LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100323\
Data File : VX038049.D
Acq On : 03 Oct 2023 20:02
Operator : JC/MD
Sample : 04566-07
Misc : 5.37g/5.0 mL/100uL/5.0mL/MSVOA_X/MeOH
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B102-08-1(5-5.5)

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\82X100323W.M
Title : SW846 8260

Signal : TIC: VX038049.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.288	516	525	536	rVB	119519	338828	5.64%	0.376%
2	5.385	689	705	710	rBV	87652	257909	4.30%	0.286%
3	5.471	711	719	727	rVV5	105036	398615	6.64%	0.442%
4	5.818	762	776	790	rBV4	136984	430443	7.17%	0.477%
5	5.958	790	799	805	rVV	100819	272225	4.53%	0.302%
6	6.031	805	811	821	rVV	113231	309168	5.15%	0.343%
7	6.153	821	831	839	rVV6	80676	298990	4.98%	0.331%
8	6.251	840	847	855	rVV3	105847	312991	5.21%	0.347%
9	6.348	855	863	875	rVB2	125593	378424	6.30%	0.420%
10	6.605	893	905	914	rBV	276688	675645	11.25%	0.749%
11	6.763	922	931	940	rBV2	223718	686232	11.43%	0.761%
12	7.373	1018	1031	1044	rBV	843984	2028580	33.79%	2.249%
13	7.769	1089	1096	1104	rVB2	117424	234803	3.91%	0.260%
14	7.952	1120	1126	1143	rVB3	149353	439183	7.32%	0.487%
15	8.141	1153	1157	1161	rVV	218717	381572	6.36%	0.423%
16	8.190	1161	1165	1174	rVV3	134961	305488	5.09%	0.339%
17	8.275	1174	1179	1182	rVV	353772	599635	9.99%	0.665%
18	8.312	1182	1185	1192	rVB3	188608	317501	5.29%	0.352%
19	8.433	1200	1205	1210	rBV	243671	410961	6.85%	0.456%
20	8.501	1212	1216	1226	rVB2	222883	422565	7.04%	0.468%
21	8.598	1226	1232	1236	rBV2	526208	1020438	17.00%	1.131%
22	8.647	1236	1240	1247	rVB2	567002	1025637	17.09%	1.137%
23	8.769	1256	1260	1265	rVV	188136	304474	5.07%	0.338%
24	8.848	1270	1273	1278	rVB	174091	253532	4.22%	0.281%
25	8.909	1278	1283	1289	rBV2	699922	1116292	18.60%	1.238%
26	8.976	1289	1294	1306	rVB2	420395	818397	13.63%	0.907%
27	9.098	1306	1314	1320	rBV2	240517	460867	7.68%	0.511%
28	9.159	1320	1324	1329	rBV2	259133	392064	6.53%	0.435%
29	9.384	1356	1361	1365	rBV2	246240	355997	5.93%	0.395%
30	9.506	1376	1381	1386	rBV3	310094	620538	10.34%	0.688%
31	9.561	1386	1390	1394	rBV	597960	783836	13.06%	0.869%
32	9.616	1394	1399	1403	rBV2	606666	955035	15.91%	1.059%
33	9.659	1403	1406	1417	rVB5	150689	342545	5.71%	0.380%
34	9.762	1417	1423	1428	rBV3	307711	534212	8.90%	0.592%
35	9.830	1428	1434	1438	rBV3	319128	659018	10.98%	0.731%
36	9.903	1442	1446	1450	rVB	440647	565445	9.42%	0.627%

LSC Area Percent Report

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Instrument :
MSVOA_X
ClientSampleId :
B102-08-1(5-5.5)

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
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Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100323W.M
Title : SW846 8260

37	10.006	1456	1463	1467	rVB	492060	689847	11.49%	0.765%
38	10.055	1467	1471	1474	rBV	513737	730990	12.18%	0.810%
39	10.195	1486	1494	1499	rVB	1231655	1737737	28.95%	1.926%
40	10.305	1499	1512	1517	rBV2	1576779	2964942	49.39%	3.287%
41	10.360	1517	1521	1532	rVB2	746011	1236073	20.59%	1.370%
42	10.457	1532	1537	1545	rBV3	211616	397619	6.62%	0.441%
43	10.586	1552	1558	1564	rBV3	495300	897262	14.95%	0.995%
44	10.640	1564	1567	1571	rBV	433949	569630	9.49%	0.632%
45	10.781	1582	1590	1594	rBV3	894086	1484540	24.73%	1.646%
46	10.854	1598	1602	1606	rVV3	657684	1004934	16.74%	1.114%
47	10.890	1606	1608	1614	rVB	246913	347304	5.79%	0.385%
48	10.963	1614	1620	1623	rBV	426971	657183	10.95%	0.729%
49	11.085	1634	1640	1643	rBV2	756874	1154331	19.23%	1.280%
50	11.189	1650	1657	1659	rBV2	372479	534236	8.90%	0.592%
51	11.220	1659	1662	1666	rVB4	429664	581180	9.68%	0.644%
52	11.305	1672	1676	1679	rBV	795521	966401	16.10%	1.071%
53	11.396	1686	1691	1695	rBV	1182556	1696332	28.26%	1.881%
54	11.451	1695	1700	1703	rBV	1456631	1847463	30.78%	2.048%
55	11.616	1722	1727	1734	rVV	986200	1499612	24.98%	1.663%
56	11.713	1740	1743	1746	rVV	438839	564037	9.40%	0.625%
57	11.756	1746	1750	1760	rVB	4644057	6003089	100.00%	6.655%
58	11.866	1760	1768	1769	rBV3	207942	373074	6.21%	0.414%
59	11.890	1769	1772	1777	rVV	425950	607457	10.12%	0.673%
60	11.945	1777	1781	1783	rVV	477899	630609	10.50%	0.699%
61	11.969	1783	1785	1789	rVV2	381276	490406	8.17%	0.544%
62	12.024	1789	1794	1800	rVV4	817848	1675642	27.91%	1.858%
63	12.085	1800	1804	1810	rVV	1584137	2409169	40.13%	2.671%
64	12.171	1813	1818	1825	rVB4	261146	589368	9.82%	0.653%
65	12.244	1825	1830	1837	rBV	1618474	2278806	37.96%	2.526%
66	12.317	1837	1842	1849	rVB4	1939880	3175589	52.90%	3.521%
67	12.421	1849	1859	1863	rBV3	504852	1057408	17.61%	1.172%
68	12.463	1863	1866	1872	rVB	670434	910036	15.16%	1.009%
69	12.530	1873	1877	1879	rVV	833573	1055902	17.59%	1.171%
70	12.555	1879	1881	1885	rVV	1008434	1214363	20.23%	1.346%
71	12.616	1887	1891	1894	rVV	1498786	1925563	32.08%	2.135%
72	12.646	1894	1896	1899	rVV	320994	360244	6.00%	0.399%
73	12.701	1899	1905	1913	rVV3	1204370	2366586	39.42%	2.624%
74	12.817	1917	1924	1926	rVV5	323052	740615	12.34%	0.821%
75	12.847	1926	1929	1932	rVV2	410907	519186	8.65%	0.576%
76	12.890	1932	1936	1938	rVV2	331393	527105	8.78%	0.584%

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ClientSampleId :
B102-08-1(5-5.5)

Integration Parameters: RTEINT.P
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Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100323W.M
Title : SW846 8260

77	12.920	1938	1941	1944	rVV	921771	1172435	19.53%	1.300%
78	12.963	1944	1948	1955	rVB	1171837	1699178	28.31%	1.884%
79	13.164	1978	1981	1985	rVB3	720414	897212	14.95%	0.995%
80	13.213	1985	1989	1992	rVB2	291540	367982	6.13%	0.408%
81	13.256	1992	1996	1998	rBV2	398459	625147	10.41%	0.693%
82	13.292	1998	2002	2007	rVV2	2169383	3032357	50.51%	3.362%
83	13.372	2012	2015	2019	rVV2	235216	320660	5.34%	0.355%
84	13.420	2019	2023	2032	rVB4	604298	1202842	20.04%	1.334%
85	13.506	2032	2037	2040	rBV2	303314	452478	7.54%	0.502%
86	13.542	2040	2043	2046	rBV	486508	576942	9.61%	0.640%
87	13.579	2046	2049	2055	rVV3	528667	925274	15.41%	1.026%
88	13.640	2056	2059	2060	rVV3	199247	242106	4.03%	0.268%
89	13.664	2060	2063	2068	rVB2	549184	800472	13.33%	0.887%
90	13.774	2077	2081	2088	rVB	1668361	2236331	37.25%	2.479%
91	13.847	2090	2093	2099	rVB5	164878	349417	5.82%	0.387%
92	13.926	2099	2106	2110	rBV5	299086	591727	9.86%	0.656%
93	14.073	2126	2130	2133	rBV3	288055	370125	6.17%	0.410%
94	14.213	2149	2153	2163	rVV2	369589	640582	10.67%	0.710%
95	14.371	2176	2179	2183	rVB	226772	263843	4.40%	0.293%
96	14.481	2193	2197	2201	rBV2	164407	271375	4.52%	0.301%
97	14.634	2215	2222	2226	rBV	1143920	1467547	24.45%	1.627%
98	14.780	2241	2246	2251	rVV	720857	951804	15.86%	1.055%
99	15.475	2355	2360	2368	rVB	183613	268618	4.47%	0.298%
100	15.609	2377	2382	2386	rBV	180178	295384	4.92%	0.327%

Sum of corrected areas: 90201793

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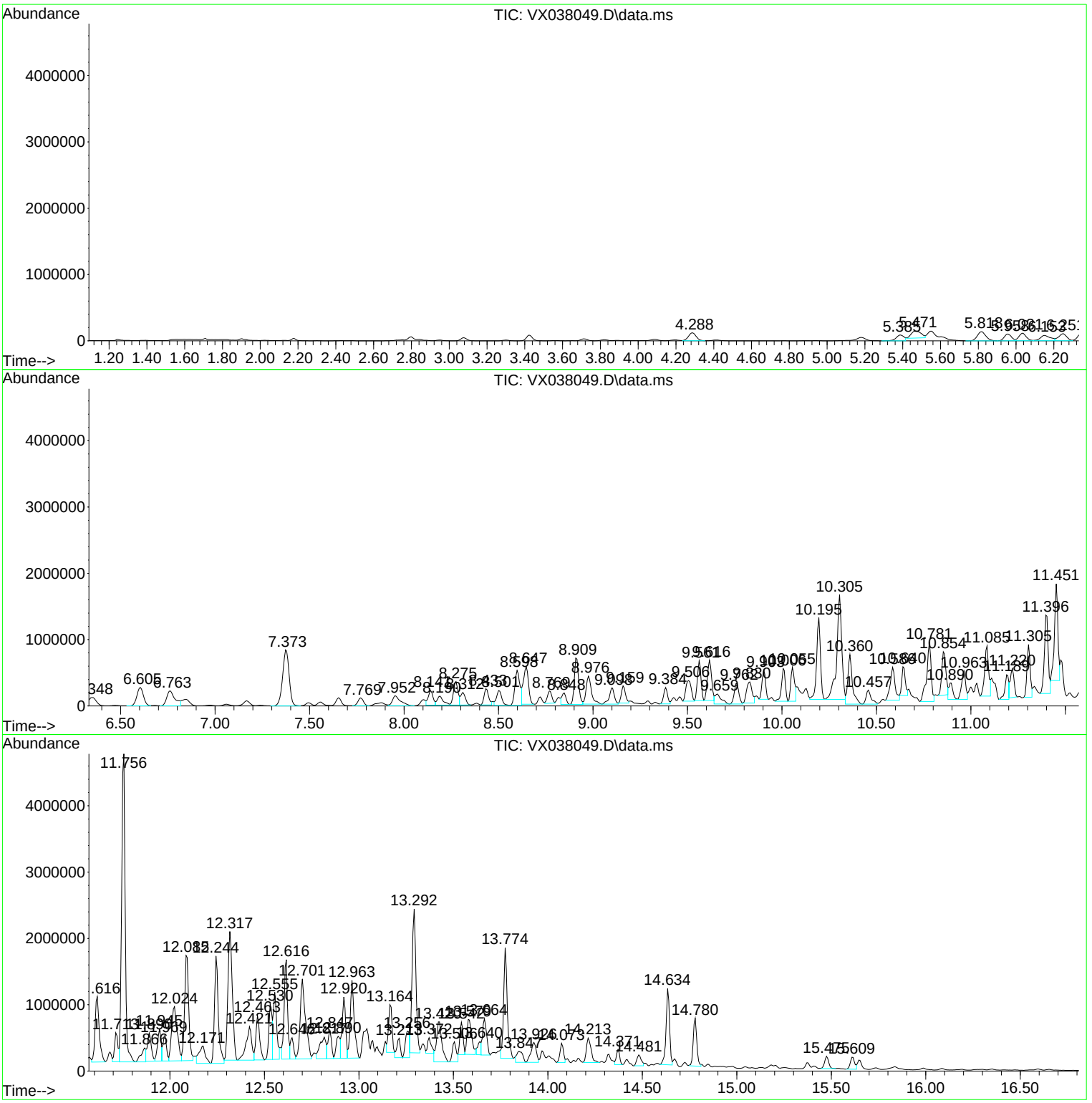
Instrument :
MSVOA_X
ClientSampleId :
B102-08-1(5-5.5)

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100323W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100323\
Data File : VX038049.D
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Instrument :
MSVOA_X
ClientSampleId :
B102-08-1(5-5.5)

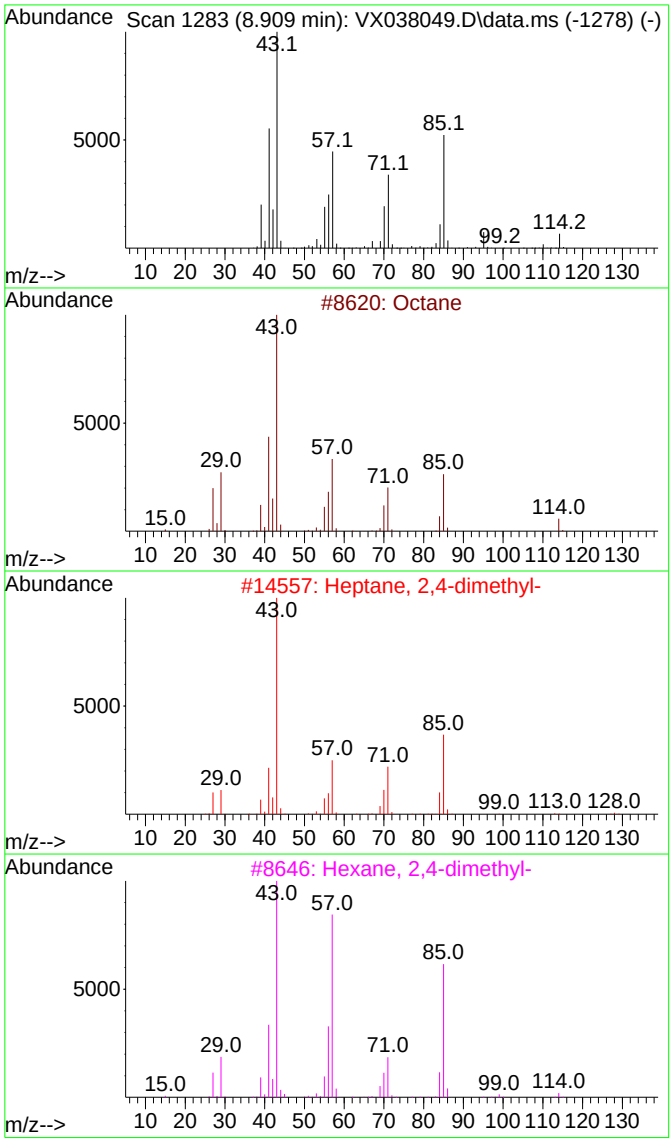
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100323W.M
Quant Title : SW846 8260

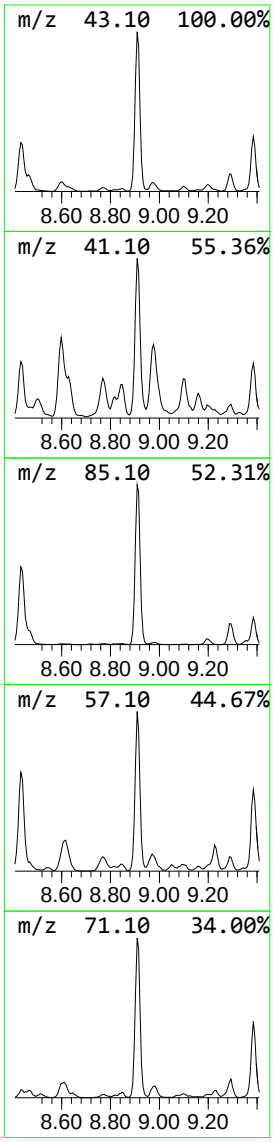
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Octane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.909	76.35 ug/l	1116290	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane	114	C8H18	000111-65-9	94
2	Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	72
3	Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	72
4	Hexane, 2-methyl-	100	C7H16	000591-76-4	53
5	Oxalic acid, diisohexyl ester	258	C14H26O4	1010309-32-9	53





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ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B102-08-1(5-5.5)

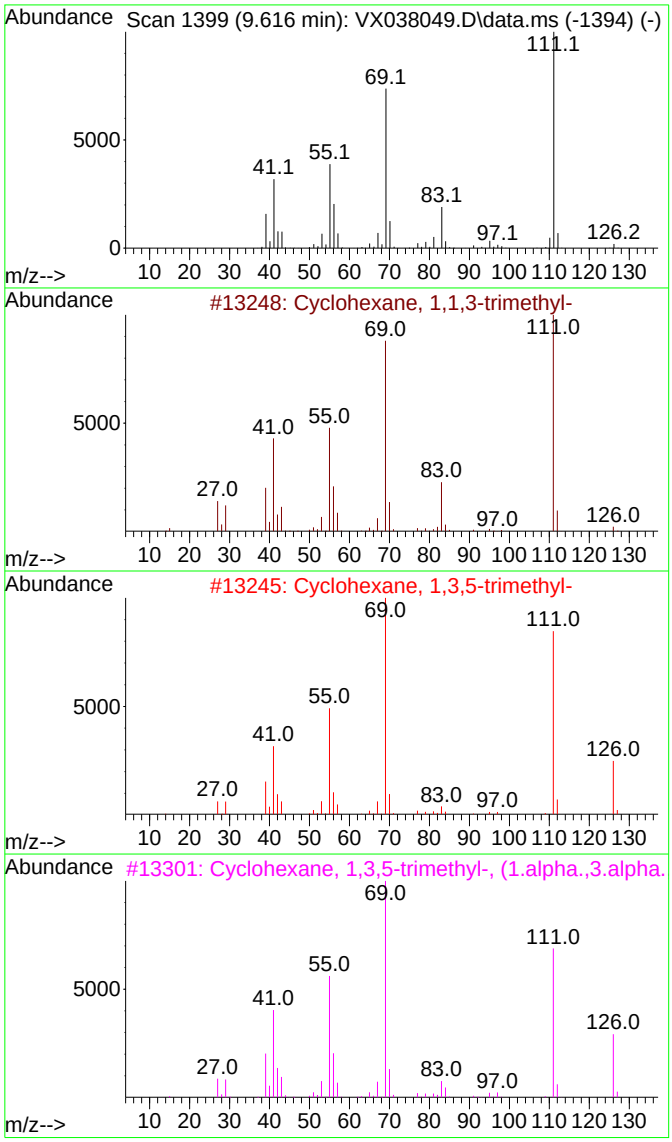
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100323W.M
Quant Title : SW846 8260

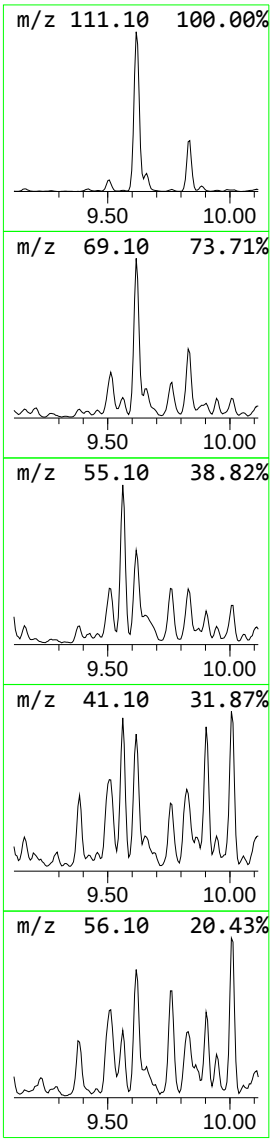
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Cyclohexane, 1,1,3-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.616	65.32 ug/l	955035	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	95
2	Cyclohexane, 1,3,5-trimethyl-	126	C9H18	001839-63-0	72
3	Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-26-2	53
4	Cyclooctane, (1-methylpropyl)-	168	C12H24	016538-89-9	50
5	3-Hexene, 2,5-dimethyl-	112	C8H16	015910-22-2	38





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Instrument :
MSVOA_X
ClientSampleId :
B102-08-1(5-5.5)

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100323W.M
Quant Title : SW846 8260

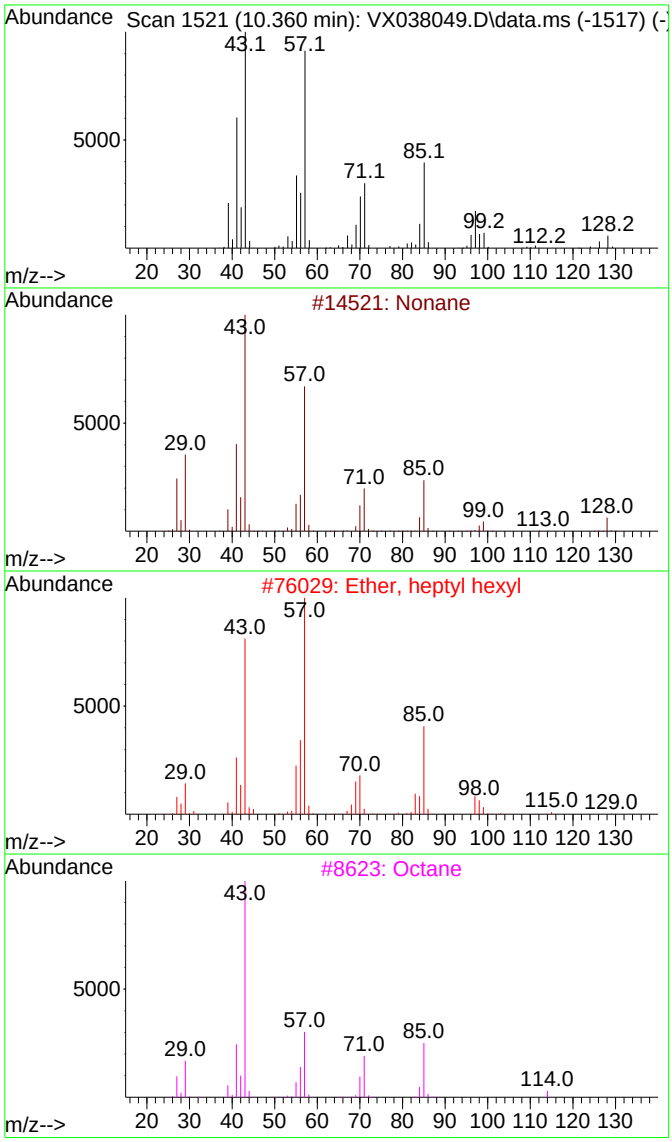
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

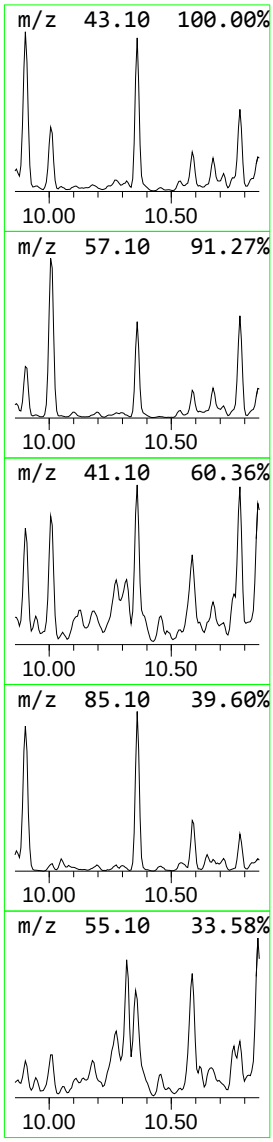
****

Peak Number 3 Nonane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.360	84.55 ug/l	1236070	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane	128	C9H20	000111-84-2	90
2	Ether, heptyl hexyl	200	C13H28O	007289-40-9	59
3	Octane	114	C8H18	000111-65-9	53
4	Hexane, 1-(hexyloxy)-4-methyl-	200	C13H28O	074421-20-8	50
5	Undecane, 2,7-dimethyl-	184	C13H28	017301-24-5	50





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ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B102-08-1(5-5.5)

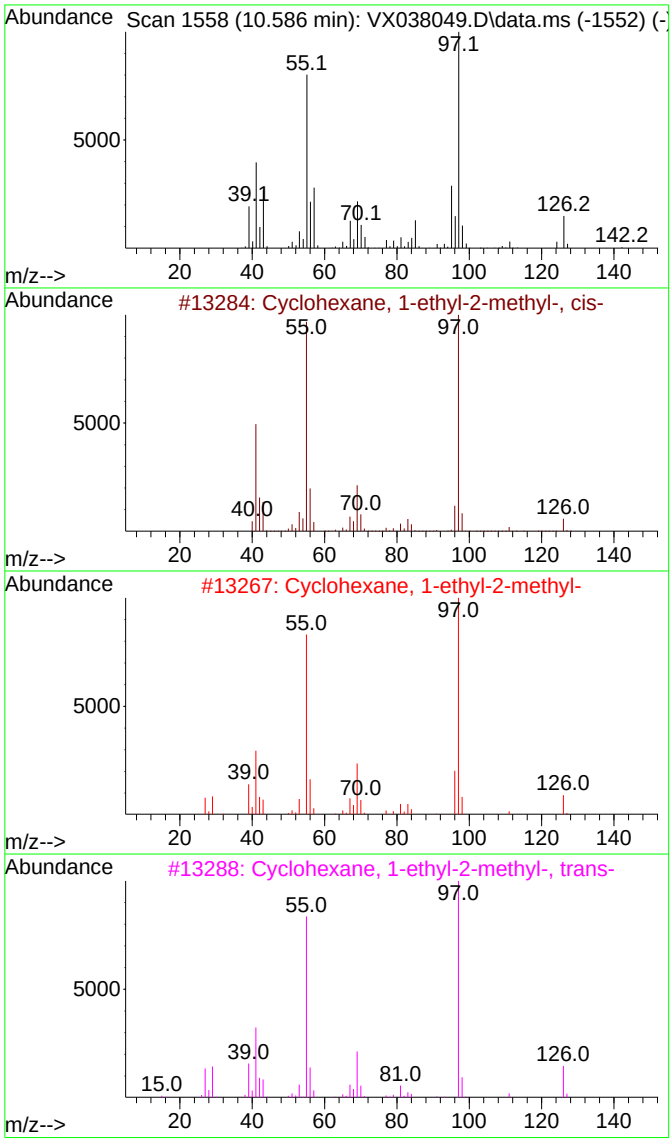
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100323W.M
Quant Title : SW846 8260

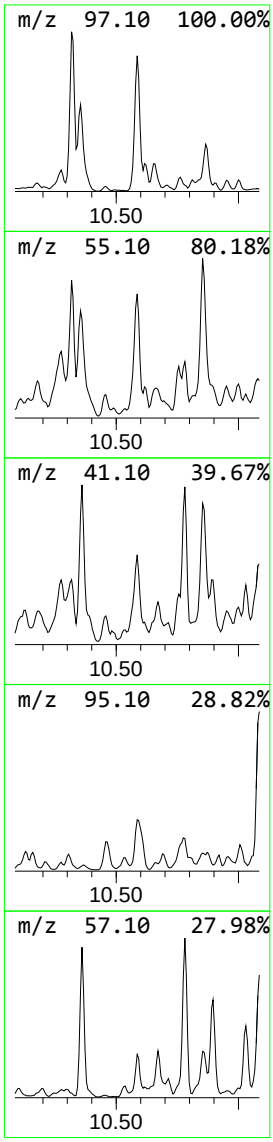
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Cyclohexane, 1-ethyl-2-meth... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.586	61.37 ug/l	897262	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-77-7	87
2	Cyclohexane, 1-ethyl-2-methyl-	126	C9H18	003728-54-9	81
3	Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-78-8	70
4	cis-1-Ethyl-3-methyl-cyclohexane	126	C9H18	019489-10-2	62
5	Cyclohexane, 1,3-dimethyl-	112	C8H16	000591-21-9	59





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Data File : VX038049.D
Acq On : 03 Oct 2023 20:02
Operator : JC/MD
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ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B102-08-1(5-5.5)

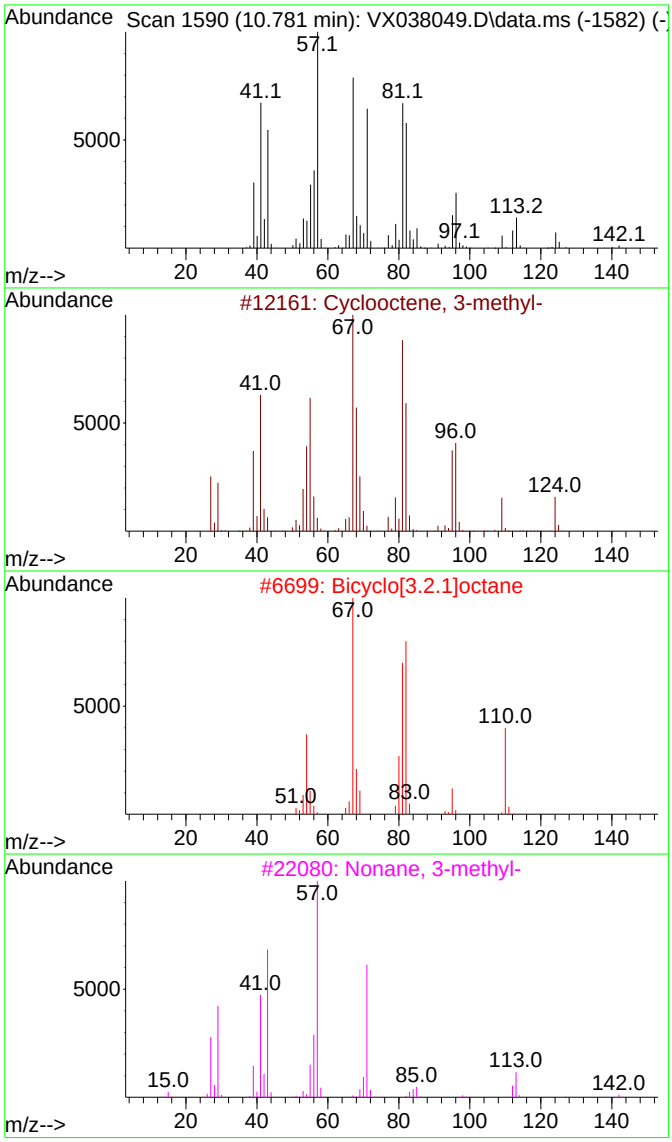
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100323W.M
Quant Title : SW846 8260

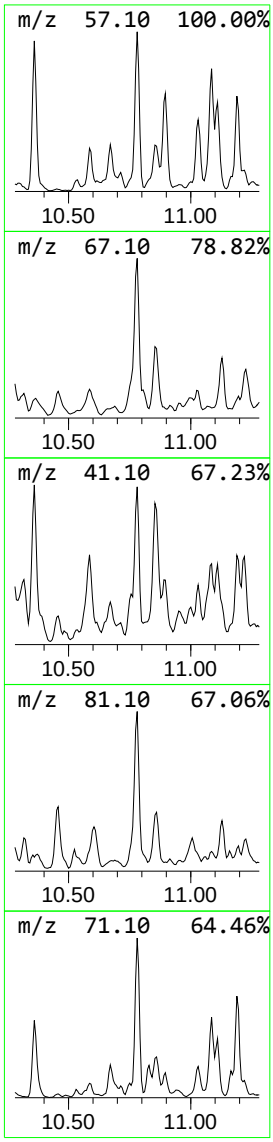
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 unknown10.781 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.781	101.54 ug/l	1484540	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclooctene, 3-methyl-	124	C9H16	013152-05-1	46
2	Bicyclo[3.2.1]octane	110	C8H14	006221-55-2	43
3	Nonane, 3-methyl-	142	C10H22	005911-04-6	42
4	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	42
5	2,4-Hexadiene, (E,Z)-	82	C6H10	005194-50-3	38





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100323\
Data File : VX038049.D
Acq On : 03 Oct 2023 20:02
Operator : JC/MD
Sample : 04566-07
Misc : 5.37g/5.0 mL/100uL/5.0mL/MSVOA_X/MeOH
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B102-08-1(5-5.5)

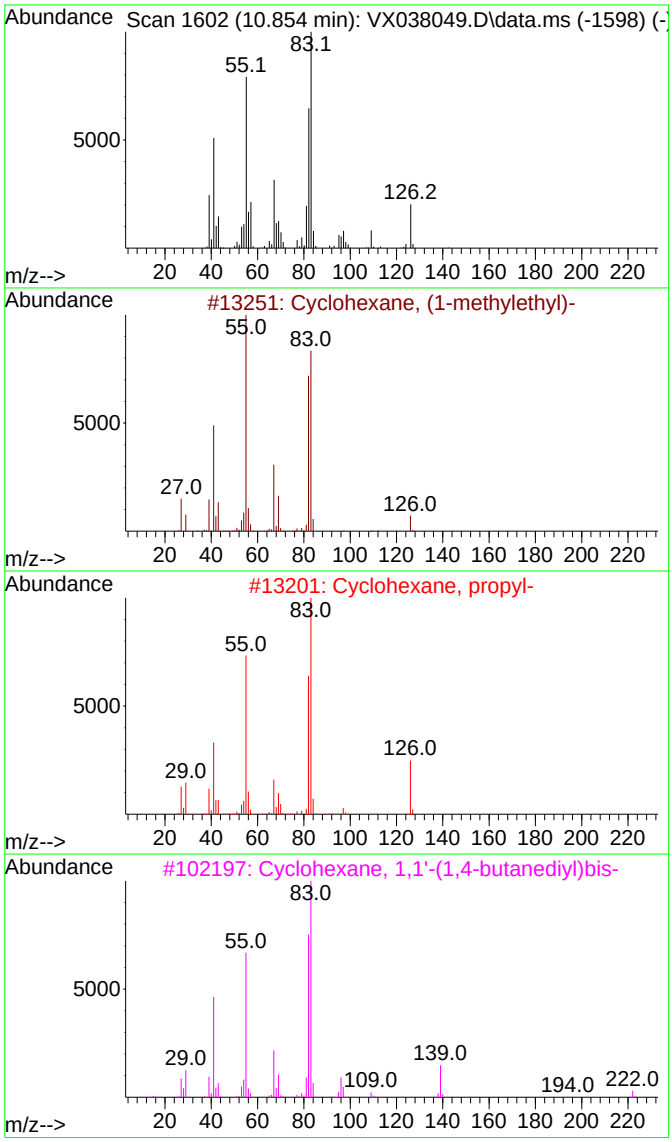
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100323W.M
Quant Title : SW846 8260

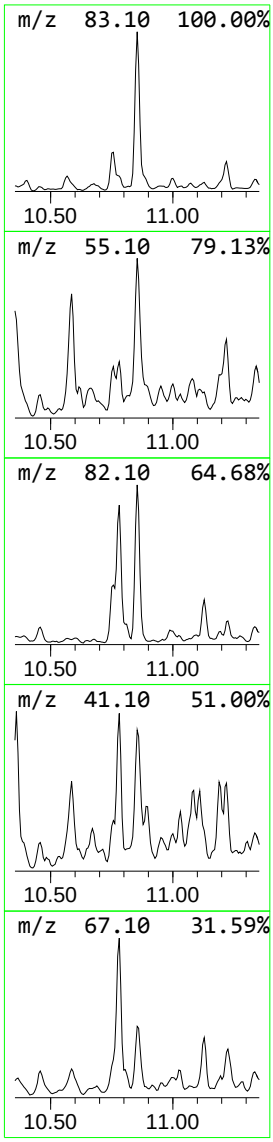
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Cyclohexane, (1-methylethyl)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.854	68.74 ug/l	1004930	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	90
2	Cyclohexane, propyl-	126	C9H18	001678-92-8	87
3	Cyclohexane, 1,1'-(1,4-butanediy...	222	C16H30	006165-44-2	64
4	Cyclohexane, butyl-	140	C10H20	001678-93-9	64
5	Cyclohexane, ethyl-	112	C8H16	001678-91-7	59





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100323\
Data File : VX038049.D
Acq On : 03 Oct 2023 20:02
Operator : JC/MD
Sample : 04566-07
Misc : 5.37g/5.0 mL/100uL/5.0mL/MSVOA_X/MeOH
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B102-08-1(5-5.5)

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100323W.M
Quant Title : SW846 8260

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

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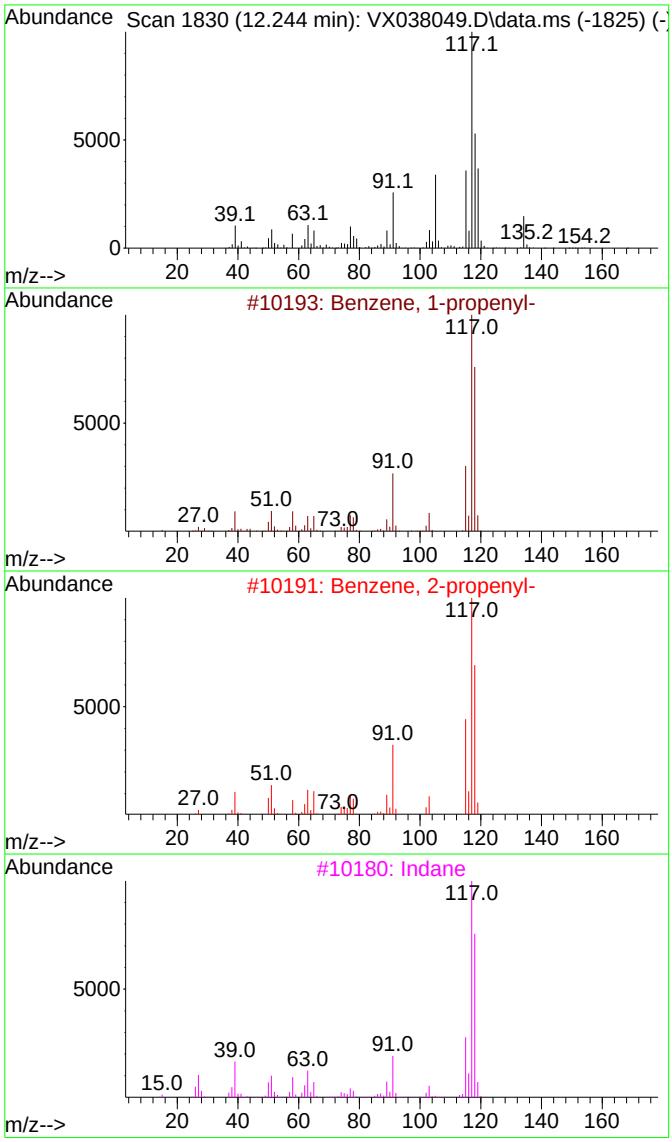
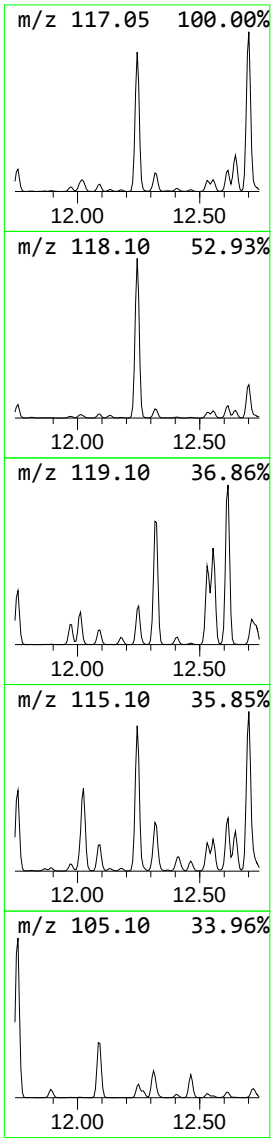
Peak Number 7 Benzene, 1-propenyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	68.00 ug/l	2278810	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-propenyl-	118	C9H10	000637-50-3	91
2	Benzene, 2-propenyl-	118	C9H10	000300-57-2	64
3	Indane	118	C9H10	000496-11-7	64
4	Benzene, cyclopropyl-	118	C9H10	000873-49-4	64
5	Deltacyclene	118	C9H10	007785-10-6	60





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100323\
Data File : VX038049.D
Acq On : 03 Oct 2023 20:02
Operator : JC/MD
Sample : 04566-07
Misc : 5.37g/5.0 mL/100uL/5.0mL/MSVOA_X/MeOH
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B102-08-1(5-5.5)

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100323W.M
Quant Title : SW846 8260

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

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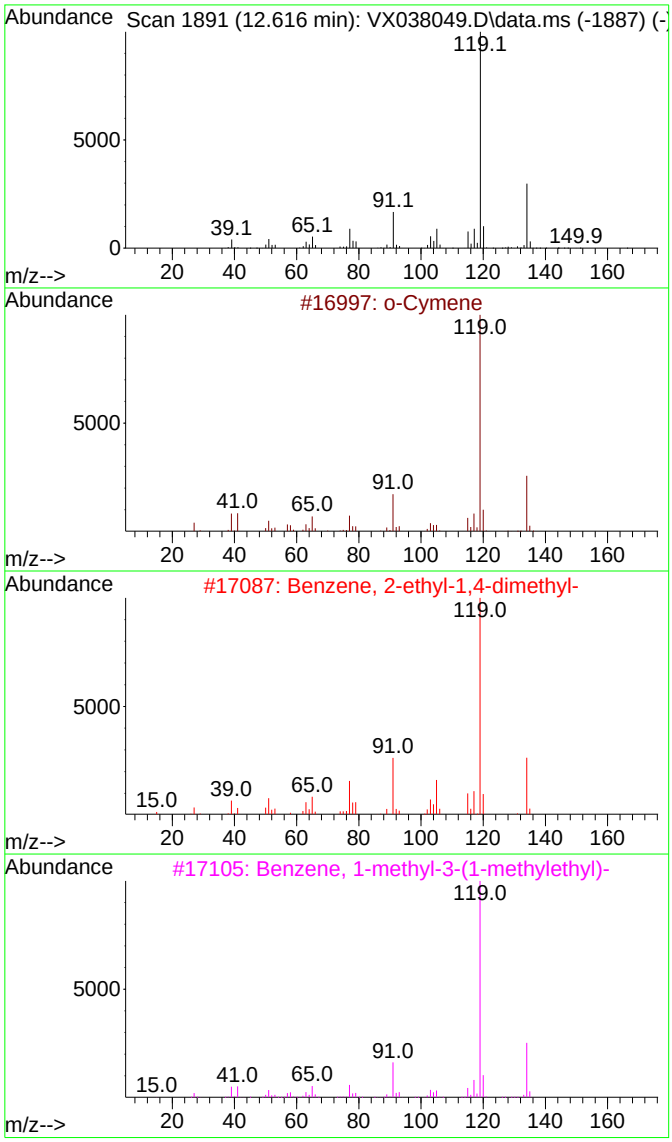
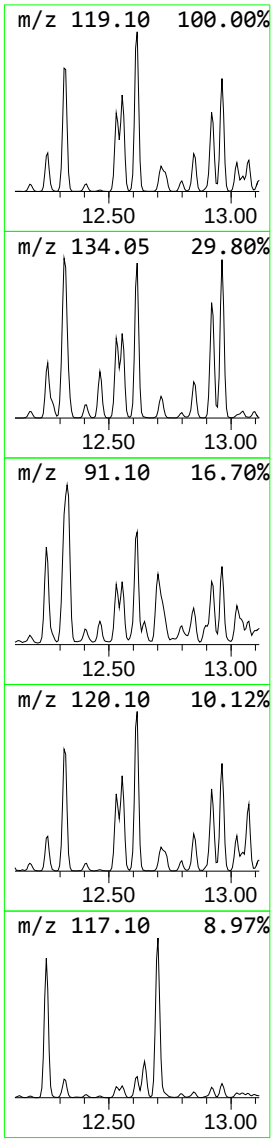
Peak Number 8 o-Cymene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.616	57.46 ug/l	1925560	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	97
2	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
3	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
4	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
5	p-Cymene	134	C10H14	000099-87-6	95





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100323\
Data File : VX038049.D
Acq On : 03 Oct 2023 20:02
Operator : JC/MD
Sample : 04566-07
Misc : 5.37g/5.0 mL/100uL/5.0mL/MSVOA_X/MeOH
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B102-08-1(5-5.5)

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100323W.M
Quant Title : SW846 8260

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

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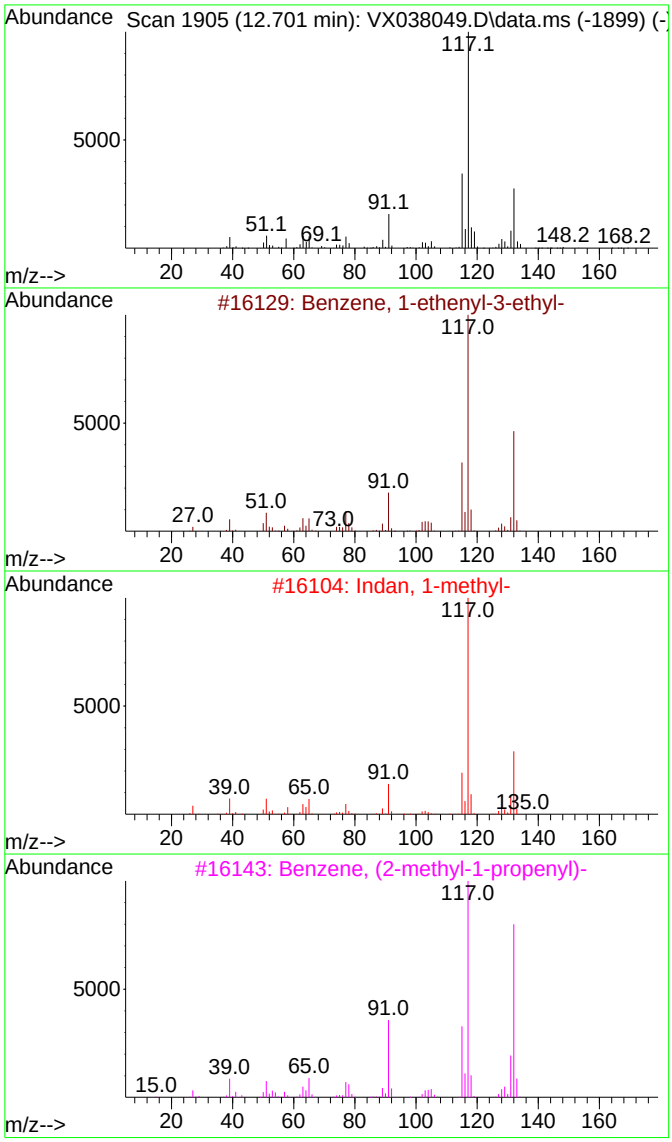
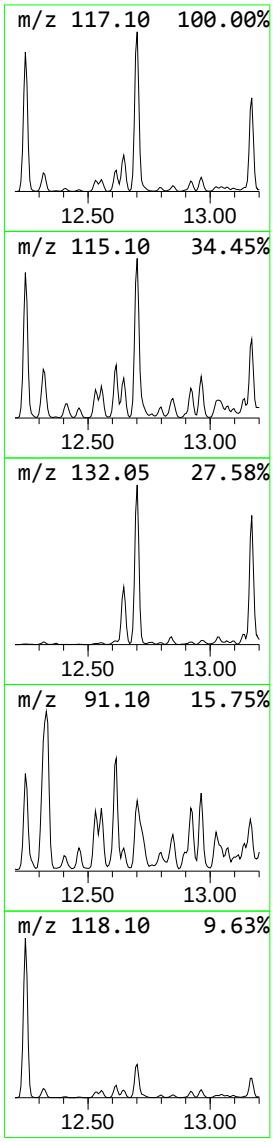
Peak Number 9 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	70.62 ug/l	2366590	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90
2	Indan, 1-methyl-	132	C10H12	000767-58-8	87
3	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	83
4	Benzene, 2-butenyl-	132	C10H12	001560-06-1	83
5	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	83





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100323\
Data File : VX038049.D
Acq On : 03 Oct 2023 20:02
Operator : JC/MD
Sample : 04566-07
Misc : 5.37g/5.0 mL/100uL/5.0mL/MSVOA_X/MeOH
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B102-08-1(5-5.5)

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100323W.M
Quant Title : SW846 8260

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

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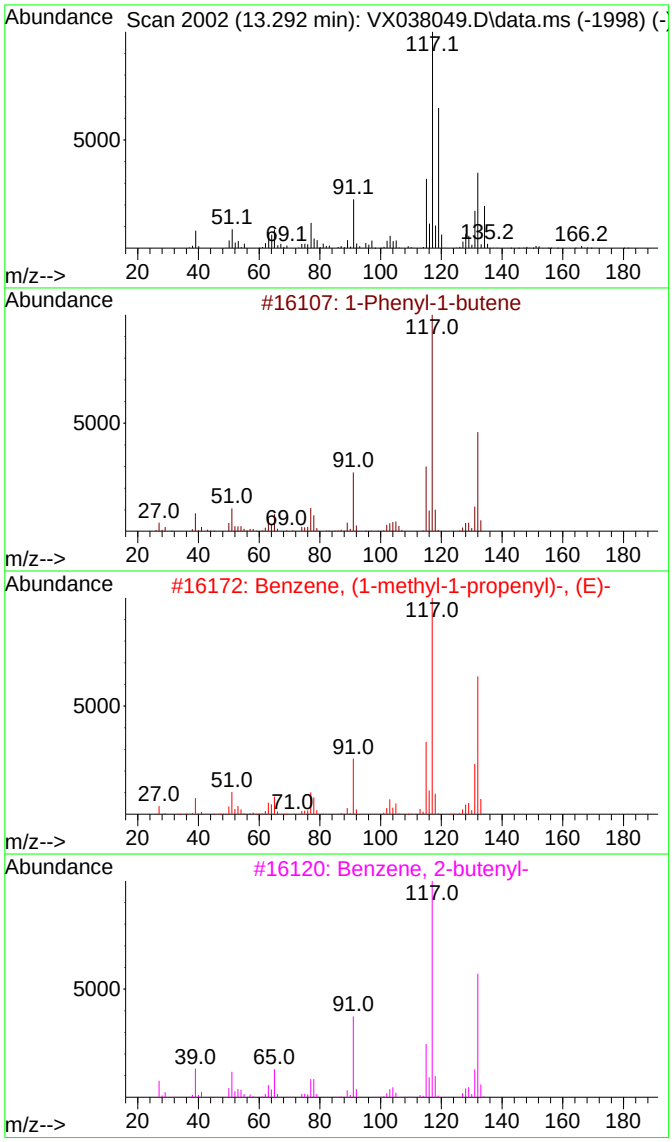
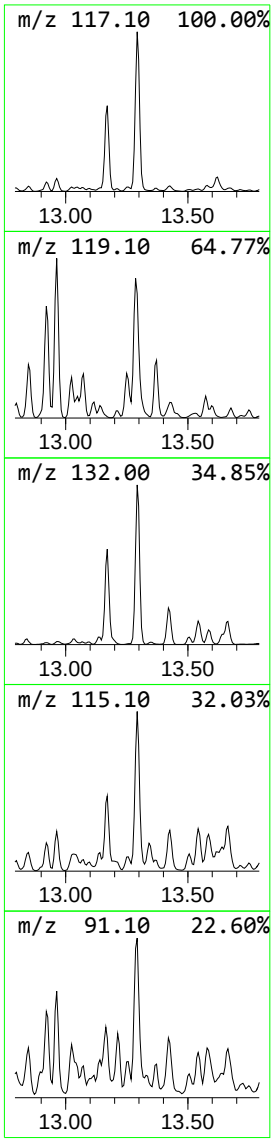
Peak Number 10 1-Phenyl-1-butene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	90.48 ug/l	3032360	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenyl-1-butene	132	C10H12	000824-90-8	86
2	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	70
3	Benzene, 2-butenyl-	132	C10H12	001560-06-1	70
4	3-Phenylbut-1-ene	132	C10H12	000934-10-1	70
5	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	64





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100323\
Data File : VX038049.D
Acq On : 03 Oct 2023 20:02
Operator : JC/MD
Sample : 04566-07
Misc : 5.37g/5.0 mL/100uL/5.0mL/MSVOA_X/MeOH
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
B102-08-1(5-5.5)

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100323W.M
Quant Title : SW846 8260

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Octane	8.909	76.3	ug/l	1116290	3	10.055	730990	50.0
Cyclohexane, 1,...	9.616	65.3	ug/l	955035	3	10.055	730990	50.0
Nonane	10.360	84.5	ug/l	1236070	3	10.055	730990	50.0
Cyclohexane, 1-...	10.586	61.4	ug/l	897262	3	10.055	730990	50.0
unknown10.781	10.781	101.5	ug/l	1484540	3	10.055	730990	50.0
Cyclohexane, (1...	10.854	68.7	ug/l	1004930	3	10.055	730990	50.0
Benzene, 1-prop...	12.244	68.0	ug/l	2278810	4	12.024	1675640	50.0
o-Cymene	12.616	57.5	ug/l	1925560	4	12.024	1675640	50.0
Benzene, 1-ethe...	12.701	70.6	ug/l	2366590	4	12.024	1675640	50.0
1-Phenyl-1-butene	13.292	90.5	ug/l	3032360	4	12.024	1675640	50.0