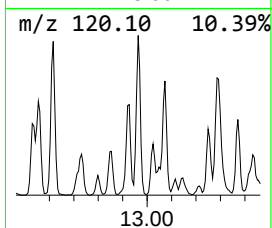
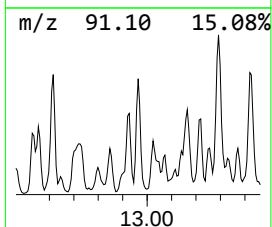
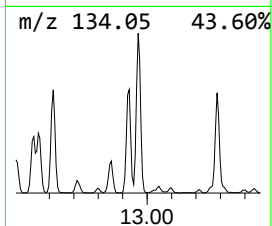
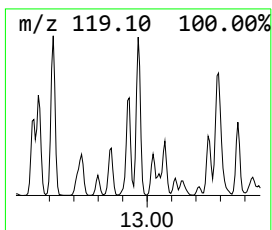
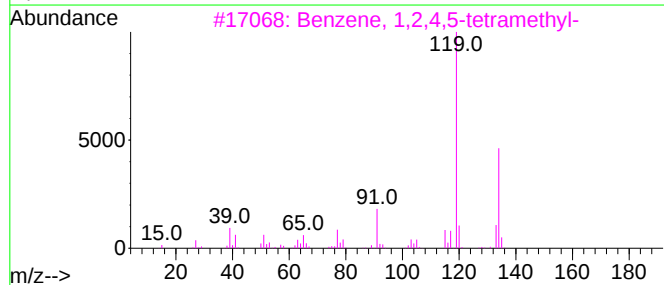
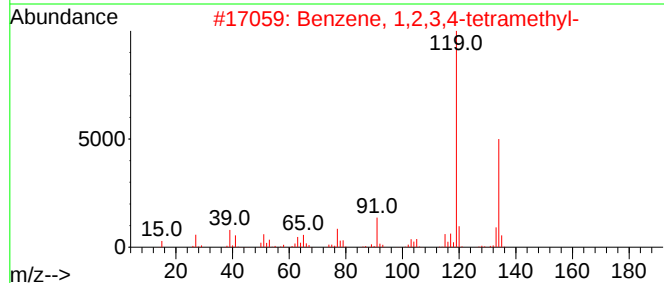
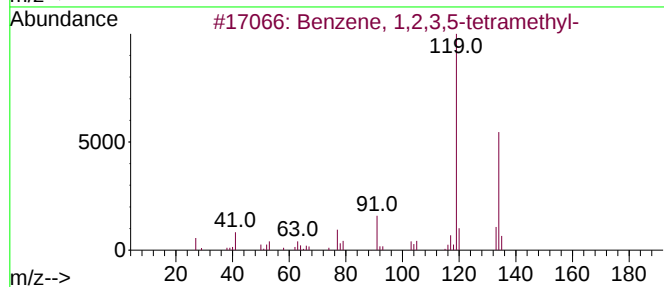
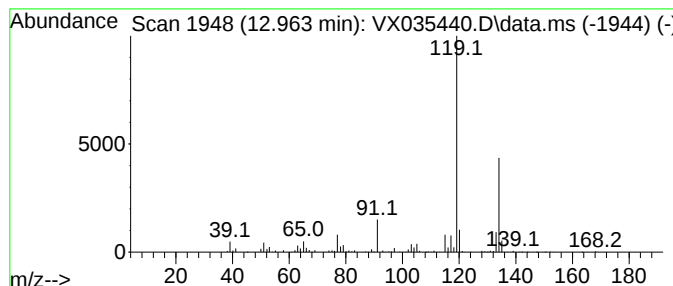
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X042523W.M
Quant Title : SW846 8260

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

Peak Number 2 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.963	67.58 ug/l	1764930	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	95
3			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
4			o-Cymene	134	C10H14	000527-84-4	94
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX050123\
Data File : VX035440.D
Acq On : 01 May 2023 15:08
Operator : JC/MD
Sample : 02558-05
Misc : 4.49g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
P001-SS012-0006-01

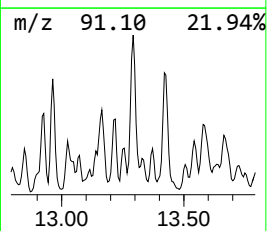
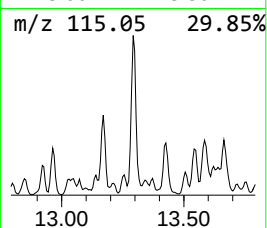
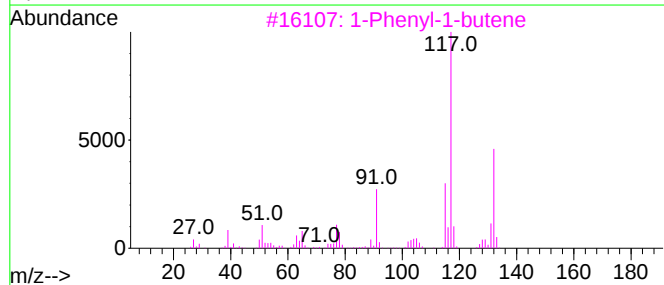
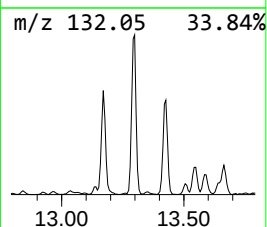
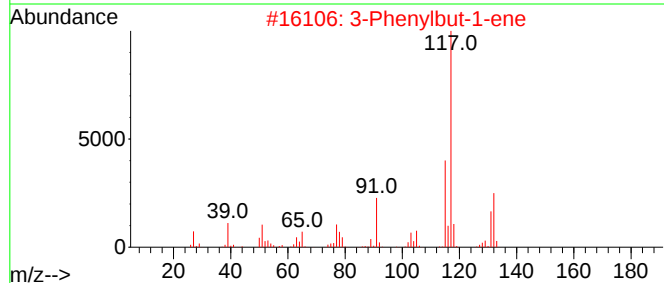
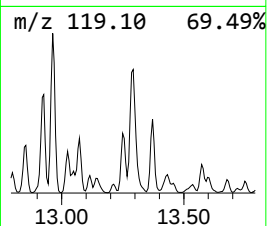
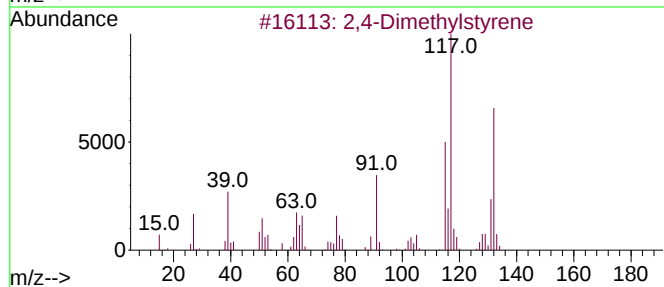
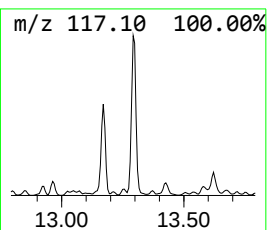
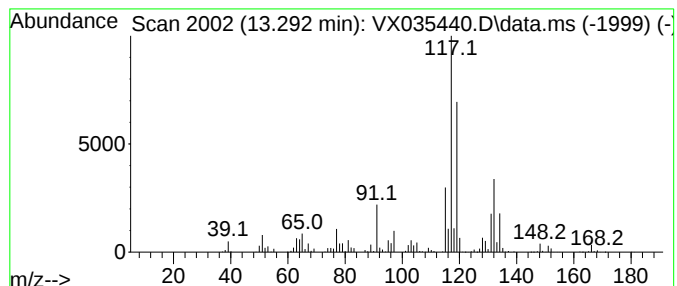
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X042523W.M
Quant Title : SW846 8260

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

Peak Number 3 2,4-Dimethylstyrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	91.56 ug/l	2391340	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,4-Dimethylstyrene	132	C10H12	002234-20-0	95
2			3-Phenylbut-1-ene	132	C10H12	000934-10-1	93
3			1-Phenyl-1-butene	132	C10H12	000824-90-8	92
4			Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	91
5			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX050123\
Data File : VX035440.D
Acq On : 01 May 2023 15:08
Operator : JC/MD
Sample : 02558-05
Misc : 4.49g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
P001-SS012-0006-01

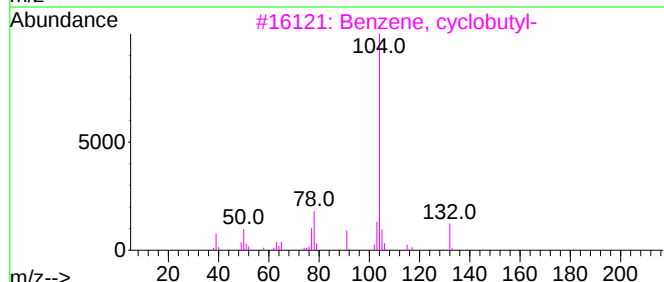
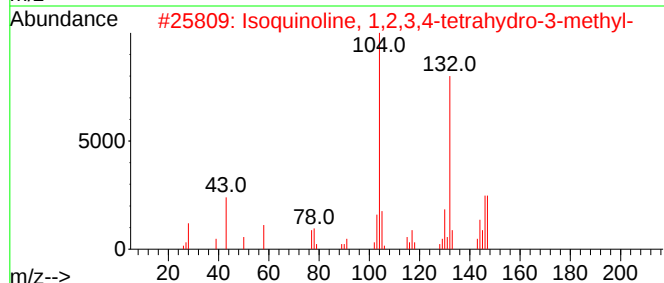
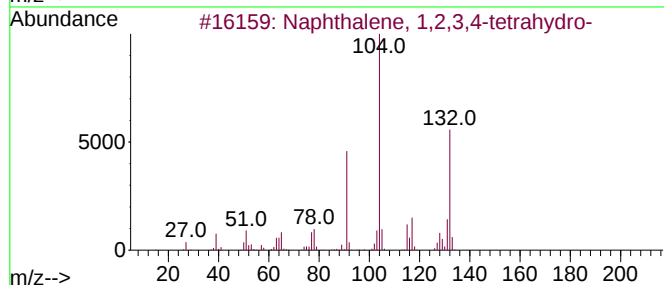
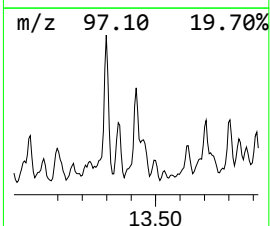
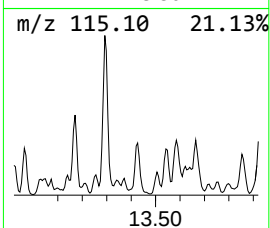
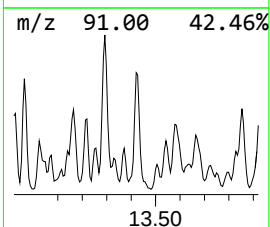
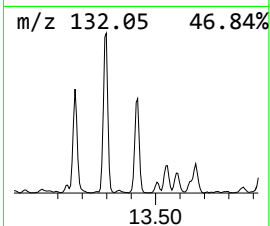
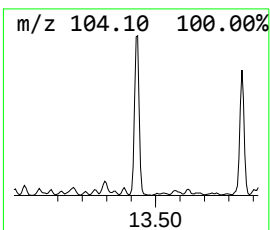
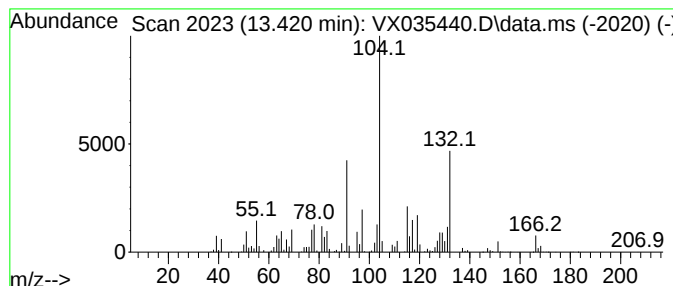
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X042523W.M
Quant Title : SW846 8260

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

Peak Number 4 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.420	74.20 ug/l	1937870	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	89
2			Isoquinoline, 1,2,3,4-tetrahydro...	147	C10H13N	029726-60-1	40
3			Benzene, cyclobutyl-	132	C10H12	004392-30-7	38
4			Pentafluoropropionic acid, 2-phe...	268	C11H9F5O2	081089-48-7	35
5			3(2H)-Furanone, dihydro-2,2-dime...	190	C12H14O2	063678-00-2	27



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX050123\
Data File : VX035440.D
Acq On : 01 May 2023 15:08
Operator : JC/MD
Sample : 02558-05
Misc : 4.49g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
P001-SS012-0006-01

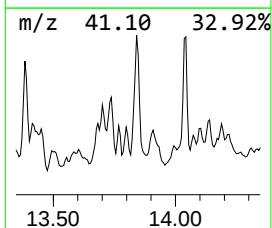
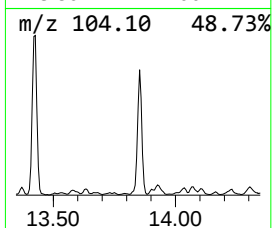
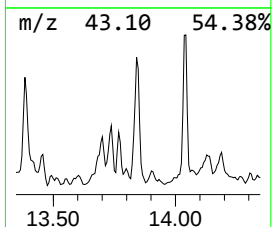
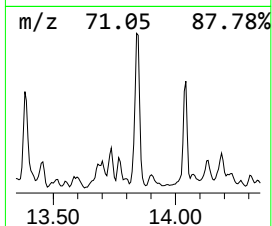
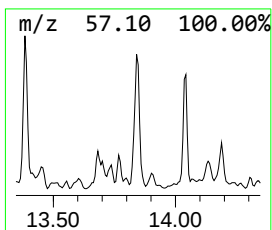
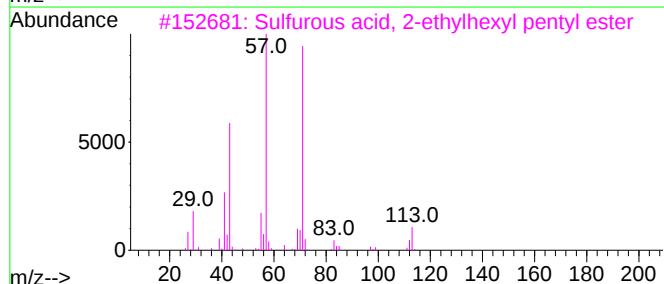
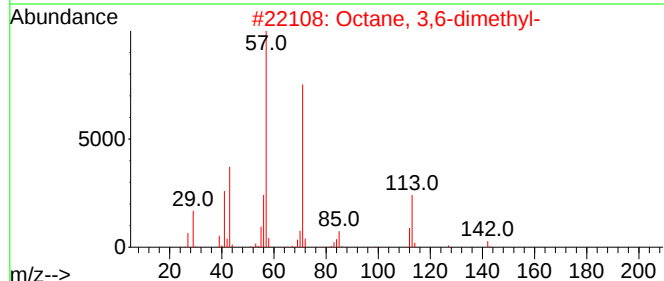
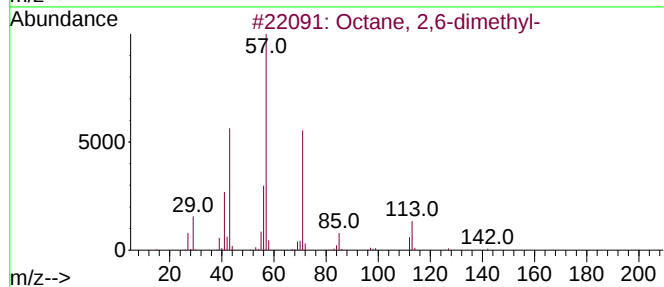
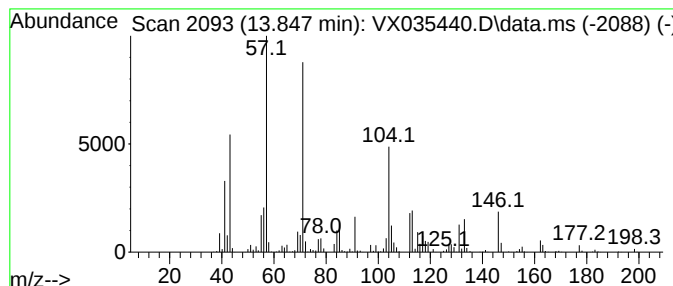
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X042523W.M
Quant Title : SW846 8260

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

Peak Number 5 Octane, 2,6-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.847	70.86 ug/l	1850500	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	60
2			Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	49
3			Sulfurous acid, 2-ethylhexyl pen...	264	C13H28O3S	1000309-18-9	47
4			Undecane, 6,6-dimethyl-	184	C13H28	017312-76-4	47
5			Sulfurous acid, 2-ethylhexyl und...	348	C19H40O3S	1000309-19-4	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX050123\
Data File : VX035440.D
Acq On : 01 May 2023 15:08
Operator : JC/MD
Sample : 02558-05
Misc : 4.49g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
P001-SS012-0006-01

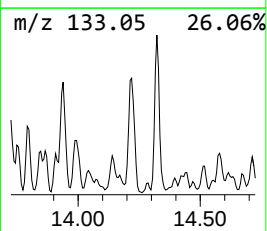
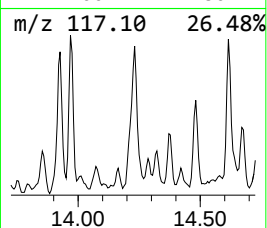
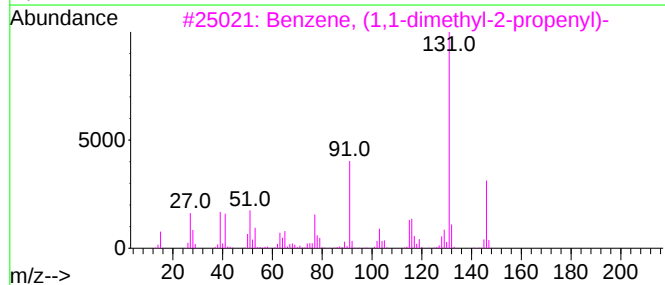
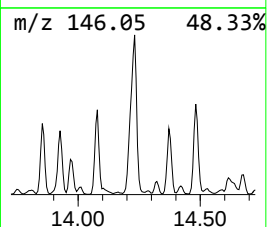
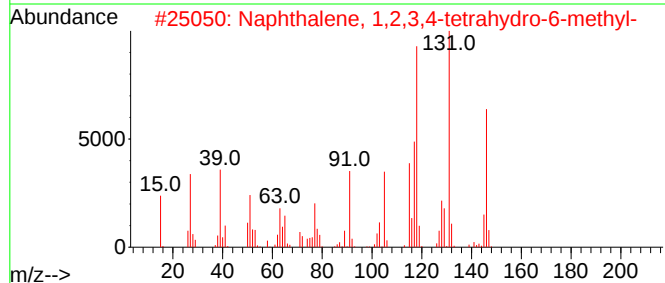
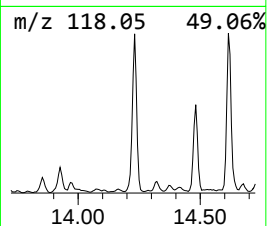
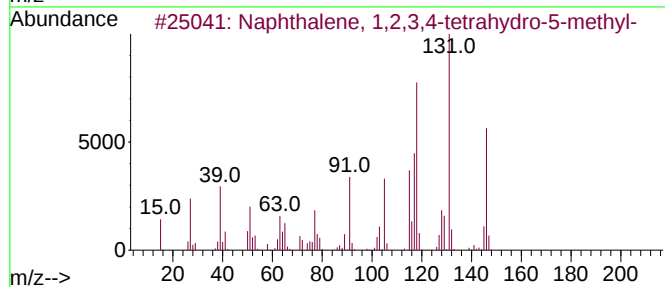
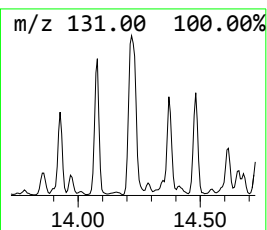
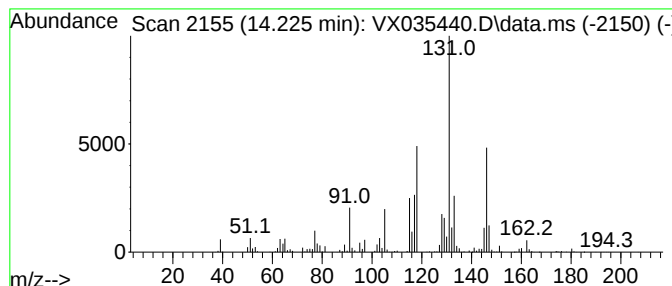
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X042523W.M
Quant Title : SW846 8260

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

Peak Number 6 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.225	79.87 ug/l	2085810	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	95
2			Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	91
3			Benzene, (1,1-dimethyl-2-propenyl)-	146	C11H14	018321-36-3	64
4			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	64
5			Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	58



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX050123\
Data File : VX035440.D
Acq On : 01 May 2023 15:08
Operator : JC/MD
Sample : 02558-05
Misc : 4.49g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
P001-SS012-0006-01

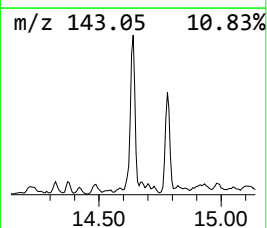
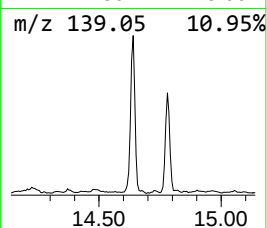
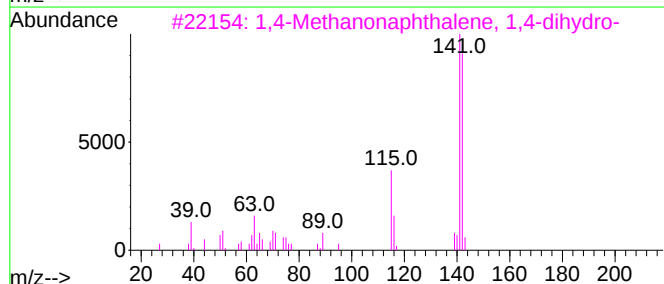
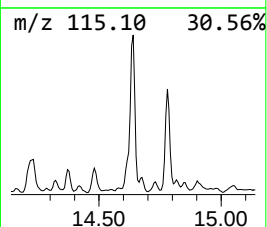
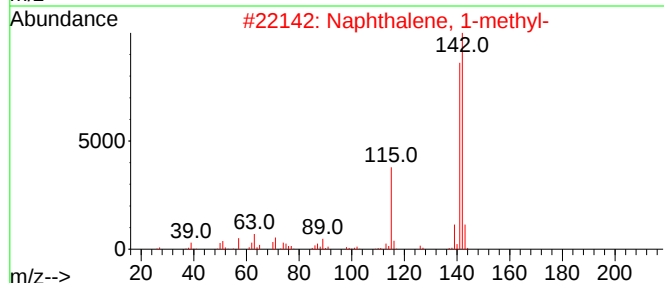
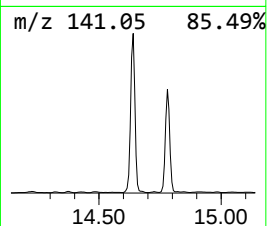
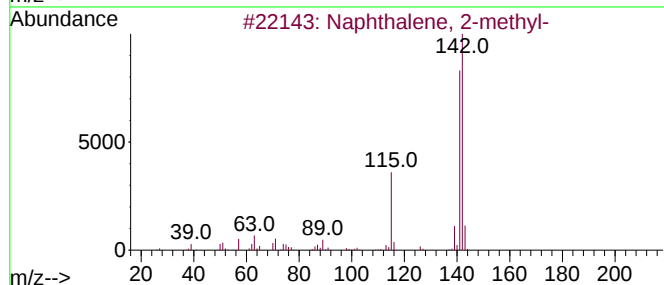
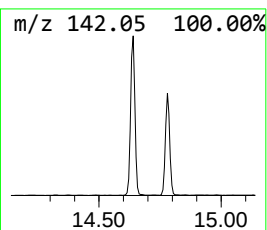
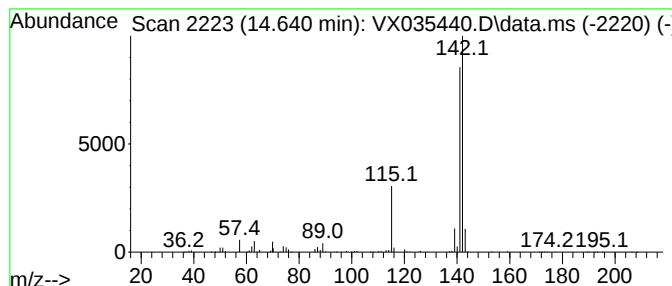
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X042523W.M
Quant Title : SW846 8260

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

Peak Number 7 1,4-Methanonaphthalene, 1,4... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.640	107.96 ug/l	2819460	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	91
2			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	91
3			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
4			Benzocycloheptatriene	142	C11H10	000264-09-5	91
5			3-Indolecarbonitrile	142	C9H6N2	005457-28-3	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX050123\
Data File : VX035440.D
Acq On : 01 May 2023 15:08
Operator : JC/MD
Sample : 02558-05
Misc : 4.49g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
P001-SS012-0006-01

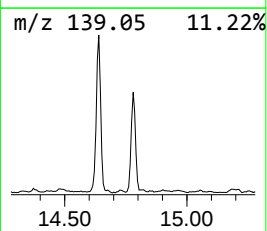
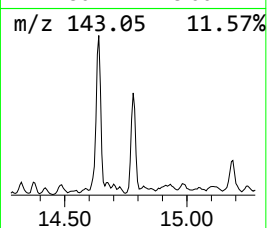
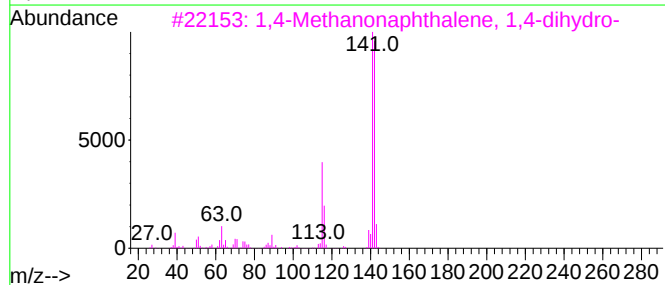
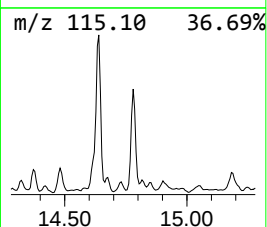
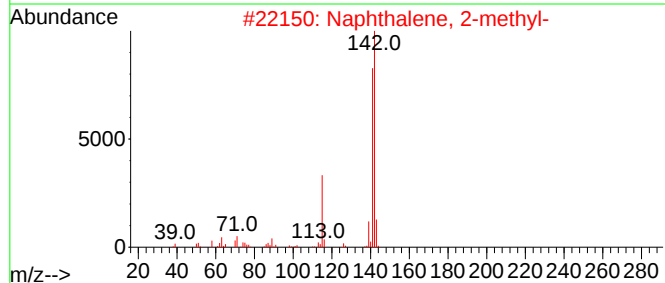
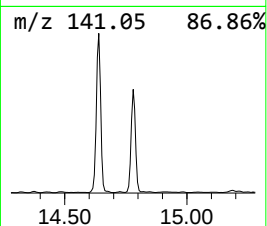
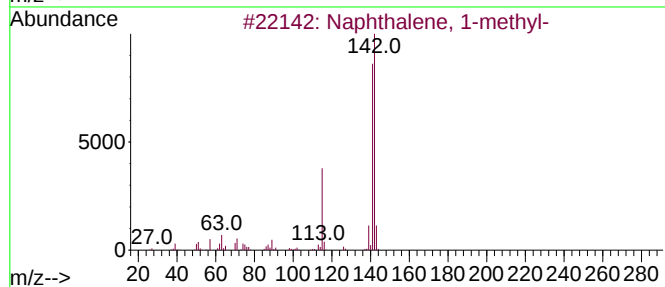
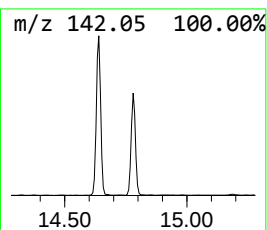
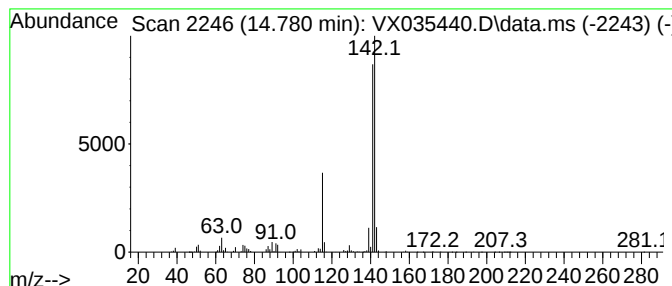
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X042523W.M
Quant Title : SW846 8260

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

Peak Number 8 Benzocycloheptatriene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.780	68.38 ug/l	1785880	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
2			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	94
3			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
4			Benzocycloheptatriene	142	C11H10	000264-09-5	90
5			Bicyclo[4.4.1]undeca-1,3,5,7,9-p...	142	C11H10	002443-46-1	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX050123\
Data File : VX035440.D
Acq On : 01 May 2023 15:08
Operator : JC/MD
Sample : 02558-05
Misc : 4.49g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
P001-SS012-0006-01

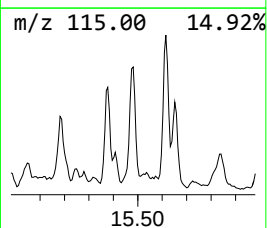
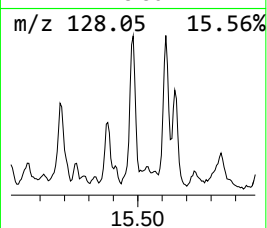
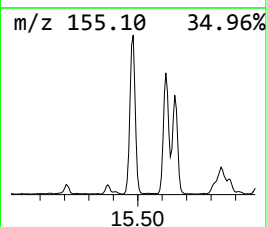
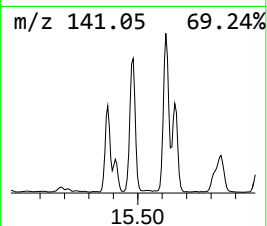
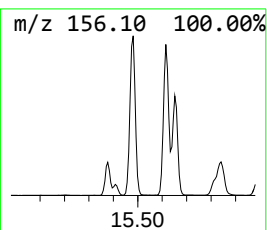
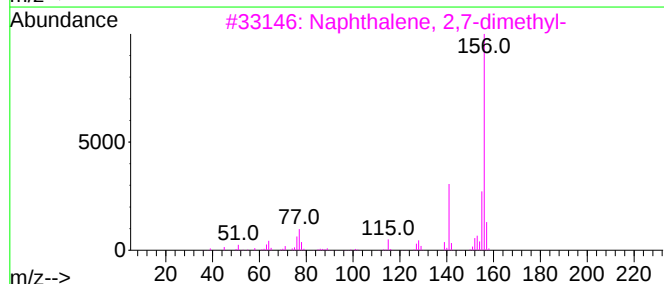
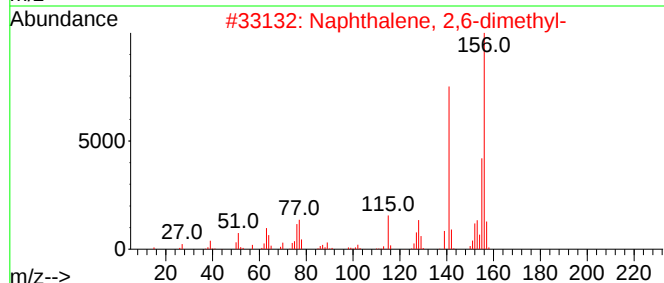
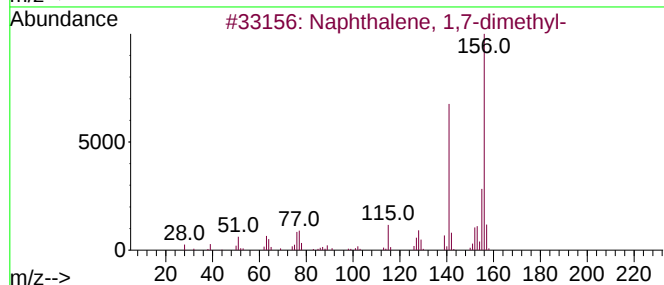
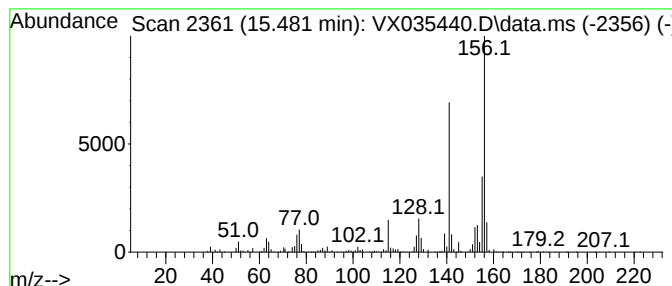
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X042523W.M
Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1,7-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.481	78.53 ug/l	2050970	1,4-Dichlorobenzene-d4	12.024

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
2			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
3			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96
4			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96
5			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX050123\
Data File : VX035440.D
Acq On : 01 May 2023 15:08
Operator : JC/MD
Sample : 02558-05
Misc : 4.49g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
P001-SS012-0006-01

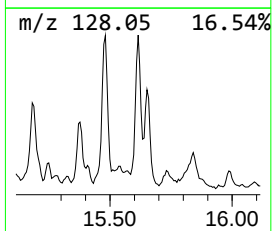
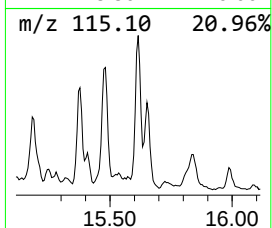
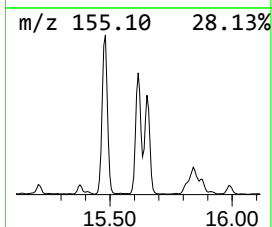
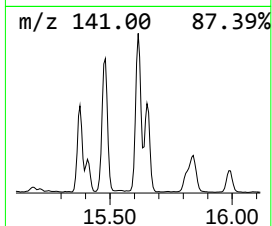
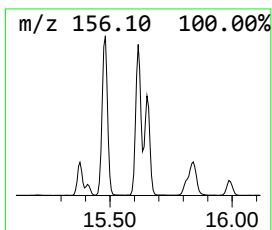
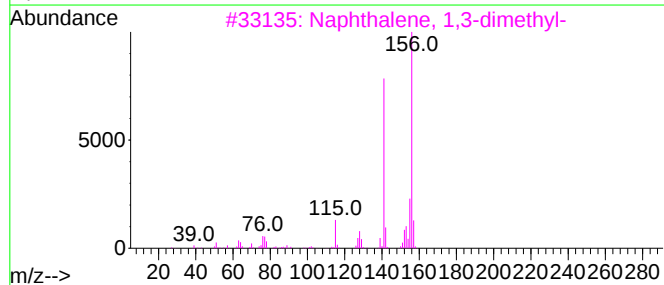
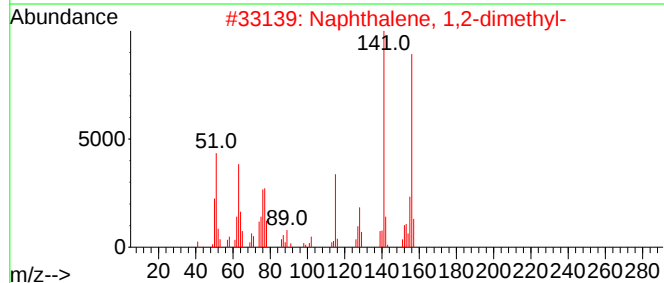
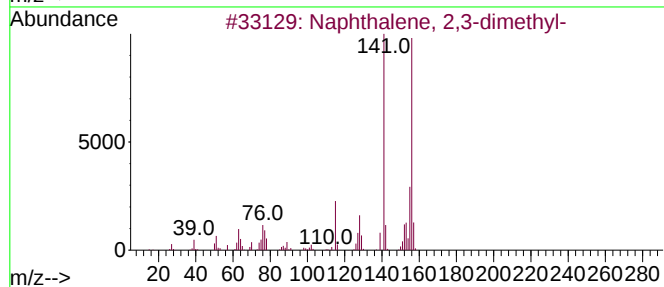
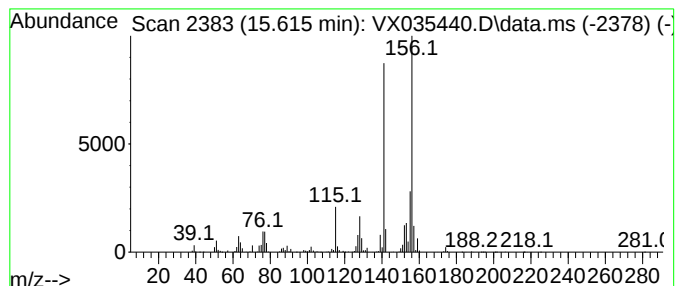
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X042523W.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 2,3-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.615	75.28 ug/l	1965960	1,4-Dichlorobenzene-d4	12.024

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2		Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	97
3		Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97
4		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
5		Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX050123\
Data File : VX035440.D
Acq On : 01 May 2023 15:08
Operator : JC/MD
Sample : 02558-05
Misc : 4.49g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
P001-SS012-0006-01

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

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1-ethy...	12.317	67.9	ug/l	1772440	4	12.024	1305820	50.0
Benzene, 1,2,3,...	12.963	67.6	ug/l	1764930	4	12.024	1305820	50.0
2,4-Dimethylsty...	13.292	91.6	ug/l	2391340	4	12.024	1305820	50.0
Naphthalene, 1,...	13.420	74.2	ug/l	1937870	4	12.024	1305820	50.0
Octane, 2,6-dim...	13.847	70.9	ug/l	1850500	4	12.024	1305820	50.0
Naphthalene, 1,...	14.225	79.9	ug/l	2085810	4	12.024	1305820	50.0
1,4-Methanonaph...	14.640	108.0	ug/l	2819460	4	12.024	1305820	50.0
Benzocyclohepta...	14.780	68.4	ug/l	1785880	4	12.024	1305820	50.0
Naphthalene, 1,...	15.481	78.5	ug/l	2050970	4	12.024	1305820	50.0
Naphthalene, 2,...	15.615	75.3	ug/l	1965960	4	12.024	1305820	50.0