LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX021224\
Data File : VX040191.D
Acq On : 12 Feb 2024 11:57
Operator : JC/MD
Sample : P1411-01
Misc : 5.00g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
72-11929

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\82X020924W.M
Title : SW846 8260

Signal : TIC: VX040191.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	721	rBV	50654	154732	3.11%	0.205%
2	5.550	721	732	749	rVB	97540	281714	5.66%	0.374%
3	5.958	787	799	815	rBV2	58540	165216	3.32%	0.219%
4	6.757	920	930	947	rBV	149143	373347	7.50%	0.495%
5	8.647	1235	1240	1251	rVB	328646	523672	10.52%	0.694%
6	9.616	1394	1399	1403	rBV	68921	109242	2.19%	0.145%
7	9.829	1428	1434	1438	rBV2	41215	81714	1.64%	0.108%
8	10.006	1456	1463	1467	rBV	64616	87280	1.75%	0.116%
9	10.055	1467	1471	1476	rBV	312081	429093	8.62%	0.569%
10	10.274	1500	1507	1510	rBV3	58983	117279	2.36%	0.156%
11	10.317	1510	1514	1517	rVV	79848	124131	2.49%	0.165%
12	10.354	1517	1520	1532	rVB4	55375	122736	2.47%	0.163%
13	10.457	1532	1537	1545	rBV3	35032	75684	1.52%	0.100%
14	10.585	1551	1558	1566	rVB3	112436	215431	4.33%	0.286%
15	10.671	1566	1572	1577	rBV5	47121	103376	2.08%	0.137%
16	10.780	1582	1590	1594	rBV2	252266	404870	8.13%	0.537%
17	10.854	1598	1602	1606	rVV3	191405	312228	6.27%	0.414%
18	10.890	1606	1608	1613	rVB	84077	110988	2.23%	0.147%
19	10.957	1614	1619	1622	rBV3	48319	89713	1.80%	0.119%
20	11.000	1622	1626	1628	rVV3	55997	81643	1.64%	0.108%
21	11.030	1628	1631	1634	rVV2	64477	75703	1.52%	0.100%
22	11.085	1634	1640	1650	rVB3	438840	817623	16.42%	1.084%
23	11.189	1653	1657	1659	rBV	148528	191497	3.85%	0.254%
24	11.219	1659	1662	1666	rVB2	178475	227698	4.57%	0.302%
25	11.335	1678	1681	1686	rVV3	66502	116008	2.33%	0.154%
26	11.390	1686	1690	1693	rVV3	85324	130151	2.61%	0.173%
27	11.433	1693	1697	1701	rVV	238823	347160	6.97%	0.460%
28	11.481	1701	1705	1709	rVB3	121722	205805	4.13%	0.273%
29	11.573	1716	1720	1722	rBV3	61476	87155	1.75%	0.116%
30	11.628	1722	1729	1734	rVB5	92695	244120	4.90%	0.324%
31	11.683	1734	1738	1740	rBV3	100935	128373	2.58%	0.170%
32	11.713	1740	1743	1746	rBV	314005	338392	6.80%	0.449%
33	11.841	1760	1764	1765	rBV	73233	91286	1.83%	0.121%
34	11.890	1766	1772	1776	rVB3	272368	424676	8.53%	0.563%
35	11.945	1776	1781	1784	rBV2	405056	557565	11.20%	0.739%
36	12.024	1789	1794	1800	rBV2	463861	840308	16.88%	1.114%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX021224\
Data File : VX040191.D
Acq On : 12 Feb 2024 11:57
Operator : JC/MD
Sample : P1411-01
Misc : 5.00g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
72-11929

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\82X020924W.M
Title : SW846 8260

37	12.073	1800	1802	1805	rVB2	163065	145415	2.92%	0.193%
38	12.103	1805	1807	1811	rBV4	72623	75010	1.51%	0.099%
39	12.170	1812	1818	1826	rVB5	215106	526343	10.57%	0.698%
40	12.243	1826	1830	1837	rBV2	364708	615120	12.35%	0.816%
41	12.311	1837	1841	1848	rBV3	486660	968702	19.46%	1.285%
42	12.420	1848	1859	1864	rBV3	328224	1050647	21.10%	1.393%
43	12.469	1864	1867	1872	rVB4	273092	442903	8.90%	0.587%
44	12.554	1875	1881	1883	rBV6	270419	472873	9.50%	0.627%
45	12.585	1883	1886	1888	rVV2	147658	231205	4.64%	0.307%
46	12.615	1888	1891	1894	rVB	379644	409960	8.23%	0.544%
47	12.695	1899	1904	1913	rBV5	611727	1607593	32.29%	2.132%
48	12.762	1913	1915	1917	rBV3	118303	135013	2.71%	0.179%
49	12.817	1920	1924	1930	rVB3	702828	1129821	22.69%	1.498%
50	12.890	1930	1936	1939	rBV2	936022	1693987	34.02%	2.246%
51	12.981	1948	1951	1955	rBV2	574626	723329	14.53%	0.959%
52	13.030	1955	1959	1965	rVB4	499998	922098	18.52%	1.223%
53	13.097	1966	1970	1975	rBV3	359888	673553	13.53%	0.893%
54	13.140	1975	1977	1979	rBV2	185961	209442	4.21%	0.278%
55	13.170	1979	1982	1987	rVV2	319914	379717	7.63%	0.504%
56	13.262	1994	1997	1998	rBV3	203031	239616	4.81%	0.318%
57	13.292	1998	2002	2008	rVB4	2476017	3931133	78.96%	5.213%
58	13.347	2008	2011	2013	rBV2	346370	452973	9.10%	0.601%
59	13.384	2013	2017	2020	rVV2	1006449	1572415	31.58%	2.085%
60	13.420	2020	2023	2032	rVB7	764209	2062207	41.42%	2.735%
61	13.505	2032	2037	2040	rBV3	565533	960351	19.29%	1.274%
62	13.548	2040	2044	2047	rBV	1062904	1409927	28.32%	1.870%
63	13.585	2047	2050	2054	rVV4	764575	963994	19.36%	1.278%
64	13.664	2054	2063	2068	rBV3	1462873	3205520	64.38%	4.251%
65	13.737	2072	2075	2082	rVB4	793627	1526724	30.66%	2.025%
66	13.798	2082	2085	2089	rBV3	886353	1377034	27.66%	1.826%
67	13.847	2089	2093	2099	rVB3	1148674	1651981	33.18%	2.191%
68	13.926	2099	2106	2111	rBV6	1094949	2360641	47.41%	3.131%
69	13.975	2111	2114	2117	rBV3	705419	1074147	21.57%	1.424%
70	14.072	2126	2130	2133	rBV2	1057692	1466959	29.46%	1.945%
71	14.103	2133	2135	2138	rVB2	428944	406506	8.16%	0.539%
72	14.140	2138	2141	2143	rBV3	293444	337855	6.79%	0.448%
73	14.164	2143	2145	2147	rVV	334436	302697	6.08%	0.401%
74	14.213	2150	2153	2158	rVB3	1641379	2335478	46.91%	3.097%
75	14.322	2168	2171	2176	rBV2	810557	1165378	23.41%	1.545%
76	14.371	2176	2179	2183	rVB	1629534	2097362	42.13%	2.781%

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Instrument :
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ClientSampleId :
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Integration Parameters: RTEINT.P
Integrator: RTE
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Title : SW846 8260

77	14.420	2183	2187	2193	rVB4	680483	1228486	24.67%	1.629%
78	14.487	2193	2198	2202	rBV3	1179400	2244817	45.09%	2.977%
79	14.554	2206	2209	2210	rBV2	207546	208025	4.18%	0.276%
80	14.578	2210	2213	2216	rVV3	399152	447854	9.00%	0.594%
81	14.615	2216	2219	2225	rVV6	1095201	1561578	31.36%	2.071%
82	14.676	2225	2229	2233	rVB2	1331579	1767291	35.50%	2.344%
83	14.725	2233	2237	2240	rBV2	666894	943456	18.95%	1.251%
84	14.780	2242	2246	2251	rVV	3383426	4978770	100.00%	6.603%
85	14.853	2255	2258	2261	rVB3	387545	445031	8.94%	0.590%
86	14.883	2261	2263	2269	rBV5	165384	222727	4.47%	0.295%
87	15.048	2287	2290	2293	rBV3	234749	281226	5.65%	0.373%
88	15.182	2307	2312	2319	rVB3	572297	1017846	20.44%	1.350%
89	15.377	2339	2344	2347	rBV	745745	1033898	20.77%	1.371%
90	15.408	2347	2349	2353	rVB2	397546	494298	9.93%	0.656%
91	15.481	2356	2361	2365	rBV	1080081	1583606	31.81%	2.100%
92	15.615	2378	2383	2386	rBV	1065484	1652433	33.19%	2.191%
93	15.651	2386	2389	2397	rVB	799293	1236193	24.83%	1.639%
94	15.731	2397	2402	2409	rVB4	174912	330491	6.64%	0.438%
95	15.834	2411	2419	2429	rVB3	269303	876432	17.60%	1.162%
96	15.987	2439	2444	2455	rVB	155969	302995	6.09%	0.402%
97	16.353	2499	2504	2508	rBV	63367	106036	2.13%	0.141%
98	16.554	2530	2537	2540	rVV5	29382	70812	1.42%	0.094%
99	16.596	2540	2544	2549	rVV2	66064	128608	2.58%	0.171%
100	16.657	2549	2554	2572	rVB2	51051	142237	2.86%	0.189%

Sum of corrected areas: 75406384

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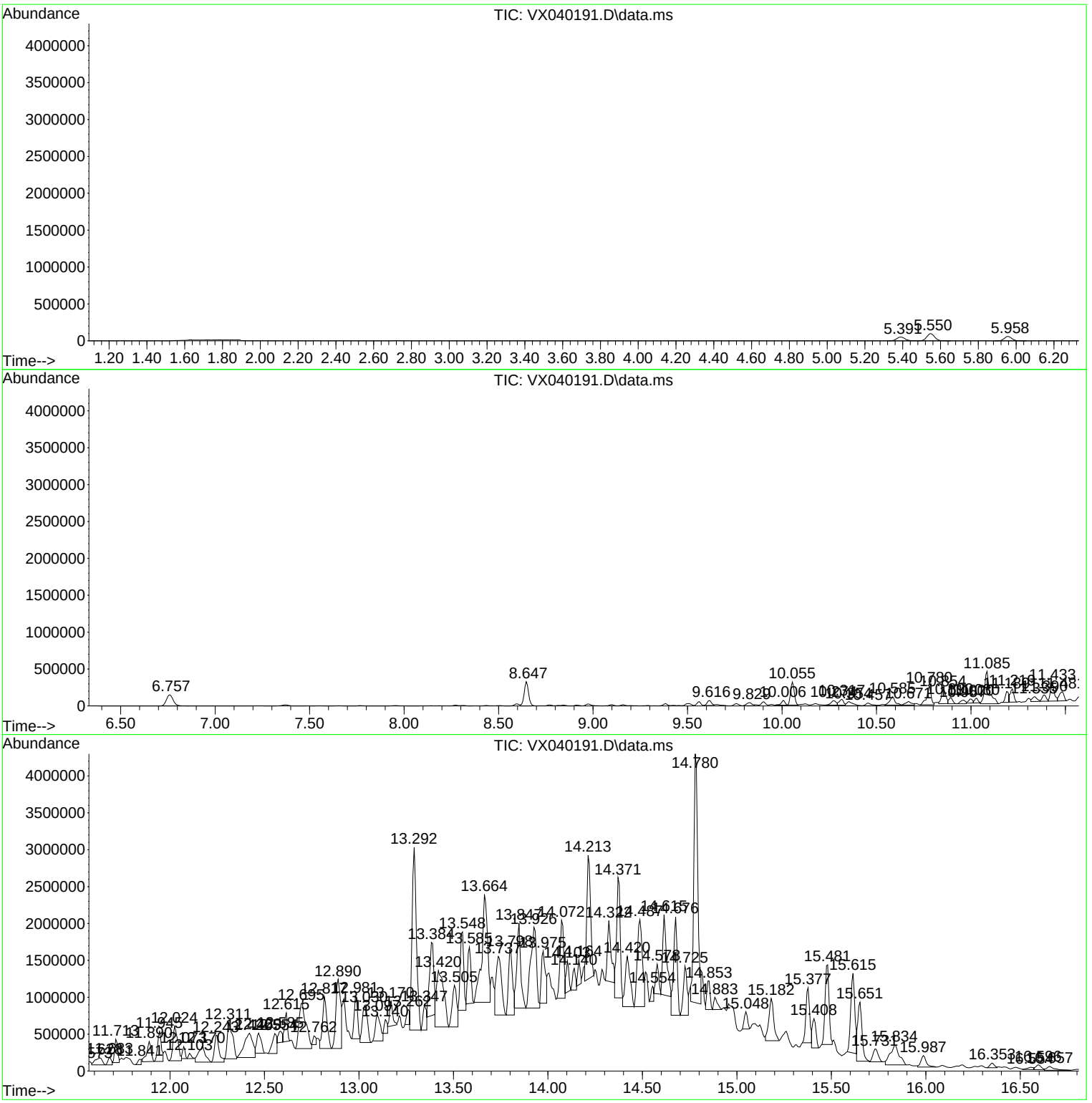
Instrument :
MSVOA_X
ClientSampleId :
72-11929

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X020924W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX021224\
Data File : VX040191.D
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Instrument :
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ClientSampleId :
72-11929

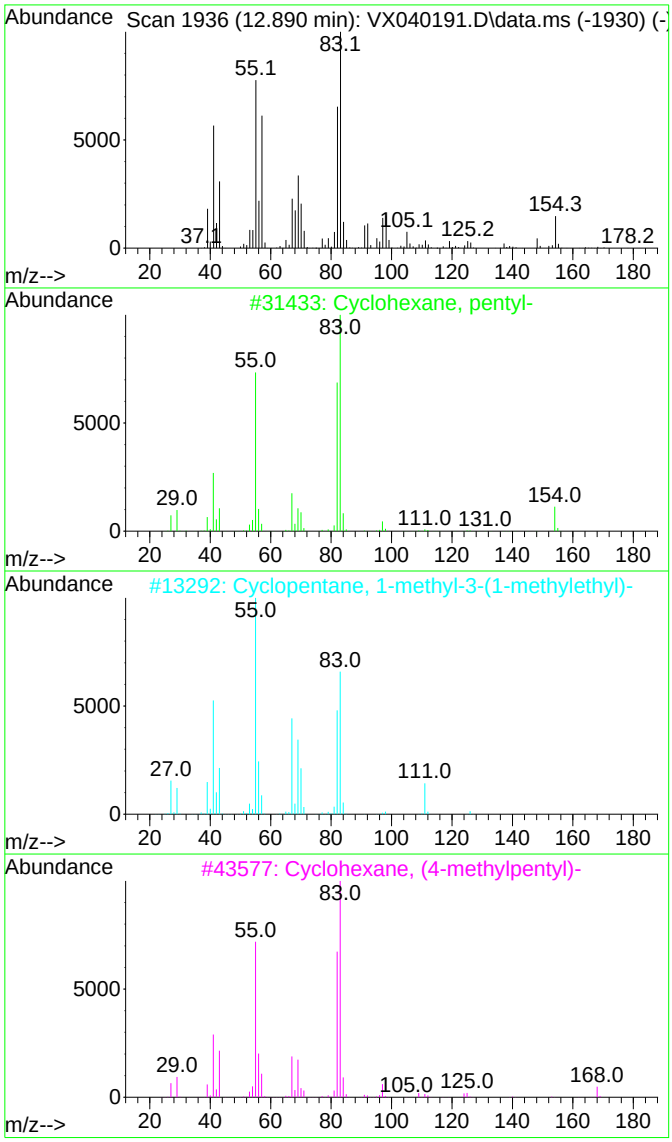
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X020924W.M
Quant Title : SW846 8260

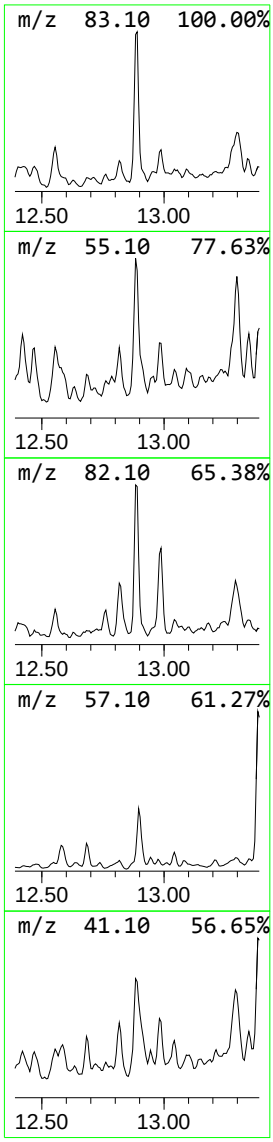
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Cyclohexane, pentyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.890	100.80 ug/l	1693990	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, pentyl-	154	C11H22	004292-92-6	72
2	Cyclopentane, 1-methyl-3-(1-meth...	126	C9H18	053771-88-3	68
3	Cyclohexane, (4-methylpentyl)-	168	C12H24	061142-20-9	59
4	Cyclohexane, methyl-	98	C7H14	000108-87-2	52
5	Cyclopentane, 1-ethyl-1-methyl-	112	C8H16	016747-50-5	46





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ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
72-11929

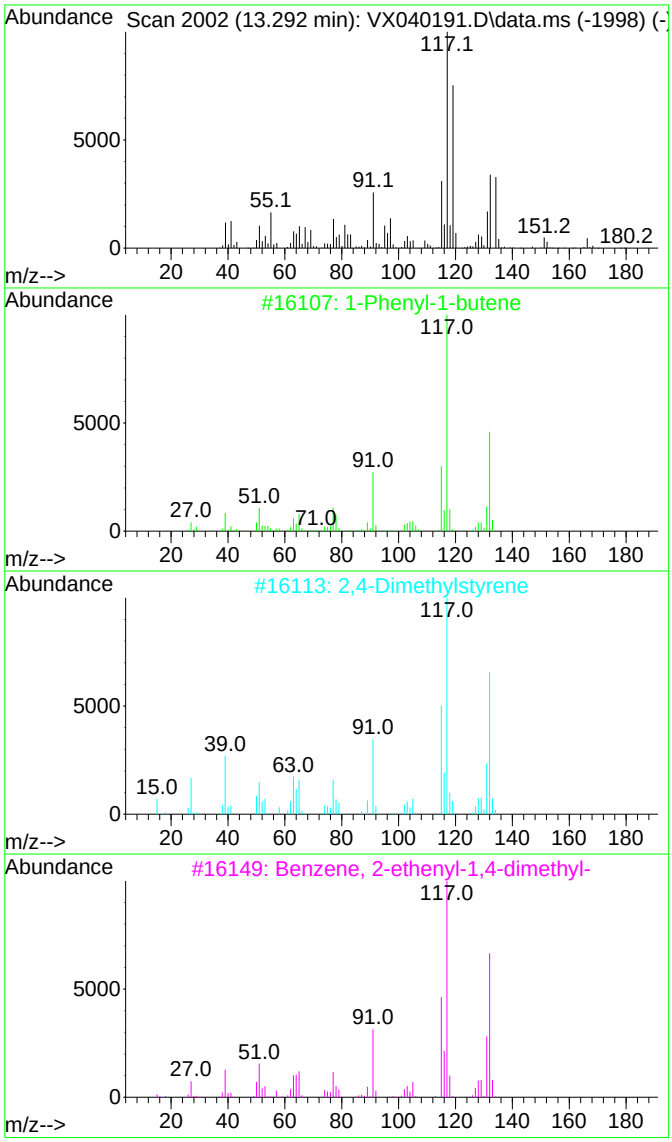
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X020924W.M
Quant Title : SW846 8260

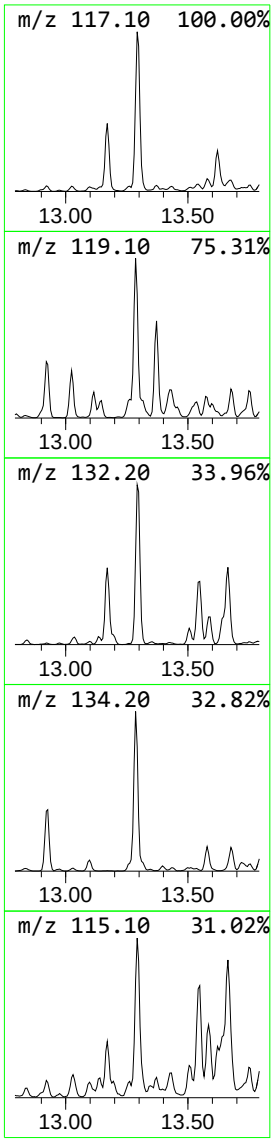
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	233.91 ug/l	3931130	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	91
2	2,4-Dimethylstyrene	132	C10H12	002234-20-0	86
3	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	83
4	Benzene, 2-butenyl-	132	C10H12	001560-06-1	80
5	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	55





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Misc : 5.00g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
72-11929

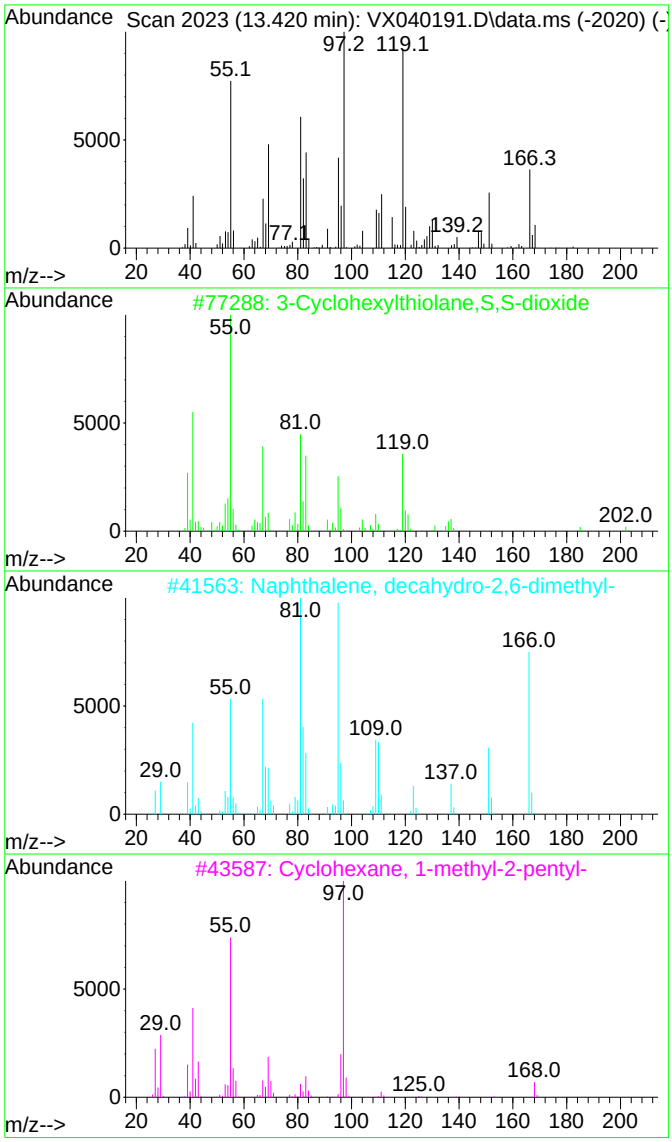
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X020924W.M
Quant Title : SW846 8260

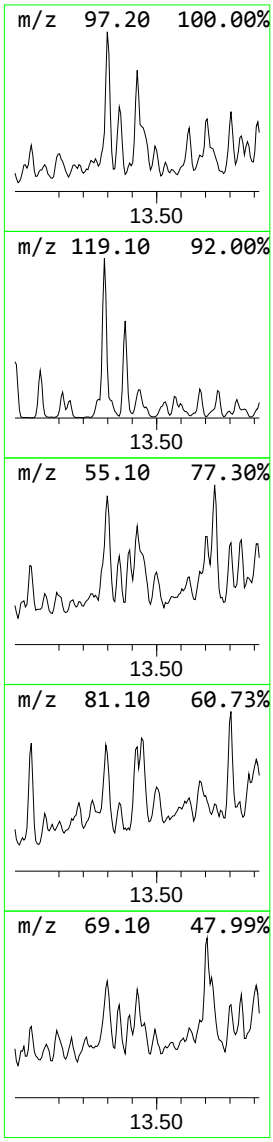
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 unknown13.420 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.420	122.71 ug/l	2062210	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Cyclohexylthiolane,S,S-dioxide	202	C10H18O2S	071053-08-2	37
2	Naphthalene, decahydro-2,6-dimet...	166	C12H22	001618-22-0	25
3	Cyclohexane, 1-methyl-2-pentyl-	168	C12H24	054411-01-7	22
4	Cyclohexane, 1-methyl-3-pentyl-	168	C12H24	054411-02-8	14
5	Pentanoic acid, 10-undecenyl ester	254	C16H30O2	1000159-93-4	12





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ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
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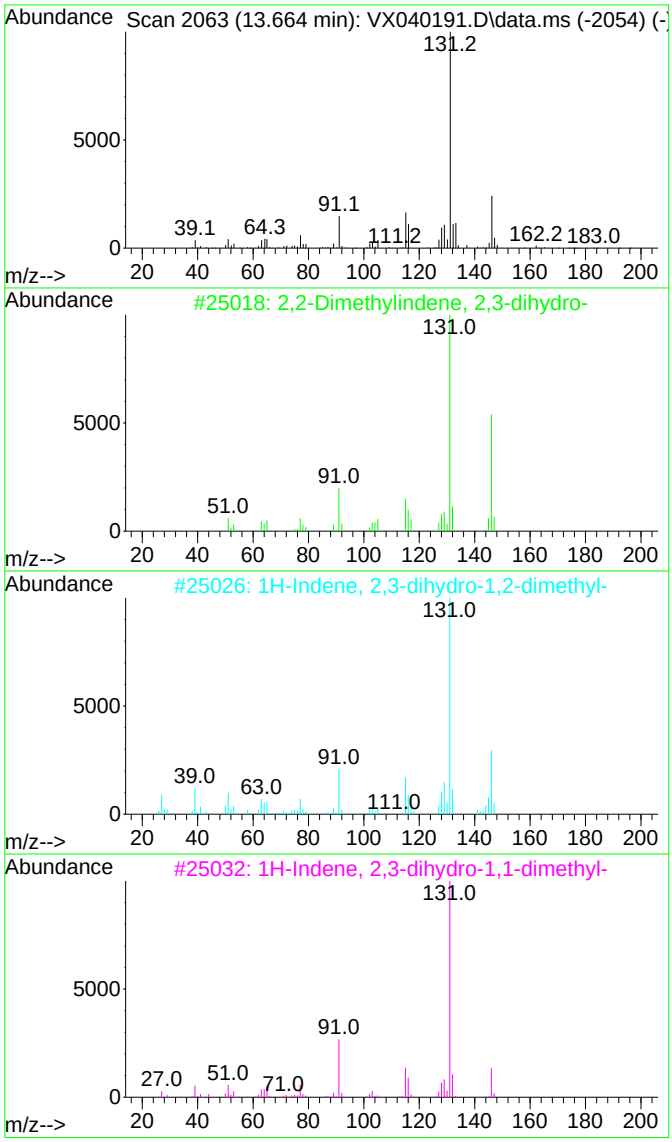
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X020924W.M
Quant Title : SW846 8260

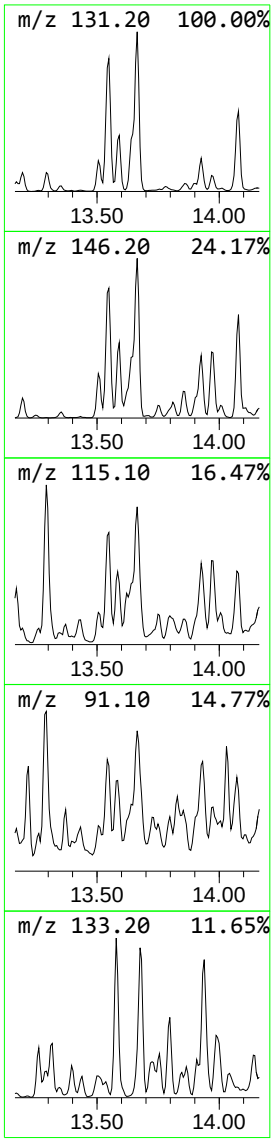
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 2,2-Dimethylindene, 2,3-dih... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.664	190.74 ug/l	3205520	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	90
2	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	90
3	1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	90
4	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	90
5	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	90





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Instrument :
MSVOA_X
ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X020924W.M
Quant Title : SW846 8260

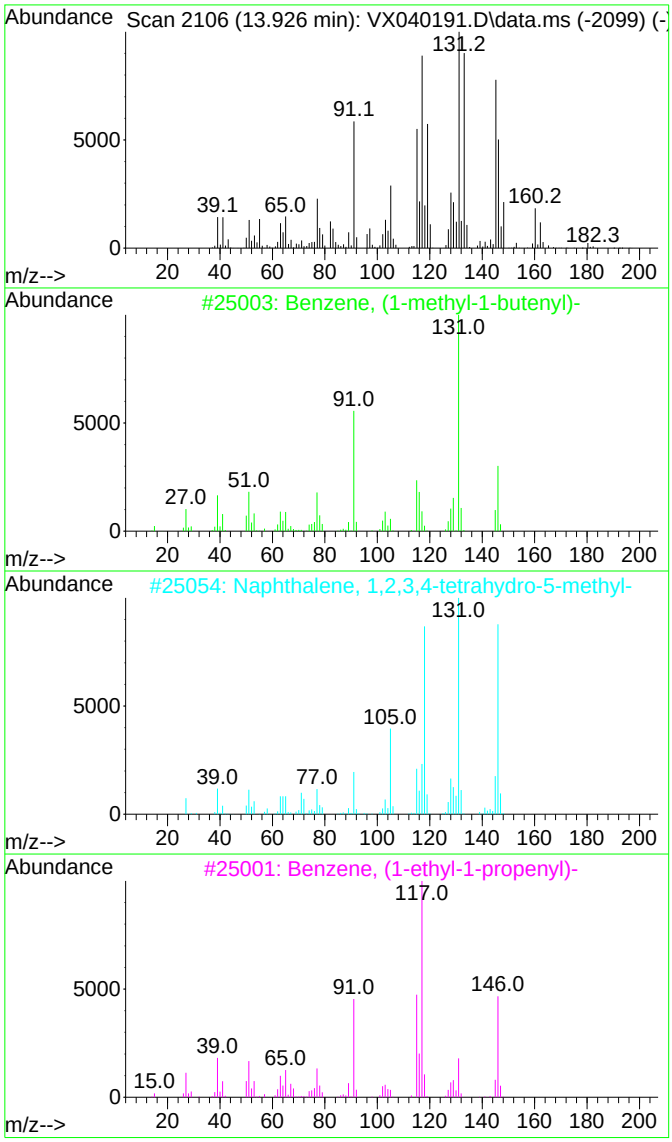
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

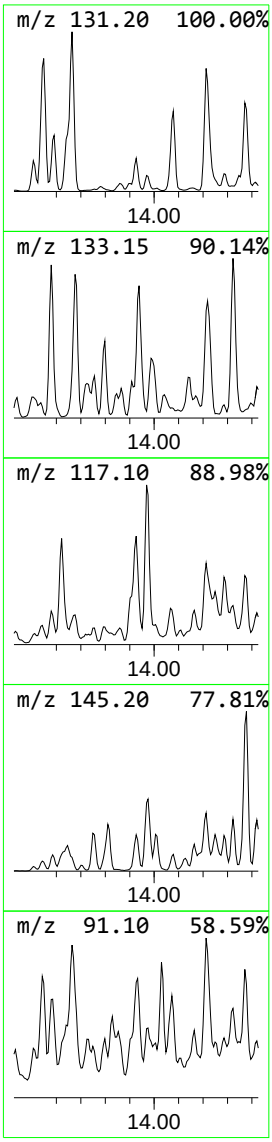
****

Peak Number 5 unknown13.926 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.926	140.46 ug/l	2360640	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	38
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	38
3	Benzene, (1-ethyl-1-propenyl)-	146	C11H14	004701-36-4	35
4	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	35
5	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	30





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Data File : VX040191.D
Acq On : 12 Feb 2024 11:57
Operator : JC/MD
Sample : P1411-01
Misc : 5.00g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
72-11929

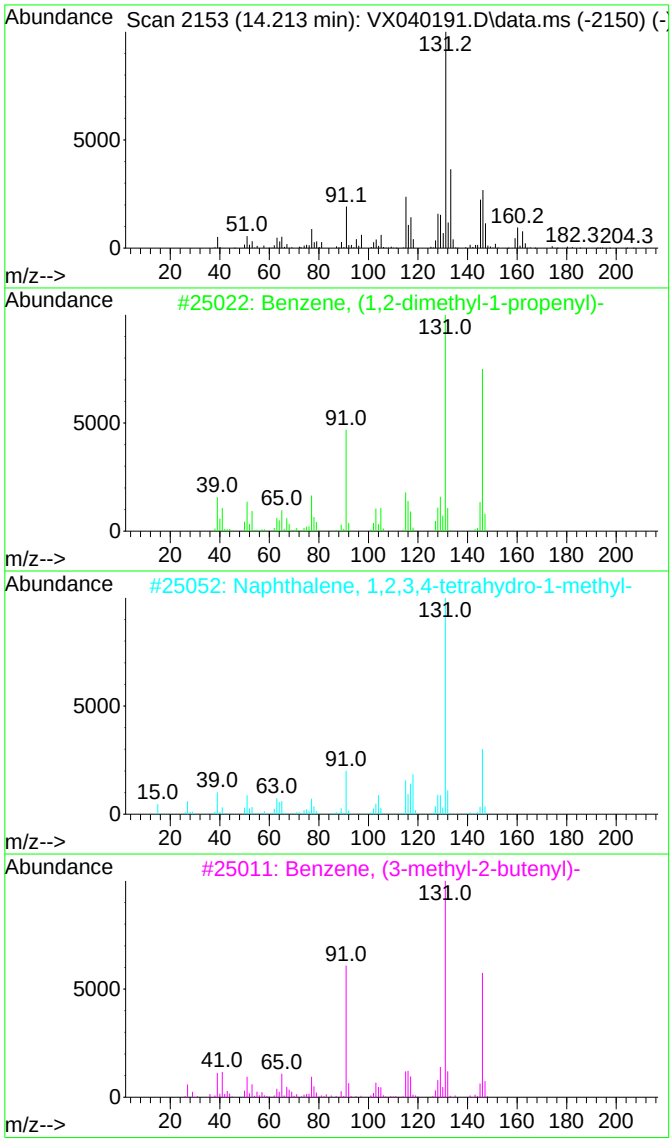
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X020924W.M
Quant Title : SW846 8260

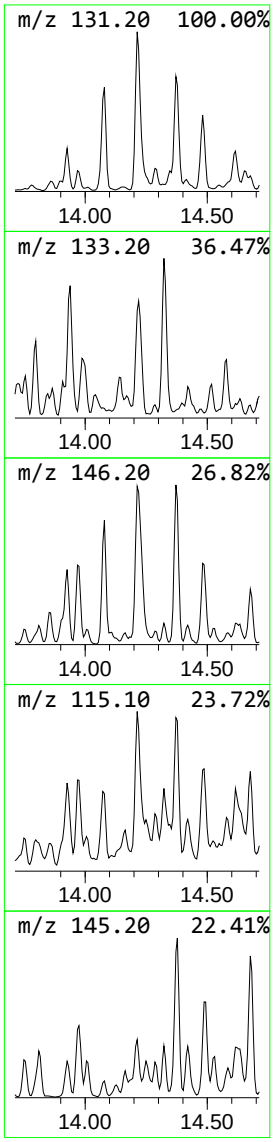
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, (1,2-dimethyl-1-pr... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.213	138.97 ug/l	2335480	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	92
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	70
3	Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	70
4	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	70
5	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	64





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Data File : VX040191.D
Acq On : 12 Feb 2024 11:57
Operator : JC/MD
Sample : P1411-01
Misc : 5.00g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
72-11929

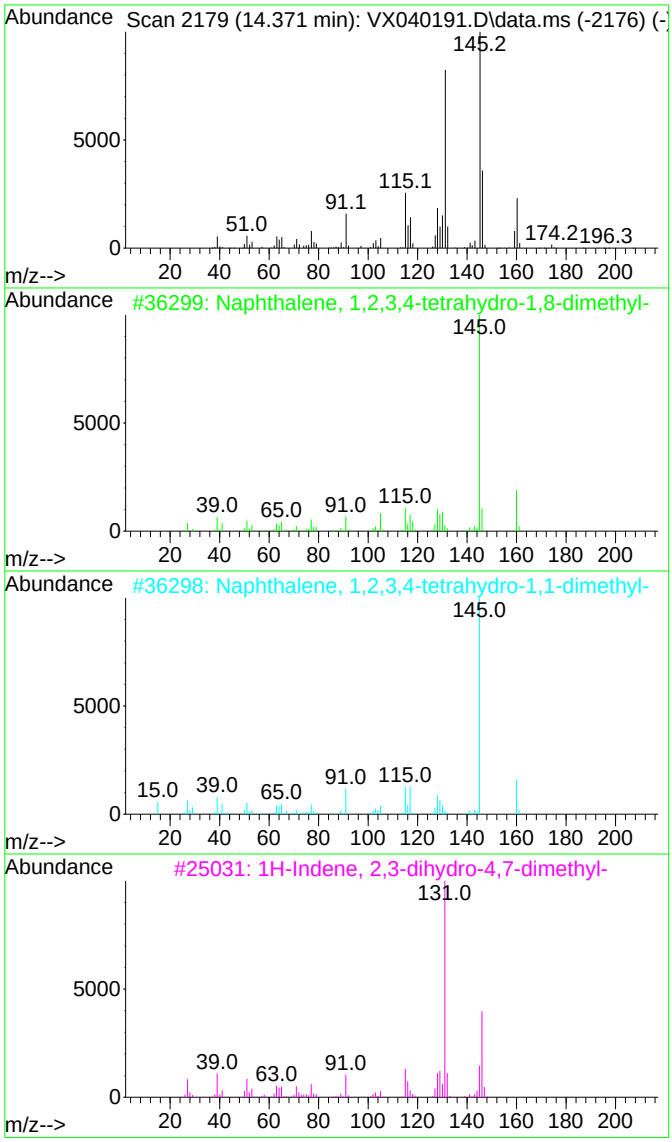
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X020924W.M
Quant Title : SW846 8260

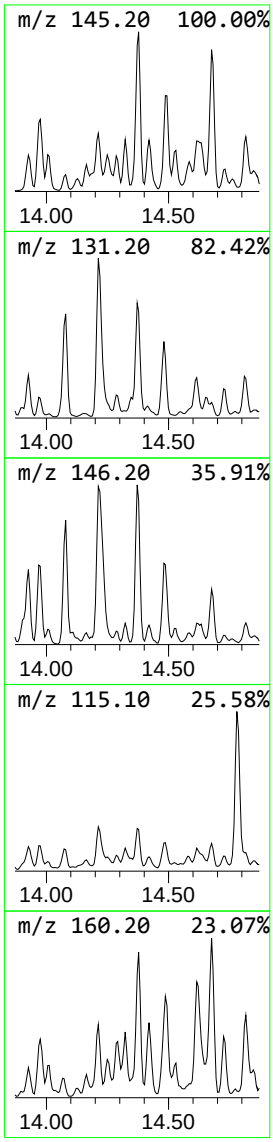
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.371	124.80 ug/l	2097360	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	025419-33-4	55
2	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	001985-59-7	55
3	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	53
4	Naphthalene, 6-ethyl-1,2,3,4-tet...	160	C12H16	022531-20-0	49
5	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	49





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX021224\
Data File : VX040191.D
Acq On : 12 Feb 2024 11:57
Operator : JC/MD
Sample : P1411-01
Misc : 5.00g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
72-11929

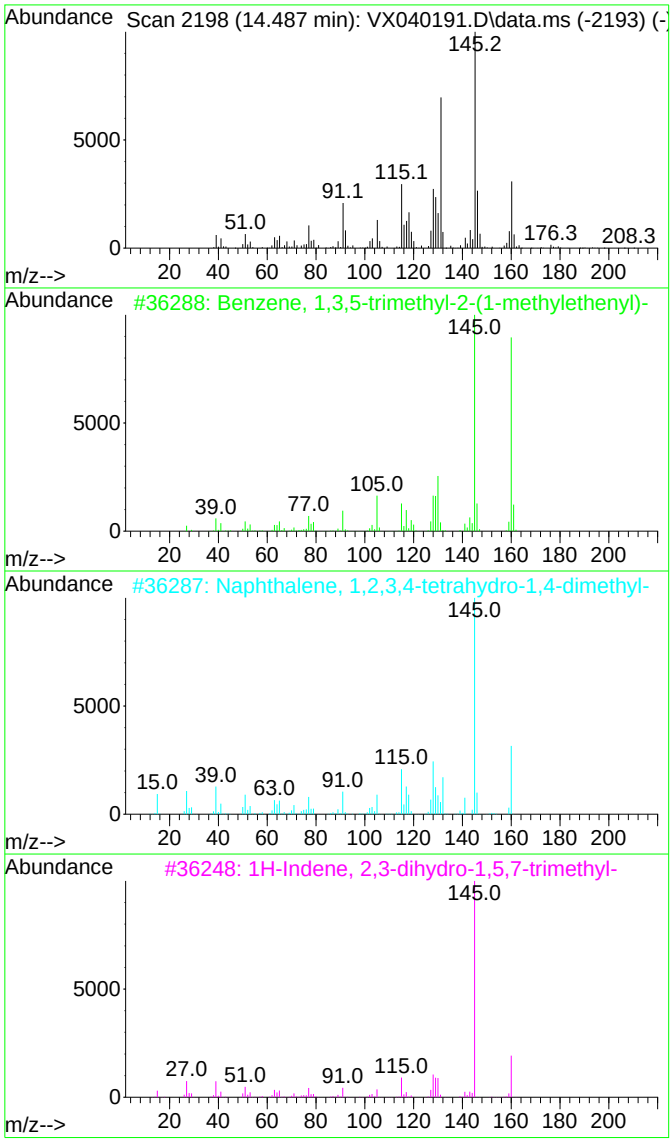
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X020924W.M
Quant Title : SW846 8260

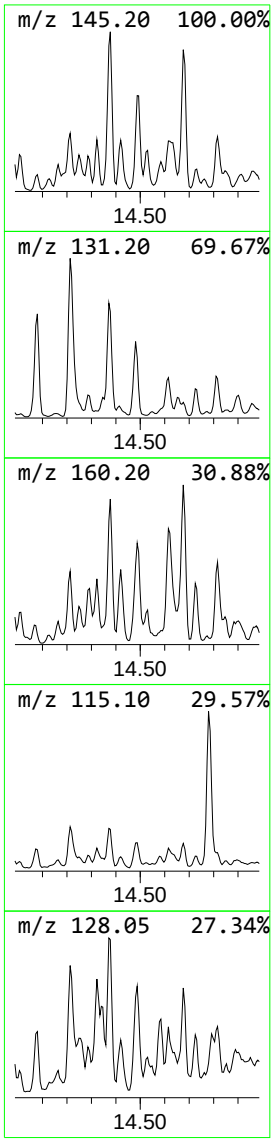
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 1,3,5-trimethyl-2-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.487	133.57 ug/l	2244820	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	83
2	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	68
3	1H-Indene, 2,3-dihydro-1,5,7-tri...	160	C12H16	054340-88-4	62
4	1H-Pyrrolo[2,3-b]pyridine, 2-ethyl-	146	C9H10N2	023612-49-9	58
5	1H-Indene, 2,3-dihydro-1,1,3-tri...	160	C12H16	002613-76-5	58





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Data File : VX040191.D
Acq On : 12 Feb 2024 11:57
Operator : JC/MD
Sample : P1411-01
Misc : 5.00g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
72-11929

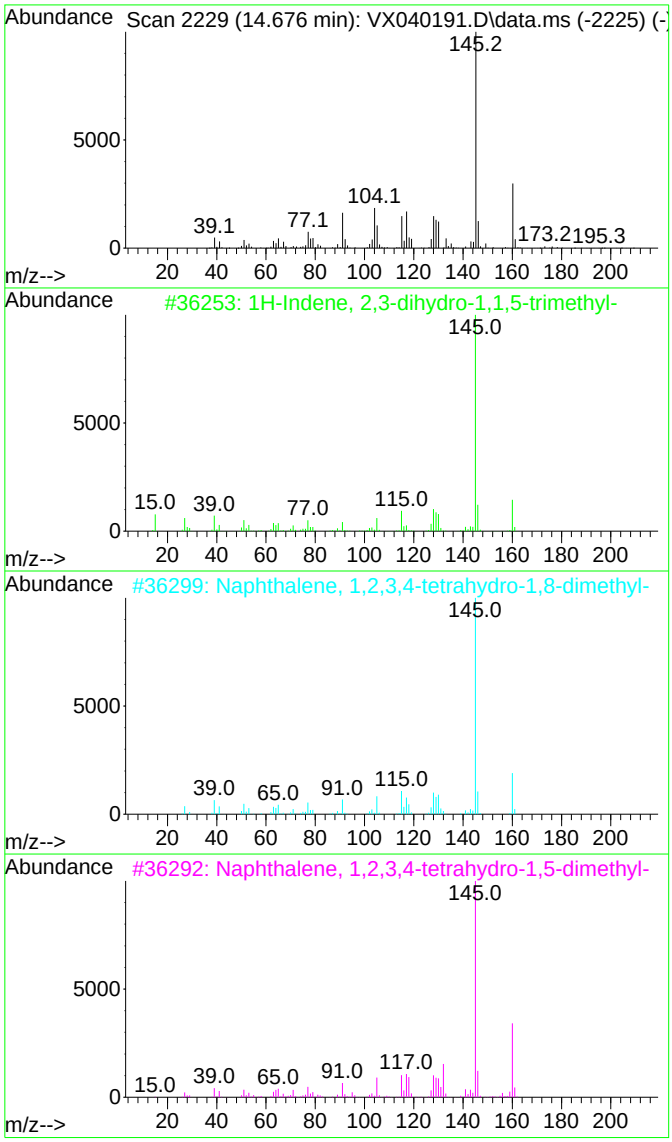
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X020924W.M
Quant Title : SW846 8260

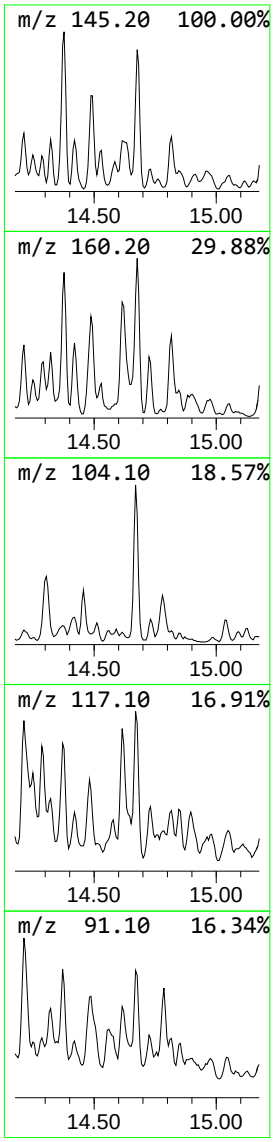
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.676	105.16 ug/l	1767290	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,1,5-tri...	160	C12H16	040650-41-7	93
2	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	025419-33-4	90
3	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	021564-91-0	87
4	Benzene, (1,3-dimethyl-2-butenyl)-	160	C12H16	050704-01-3	87
5	Benzene, 1-(1-methylethenyl)-3-(...	160	C12H16	001129-29-9	87





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Data File : VX040191.D
Acq On : 12 Feb 2024 11:57
Operator : JC/MD
Sample : P1411-01
Misc : 5.00g/10mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
72-11929

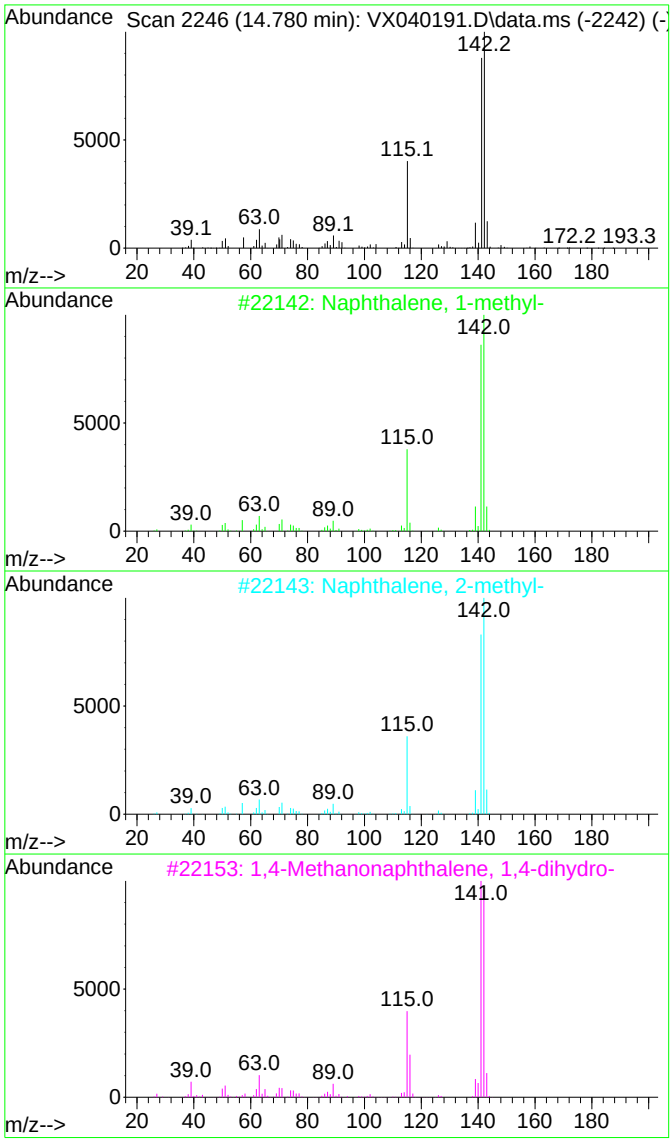
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X020924W.M
Quant Title : SW846 8260

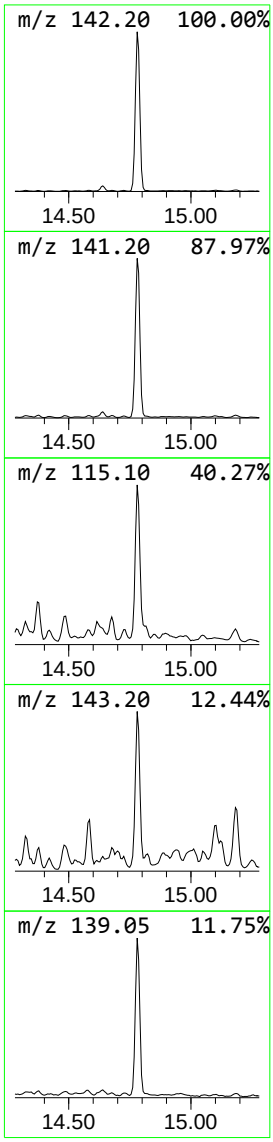
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.780	296.25 ug/l	4978770	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	93
4	Benzocycloheptatriene	142	C11H10	000264-09-5	91
5	1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	64





Tentatively Identified Compound (LSC) summary

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Data File : VX040191.D
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Instrument :
MSVOA_X
ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X020924W.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, pe...	12.890	100.8	ug/l	1693990	4	12.024	840308	50.0
1-Phenyl-1-butene	13.292	233.9	ug/l	3931130	4	12.024	840308	50.0
unknown13.420	13.420	122.7	ug/l	2062210	4	12.024	840308	50.0
2,2-Dimethylind...	13.664	190.7	ug/l	3205520	4	12.024	840308	50.0
unknown13.926	13.926	140.5	ug/l	2360640	4	12.024	840308	50.0
Benzene, (1,2-d...	14.213	139.0	ug/l	2335480	4	12.024	840308	50.0
Naphthalene, 1,...	14.371	124.8	ug/l	2097360	4	12.024	840308	50.0
Benzene, 1,3,5-...	14.487	133.6	ug/l	2244820	4	12.024	840308	50.0
1H-Indene, 2,3-...	14.676	105.2	ug/l	1767290	4	12.024	840308	50.0
Naphthalene, 1-...	14.780	296.3	ug/l	4978770	4	12.024	840308	50.0