

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX082221\
 Data File : VX023880.D
 Acq On : 20 Aug 2021 16:23
 Operator : JC/MD
 Sample : M3414-01
 Misc : 5.02g/10mL/100uL/5.00mL/MSVOA_X/MEOH
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 B-1

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: VX023880.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.794	275	280	292	rVB	124635	249451	2.54%	0.228%
2	5.397	690	707	714	rBV	79408	254640	2.59%	0.232%
3	5.495	714	723	729	rBV	210550	616910	6.27%	0.563%
4	5.818	763	776	790	rBV	279113	810120	8.24%	0.739%
5	5.964	790	800	816	rBV	96195	258894	2.63%	0.236%
6	6.153	816	831	838	rBV3	58865	216127	2.20%	0.197%
7	6.245	838	846	856	rVV2	187647	582481	5.92%	0.532%
8	6.604	894	905	915	rBV	251943	603270	6.13%	0.551%
9	6.769	923	932	939	rBV	232218	583590	5.93%	0.533%
10	7.379	1021	1032	1043	rVB2	270813	701697	7.13%	0.641%
11	7.562	1056	1062	1072	rVV	143409	318604	3.24%	0.291%
12	7.982	1121	1131	1143	rVB2	125148	351189	3.57%	0.321%
13	8.104	1143	1151	1155	rBV3	103377	220394	2.24%	0.201%
14	8.275	1174	1179	1183	rBV	497318	833126	8.47%	0.760%
15	8.311	1183	1185	1192	rVB	235746	346022	3.52%	0.316%
16	8.439	1201	1206	1215	rBV	541671	970555	9.87%	0.886%
17	8.604	1227	1233	1236	rBV3	240564	472825	4.81%	0.432%
18	8.653	1236	1241	1249	rVB	514608	934820	9.50%	0.853%
19	8.775	1249	1261	1265	rBV2	128145	281405	2.86%	0.257%
20	8.915	1277	1284	1290	rBV	566516	884015	8.99%	0.807%
21	8.982	1290	1295	1302	rVB3	122032	232595	2.36%	0.212%
22	9.104	1307	1315	1321	rBV2	133344	278518	2.83%	0.254%
23	9.384	1355	1361	1371	rBV2	177913	319516	3.25%	0.292%
24	9.500	1375	1380	1387	rBV2	338060	684422	6.96%	0.625%
25	9.561	1387	1390	1395	rVB	171350	244382	2.48%	0.223%
26	9.823	1428	1433	1437	rVV2	163208	325906	3.31%	0.297%
27	9.866	1437	1440	1443	rVV2	157319	272573	2.77%	0.249%
28	9.909	1443	1447	1452	rVB	1032775	1453764	14.78%	1.327%
29	10.012	1457	1464	1468	rBV	804084	1099527	11.18%	1.004%
30	10.061	1468	1472	1476	rVV	472808	623813	6.34%	0.569%
31	10.201	1489	1495	1500	rVB	966218	1352209	13.75%	1.234%
32	10.305	1500	1512	1518	rVV	3310601	4970562	50.54%	4.537%
33	10.366	1518	1522	1532	rVB	873165	1322535	13.45%	1.207%
34	10.591	1554	1559	1563	rVV	279463	382907	3.89%	0.350%
35	10.677	1569	1573	1577	rBV	210190	279335	2.84%	0.255%
36	10.787	1583	1591	1594	rBV	335538	551663	5.61%	0.504%

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX082221\
 Data File : VX023880.D
 Acq On : 20 Aug 2021 16:23
 Operator : JC/MD
 Sample : M3414-01
 Misc : 5.02g/10mL/100uL/5.00mL/MSVOA_X/MEOH
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 B-1

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

37	10.860	1599	1603	1607	rVB3	243283	332987	3.39%	0.304%
38	10.969	1614	1621	1625	rBV	327360	494365	5.03%	0.451%
39	11.030	1627	1631	1635	rVV	294536	416558	4.24%	0.380%
40	11.085	1635	1640	1643	rVV2	935271	1588786	16.15%	1.450%
41	11.110	1643	1644	1651	rVV	689980	686773	6.98%	0.627%
42	11.195	1654	1658	1665	rVB	728845	932849	9.48%	0.852%
43	11.311	1672	1677	1684	rVV	926154	1166520	11.86%	1.065%
44	11.402	1684	1692	1696	rVV	1710568	3135716	31.88%	2.862%
45	11.457	1696	1701	1704	rVV	2293453	3260672	33.15%	2.976%
46	11.482	1704	1705	1712	rVB	970398	741444	7.54%	0.677%
47	11.616	1723	1727	1735	rBV	1076046	1506830	15.32%	1.375%
48	11.689	1735	1739	1741	rVV	177825	216724	2.20%	0.198%
49	11.719	1741	1744	1746	rVV	225760	261044	2.65%	0.238%
50	11.756	1746	1750	1759	rVB	7583202	9835176	100.00%	8.978%
51	11.841	1759	1764	1766	rBV	218642	278251	2.83%	0.254%
52	11.872	1766	1769	1770	rVV	202822	235841	2.40%	0.215%
53	11.896	1770	1773	1777	rVB	301729	370413	3.77%	0.338%
54	11.975	1783	1786	1789	rVV	422328	459443	4.67%	0.419%
55	12.024	1789	1794	1800	rVV3	650769	1363339	13.86%	1.244%
56	12.091	1800	1805	1811	rVB	2184642	3245725	33.00%	2.963%
57	12.177	1816	1819	1826	rVB2	537245	628626	6.39%	0.574%
58	12.250	1826	1831	1833	rBV3	1755794	2448491	24.90%	2.235%
59	12.274	1833	1835	1838	rVV	1417528	1452097	14.76%	1.326%
60	12.323	1838	1843	1849	rVB3	2886928	4387135	44.61%	4.005%
61	12.426	1854	1860	1863	rVV2	794016	1048205	10.66%	0.957%
62	12.469	1863	1867	1873	rVB	685115	914811	9.30%	0.835%
63	12.536	1873	1878	1880	rBV	1491540	1988868	20.22%	1.815%
64	12.561	1880	1882	1887	rVV	1408730	1630676	16.58%	1.489%
65	12.615	1887	1891	1894	rVV	2470674	2934378	29.84%	2.679%
66	12.646	1894	1896	1900	rVV2	416979	500257	5.09%	0.457%
67	12.707	1900	1906	1915	rVV2	898068	2050349	20.85%	1.872%
68	12.798	1916	1921	1926	rVV4	190789	387150	3.94%	0.353%
69	12.853	1926	1930	1934	rVV	553048	711211	7.23%	0.649%
70	12.926	1934	1942	1945	rVV	1507313	2387294	24.27%	2.179%
71	12.969	1945	1949	1954	rVV	2159095	2769695	28.16%	2.528%
72	13.030	1954	1959	1960	rVV	485754	649175	6.60%	0.593%
73	13.048	1960	1962	1964	rVV2	655466	734469	7.47%	0.670%
74	13.073	1964	1966	1969	rVV	496402	561611	5.71%	0.513%
75	13.115	1970	1973	1976	rVV	434375	588289	5.98%	0.537%
76	13.170	1976	1982	1986	rVV3	1471713	2232706	22.70%	2.038%

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX082221\
 Data File : VX023880.D
 Acq On : 20 Aug 2021 16:23
 Operator : JC/MD
 Sample : M3414-01
 Misc : 5.02g/10mL/100uL/5.00mL/MSVOA_X/MEOH
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 B-1

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

77	13.213	1986	1989	1992	rVV2	519001	696130	7.08%	0.635%
78	13.268	1992	1998	1999	rVV5	804159	1407886	14.31%	1.285%
79	13.298	1999	2003	2008	rVV3	2625910	4028005	40.96%	3.677%
80	13.371	2012	2015	2020	rVV	607727	923116	9.39%	0.843%
81	13.426	2020	2024	2032	rVB5	339063	680000	6.91%	0.621%
82	13.506	2032	2037	2040	rBV2	385698	579241	5.89%	0.529%
83	13.548	2040	2044	2047	rVV	628516	961209	9.77%	0.877%
84	13.585	2047	2050	2057	rVV3	1003368	1948426	19.81%	1.779%
85	13.646	2057	2060	2061	rVV	235892	308216	3.13%	0.281%
86	13.670	2061	2064	2069	rVV2	610894	1142388	11.62%	1.043%
87	13.719	2069	2072	2077	rVV2	169437	365190	3.71%	0.333%
88	13.780	2077	2082	2088	rVV	1623877	2177756	22.14%	1.988%
89	13.835	2088	2091	2094	rVV4	154085	262681	2.67%	0.240%
90	13.932	2100	2107	2111	rBV4	370021	757947	7.71%	0.692%
91	13.975	2111	2114	2117	rVV	216066	267917	2.72%	0.245%
92	14.042	2121	2125	2127	rVB3	238815	281480	2.86%	0.257%
93	14.079	2127	2131	2135	rBV	570436	652786	6.64%	0.596%
94	14.219	2150	2154	2164	rVB	544028	856249	8.71%	0.782%
95	14.322	2167	2171	2176	rBV	287678	395354	4.02%	0.361%
96	14.377	2176	2180	2184	rVB	358335	422553	4.30%	0.386%
97	14.639	2215	2223	2228	rBV	1917366	2407826	24.48%	2.198%
98	14.780	2240	2246	2251	rBV	835397	1045445	10.63%	0.954%
99	15.481	2355	2361	2371	rVB	177144	289778	2.95%	0.265%
100	15.615	2377	2383	2387	rBV	171521	273495	2.78%	0.250%

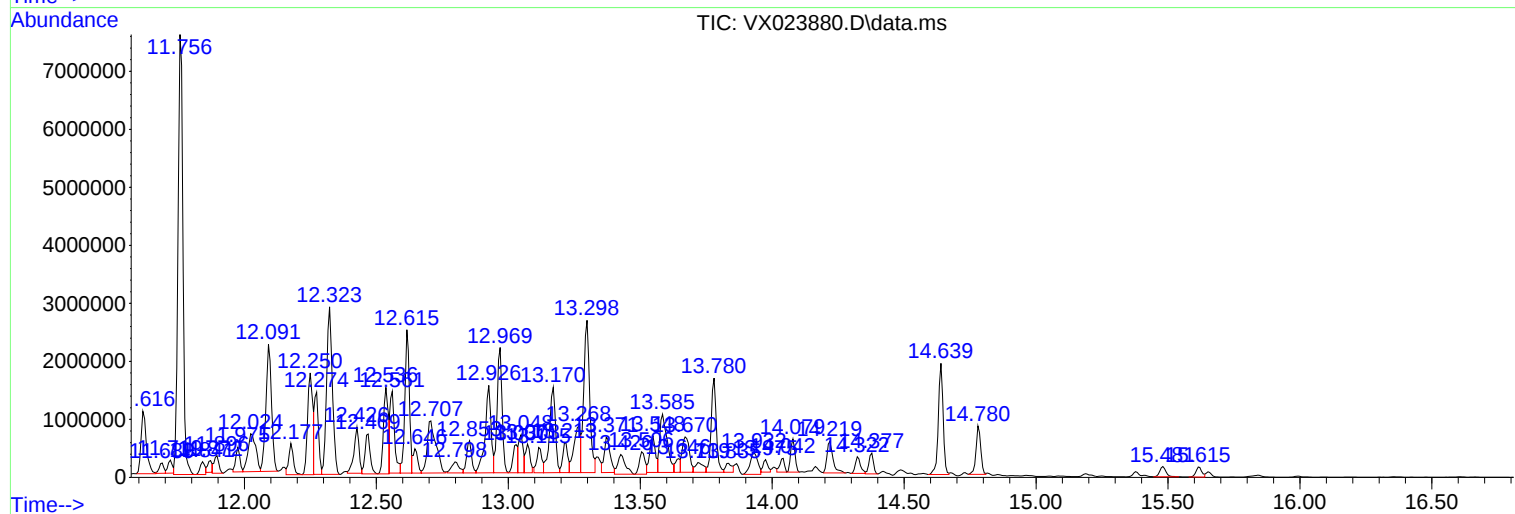
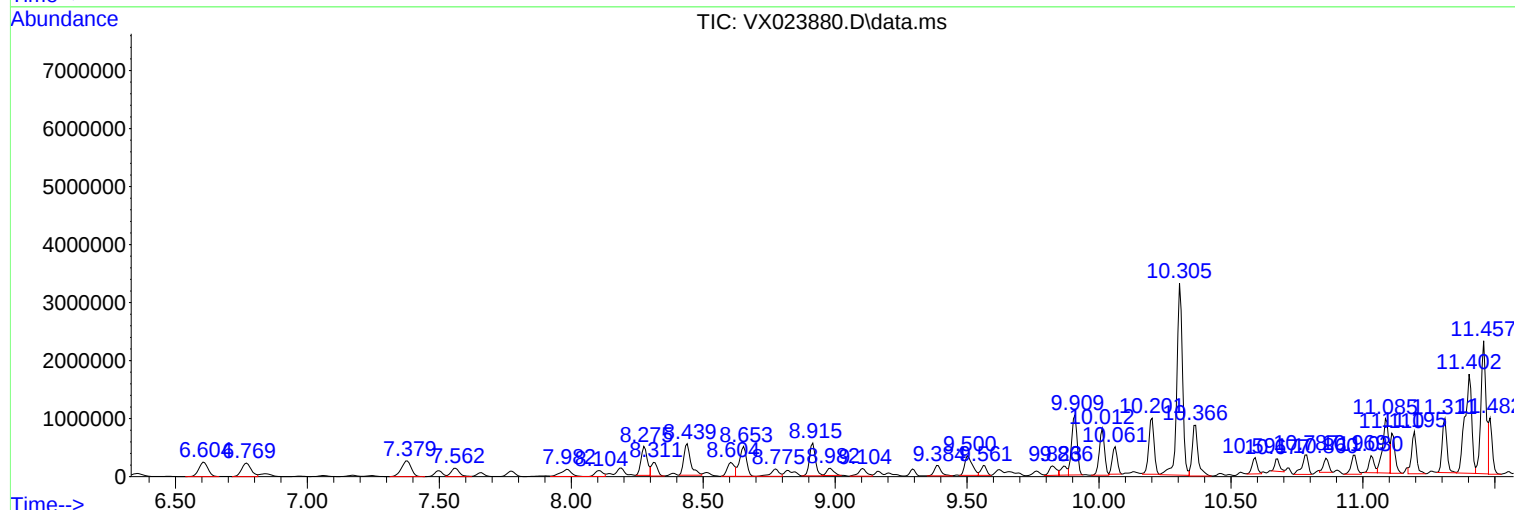
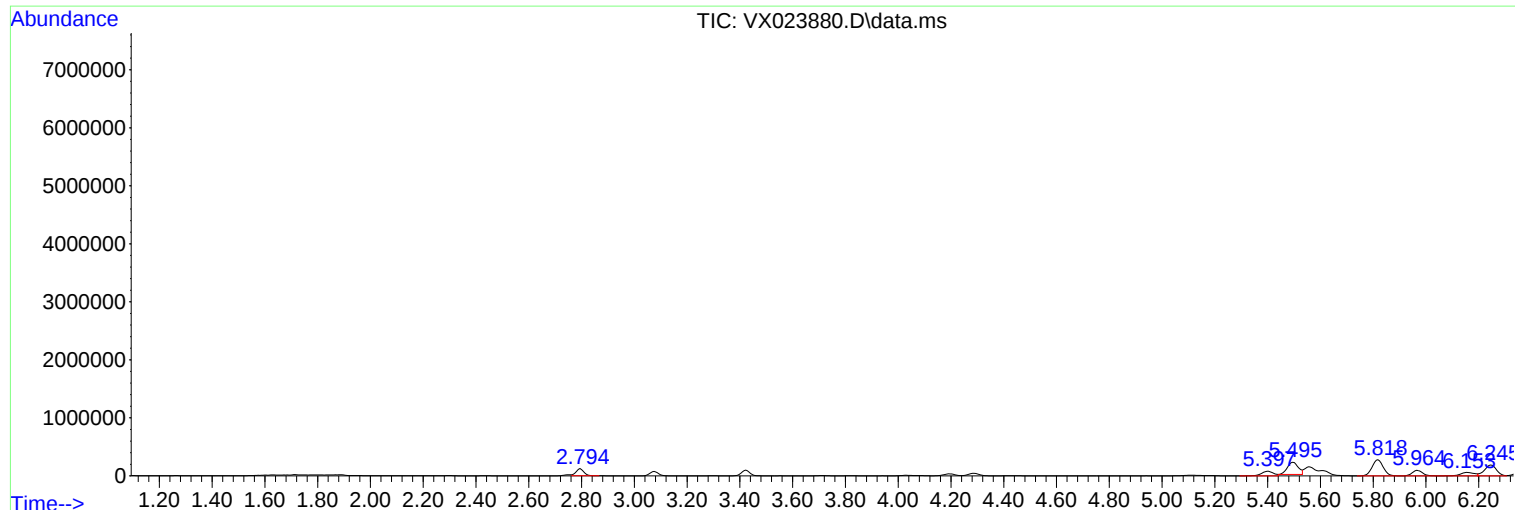
Sum of corrected areas: 109550385

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX082221\
Data File : VX023880.D
Acq On : 20 Aug 2021 16:23
Operator : JC/MD
Sample : M3414-01
Misc : 5.02g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B-1

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX082221\
Data File : VX023880.D
Acq On : 20 Aug 2021 16:23
Operator : JC/MD
Sample : M3414-01
Misc : 5.02g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B-1

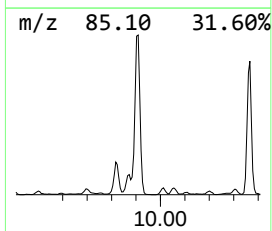
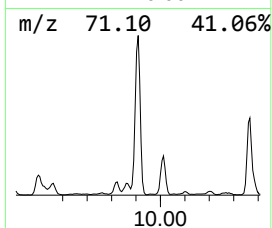
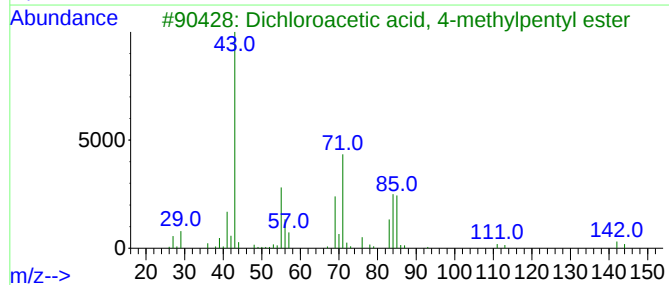
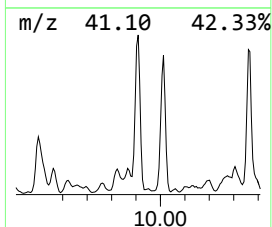
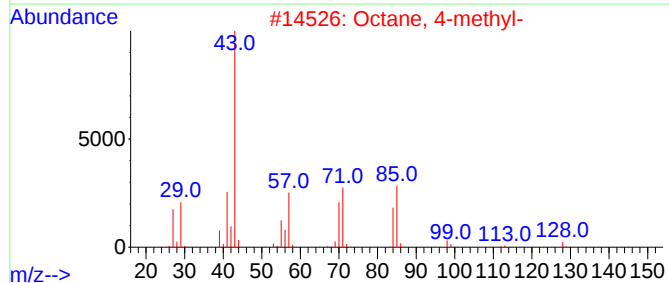
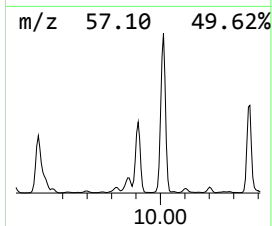
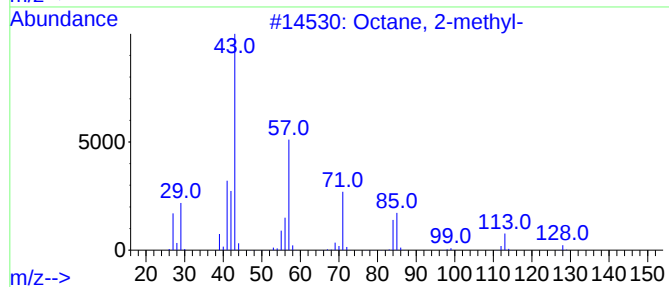
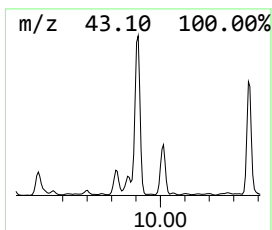
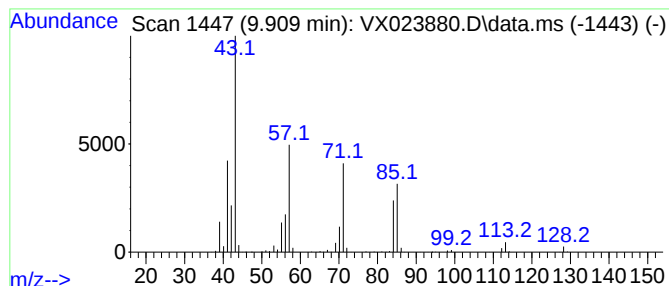
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081721W.M
Quant Title : SW846 8260

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

Peak Number 1 Octane, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.909	116.52 ug/l	1453760	Chlorobenzene-d5	10.061

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 2-methyl-	128	C9H20	003221-61-2	80
2			Octane, 4-methyl-	128	C9H20	002216-34-4	72
3			Dichloroacetic acid, 4-methylpen...	212	C8H14Cl2O2	1000282-43-7	64
4			Dodecane, 5-methyl-	184	C13H28	017453-93-9	59
5			Octane	114	C8H18	000111-65-9	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX082221\
Data File : VX023880.D
Acq On : 20 Aug 2021 16:23
Operator : JC/MD
Sample : M3414-01
Misc : 5.02g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B-1

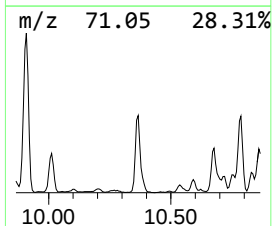
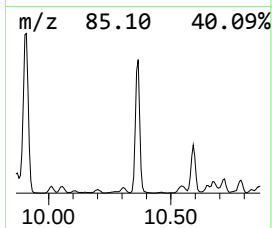
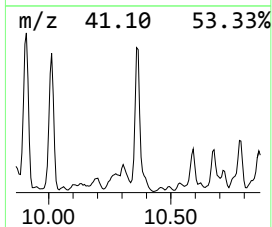
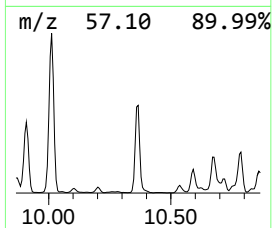
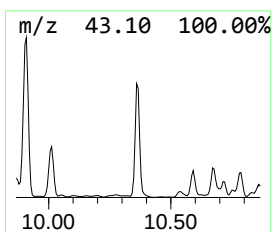
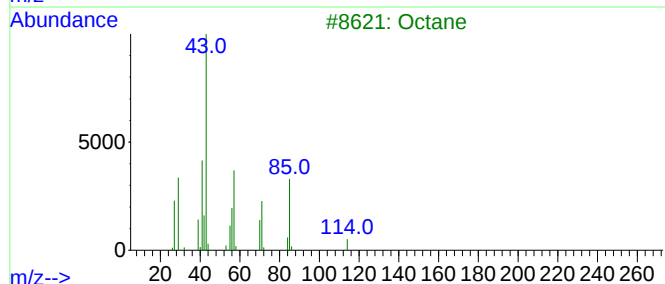
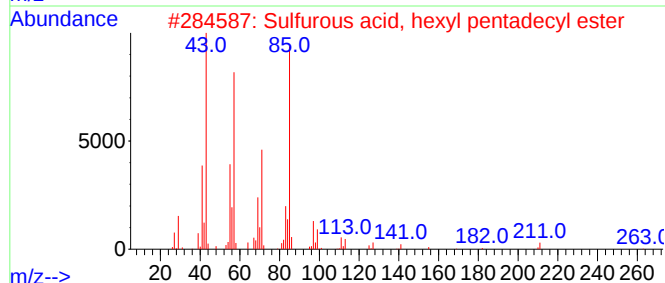
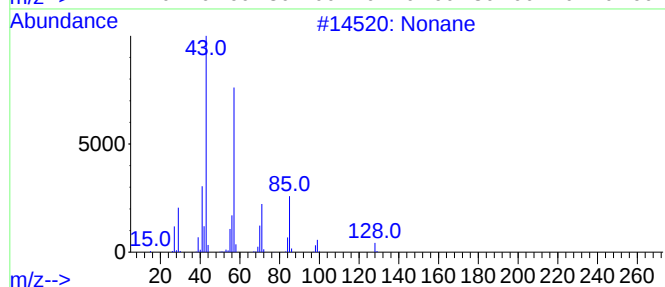
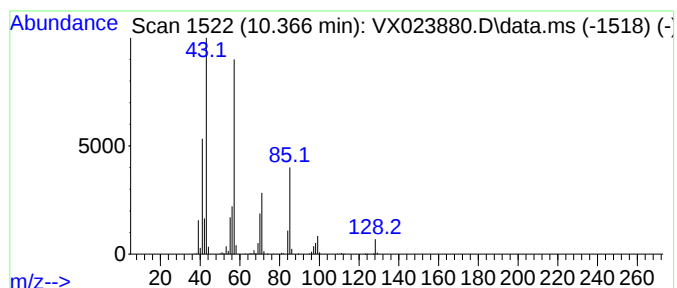
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081721W.M
Quant Title : SW846 8260

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

Peak Number 2 Nonane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.366	106.00 ug/l	1322540	Chlorobenzene-d5	10.061

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane	128	C9H20	000111-84-2	94
2			Sulfurous acid, hexyl pentadecyl...	376	C21H44O3S	1000309-13-7	72
3			Octane	114	C8H18	000111-65-9	64
4			Decane, 2,4,6-trimethyl-	184	C13H28	062108-27-4	53
5			4,4-Dimethyl octane	142	C10H22	015869-95-1	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX082221\
Data File : VX023880.D
Acq On : 20 Aug 2021 16:23
Operator : JC/MD
Sample : M3414-01
Misc : 5.02g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B-1

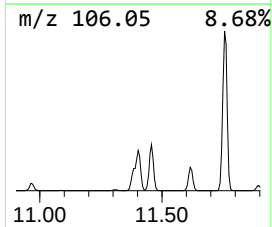
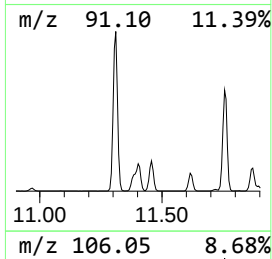
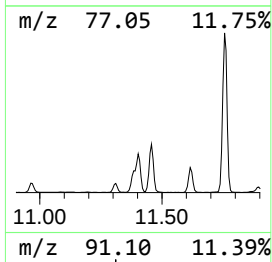
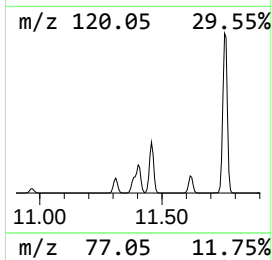
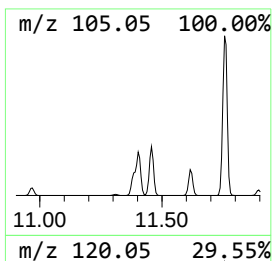
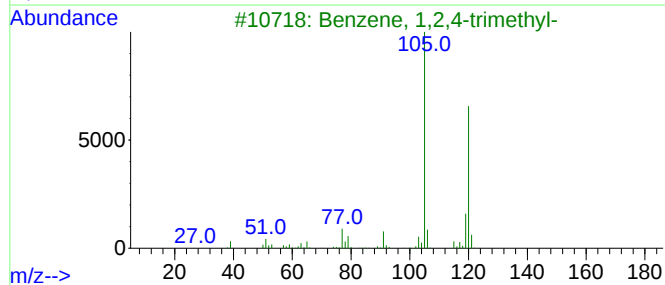
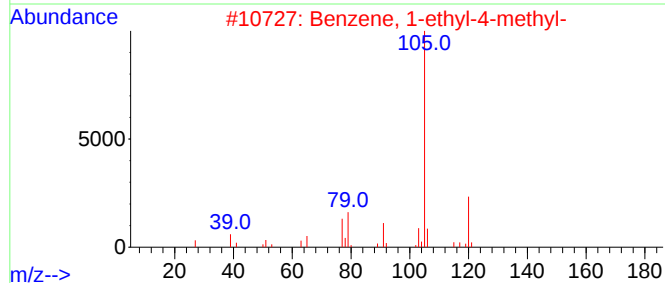
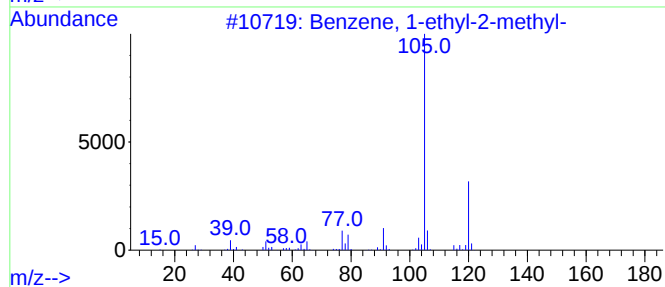
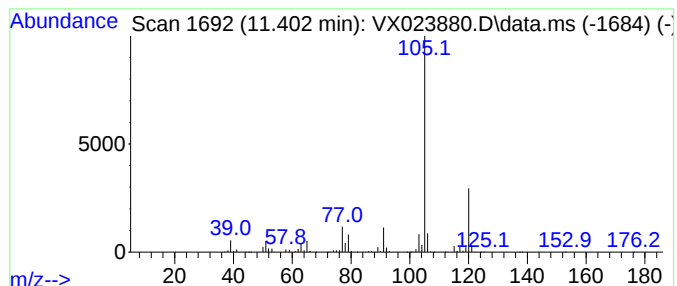
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081721W.M
Quant Title : SW846 8260

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

Peak Number 3 Benzene, 1-ethyl-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.402	115.00 ug/l	3135720	1,4-Dichlorobenzene-d4	12.030

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	94
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4			Mesitylene	120	C9H12	000108-67-8	91
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX082221\
Data File : VX023880.D
Acq On : 20 Aug 2021 16:23
Operator : JC/MD
Sample : M3414-01
Misc : 5.02g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B-1

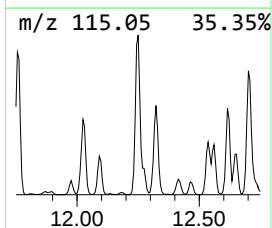
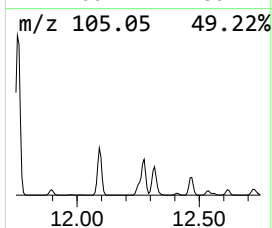
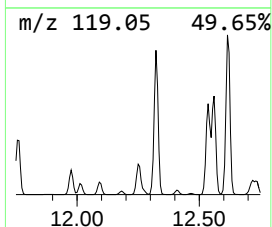
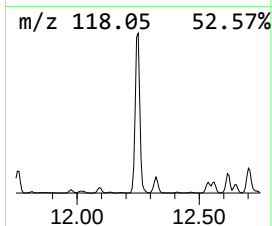
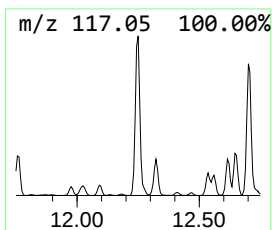
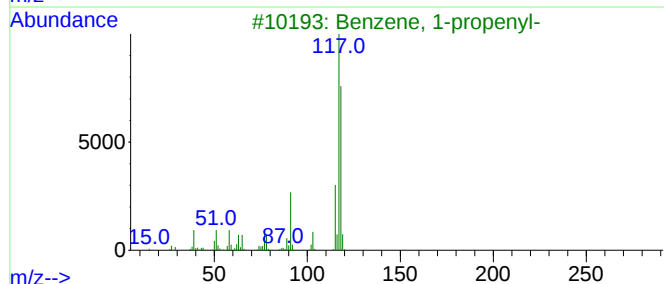
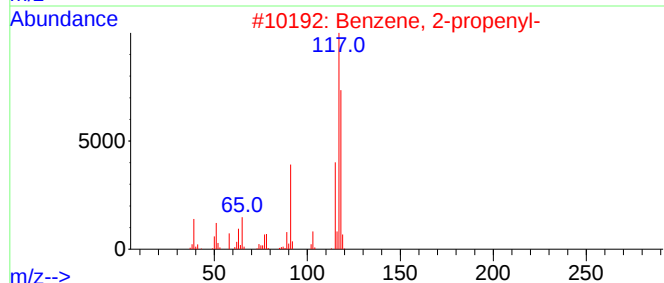
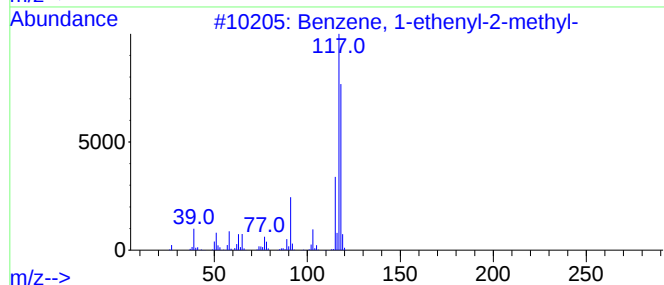
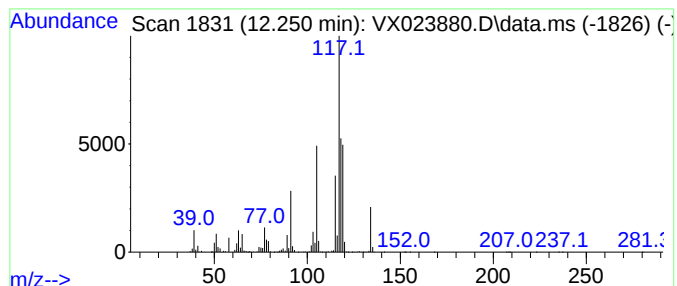
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081721W.M
Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1-ethenyl-2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.250	89.80 ug/l	2448490	1,4-Dichlorobenzene-d4	12.030

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	60
2			Benzene, 2-propenyl-	118	C9H10	000300-57-2	60
3			Benzene, 1-propenyl-	118	C9H10	000637-50-3	60
4			Benzene, cyclopropyl-	118	C9H10	000873-49-4	60
5			Deltacyclene	118	C9H10	007785-10-6	55



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX082221\
Data File : VX023880.D
Acq On : 20 Aug 2021 16:23
Operator : JC/MD
Sample : M3414-01
Misc : 5.02g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B-1

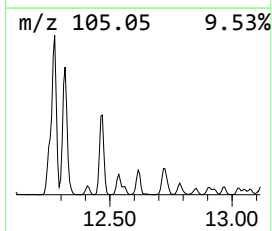
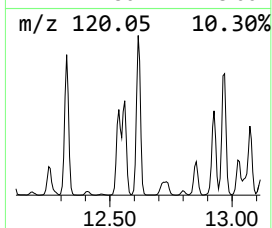
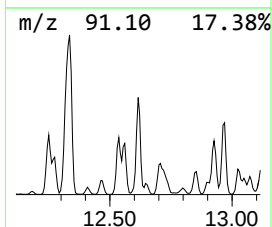
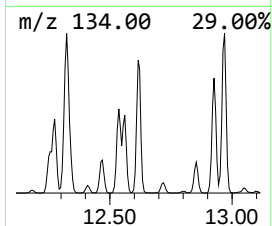
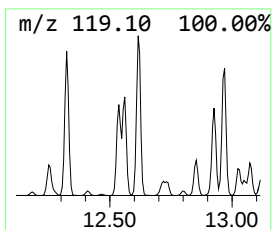
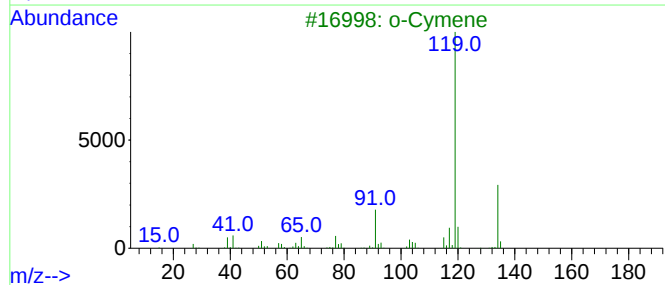
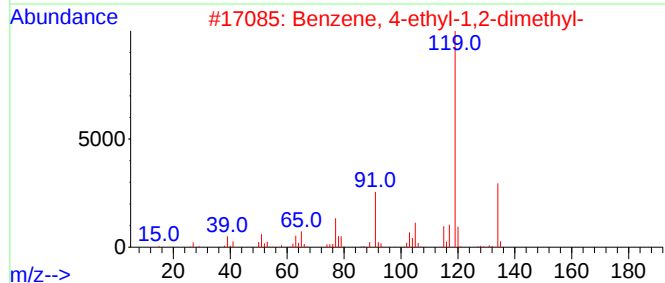
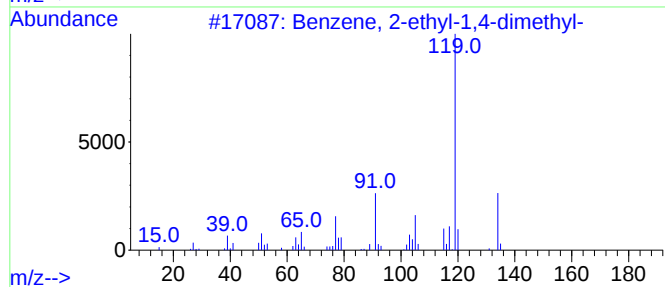
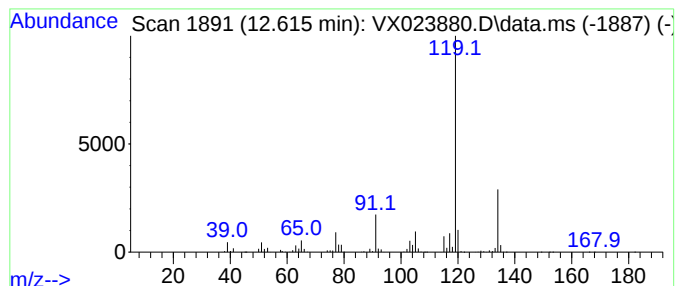
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081721W.M
Quant Title : SW846 8260

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

Peak Number 6 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.616	107.62 ug/l	2934380	1,4-Dichlorobenzene-d4	12.030

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			o-Cymene	134	C10H14	000527-84-4	95
4			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX082221\
Data File : VX023880.D
Acq On : 20 Aug 2021 16:23
Operator : JC/MD
Sample : M3414-01
Misc : 5.02g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampled :
B-1

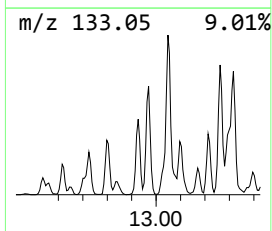
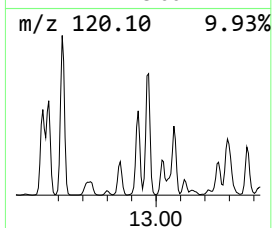
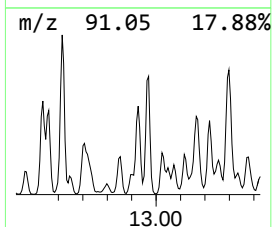
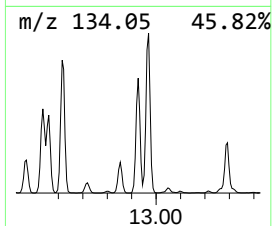
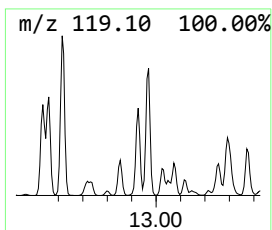
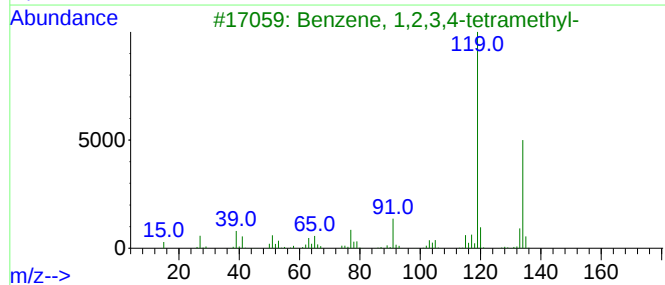
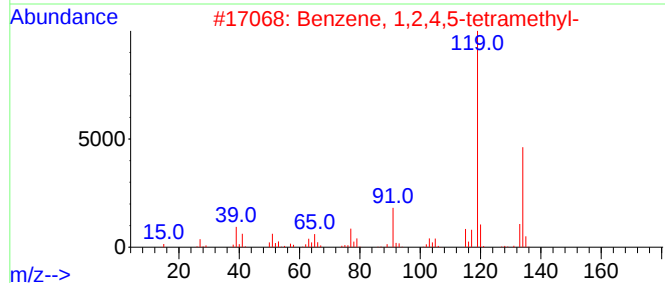
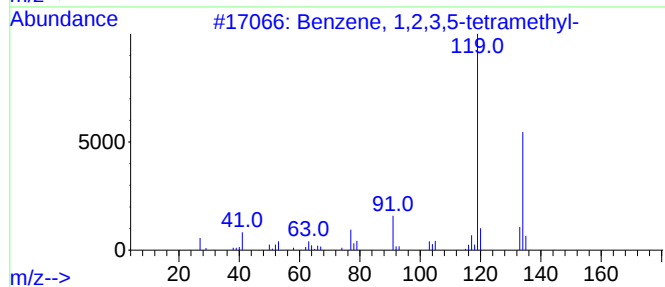
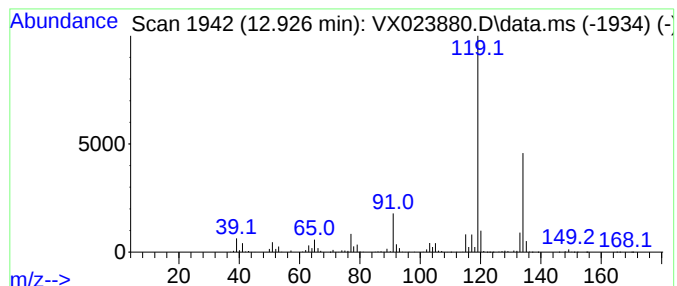
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081721W.M
Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.926	87.55 ug/l	2387290	1,4-Dichlorobenzene-d4	12.030

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX082221\
Data File : VX023880.D
Acq On : 20 Aug 2021 16:23
Operator : JC/MD
Sample : M3414-01
Misc : 5.02g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B-1

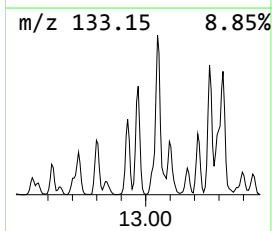
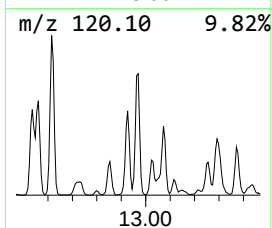
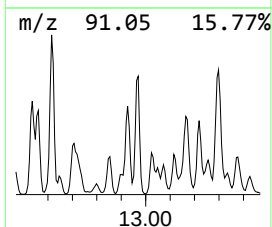
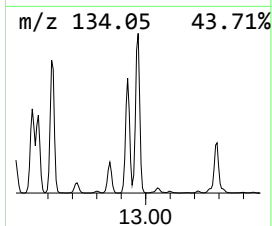
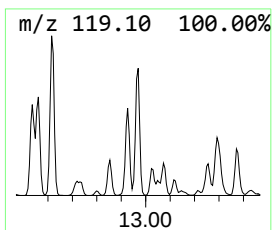
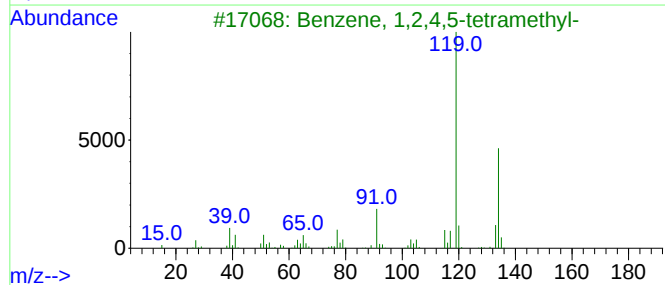
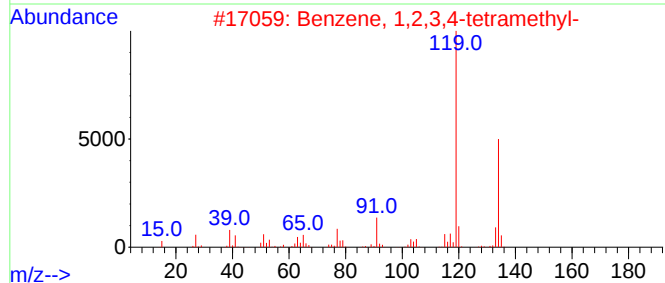
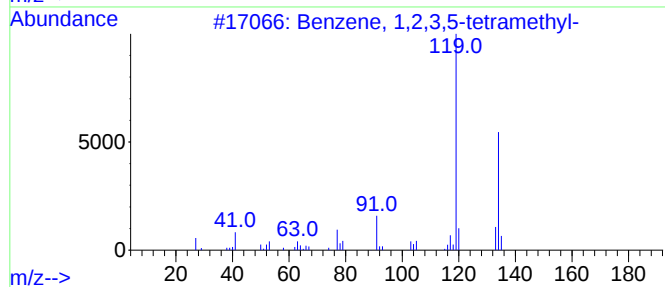
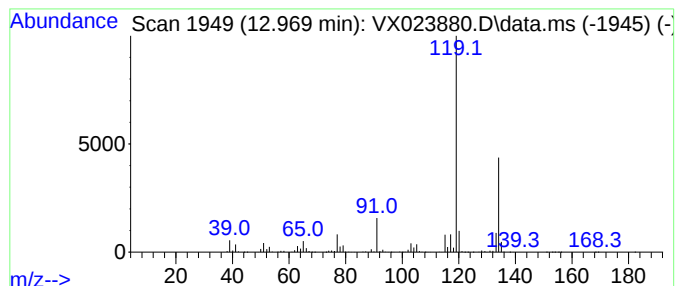
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081721W.M
Quant Title : SW846 8260

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

Peak Number 8 o-Cymene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.969	101.58 ug/l	2769700	1,4-Dichlorobenzene-d4	12.030

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			p-Cymene	134	C10H14	000099-87-6	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX082221\
Data File : VX023880.D
Acq On : 20 Aug 2021 16:23
Operator : JC/MD
Sample : M3414-01
Misc : 5.02g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B-1

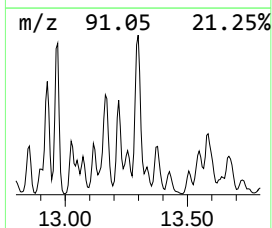
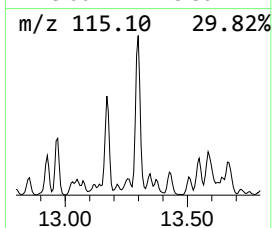
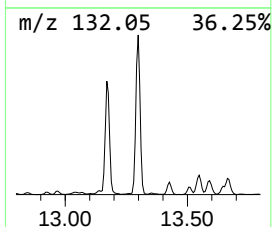
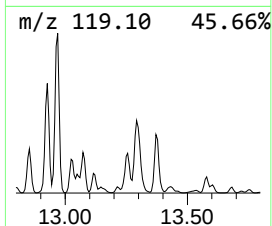
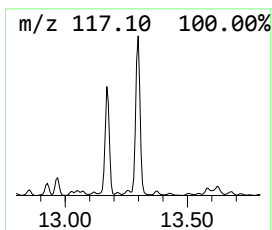
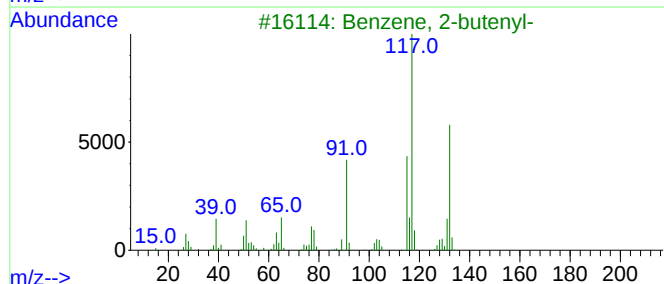
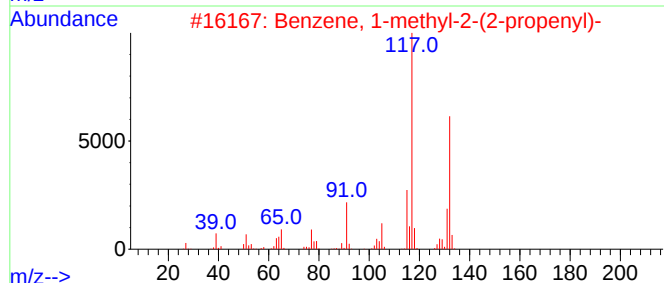
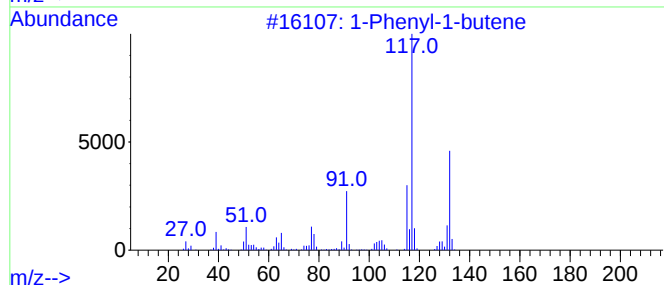
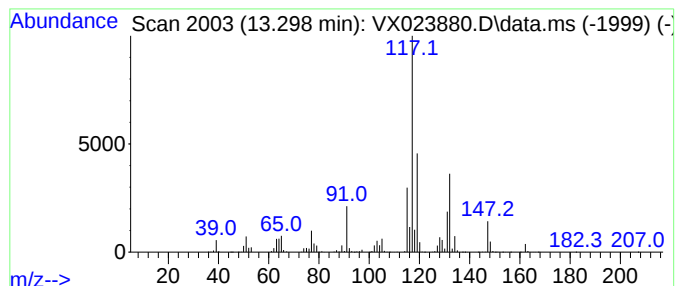
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081721W.M
Quant Title : SW846 8260

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

Peak Number 9 1-Phenyl-1-butene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.298	147.73 ug/l	4028010	1,4-Dichlorobenzene-d4	12.030

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenyl-1-butene	132	C10H12	000824-90-8	86
2			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	70
3			Benzene, 2-butenyl-	132	C10H12	001560-06-1	70
4			Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	70
5			3-Phenylbut-1-ene	132	C10H12	000934-10-1	70



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX082221\
Data File : VX023880.D
Acq On : 20 Aug 2021 16:23
Operator : JC/MD
Sample : M3414-01
Misc : 5.02g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B-1

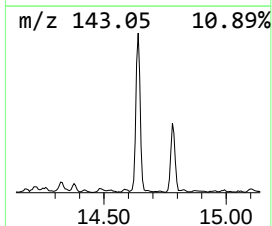
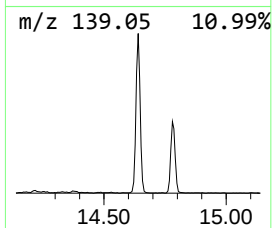
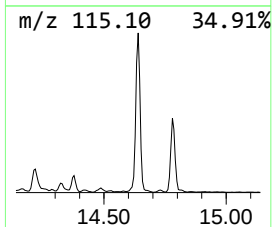
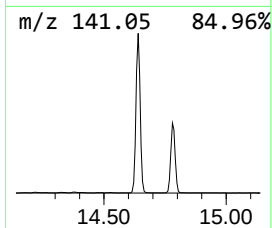
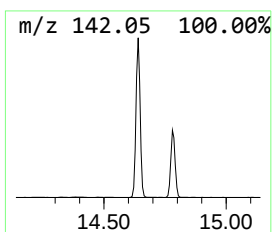
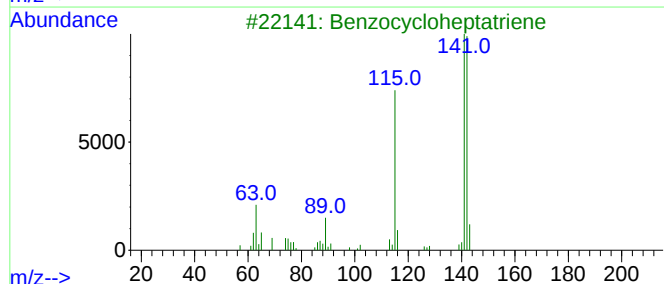
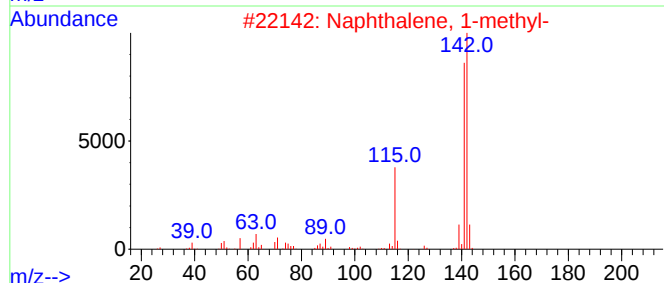
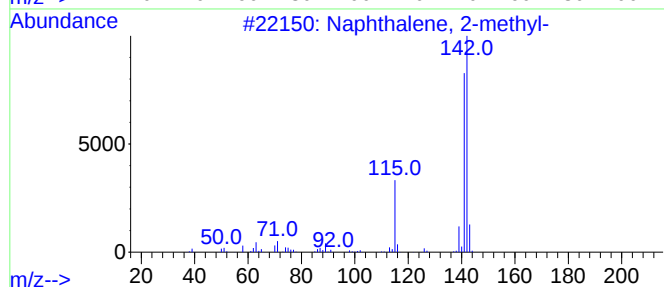
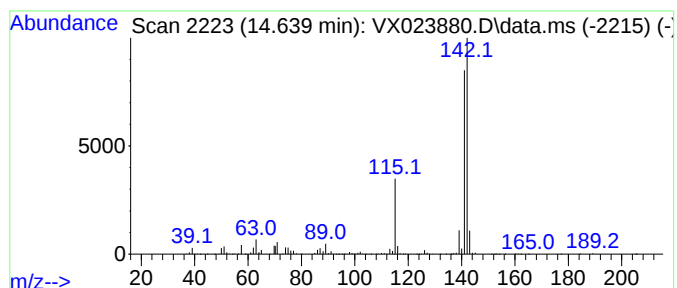
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081721W.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.640	88.31 ug/l	2407830	1,4-Dichlorobenzene-d4	12.030

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2			Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3			Benzocycloheptatriene	142	C11H10	000264-09-5	91
4			1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
5			Spiro(tricyclo[6.2.1.0(2,7)]unde...	186	C13H14O	1000210-02-3	74



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX082221\
Data File : VX023880.D
Acq On : 20 Aug 2021 16:23
Operator : JC/MD
Sample : M3414-01
Misc : 5.02g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B-1

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
Octane, 2-methyl-	9.909	116.5	ug/l	1453760	3	10.061	623813	50.0
Nonane	10.366	106.0	ug/l	1322540	3	10.061	623813	50.0
Benzene, 1-ethy...	11.402	115.0	ug/l	3135720	4	12.030	1363340	50.0
Benzene, 1-ethe...	12.250	89.8	ug/l	2448490	4	12.030	1363340	50.0
Benzene, 2-ethy...	12.616	107.6	ug/l	2934380	4	12.030	1363340	50.0
Benzene, 1,2,3,...	12.926	87.5	ug/l	2387290	4	12.030	1363340	50.0
o-Cymene	12.969	101.6	ug/l	2769700	4	12.030	1363340	50.0
1-Phenyl-1-butene	13.298	147.7	ug/l	4028010	4	12.030	1363340	50.0
Naphthalene, 1-...	14.640	88.3	ug/l	2407830	4	12.030	1363340	50.0