LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060922\
Data File : VX029349.D
Acq On : 09 Jun 2022 16:44
Operator : JC/MD
Sample : N3209-11 10X
Misc : 4.00g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
WC-14B-2-1

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\82X060722W.M
Title : SW846 8260

Signal : TIC: VX029349.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.391	695	706	716	rBV	184665	538166	31.66%	1.082%
2	5.556	723	733	741	rBV	238834	692418	40.73%	1.392%
3	5.958	790	799	811	rBV	186207	493532	29.03%	0.992%
4	6.160	821	832	842	rBV3	66552	227238	13.37%	0.457%
5	6.263	842	849	856	rVV2	60209	167403	9.85%	0.337%
6	6.361	856	865	876	rVB	105580	296177	17.42%	0.596%
7	6.763	921	931	943	rBV	364250	875145	51.48%	1.760%
8	7.379	1020	1032	1043	rVB2	447609	1109141	65.25%	2.230%
9	7.775	1088	1097	1105	rBV2	150915	304237	17.90%	0.612%
10	7.958	1120	1127	1143	rVB	173894	355082	20.89%	0.714%
11	8.604	1225	1233	1236	rBV2	442587	828076	48.71%	1.665%
12	8.647	1236	1240	1251	rVB	935012	1699897	100.00%	3.418%
13	8.775	1255	1261	1265	rBV	168043	265729	15.63%	0.534%
14	8.848	1270	1273	1280	rVB	135505	202024	11.88%	0.406%
15	8.982	1289	1295	1305	rVB2	405759	681878	40.11%	1.371%
16	9.104	1306	1315	1320	rBV	176426	295926	17.41%	0.595%
17	9.385	1353	1361	1366	rBV	178402	302052	17.77%	0.607%
18	9.506	1376	1381	1386	rBV4	246790	496714	29.22%	0.999%
19	9.567	1386	1391	1395	rVB	370197	544766	32.05%	1.095%
20	9.622	1395	1400	1404	rBV	545803	868504	51.09%	1.746%
21	9.763	1416	1423	1428	rBV	144895	229235	13.49%	0.461%
22	9.830	1428	1434	1438	rBV3	299918	605243	35.60%	1.217%
23	10.006	1456	1463	1467	rBV2	97168	161223	9.48%	0.324%
24	10.055	1467	1471	1476	rVB	793475	1035033	60.89%	2.081%
25	10.128	1476	1483	1487	rBV4	98751	203666	11.98%	0.410%
26	10.183	1487	1492	1499	rVB2	128327	215896	12.70%	0.434%
27	10.281	1499	1508	1511	rBV2	246153	508022	29.89%	1.021%
28	10.317	1511	1514	1518	rVV	312626	457115	26.89%	0.919%
29	10.354	1518	1520	1532	rVB	204330	394054	23.18%	0.792%
30	10.458	1532	1537	1546	rBV	178007	287465	16.91%	0.578%
31	10.586	1552	1558	1566	rVB3	429830	819190	48.19%	1.647%
32	10.671	1566	1572	1577	rBV3	166757	335538	19.74%	0.675%
33	10.781	1582	1590	1594	rBV2	804052	1277778	75.17%	2.569%
34	10.860	1594	1603	1606	rBV2	575805	1005238	59.14%	2.021%
35	10.896	1606	1609	1614	rVB	435827	584207	34.37%	1.175%
36	10.964	1614	1620	1623	rBV2	217490	364342	21.43%	0.733%

LSC Area Percent Report

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Instrument :
MSVOA_X
ClientSampleId :
WC-14B-2-1

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

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Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X060722W.M
Title : SW846 8260

37	11.006	1623	1627	1629	rBV2	138654	214926	12.64%	0.432%
38	11.031	1629	1631	1635	rVB2	243211	274142	16.13%	0.551%
39	11.085	1635	1640	1644	rVB	919993	1196459	70.38%	2.406%
40	11.220	1655	1662	1667	rBV2	565168	841944	49.53%	1.693%
41	11.341	1678	1682	1686	rVV3	175951	297886	17.52%	0.599%
42	11.390	1686	1690	1693	rVV3	198330	299645	17.63%	0.603%
43	11.433	1693	1697	1701	rVV	434753	553044	32.53%	1.112%
44	11.482	1701	1705	1709	rVB3	253379	390557	22.98%	0.785%
45	11.634	1723	1730	1734	rVB5	199140	452272	26.61%	0.909%
46	11.683	1734	1738	1741	rBV3	140280	192950	11.35%	0.388%
47	11.719	1741	1744	1746	rBV	274260	300354	17.67%	0.604%
48	11.841	1759	1764	1765	rBV2	155586	195310	11.49%	0.393%
49	11.890	1766	1772	1777	rVV4	511952	1057647	62.22%	2.127%
50	11.945	1777	1781	1784	rVV	482690	687257	40.43%	1.382%
51	11.976	1784	1786	1789	rVV2	200508	258098	15.18%	0.519%
52	12.024	1789	1794	1800	rVV	1052425	1460034	85.89%	2.936%
53	12.158	1812	1816	1826	rVB5	227958	563443	33.15%	1.133%
54	12.250	1826	1831	1836	rBV2	347005	560894	33.00%	1.128%
55	12.311	1837	1841	1848	rBV4	372248	777579	45.74%	1.563%
56	12.408	1853	1857	1858	rBV3	161239	209644	12.33%	0.422%
57	12.469	1864	1867	1873	rVB3	307833	533969	31.41%	1.074%
58	12.555	1876	1881	1883	rBV5	206316	344015	20.24%	0.692%
59	12.616	1887	1891	1894	rVB	465144	558097	32.83%	1.122%
60	12.701	1899	1905	1913	rVV5	291023	842652	49.57%	1.694%
61	12.823	1921	1925	1931	rVB3	388192	642547	37.80%	1.292%
62	12.890	1931	1936	1939	rBV2	410666	715730	42.10%	1.439%
63	12.920	1939	1941	1945	rVB	475656	535764	31.52%	1.077%
64	12.988	1948	1952	1955	rBV2	303768	343822	20.23%	0.691%
65	13.024	1955	1958	1966	rVB2	216045	372509	21.91%	0.749%
66	13.097	1966	1970	1974	rBV	223027	307164	18.07%	0.618%
67	13.140	1974	1977	1980	rBV2	219997	274846	16.17%	0.553%
68	13.292	1998	2002	2008	rVB4	685358	1164764	68.52%	2.342%
69	13.372	2013	2015	2020	rVV2	191811	339456	19.97%	0.683%
70	13.420	2020	2023	2032	rVB7	337690	868449	51.09%	1.746%
71	13.506	2033	2037	2040	rBV2	179139	267871	15.76%	0.539%
72	13.542	2040	2043	2046	rVB2	141090	166038	9.77%	0.334%
73	13.585	2046	2050	2053	rBV2	336031	457357	26.90%	0.920%
74	13.670	2061	2064	2068	rVB3	232567	342051	20.12%	0.688%
75	13.744	2072	2076	2081	rVB2	258088	491574	28.92%	0.988%
76	13.798	2081	2085	2089	rBV2	308116	399047	23.47%	0.802%

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ClientSampleId :
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Integration Parameters: RTEINT.P
Integrator: RTE
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Sampling : 1 Min Area: 3 % of largest Peak
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Title : SW846 8260

77	13.847	2089	2093	2099	rVB4	390765	526473	30.97%	1.059%
78	13.926	2099	2106	2111	rVB9	287817	651355	38.32%	1.310%
79	13.993	2111	2117	2118	rBV3	150522	290558	17.09%	0.584%
80	14.073	2126	2130	2133	rBV3	196049	276263	16.25%	0.555%
81	14.103	2133	2135	2138	rVV3	175281	192538	11.33%	0.387%
82	14.140	2138	2141	2143	rVV3	149159	171272	10.08%	0.344%
83	14.213	2150	2153	2161	rVB2	378120	603559	35.51%	1.214%
84	14.323	2168	2171	2175	rVB	184065	220119	12.95%	0.443%
85	14.371	2176	2179	2183	rVB	195731	245356	14.43%	0.493%
86	14.420	2184	2187	2193	rVB3	210751	290679	17.10%	0.584%
87	14.487	2193	2198	2201	rBV3	270209	480276	28.25%	0.966%
88	14.615	2216	2219	2226	rBV4	393262	559920	32.94%	1.126%
89	14.676	2226	2229	2233	rVB	267698	334226	19.66%	0.672%
90	14.725	2233	2237	2240	rBV3	163052	272013	16.00%	0.547%
91	14.792	2244	2248	2254	rBV6	146196	264366	15.55%	0.532%
92	14.853	2254	2258	2261	rVB3	149307	175577	10.33%	0.353%
93	15.182	2309	2312	2315	rBV3	118078	181817	10.70%	0.366%
94	15.481	2357	2361	2368	rVB	332198	484195	28.48%	0.974%
95	15.615	2378	2383	2388	rBV	683887	1067899	62.82%	2.147%
96	15.737	2398	2403	2409	rBV5	101183	210108	12.36%	0.422%
97	15.841	2413	2420	2424	rVV	192144	407543	23.97%	0.819%
98	15.987	2441	2444	2456	rVB3	153423	354303	20.84%	0.712%
99	16.603	2540	2545	2550	rVV2	252366	516563	30.39%	1.039%
100	16.658	2550	2554	2566	rVB	235132	477362	28.08%	0.960%

Sum of corrected areas: 49733497

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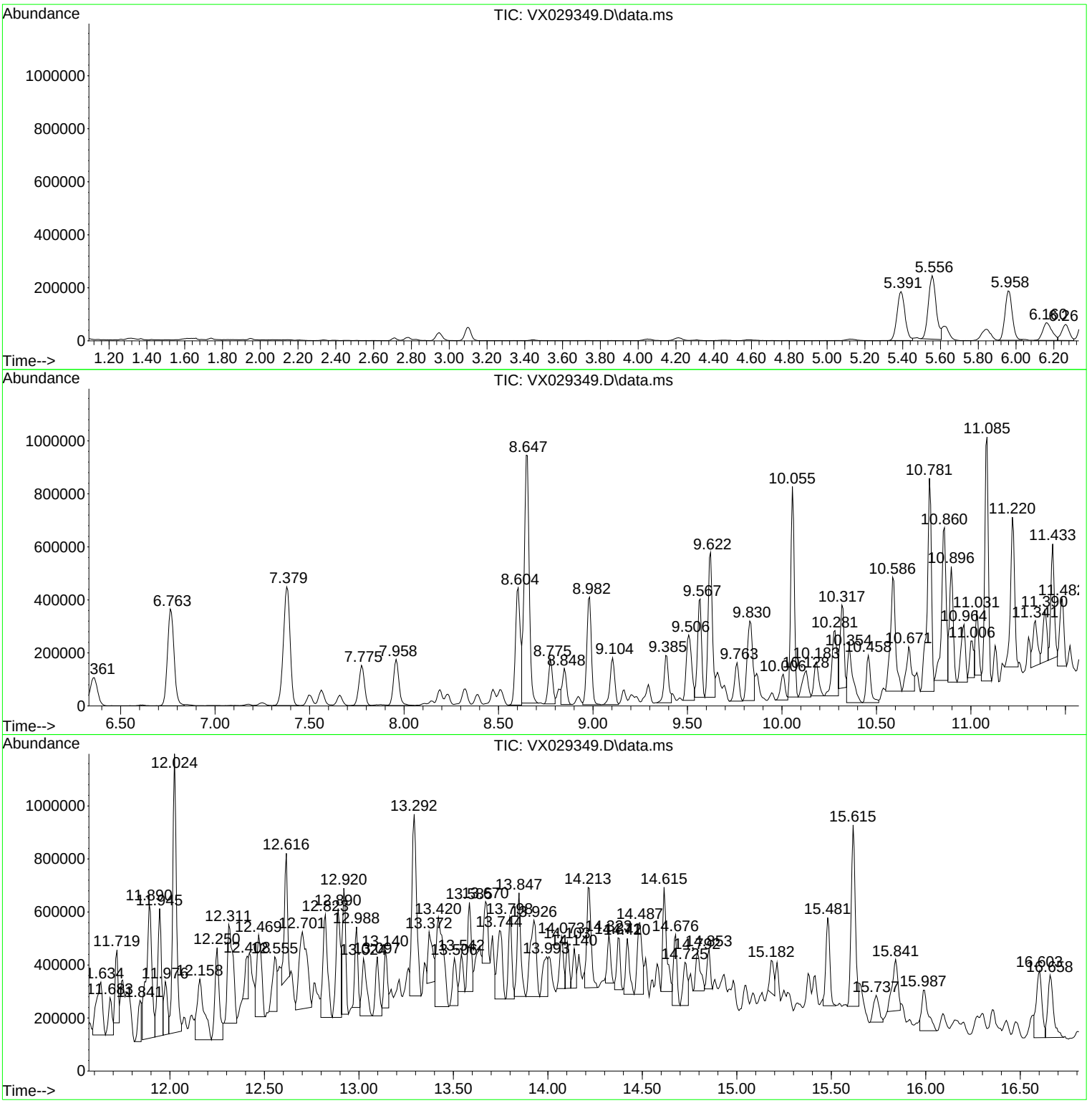
Instrument :
MSVOA_X
ClientSampleId :
WC-14B-2-1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X060722W.M
Quant Title : SW846 8260

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



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Instrument :
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ClientSampleId :
WC-14B-2-1

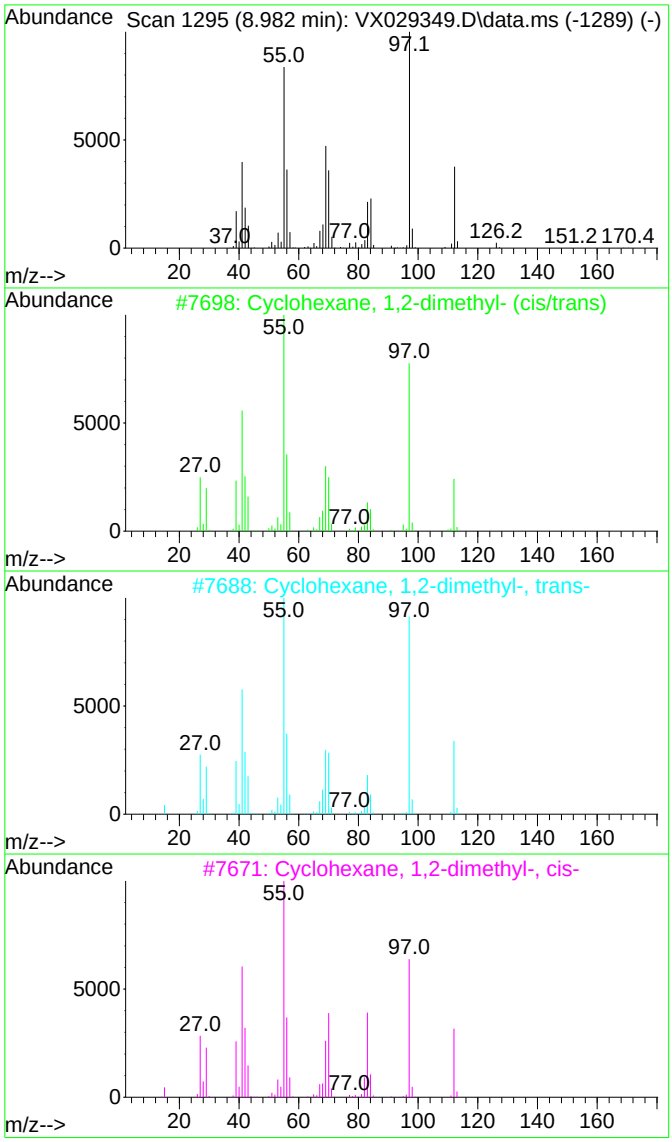
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X060722W.M
Quant Title : SW846 8260

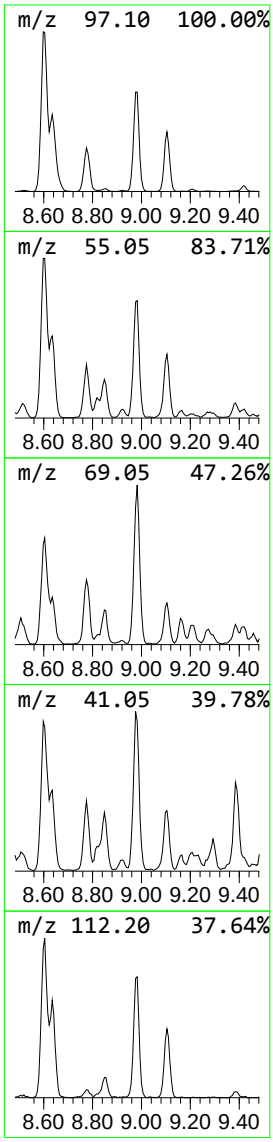
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Cyclohexane, 1,2-dimethyl- ... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.982	32.94 ug/l	681878	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,2-dimethyl- (cis/...	112	C8H16	000583-57-3	90
2	Cyclohexane, 1,2-dimethyl-, trans-	112	C8H16	006876-23-9	90
3	Cyclohexane, 1,2-dimethyl-, cis-	112	C8H16	002207-01-4	76
4	Cyclohexane, 1,3-dimethyl-, trans-	112	C8H16	002207-03-6	64
5	Cyclohexane, 1,3-dimethyl-, cis-	112	C8H16	000638-04-0	64





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Instrument :
MSVOA_X
ClientSampleId :
WC-14B-2-1

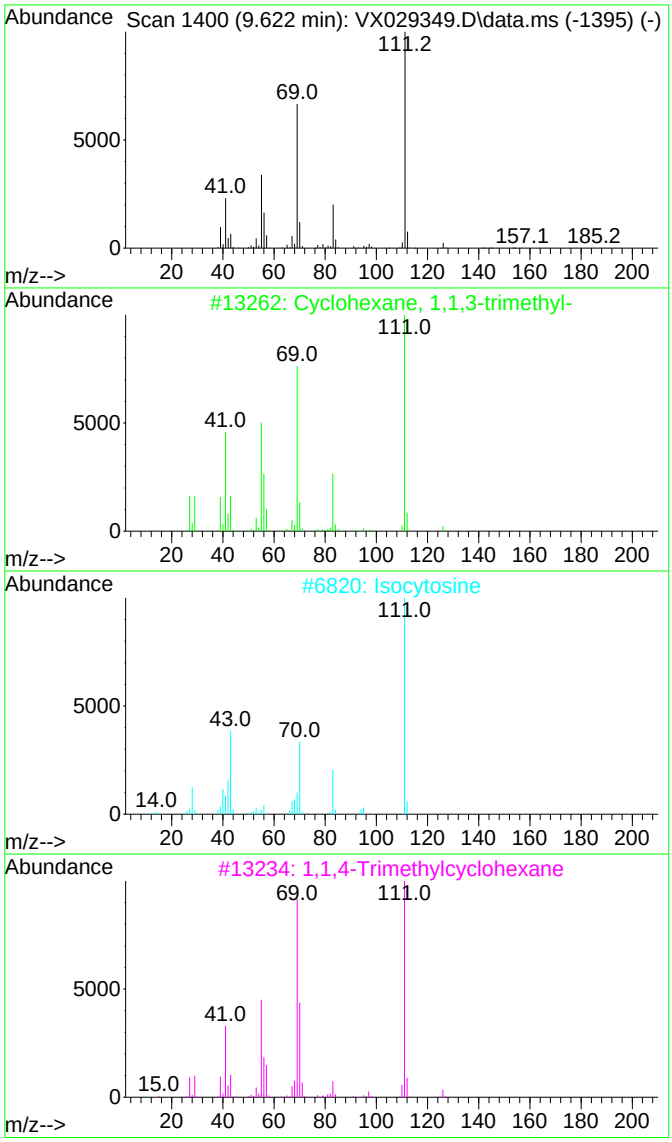
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X060722W.M
Quant Title : SW846 8260

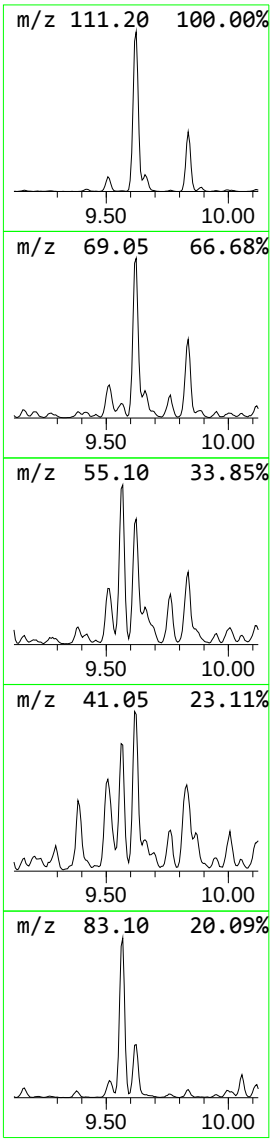
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.622	41.96 ug/l	868504	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	92
2	Isocytosine	111	C4H5N3O	000108-53-2	72
3	1,1,4-Trimethylcyclohexane	126	C9H18	007094-27-1	64
4	Cyclohexane, 1,3,5-trimethyl-	126	C9H18	001839-63-0	64
5	Cyclopentane, (2-methylbutyl)-	140	C10H20	053366-38-4	50





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Instrument :
MSVOA_X
ClientSampleId :
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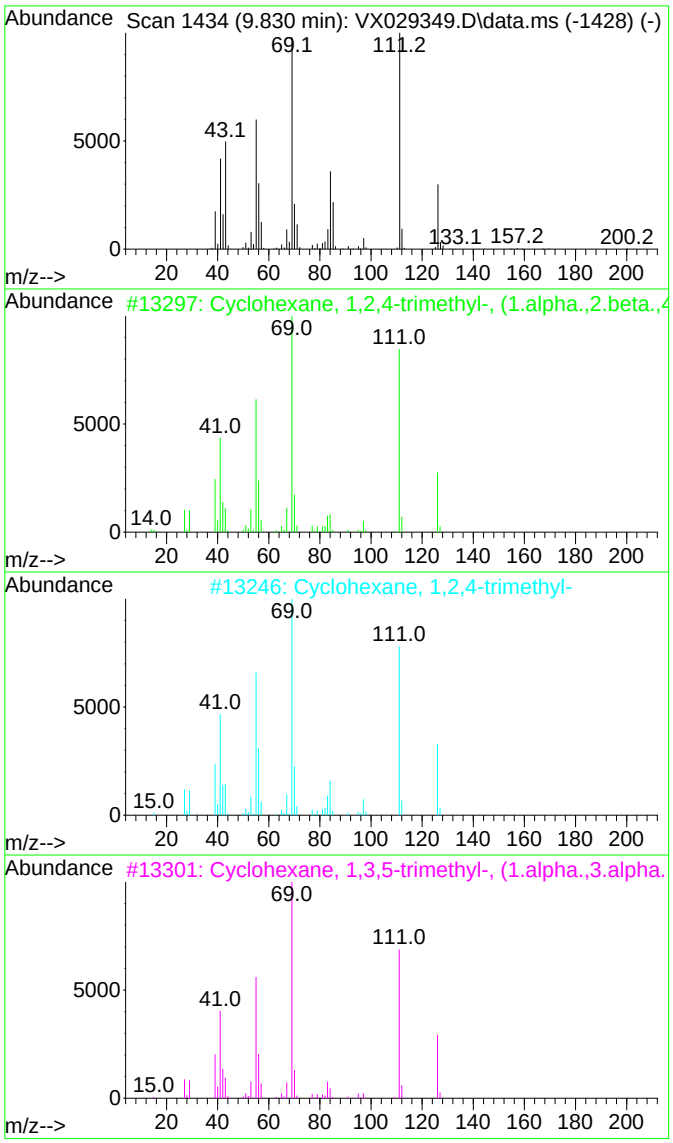
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X060722W.M
Quant Title : SW846 8260

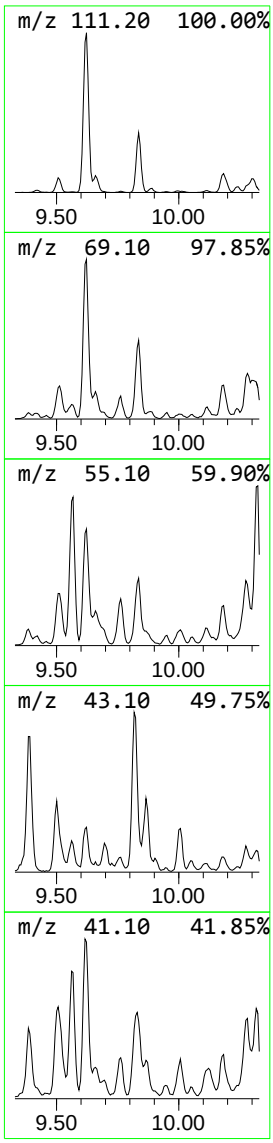
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Cyclohexane, 1,2,4-trimethy... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.830	29.24 ug/l	605243	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,2,4-trimethyl-, (...	126	C9H18	007667-60-9	91
2	Cyclohexane, 1,2,4-trimethyl-	126	C9H18	002234-75-5	91
3	Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-26-2	81
4	Cyclohexane, 1,3,5-trimethyl-	126	C9H18	001839-63-0	74
5	Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-27-3	64





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ClientSampleId :
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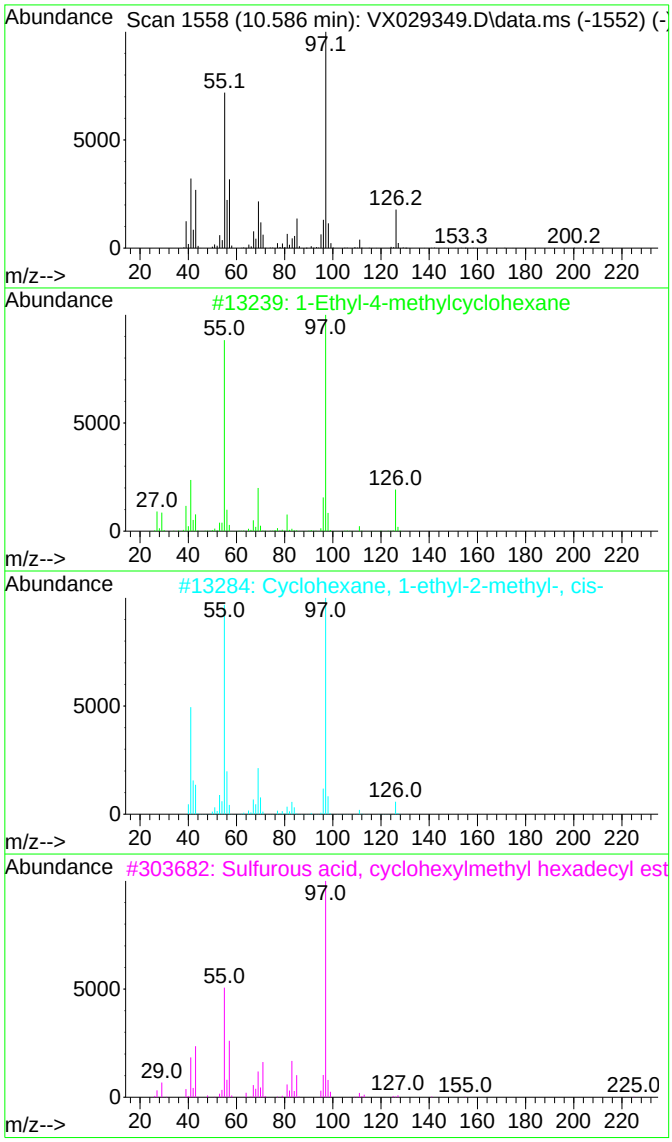
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X060722W.M
Quant Title : SW846 8260

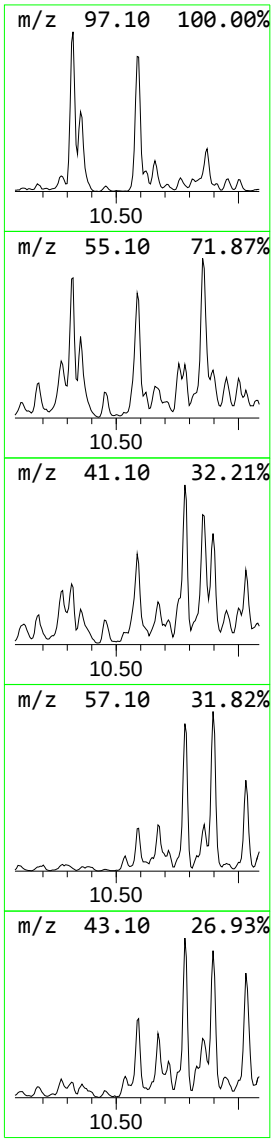
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 1-Ethyl-4-methylcyclohexane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.586	39.57 ug/l	819190	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Ethyl-4-methylcyclohexane	126	C9H18	003728-56-1	76
2	Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-77-7	72
3	Sulfurous acid, cyclohexylmethyl...	402	C23H46O3S	1000309-22-4	72
4	1-Ethyl-3-methylcyclohexane (c,t)	126	C9H18	003728-55-0	70
5	Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-78-8	70





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ClientSampleId :
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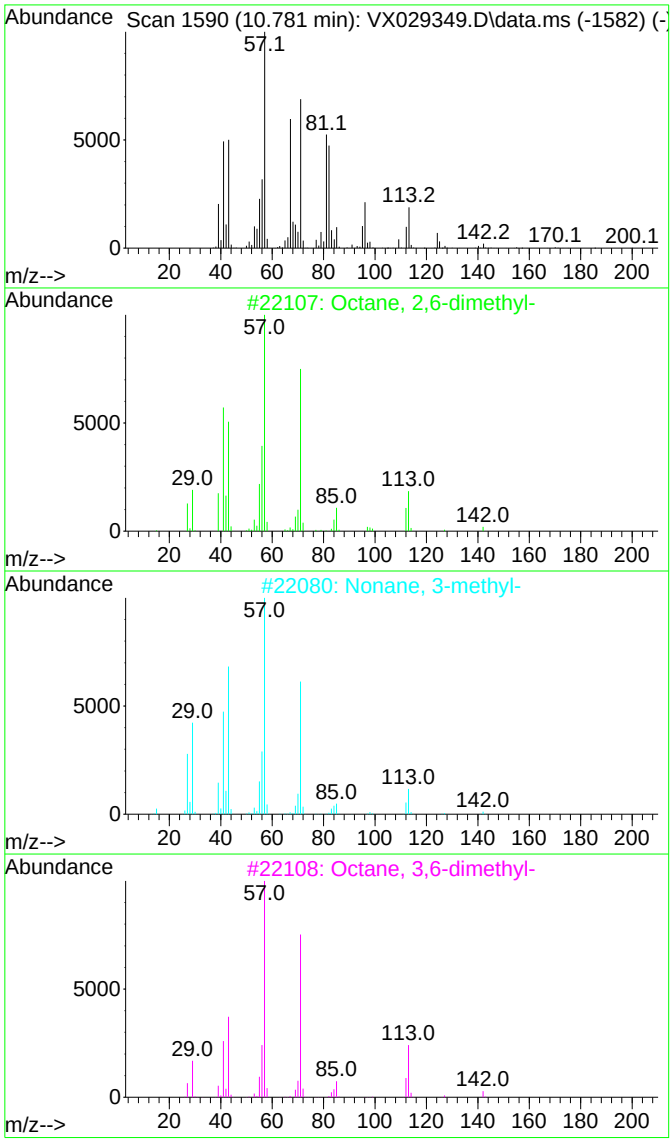
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X060722W.M
Quant Title : SW846 8260

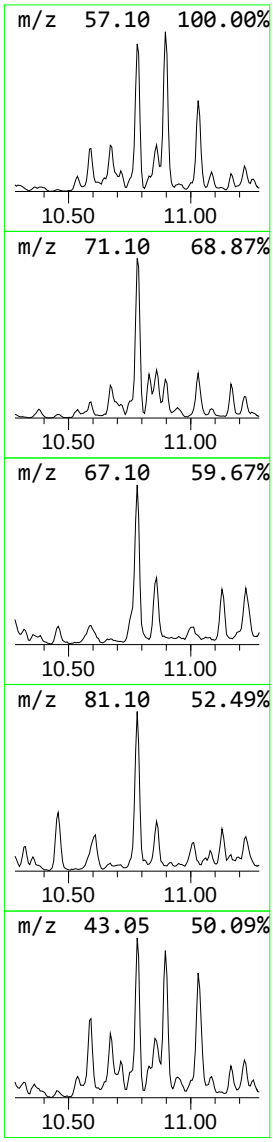
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Octane, 2,6-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.781	61.73 ug/l	1277780	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	50
2	Nonane, 3-methyl-	142	C10H22	005911-04-6	50
3	Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	46
4	Sulfurous acid, dodecyl 2-ethylh...	362	C20H42O3S	1000309-19-5	38
5	Sulfurous acid, 2-ethylhexyl hex...	418	C24H50O3S	1000309-19-9	30





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060922\
Data File : VX029349.D
Acq On : 09 Jun 2022 16:44
Operator : JC/MD
Sample : N3209-11 10X
Misc : 4.00g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
WC-14B-2-1

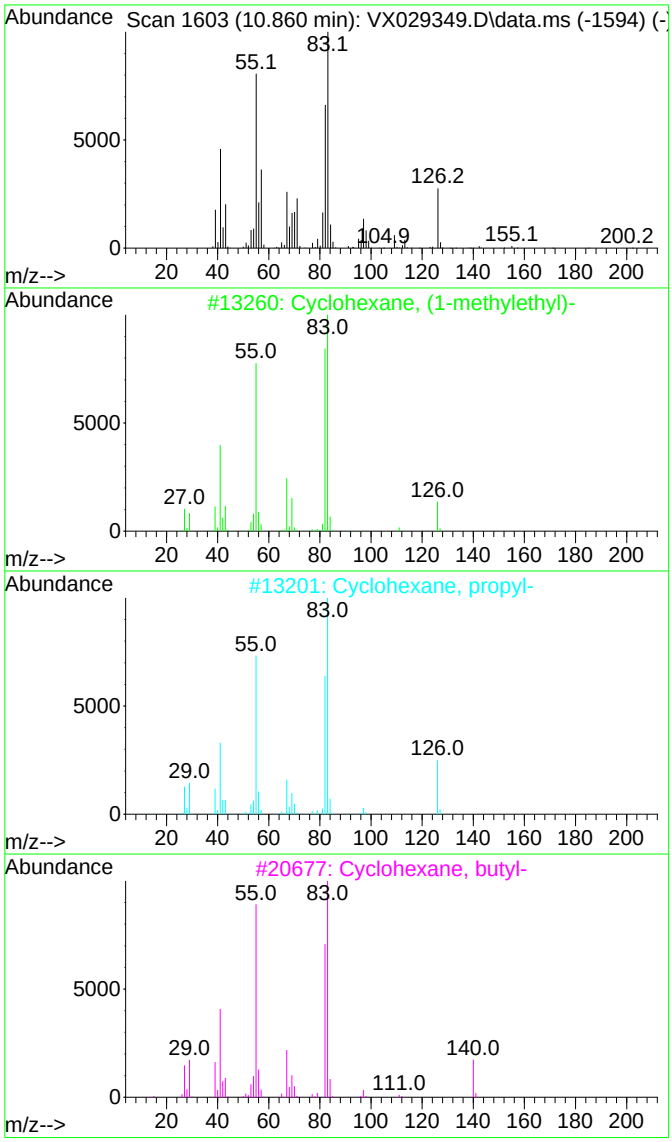
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X060722W.M
Quant Title : SW846 8260

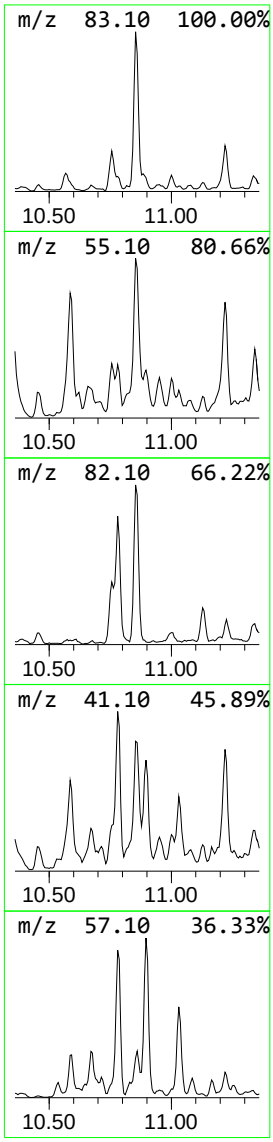
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Cyclohexane, (1-methylethyl)- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.860	48.56 ug/l	1005240	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	81
2	Cyclohexane, propyl-	126	C9H18	001678-92-8	68
3	Cyclohexane, butyl-	140	C10H20	001678-93-9	59
4	Cyclopentane, 1-ethyl-1-methyl-	112	C8H16	016747-50-5	58
5	Cyclohexane, pentyl-	154	C11H22	004292-92-6	53





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060922\
Data File : VX029349.D
Acq On : 09 Jun 2022 16:44
Operator : JC/MD
Sample : N3209-11 10X
Misc : 4.00g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
WC-14B-2-1

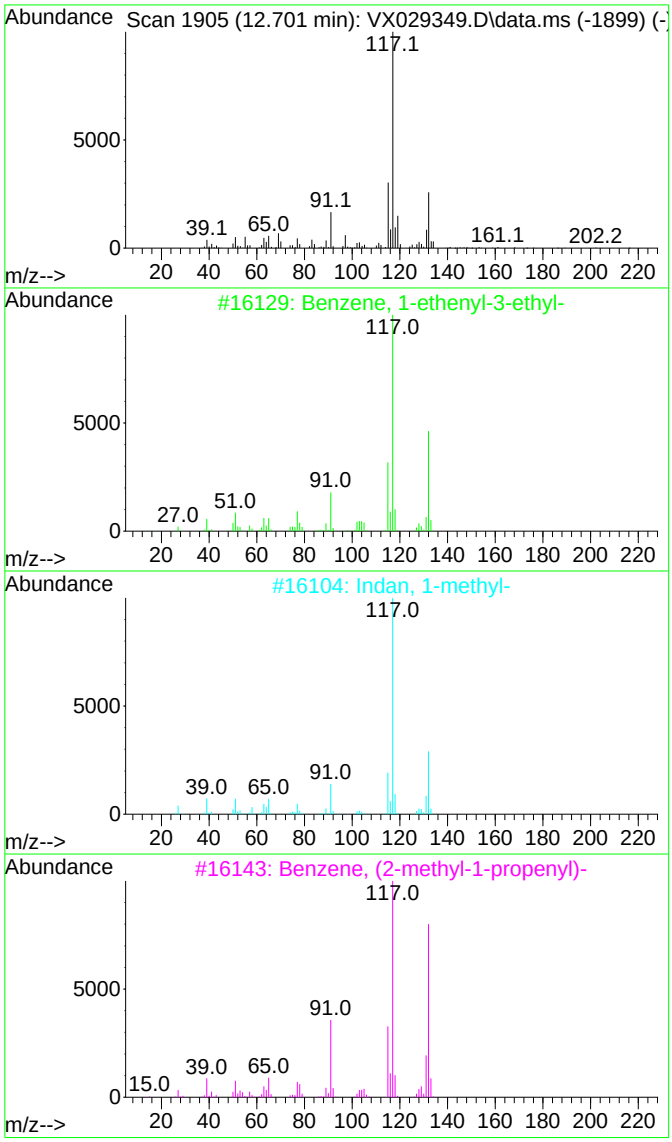
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X060722W.M
Quant Title : SW846 8260

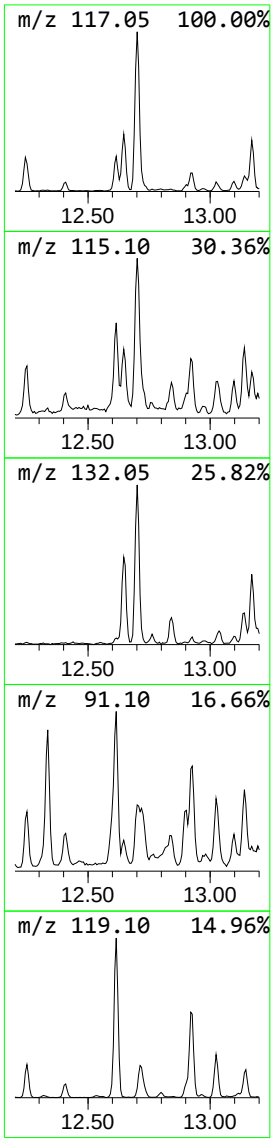
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	28.86 ug/l	842652	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	87
2	Indan, 1-methyl-	132	C10H12	000767-58-8	87
3	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	87
4	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	81
5	Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	81





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060922\
Data File : VX029349.D
Acq On : 09 Jun 2022 16:44
Operator : JC/MD
Sample : N3209-11 10X
Misc : 4.00g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
WC-14B-2-1

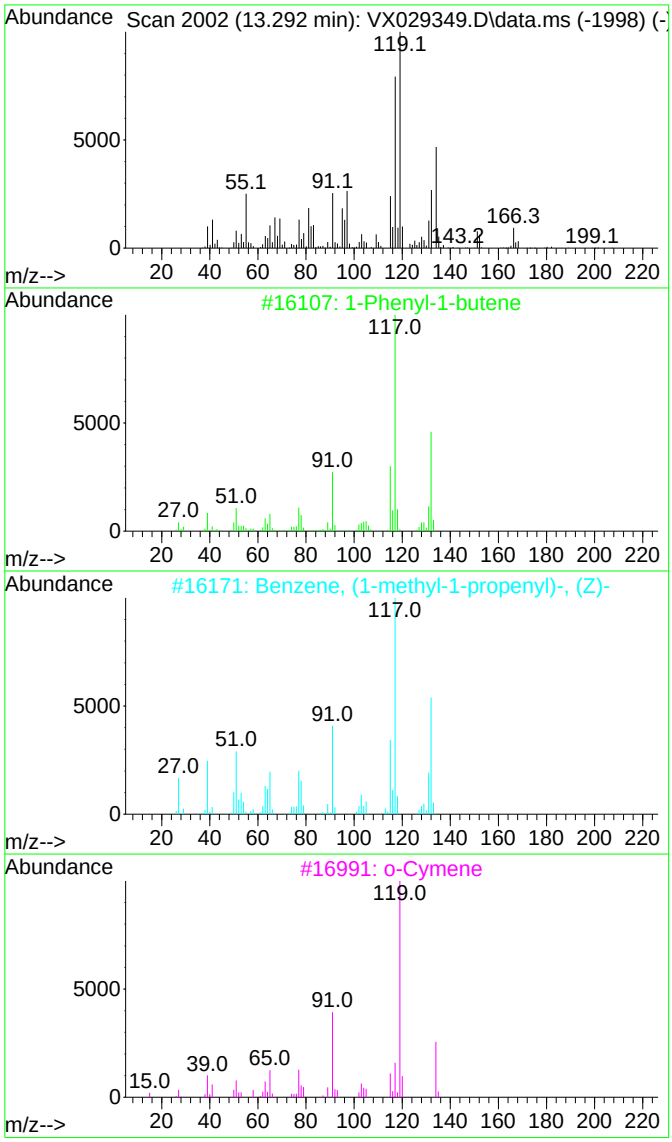
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X060722W.M
Quant Title : SW846 8260

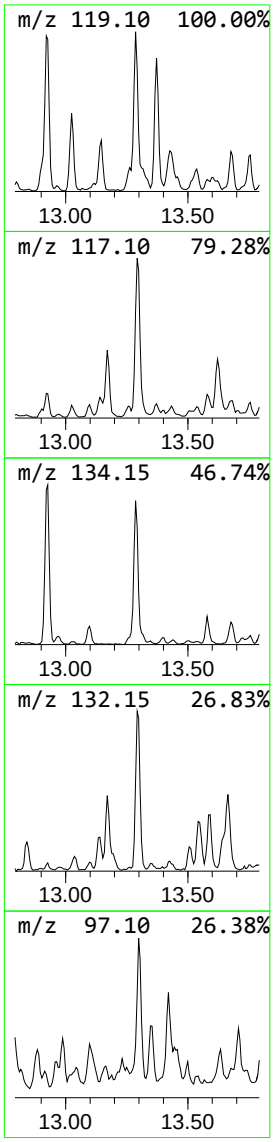
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 1-Phenyl-1-butene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	39.89 ug/l	1164760	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	83
2	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	52
3	o-Cymene	134	C10H14	000527-84-4	50
4	p-Cymene	134	C10H14	000099-87-6	50
5	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	46





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060922\
Data File : VX029349.D
Acq On : 09 Jun 2022 16:44
Operator : JC/MD
Sample : N3209-11 10X
Misc : 4.00g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
WC-14B-2-1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X060722W.M
Quant Title : SW846 8260

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

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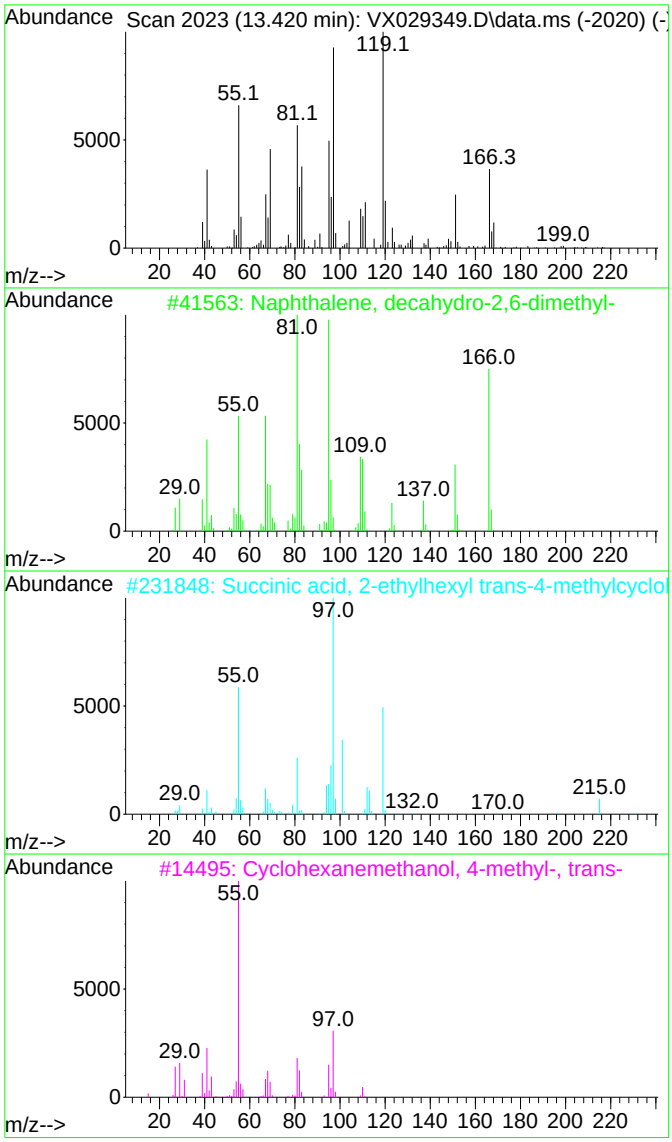
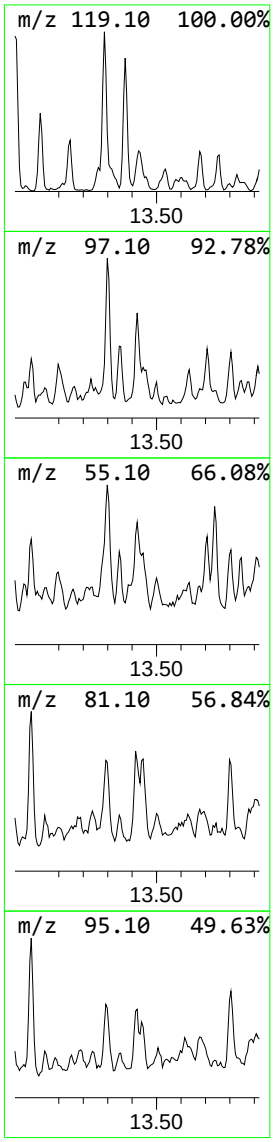
Peak Number 9 unknown13.420 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.420	29.74 ug/l	868449	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, decahydro-2,6-dimet...	166	C12H22	001618-22-0	35
2	Succinic acid, 2-ethylhexyl tran...	326	C19H34O4	1000390-07-1	35
3	Cyclohexanemethanol, 4-methyl-, ...	128	C8H16O	003937-49-3	27
4	Cyclohexane, 1-methyl-3-pentyl-	168	C12H24	054411-02-8	25
5	2-Bromopropionic acid, 1-(cyclop...	248	C10H17BrO2	1000293-39-1	22





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060922\
Data File : VX029349.D
Acq On : 09 Jun 2022 16:44
Operator : JC/MD
Sample : N3209-11 10X
Misc : 4.00g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
WC-14B-2-1

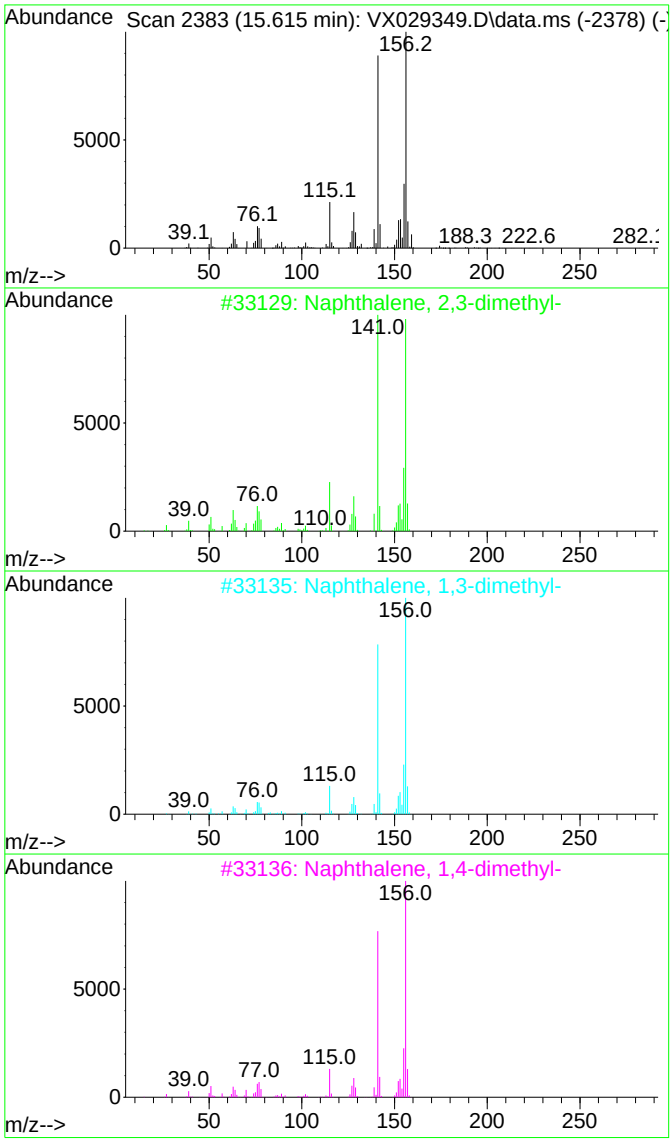
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X060722W.M
Quant Title : SW846 8260

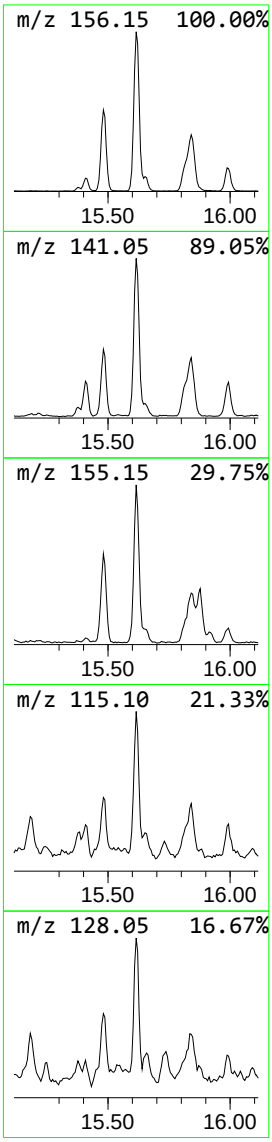
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Naphthalene, 2,3-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.615	36.57 ug/l	1067900	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, 1,3-dimethyl-	156	C12H12	000575-41-7	97
3	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97
4	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
5	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060922\
Data File : VX029349.D
Acq On : 09 Jun 2022 16:44
Operator : JC/MD
Sample : N3209-11 10X
Misc : 4.00g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
WC-14B-2-1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X060722W.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
Cyclohexane, 1,...	8.982	32.9	ug/l	681878	3	10.055	1035030	50.0
Cyclohexane, 1,...	9.622	42.0	ug/l	868504	3	10.055	1035030	50.0
Cyclohexane, 1,...	9.830	29.2	ug/l	605243	3	10.055	1035030	50.0
1-Ethyl-4-methy...	10.586	39.6	ug/l	819190	3	10.055	1035030	50.0
Octane, 2,6-dim...	10.781	61.7	ug/l	1277780	3	10.055	1035030	50.0
Cyclohexane, (1...	10.860	48.6	ug/l	1005240	3	10.055	1035030	50.0
Benzene, 1-ethe...	12.701	28.9	ug/l	842652	4	12.024	1460030	50.0
1-Phenyl-1-butene	13.292	39.9	ug/l	1164760	4	12.024	1460030	50.0
unknown13.420	13.420	29.7	ug/l	868449	4	12.024	1460030	50.0
Naphthalene, 2,...	15.615	36.6	ug/l	1067900	4	12.024	1460030	50.0