LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0 % 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\SFAMXLM122821WMA.M
Title : VOC Analysis

Signal : TIC: VX026120.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.374	41	47	52	rBV	142931	177757	3.91%	0.384%
2	1.666	90	95	102	rVB	109637	137295	3.02%	0.296%
3	2.313	194	201	211	rBV2	227515	420922	9.25%	0.909%
4	3.447	379	387	401	rBV2	44106	99174	2.18%	0.214%
5	3.617	401	415	421	rBV	64001	173708	3.82%	0.375%
6	4.495	541	559	574	rVB2	185747	658179	14.47%	1.421%
7	5.068	642	653	663	rVV	116417	367538	8.08%	0.794%
8	5.190	664	673	687	rVB5	40901	153880	3.38%	0.332%
9	5.391	693	706	716	rBV	500203	1570786	34.52%	3.392%
10	5.495	716	723	734	rVB5	39687	152104	3.34%	0.328%
11	5.623	734	744	761	rVB4	33123	135416	2.98%	0.292%
12	5.830	767	778	790	rBV2	44504	135520	2.98%	0.293%
13	5.977	790	802	810	rBV2	306045	918089	20.18%	1.982%
14	6.257	841	848	857	rVB2	58214	174173	3.83%	0.376%
15	6.367	857	866	875	rVB3	38825	109086	2.40%	0.236%
16	6.617	895	907	914	rBV	99437	243654	5.36%	0.526%
17	6.769	924	932	939	rBV	189888	475762	10.46%	1.027%
18	6.848	940	945	953	rVB5	49053	135346	2.97%	0.292%
19	7.178	992	999	1006	rVB3	46561	105919	2.33%	0.229%
20	7.312	1013	1021	1027	rBV	191013	423606	9.31%	0.915%
21	7.385	1027	1033	1044	rVB2	174068	408727	8.98%	0.883%
22	7.988	1123	1132	1142	rVB4	41380	144035	3.17%	0.311%
23	8.104	1144	1151	1154	rBV3	54135	108718	2.39%	0.235%
24	8.147	1154	1158	1162	rVV	55962	100716	2.21%	0.217%
25	8.281	1174	1180	1183	rBV2	93495	176557	3.88%	0.381%
26	8.324	1183	1187	1194	rVB2	158967	287719	6.32%	0.621%
27	8.440	1201	1206	1210	rBV	77091	125845	2.77%	0.272%
28	8.507	1210	1217	1222	rVB3	63498	119421	2.62%	0.258%
29	8.604	1227	1233	1235	rBV2	74512	124351	2.73%	0.269%
30	8.653	1235	1241	1246	rVB	508623	855007	18.79%	1.846%
31	8.720	1246	1252	1259	rBV	2916885	4549918	100.00%	9.824%
32	8.915	1277	1284	1287	rBV2	186608	315160	6.93%	0.681%
33	9.275	1338	1343	1349	rBV2	85282	130113	2.86%	0.281%
34	9.391	1356	1362	1367	rBV	310975	462384	10.16%	0.998%
35	9.500	1376	1380	1386	rVV4	66003	133899	2.94%	0.289%
36	9.567	1386	1391	1394	rVV	71798	101260	2.23%	0.219%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0 % 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\SFAMXLM122821WMA.M
Title : VOC Analysis

37	9.616	1394	1399	1403	rVV2	61694	110890	2.44%	0.239%
38	9.762	1417	1423	1429	rBV3	55396	105228	2.31%	0.227%
39	9.836	1429	1435	1438	rBV3	50126	105807	2.33%	0.228%
40	9.909	1443	1447	1451	rVV	110601	159650	3.51%	0.345%
41	10.012	1459	1464	1467	rVV	124994	166908	3.67%	0.360%
42	10.055	1467	1471	1476	rVV	414180	552959	12.15%	1.194%
43	10.195	1489	1494	1499	rVV	792153	1053067	23.14%	2.274%
44	10.305	1499	1512	1517	rVV	2536452	3699172	81.30%	7.987%
45	10.360	1517	1521	1531	rVB	509927	701595	15.42%	1.515%
46	10.457	1531	1537	1541	rBV4	56325	97389	2.14%	0.210%
47	10.586	1551	1558	1563	rBV2	122404	217543	4.78%	0.470%
48	10.646	1563	1568	1582	rVB	1027273	1483645	32.61%	3.204%
49	10.781	1582	1590	1594	rBV2	242781	394353	8.67%	0.852%
50	10.860	1594	1603	1606	rBV3	180654	305461	6.71%	0.660%
51	10.896	1606	1609	1614	rVB	127663	174378	3.83%	0.377%
52	10.963	1614	1620	1624	rBV	171866	264495	5.81%	0.571%
53	11.085	1634	1640	1642	rBV2	102370	146710	3.22%	0.317%
54	11.110	1642	1644	1651	rVB2	126864	200319	4.40%	0.433%
55	11.195	1651	1658	1661	rBV	432355	618663	13.60%	1.336%
56	11.305	1671	1676	1680	rBV	338856	425259	9.35%	0.918%
57	11.378	1683	1688	1695	rVV	1080305	1962119	43.12%	4.237%
58	11.451	1695	1700	1702	rVV	596248	850523	18.69%	1.837%
59	11.482	1702	1705	1711	rVB	637685	775808	17.05%	1.675%
60	11.616	1722	1727	1734	rVB	501484	668562	14.69%	1.444%
61	11.719	1740	1744	1746	rVV	147518	188718	4.15%	0.407%
62	11.756	1746	1750	1760	rVB	1977770	2454343	53.94%	5.300%
63	11.866	1761	1768	1770	rBV2	63442	121073	2.66%	0.261%
64	11.890	1770	1772	1777	rVV	135399	169836	3.73%	0.367%
65	11.945	1777	1781	1783	rVV2	135221	178318	3.92%	0.385%
66	11.969	1783	1785	1789	rVV	188498	234872	5.16%	0.507%
67	12.024	1789	1794	1800	rVV2	477666	765615	16.83%	1.653%
68	12.091	1800	1805	1813	rVV	656312	924745	20.32%	1.997%
69	12.177	1813	1819	1826	rVB4	73131	142297	3.13%	0.307%
70	12.244	1826	1830	1832	rBV2	349233	453220	9.96%	0.979%
71	12.268	1832	1834	1838	rVV	369146	417534	9.18%	0.902%
72	12.323	1838	1843	1849	rVB4	1009441	1591534	34.98%	3.437%
73	12.427	1850	1860	1863	rBV	335682	491362	10.80%	1.061%
74	12.463	1863	1866	1873	rVB	198247	263510	5.79%	0.569%
75	12.536	1873	1878	1879	rBV	232590	289994	6.37%	0.626%
76	12.555	1879	1881	1885	rVV	252615	279516	6.14%	0.604%

LSC Area Percent Report

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Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0 % 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\SFAMXLM122821WMA.M
Title : VOC Analysis

77	12.616	1887	1891	1894	rVV	347103	402190	8.84%	0.868%
78	12.701	1900	1905	1913	rBV4	209719	472125	10.38%	1.019%
79	12.847	1926	1929	1933	rVB2	119612	150578	3.31%	0.325%
80	12.920	1938	1941	1945	rVV	168985	197656	4.34%	0.427%
81	12.963	1945	1948	1955	rVB	332793	433410	9.53%	0.936%
82	13.170	1974	1982	1986	rBV4	173303	280627	6.17%	0.606%
83	13.292	1992	2002	2007	rBV3	481080	870945	19.14%	1.881%
84	13.426	2020	2024	2032	rVB2	138677	215279	4.73%	0.465%
85	13.548	2040	2044	2046	rBV	86262	109949	2.42%	0.237%
86	13.585	2046	2050	2056	rVV3	134547	252302	5.55%	0.545%
87	13.664	2060	2063	2069	rVB2	120198	187933	4.13%	0.406%
88	13.780	2075	2082	2089	rVB	379826	526566	11.57%	1.137%
89	13.926	2099	2106	2111	rBV4	87379	162639	3.57%	0.351%
90	14.079	2127	2131	2135	rBV2	88492	109011	2.40%	0.235%
91	14.219	2149	2154	2163	rBV4	119750	227082	4.99%	0.490%
92	14.371	2176	2179	2184	rVB	98286	125403	2.76%	0.271%
93	14.481	2193	2197	2202	rBV2	68719	117581	2.58%	0.254%
94	14.640	2215	2223	2227	rVV	684669	904869	19.89%	1.954%
95	14.780	2240	2246	2255	rVV	356888	503316	11.06%	1.087%
96	15.377	2336	2344	2347	rBV	79686	121043	2.66%	0.261%
97	15.481	2355	2361	2374	rVV	152695	279942	6.15%	0.604%
98	15.615	2377	2383	2386	rVV	169232	252131	5.54%	0.544%
99	15.652	2386	2389	2398	rVB	104891	172870	3.80%	0.373%
100	15.841	2410	2420	2430	rVB2	42730	117972	2.59%	0.255%

Sum of corrected areas: 46312153

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
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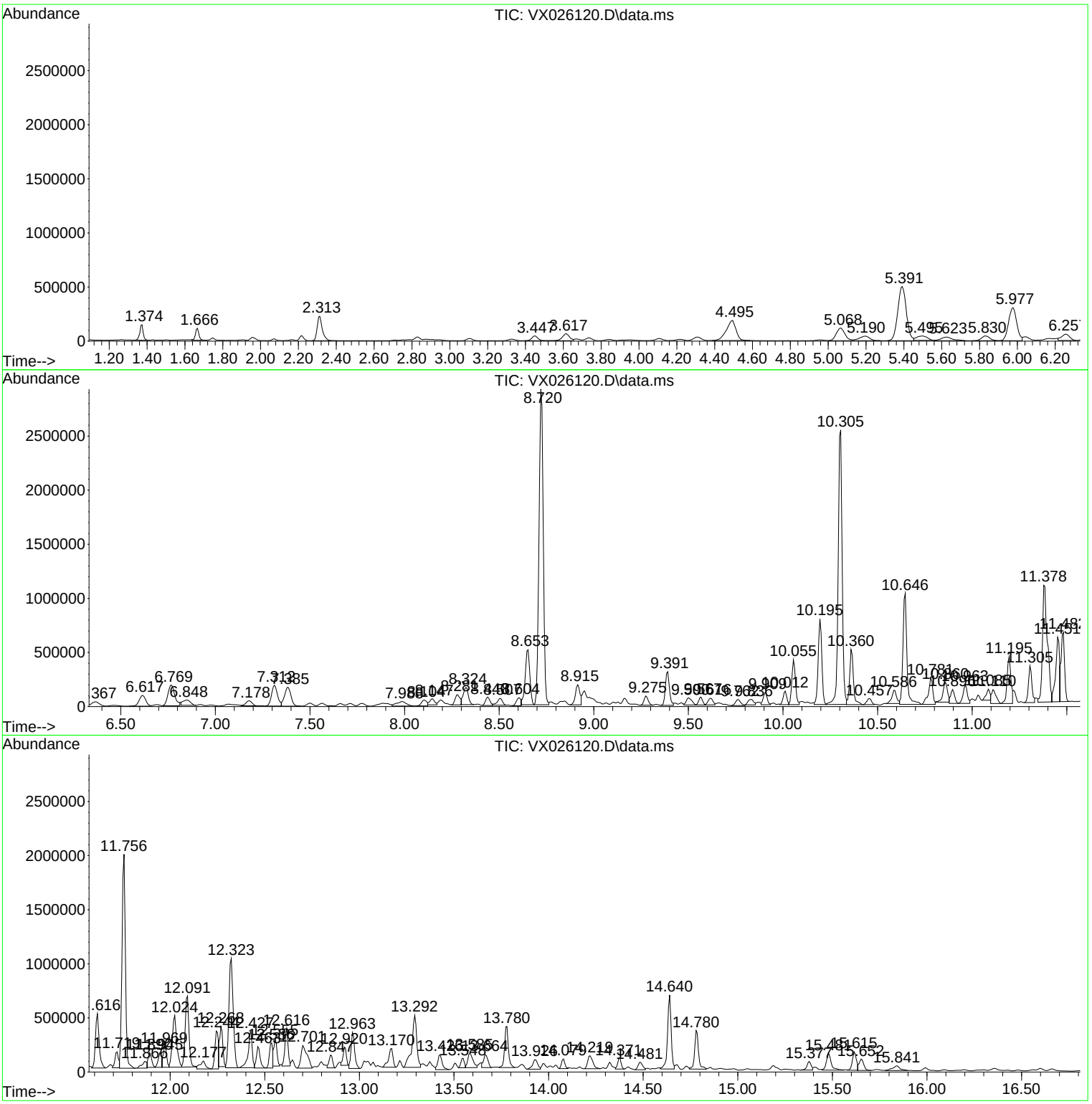
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

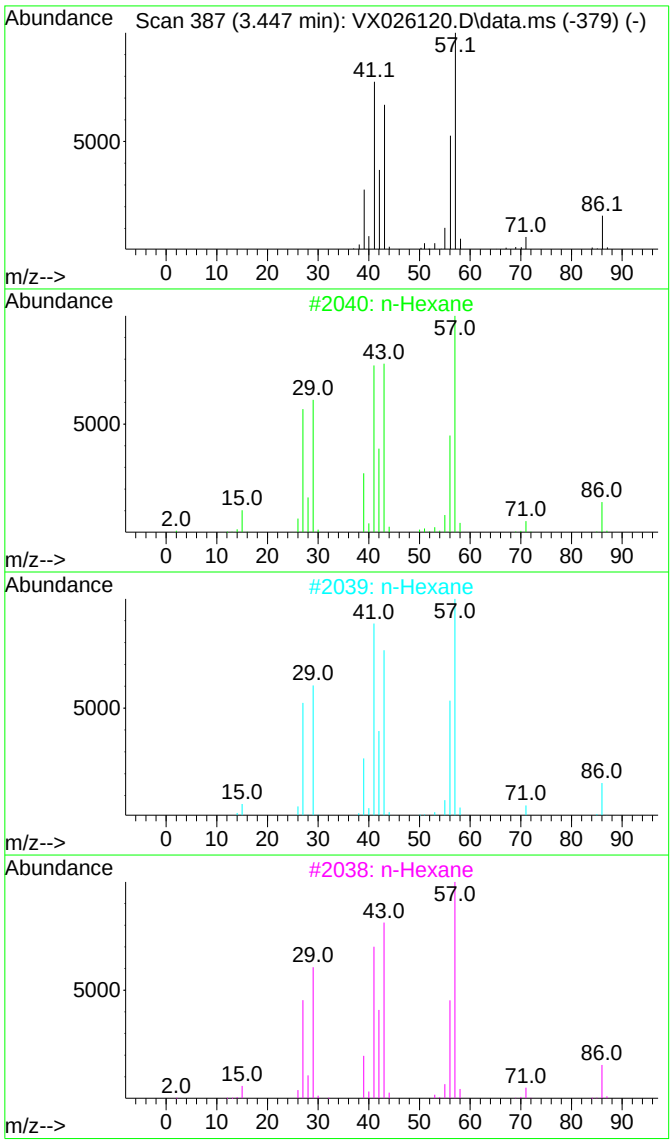
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

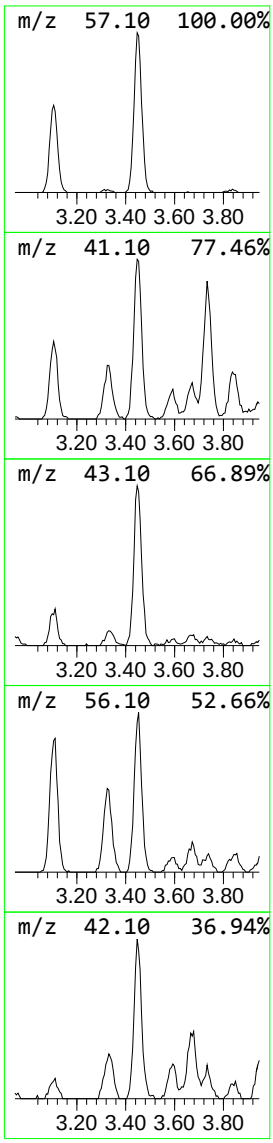
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 51

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.447	10.42 ug/L	99174	1,4-Difluorobenzene	6.763

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexane	86	C6H14	000110-54-3	91
2	n-Hexane	86	C6H14	000110-54-3	91
3	n-Hexane	86	C6H14	000110-54-3	91
4	2-Ethyl-oxetane	86	C5H10O	1010386-40-2	78
5	n-Hexane	86	C6H14	000110-54-3	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

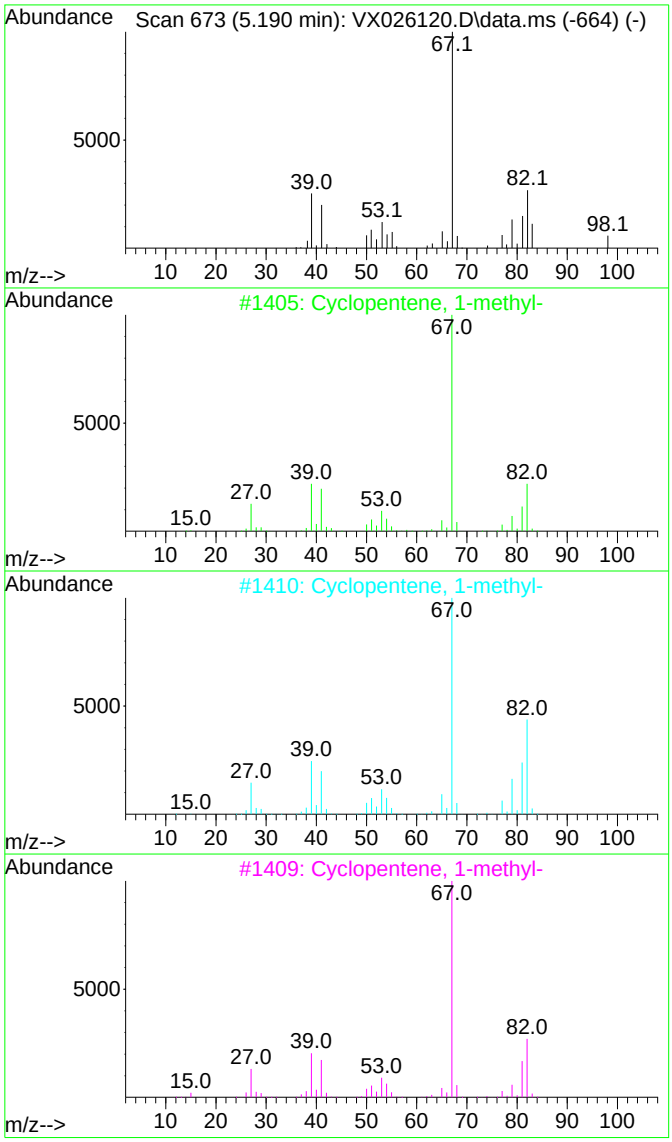
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

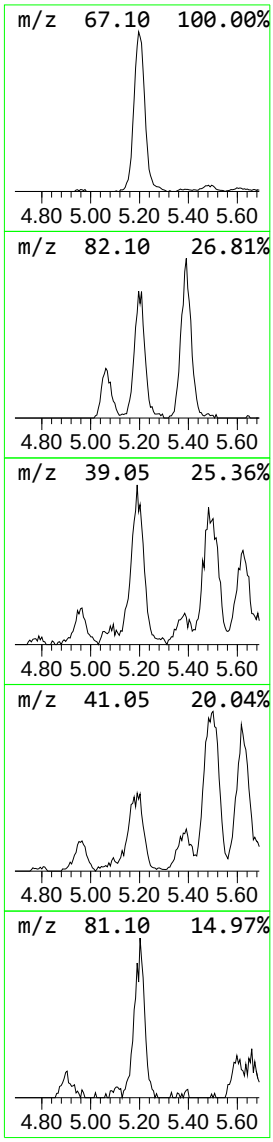
****

Peak Number 2 Cyclopentene, 1-methyl- Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.190	16.17 ug/L	153880	1,4-Difluorobenzene	6.763

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentene, 1-methyl-	82	C6H10	000693-89-0	81
2	Cyclopentene, 1-methyl-	82	C6H10	000693-89-0	81
3	Cyclopentene, 1-methyl-	82	C6H10	000693-89-0	74
4	1,4-Hexadiene	82	C6H10	000592-45-0	74
5	2,4-Hexadiene	82	C6H10	000592-46-1	74





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

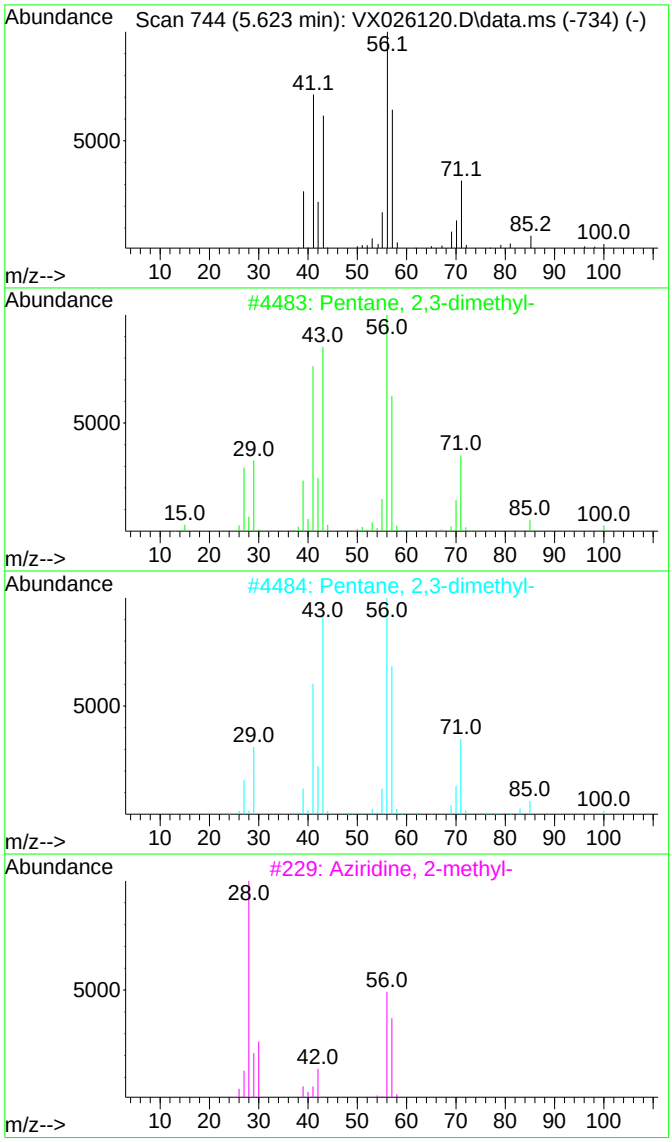
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

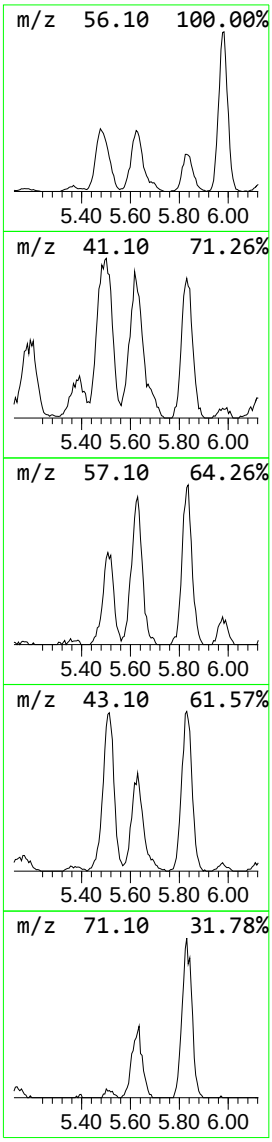
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.623	14.23 ug/L	135416	1,4-Difluorobenzene	6.763

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	91
2	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	91
3	Aziridine, 2-methyl-	57	C3H7N	000075-55-8	53
4	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	52
5	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	52





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

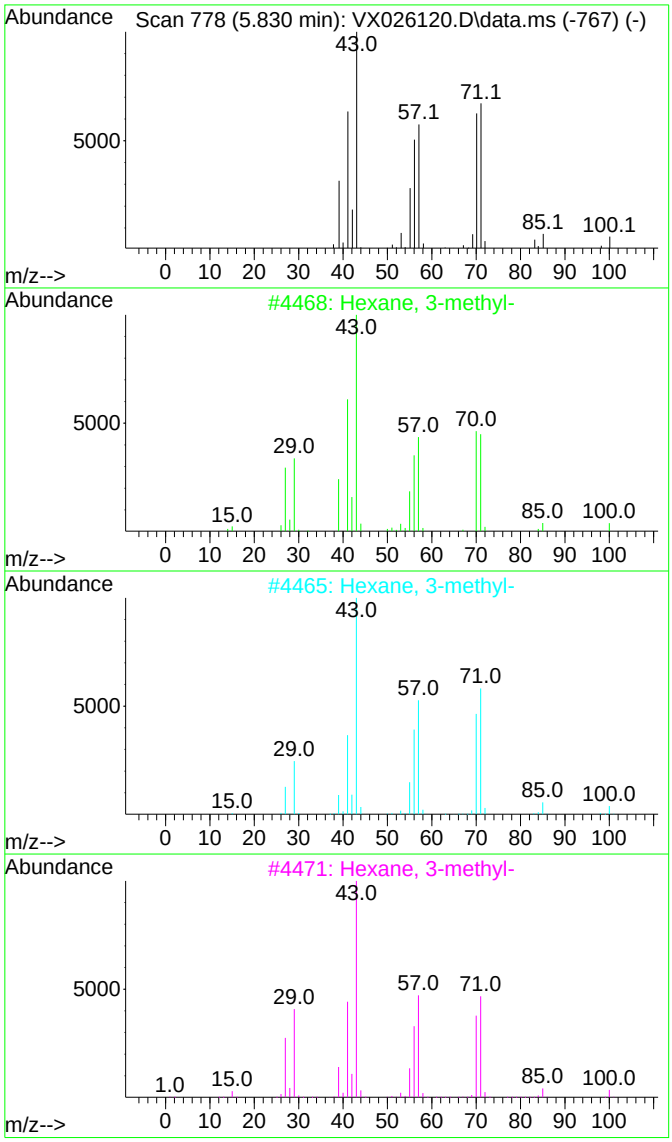
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

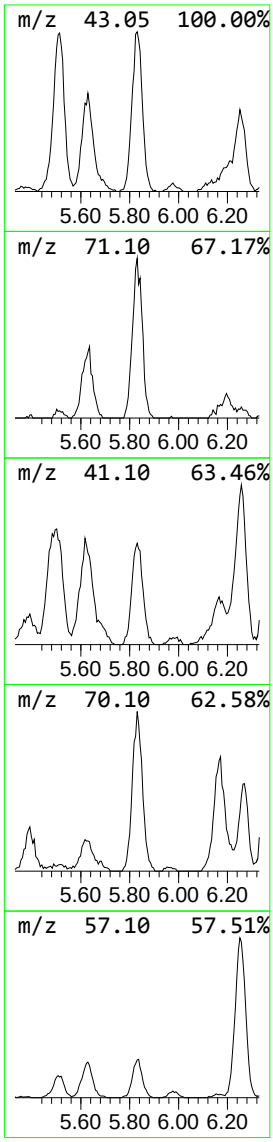
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.830	14.24 ug/L	135520	1,4-Difluorobenzene	6.763

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 3-methyl-	100	C7H16	000589-34-4	94
2	Hexane, 3-methyl-	100	C7H16	000589-34-4	91
3	Hexane, 3-methyl-	100	C7H16	000589-34-4	87
4	Heptane, 3,4-dimethyl-	128	C9H20	000922-28-1	72
5	Octane, 4,5-dimethyl-	142	C10H22	015869-96-2	59





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

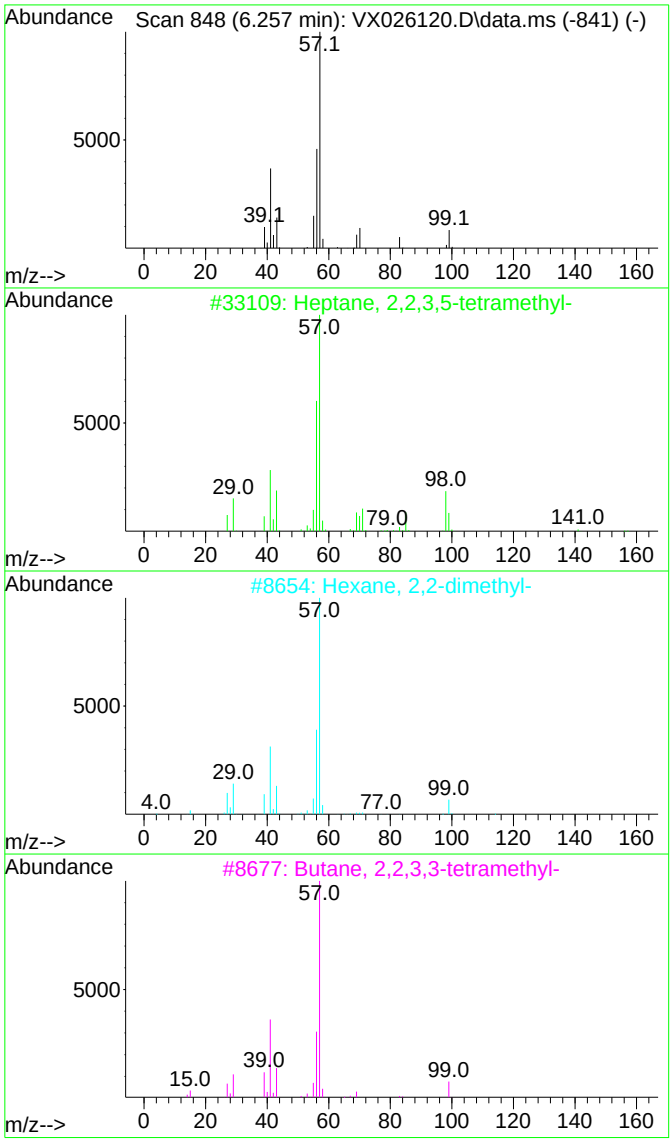
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

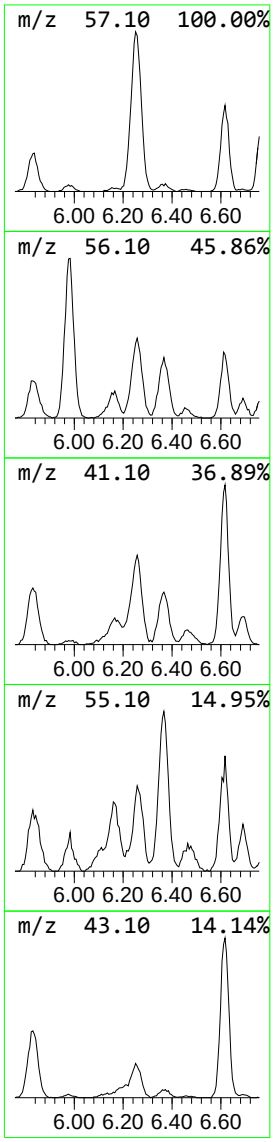
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.257	18.30 ug/L	174173	1,4-Difluorobenzene	6.763

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 2,2,3,5-tetramethyl-	156	C11H24	061868-42-6	74
2	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	59
3	Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	50
4	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	45
5	Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	42





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

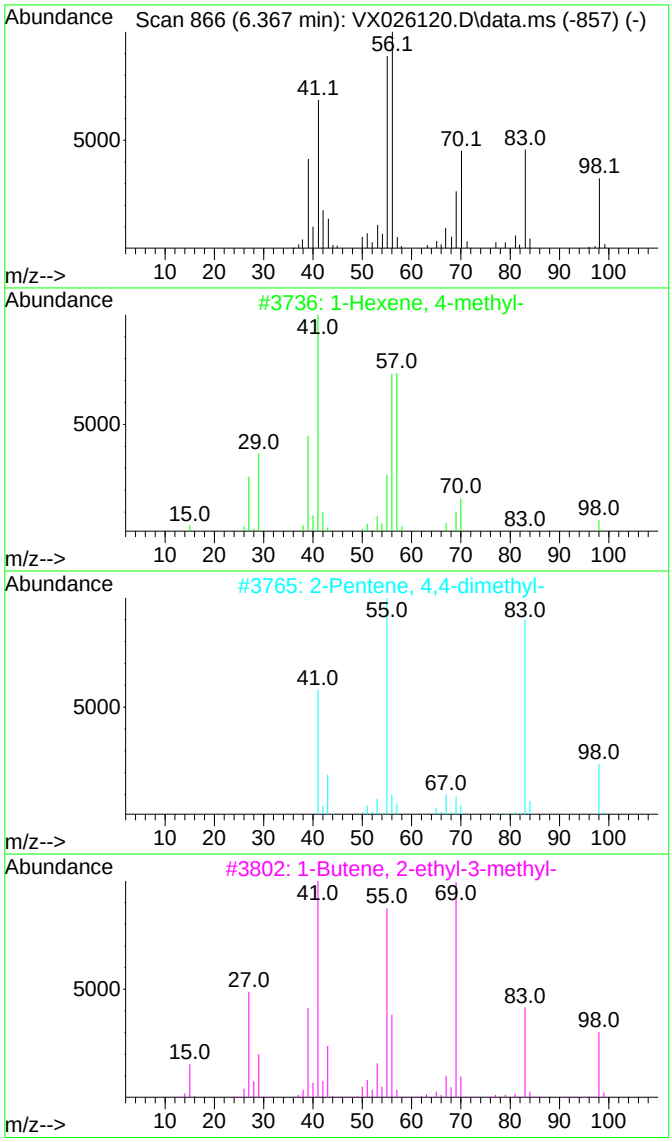
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

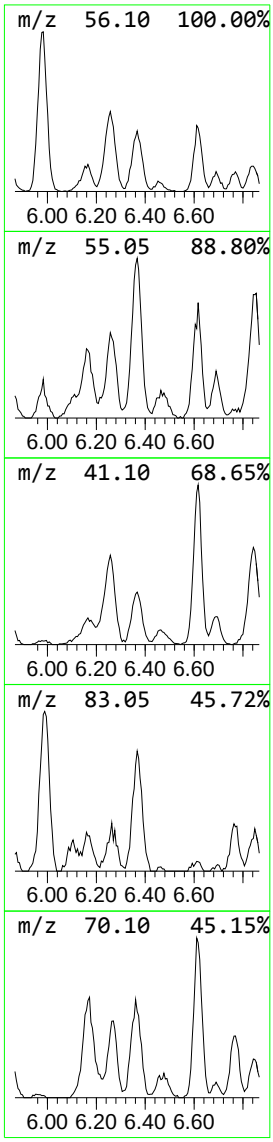
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 42

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.367	11.46 ug/L	109086	1,4-Difluorobenzene	6.763

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexene, 4-methyl-	98	C7H14	003769-23-1	68
2	2-Pentene, 4,4-dimethyl-	98	C7H14	026232-98-4	50
3	1-Butene, 2-ethyl-3-methyl-	98	C7H14	007357-93-9	46
4	(Z)-2-Heptene	98	C7H14	006443-92-1	43
5	Cyclopropane, 1,1,2,2-tetramethyl-	98	C7H14	004127-47-3	38





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

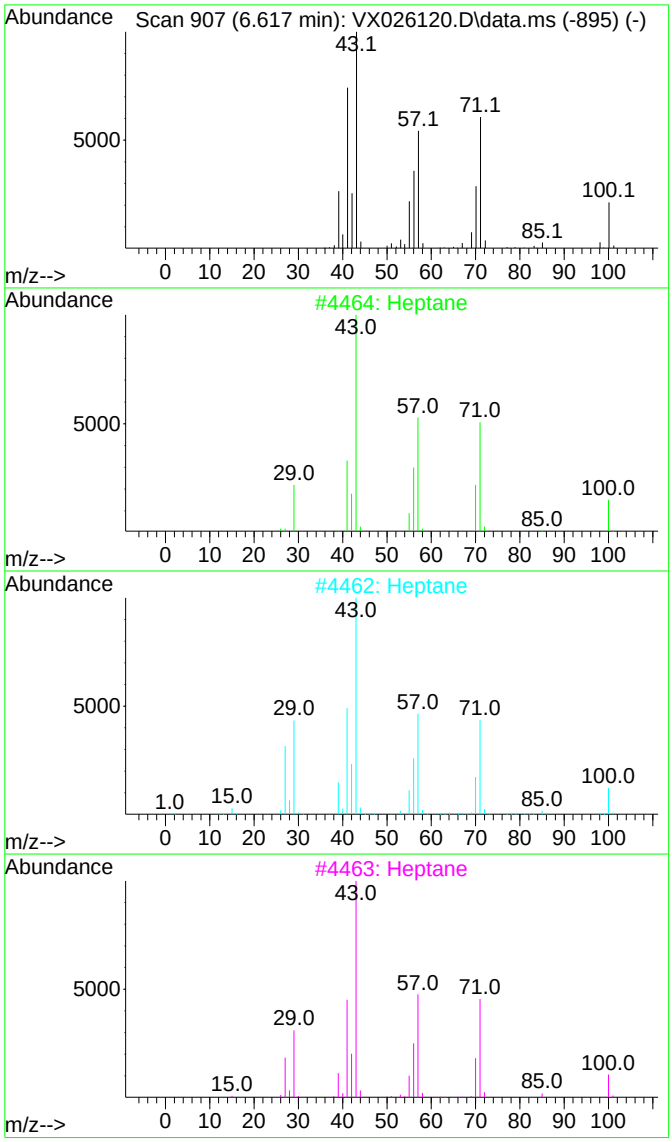
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

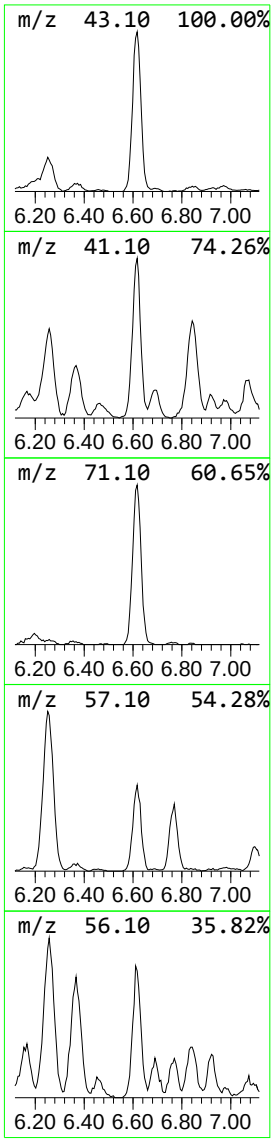
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.617	25.61 ug/L	243654	1,4-Difluorobenzene	6.763

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane	100	C7H16	000142-82-5	90
2	Heptane	100	C7H16	000142-82-5	68
3	Heptane	100	C7H16	000142-82-5	64
4	Heptane	100	C7H16	000142-82-5	64
5	Hexane, 3-methyl-	100	C7H16	000589-34-4	47





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

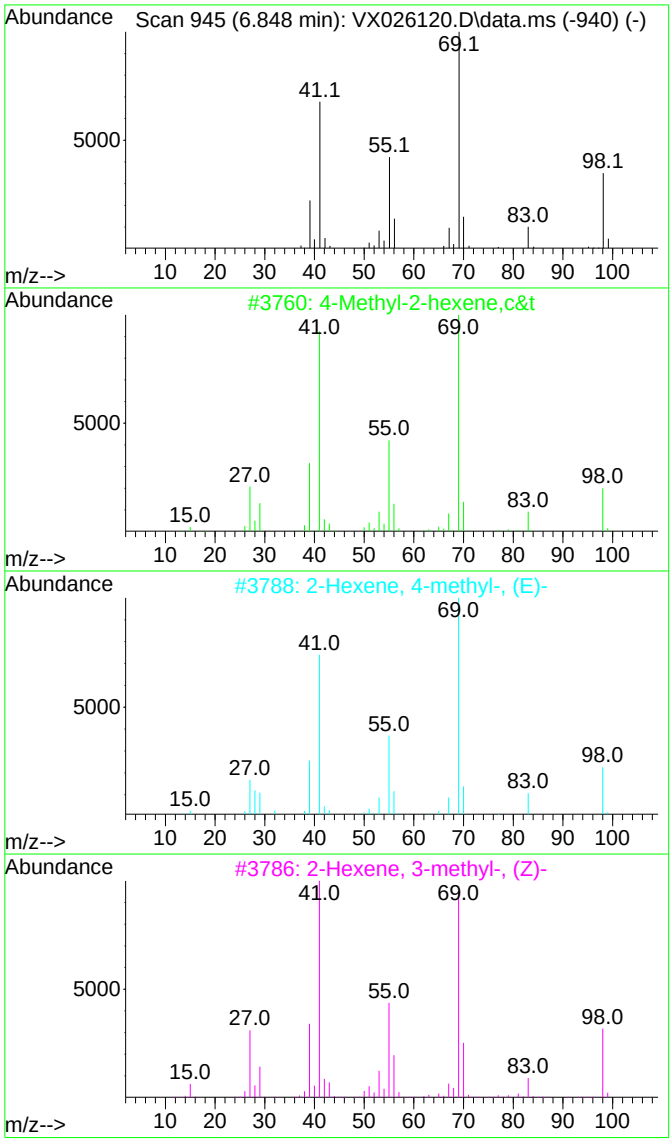
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

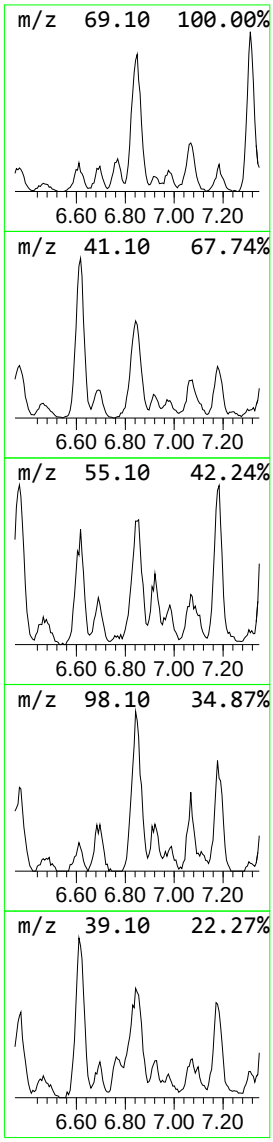
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.848	14.22 ug/L	135346	1,4-Difluorobenzene	6.763

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Methyl-2-hexene,c&t	98	C7H14	003404-55-5	91
2	2-Hexene, 4-methyl-, (E)-	98	C7H14	003683-22-5	91
3	2-Hexene, 3-methyl-, (Z)-	98	C7H14	010574-36-4	86
4	3-Hexene, 3-methyl-, (Z)-	98	C7H14	004914-89-0	86
5	(Z)-4-Methyl-2-hexene	98	C7H14	003683-19-0	83





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

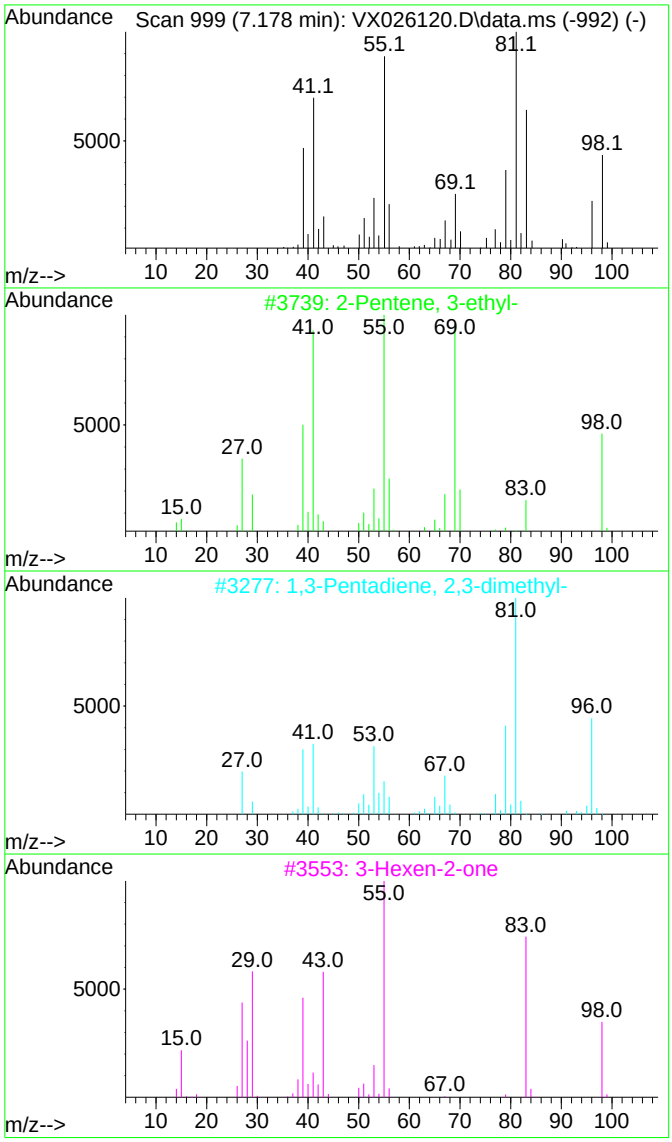
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

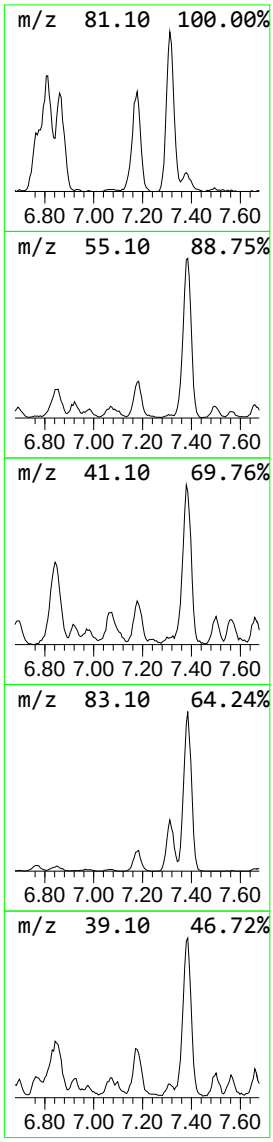
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 46

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.178	11.13 ug/L	105919	1,4-Difluorobenzene	6.763

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentene, 3-ethyl-	98	C7H14	000816-79-5	46
2	1,3-Pentadiene, 2,3-dimethyl-	96	C7H12	001113-56-0	46
3	3-Hexen-2-one	98	C6H10O	000763-93-9	43
4	2,3-Dimethyl-1,4-pentadiene	96	C7H12	000758-86-1	43
5	2-Pentyne, 4,4-dimethyl-	96	C7H12	000999-78-0	38





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

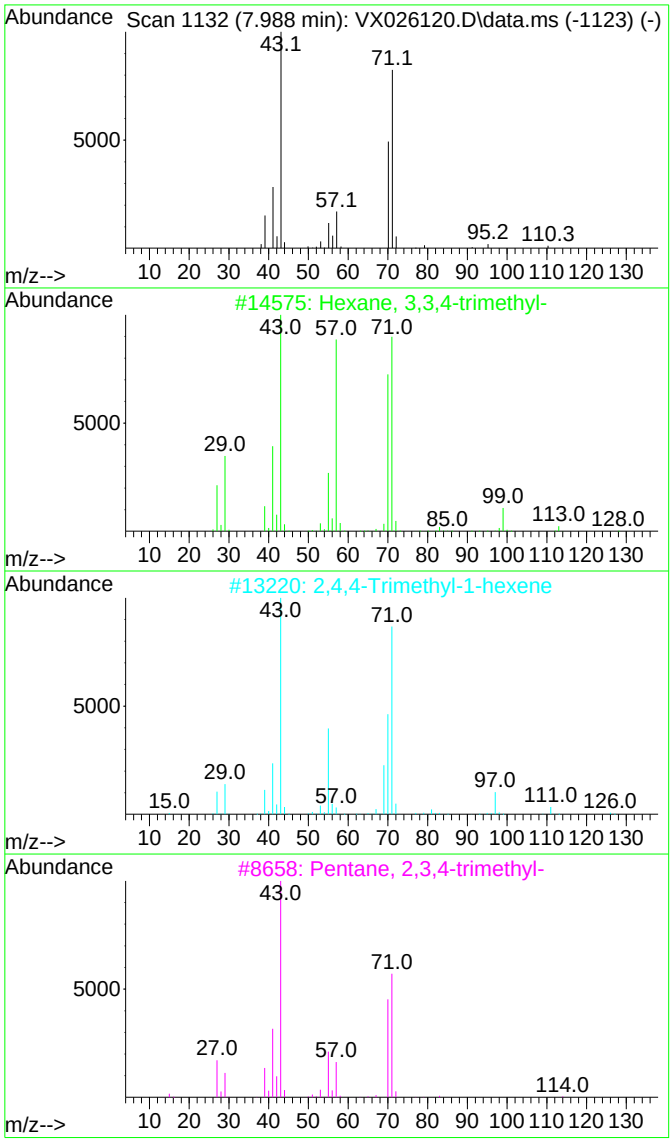
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

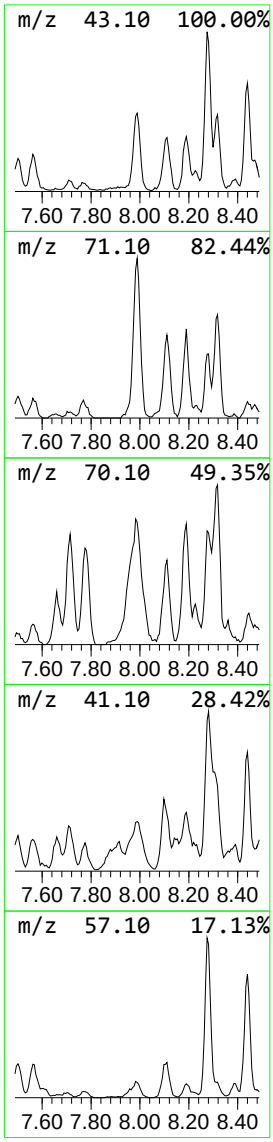
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.988	15.14 ug/L	144035	1,4-Difluorobenzene	6.763

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	78
2	2,4,4-Trimethyl-1-hexene	126	C9H18	051174-12-0	78
3	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	64
4	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	64
5	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	56





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

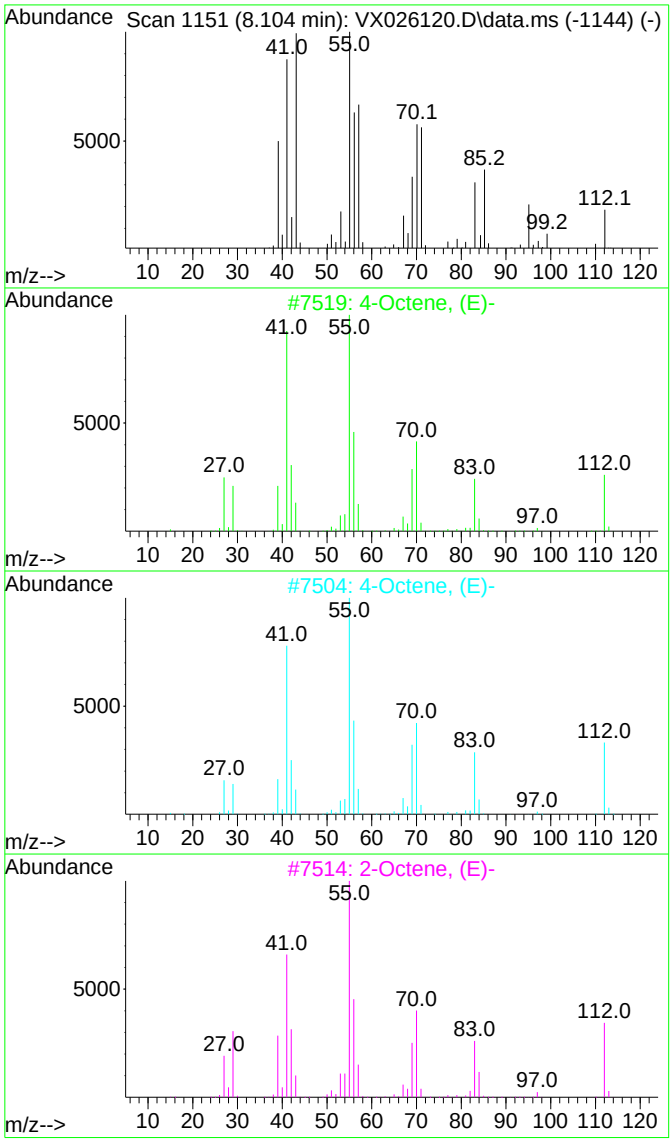
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

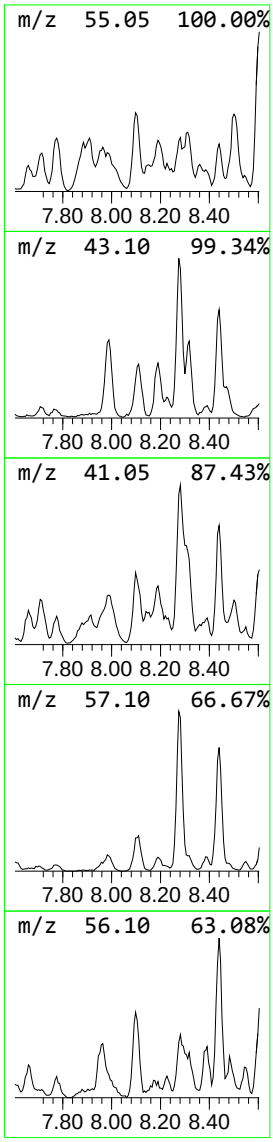
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 43

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.104	11.43 ug/L	108718	1,4-Difluorobenzene	6.763

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Octene, (E)-	112	C8H16	014850-23-8	45
2	4-Octene, (E)-	112	C8H16	014850-23-8	45
3	2-Octene, (E)-	112	C8H16	013389-42-9	43
4	4-Octene, (Z)-	112	C8H16	007642-15-1	43
5	1-Octene	112	C8H16	000111-66-0	43





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

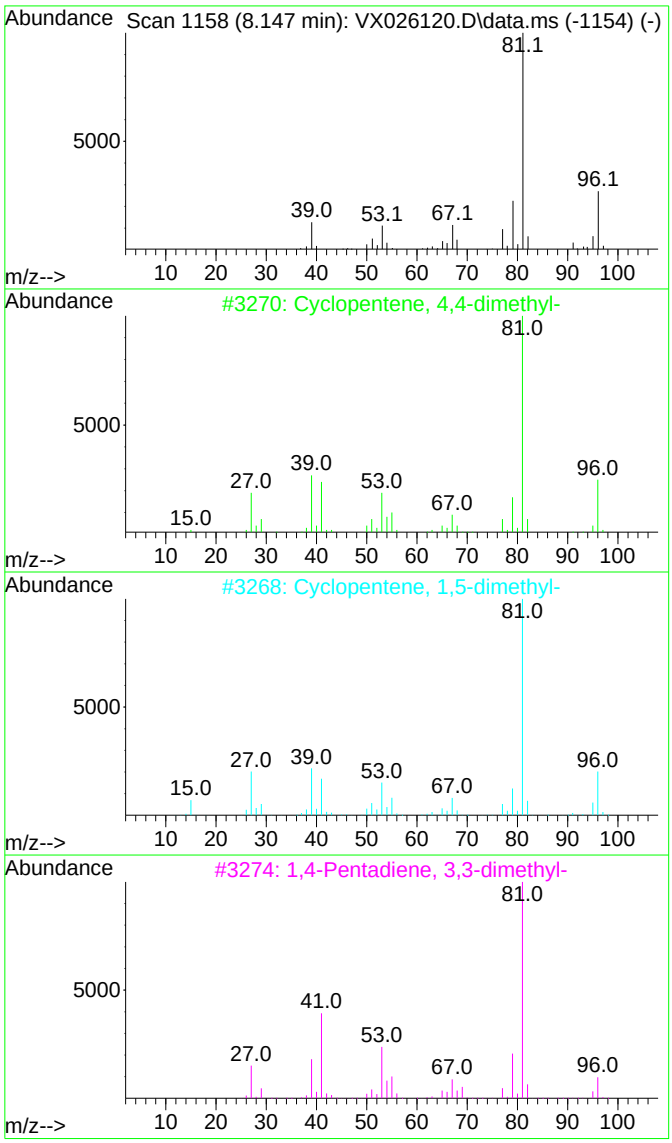
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

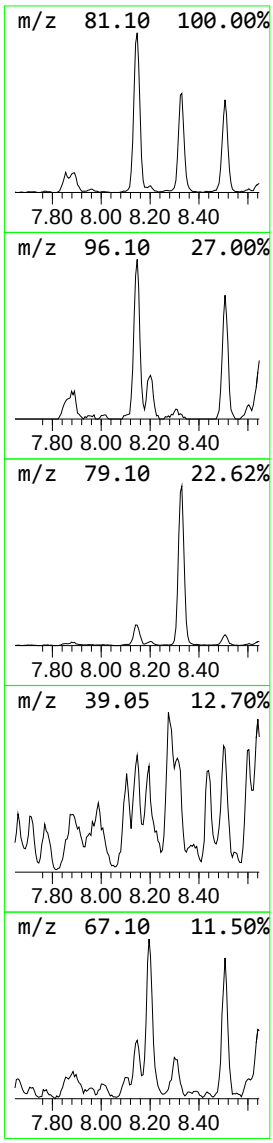
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Cyclopentene, 4,4-dimethyl- Concentration Rank 50

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.147	10.58 ug/L	100716	1,4-Difluorobenzene	6.763

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentene, 4,4-dimethyl-	96	C7H12	019037-72-0	83
2	Cyclopentene, 1,5-dimethyl-	96	C7H12	016491-15-9	80
3	1,4-Pentadiene, 3,3-dimethyl-	96	C7H12	001112-35-2	72
4	Cyclobutane, (1-methylethylidene)-	96	C7H12	001528-22-9	72
5	3,5-Dimethylcyclopentene	96	C7H12	007459-71-4	72





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

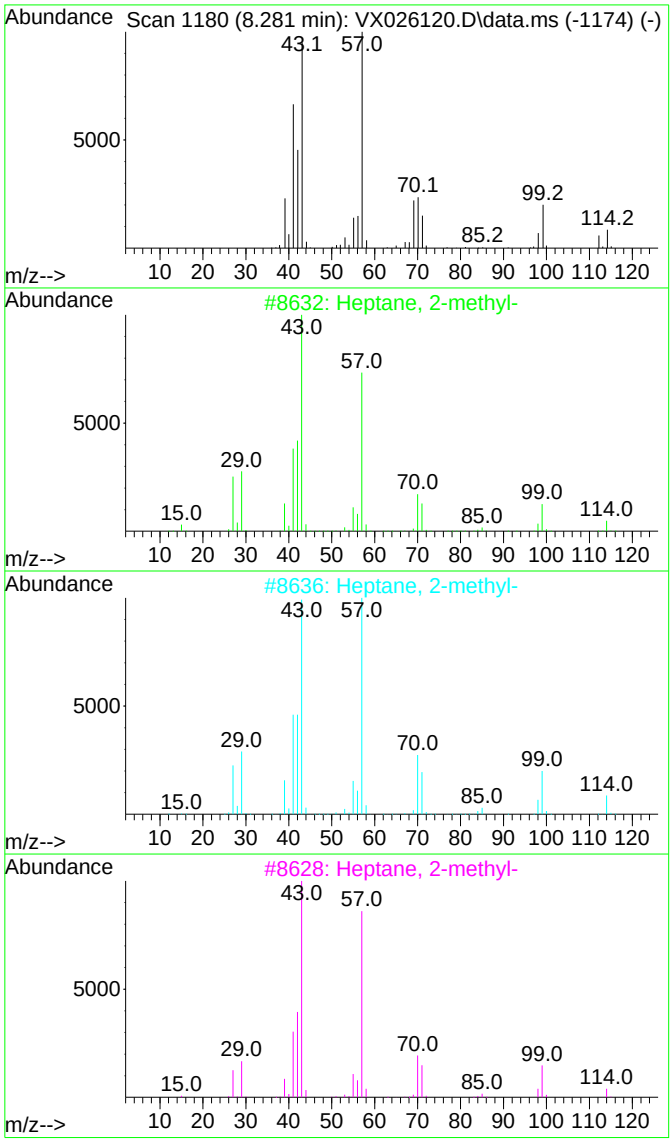
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

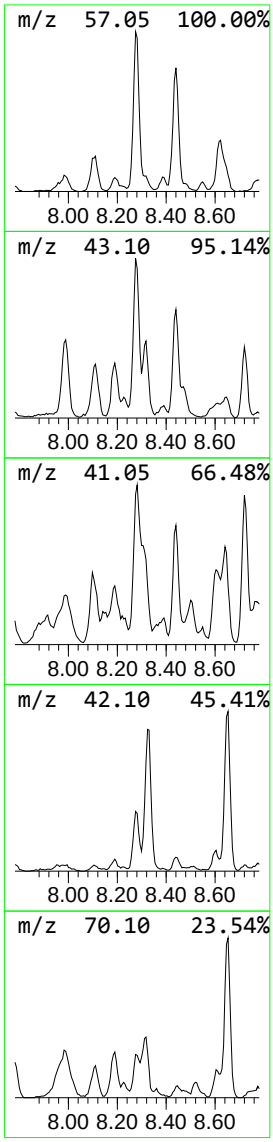
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.281	18.56 ug/L	176557	1,4-Difluorobenzene	6.763

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 2-methyl-	114	C8H18	000592-27-8	91
2	Heptane, 2-methyl-	114	C8H18	000592-27-8	91
3	Heptane, 2-methyl-	114	C8H18	000592-27-8	83
4	Heptane, 2,3,6-trimethyl-	142	C10H22	004032-93-3	43
5	Hexane, 2,5-dimethyl-	114	C8H18	000592-13-2	43





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

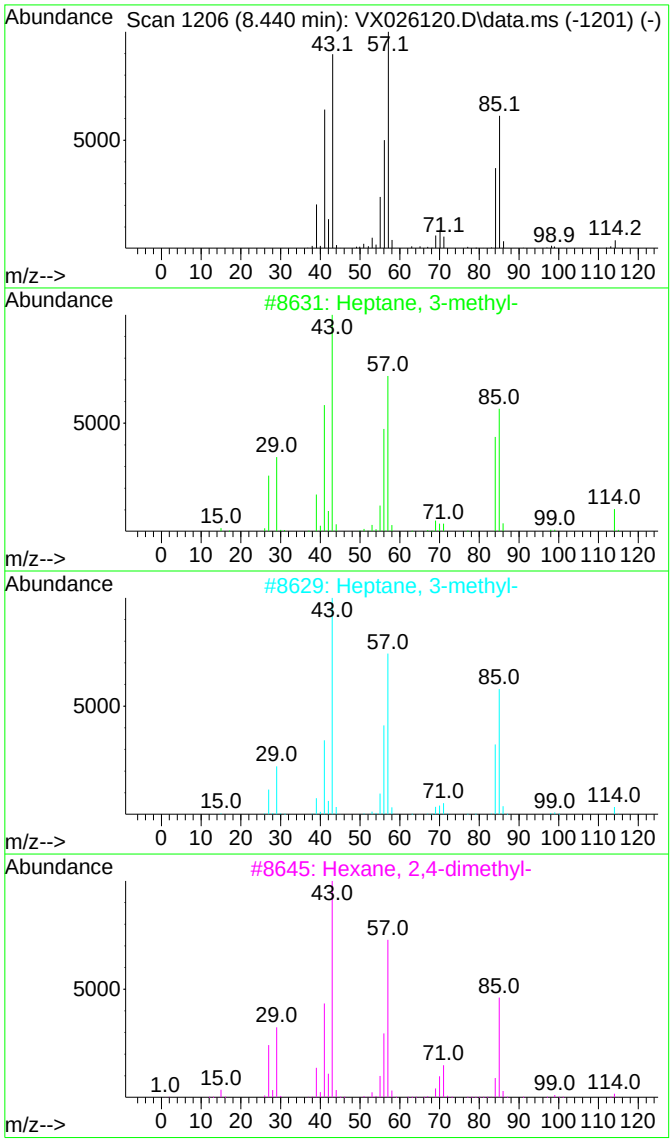
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

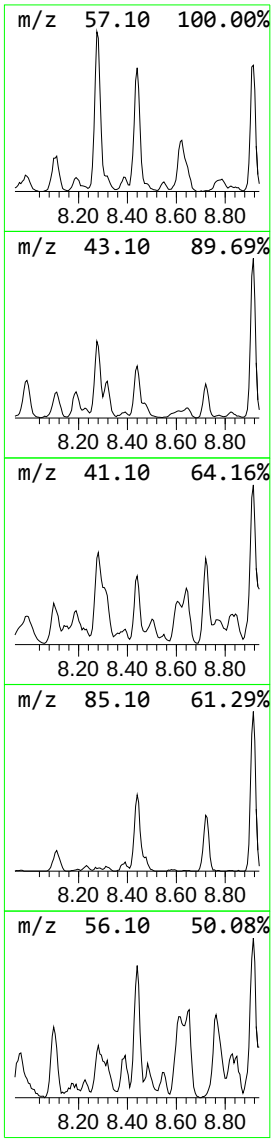
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 44

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.440	11.38 ug/L	125845	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 3-methyl-	114	C8H18	000589-81-1	78
2	Heptane, 3-methyl-	114	C8H18	000589-81-1	74
3	Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	64
4	Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	64
5	Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	59





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

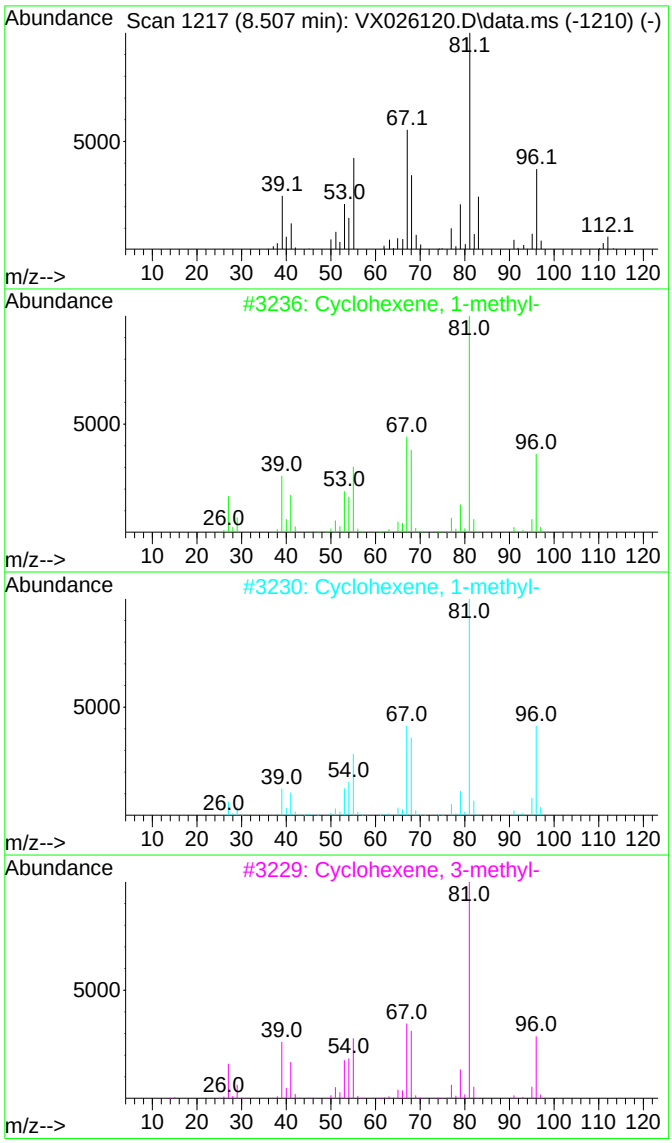
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

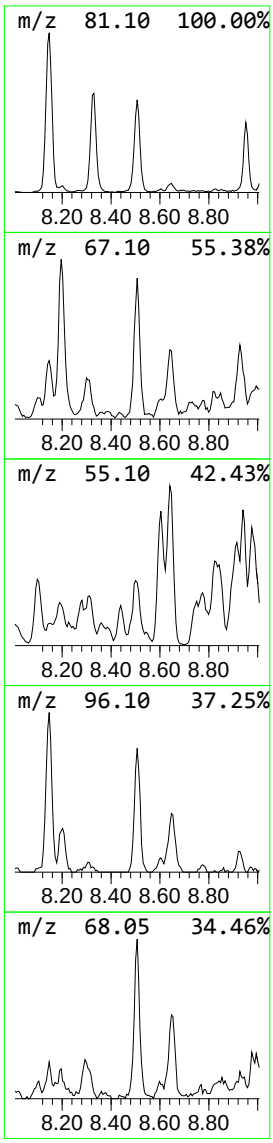
****

Peak Number 16 Cyclohexene, 1-methyl- Concentration Rank 48

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.507	10.80 ug/L	119421	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexene, 1-methyl-	96	C7H12	000591-49-1	94
2	Cyclohexene, 1-methyl-	96	C7H12	000591-49-1	93
3	Cyclohexene, 3-methyl-	96	C7H12	000591-48-0	76
4	Cyclohexene, 4-methyl-	96	C7H12	000591-47-9	74
5	2,4-Hexadiene, 3-methyl-	96	C7H12	028823-42-9	70





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

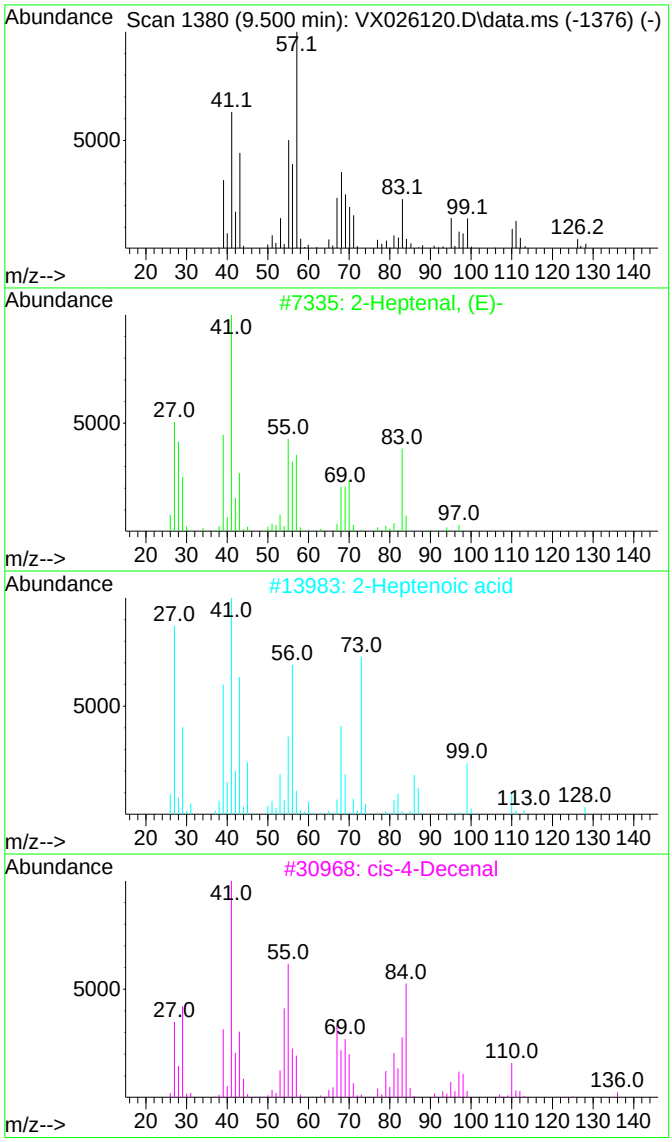
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

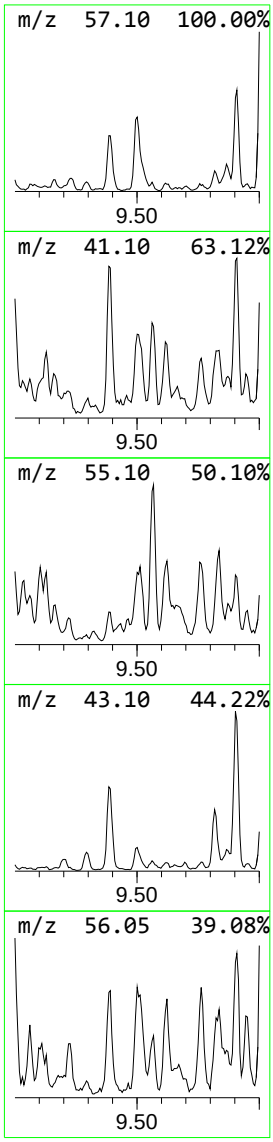
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 unknown-01 Concentration Rank 41

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.500	12.11 ug/L	133899	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Heptenal, (E)-	112	C7H12O	018829-55-5	38
2	2-Heptenoic acid	128	C7H12O2	018999-28-5	38
3	cis-4-Decenal	154	C10H18O	021662-09-9	35
4	1,10-Decanediol	174	C10H22O2	000112-47-0	35
5	2-Octen-1-ol, (E)-	128	C8H16O	018409-17-1	35





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

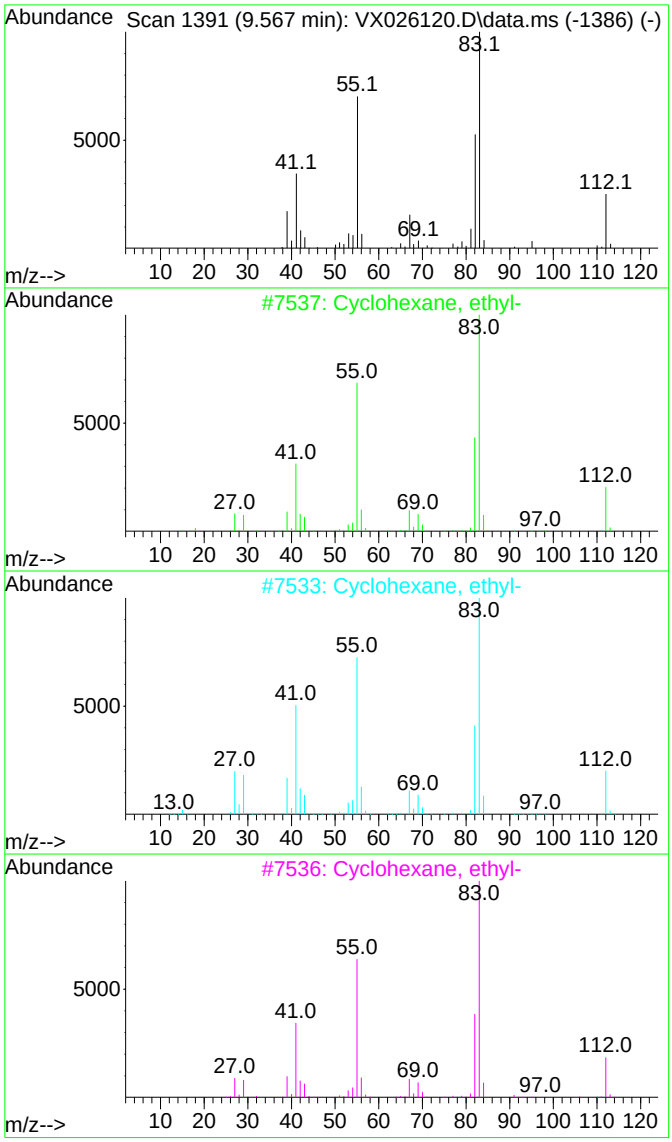
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

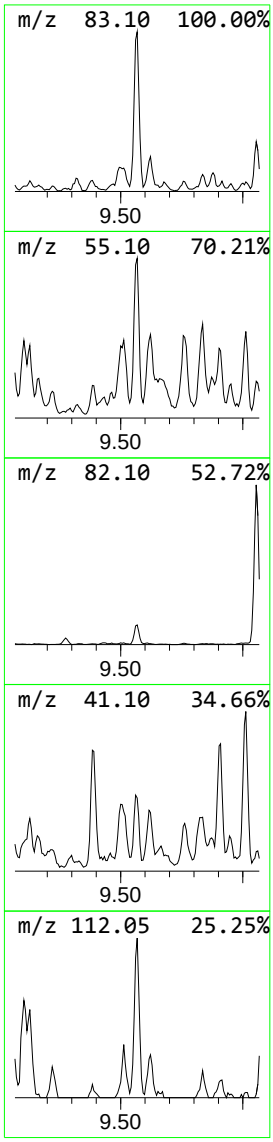
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 (DEL) Alkane: Cyclic9.567 Concentration Rank 58

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.567	9.16 ug/L	101260	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, ethyl-	112	C8H16	001678-91-7	86
2	Cyclohexane, ethyl-	112	C8H16	001678-91-7	86
3	Cyclohexane, ethyl-	112	C8H16	001678-91-7	86
4	Cyclohexane, ethyl-	112	C8H16	001678-91-7	80
5	Cyclohexane, ethyl-	112	C8H16	001678-91-7	80





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

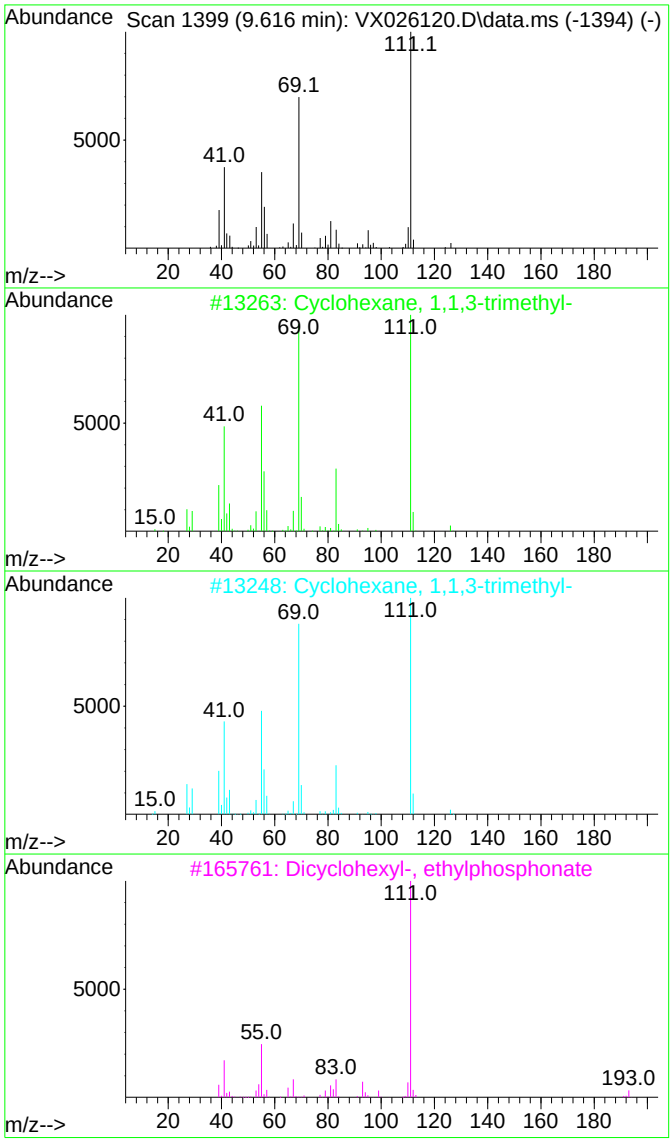
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

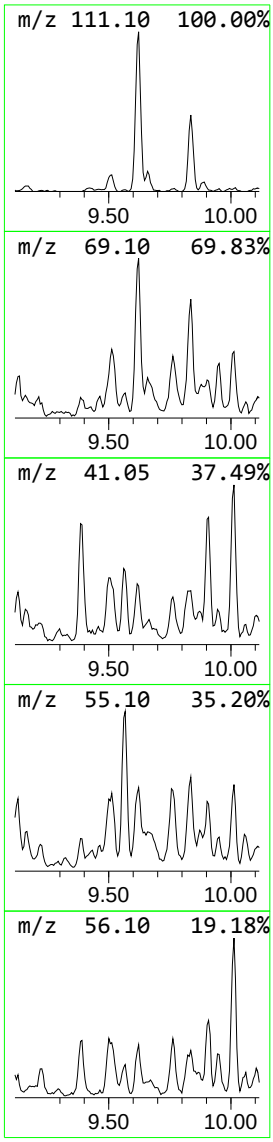
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 (DEL) Alkane: Cyclic9.616 Concentration Rank 52

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.616	10.03 ug/L	110890	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	64
2	Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	64
3	Dicyclohexyl-, ethylphosphonate	274	C14H27O3P	169739-47-3	53
4	1,1,4-Trimethylcyclohexane	126	C9H18	007094-27-1	53
5	Methoxyacetic acid, 2-ethylcyclo...	200	C11H20O3	1000282-69-7	50





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

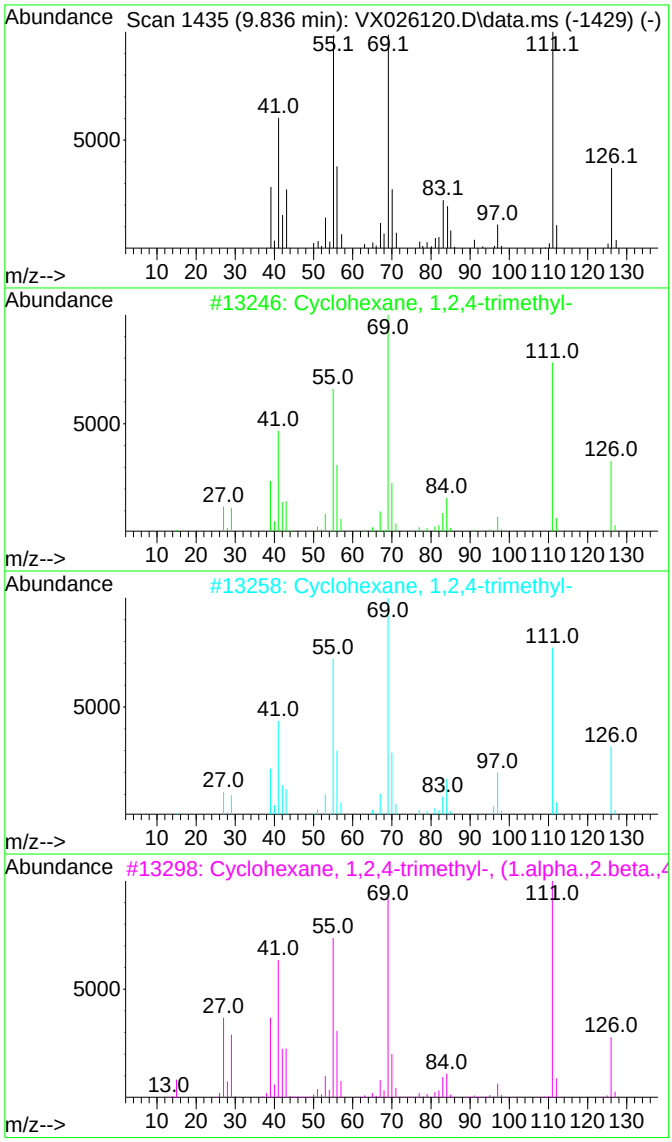
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

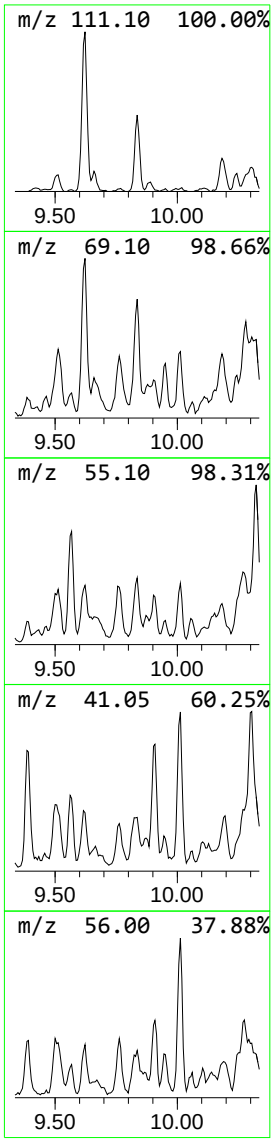
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 (DEL) Alkane: Cyclic9.836 Concentration Rank 55

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.836	9.57 ug/L	105807	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,2,4-trimethyl-	126	C9H18	002234-75-5	93
2	Cyclohexane, 1,2,4-trimethyl-	126	C9H18	002234-75-5	93
3	Cyclohexane, 1,2,4-trimethyl-, (...	126	C9H18	007667-60-9	93
4	Cyclohexane, 1,1,2-trimethyl-	126	C9H18	007094-26-0	83
5	Cyclohexane, 1,2,4-trimethyl-, (...	126	C9H18	007667-60-9	81





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

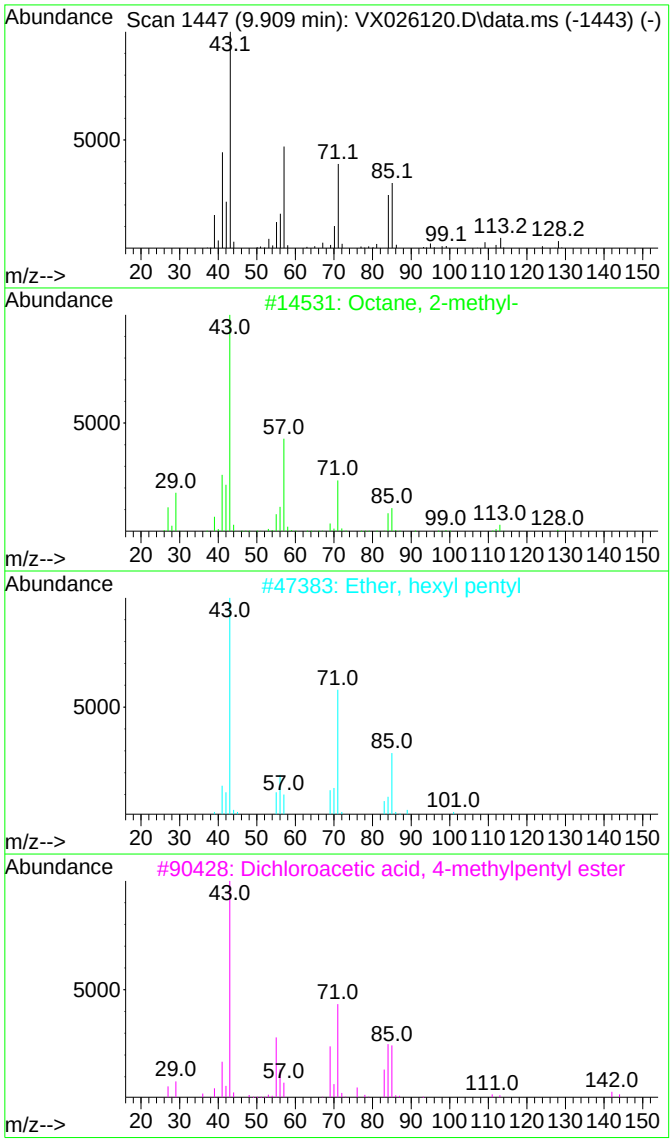
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

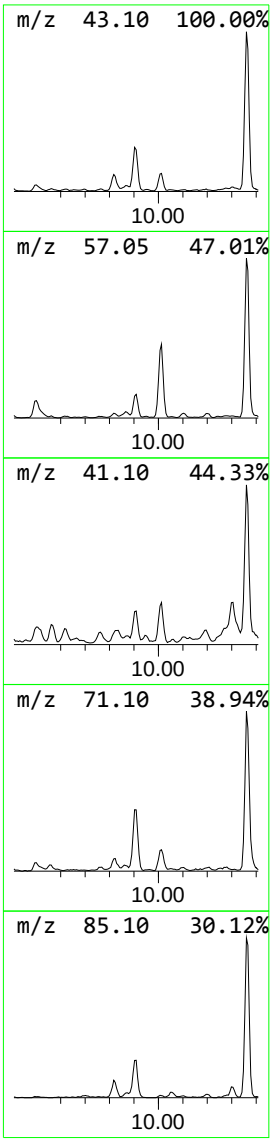
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.909	14.44 ug/L	159650	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2-methyl-	128	C9H20	003221-61-2	72
2	Ether, hexyl pentyl	172	C11H24O	032357-83-8	64
3	Dichloroacetic acid, 4-methylpen...	212	C8H14Cl2O2	1000282-43-7	59
4	Sulfurous acid, 2-propyl tetrade...	320	C17H36O3S	1010309-12-5	53
5	Pentane, 3-ethyl-2-methyl-	114	C8H18	000609-26-7	53





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

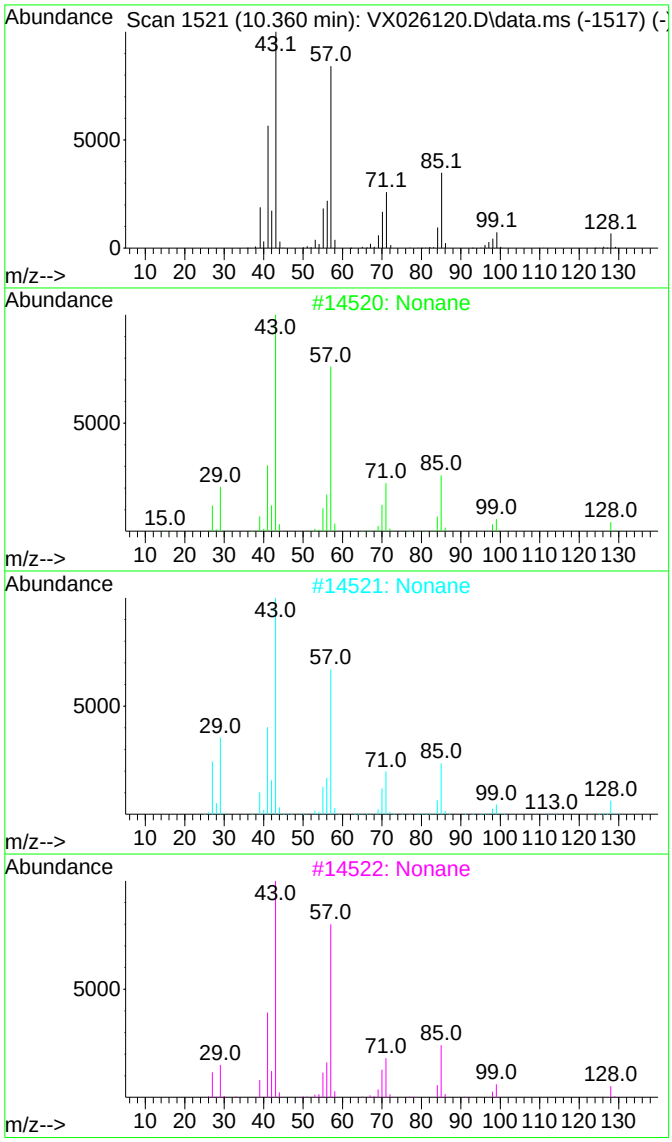
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

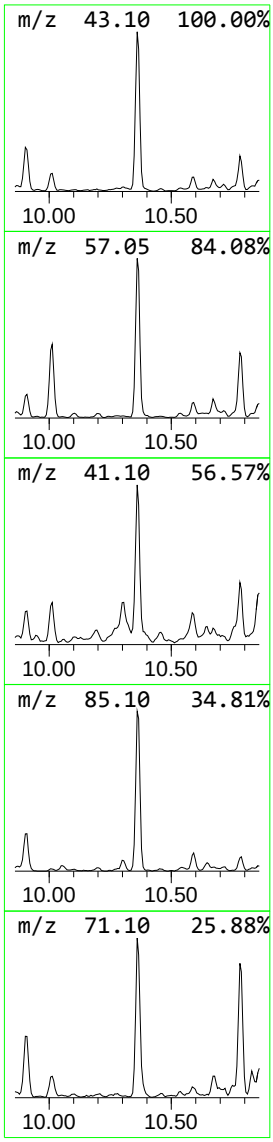
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.360	63.44 ug/L	701595	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane	128	C9H20	000111-84-2	97
2	Nonane	128	C9H20	000111-84-2	97
3	Nonane	128	C9H20	000111-84-2	94
4	Nonane	128	C9H20	000111-84-2	91
5	Sulfurous acid, hexyl pentadecyl...	376	C21H44O3S	1000309-13-7	72





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

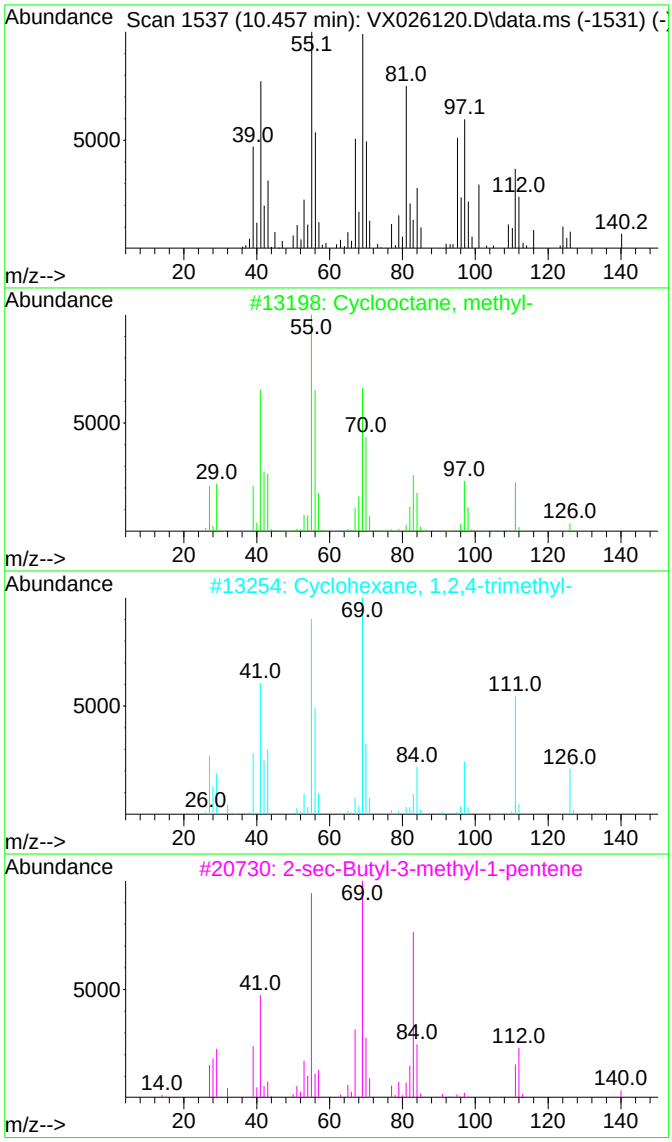
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

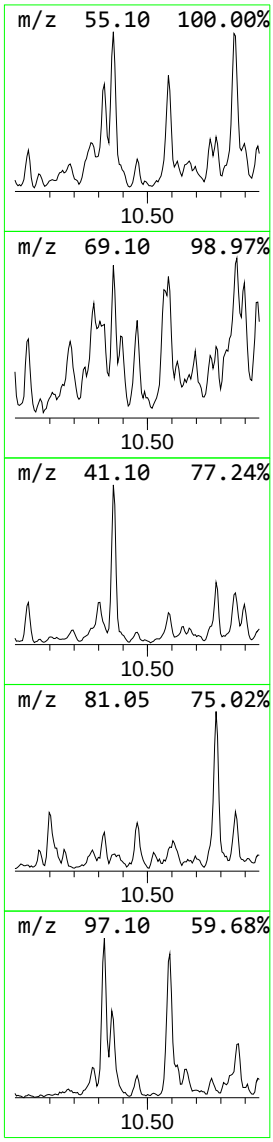
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 (DEL) Alkane: Cyclic10.458 Concentration Rank 59

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.458	8.81 ug/L	97389	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclooctane, methyl-	126	C9H18	001502-38-1	46
2	Cyclohexane, 1,2,4-trimethyl-	126	C9H18	002234-75-5	42
3	2-sec-Butyl-3-methyl-1-pentene	140	C10H20	075144-24-0	38
4	1,4-Hexadiene, 4-methyl-	96	C7H12	001116-90-1	35
5	Cyclooctane, methyl-	126	C9H18	001502-38-1	30





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

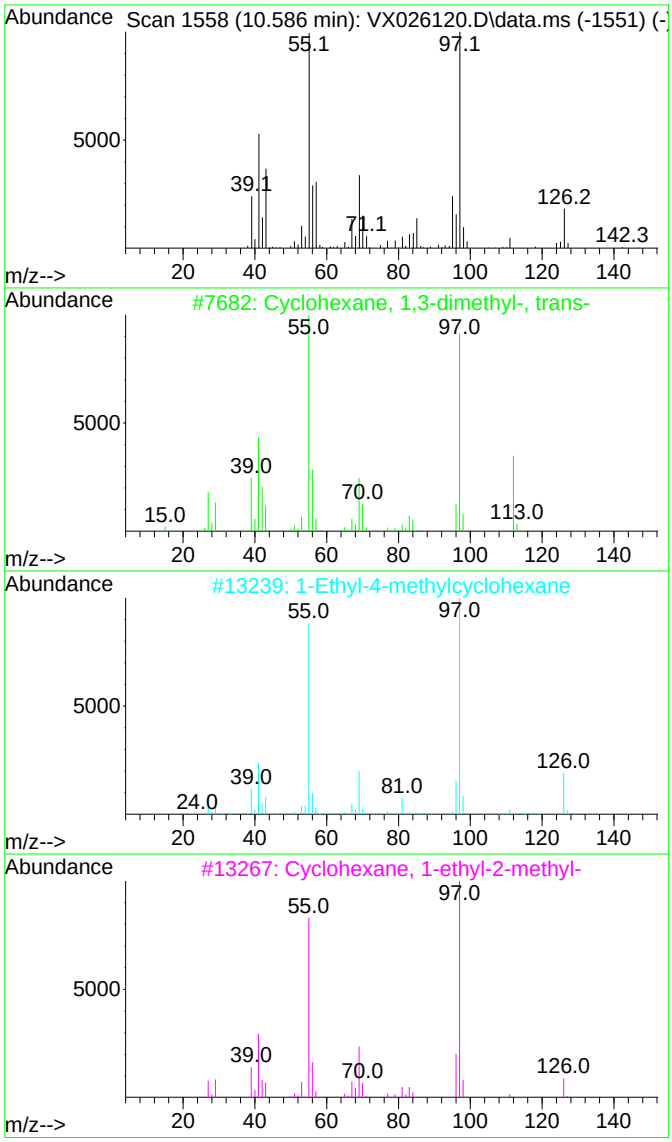
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

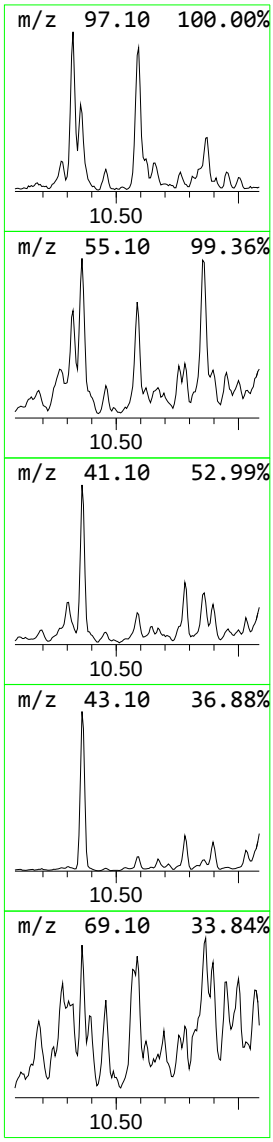
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 25 (DEL) Alkane: Cyclic10.586 Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.586	19.67 ug/L	217543	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,3-dimethyl-, trans-	112	C8H16	002207-03-6	64
2	1-Ethyl-4-methylcyclohexane	126	C9H18	003728-56-1	64
3	Cyclohexane, 1-ethyl-2-methyl-	126	C9H18	003728-54-9	59
4	Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-78-8	58
5	Cyclohexane, 1,3-dimethyl-	112	C8H16	000591-21-9	58





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

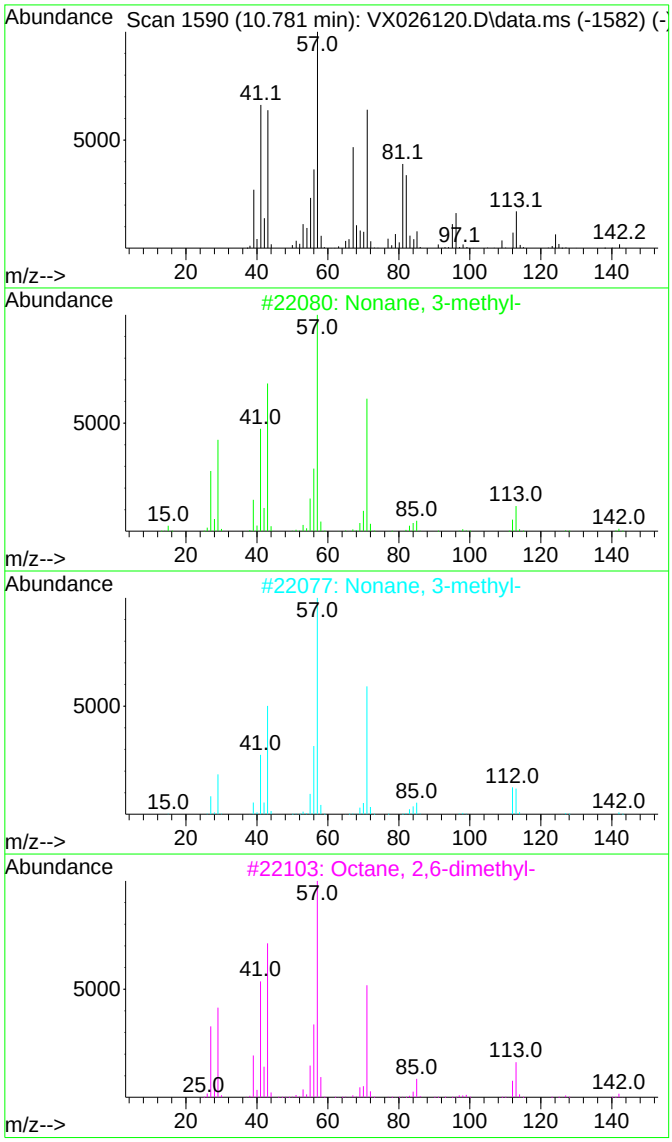
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

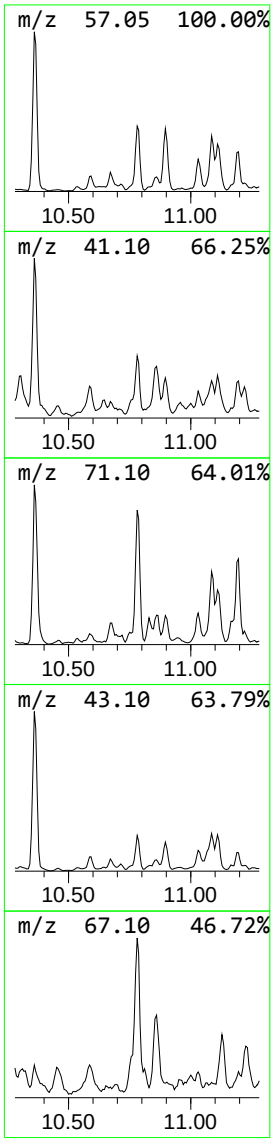
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 26 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.781	35.66 ug/L	394353	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane, 3-methyl-	142	C10H22	005911-04-6	50
2	Nonane, 3-methyl-	142	C10H22	005911-04-6	50
3	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	50
4	Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	46
5	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	43





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

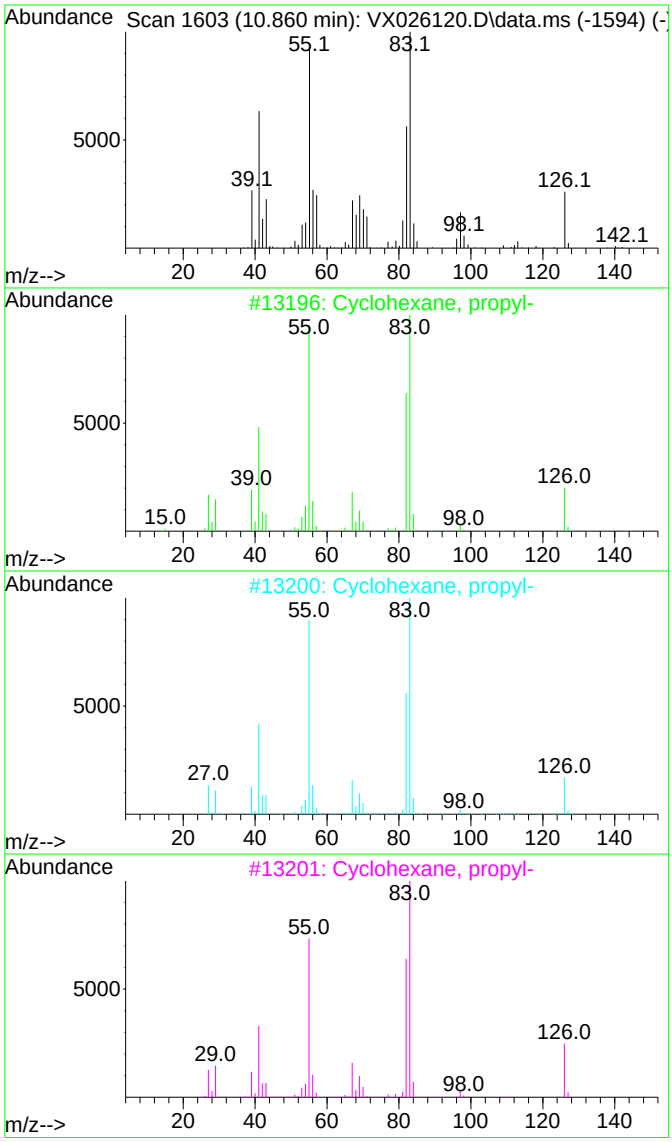
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

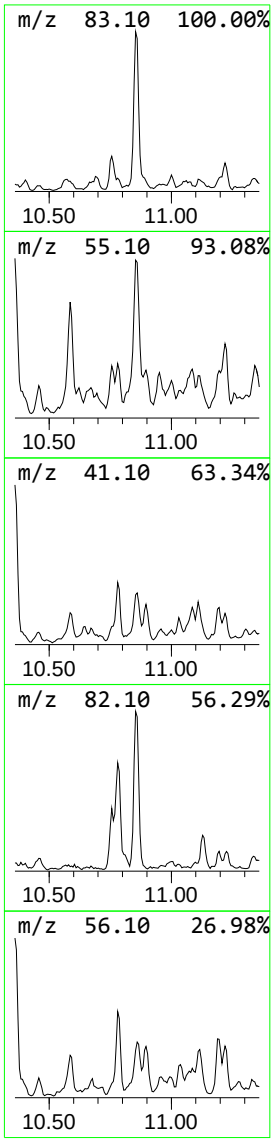
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 27 (DEL) Alkane: Cyclic10.860 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.860	27.62 ug/L	305461	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, propyl-	126	C9H18	001678-92-8	87
2	Cyclohexane, propyl-	126	C9H18	001678-92-8	80
3	Cyclohexane, propyl-	126	C9H18	001678-92-8	72
4	Cyclohexane, butyl-	140	C10H20	001678-93-9	64
5	Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

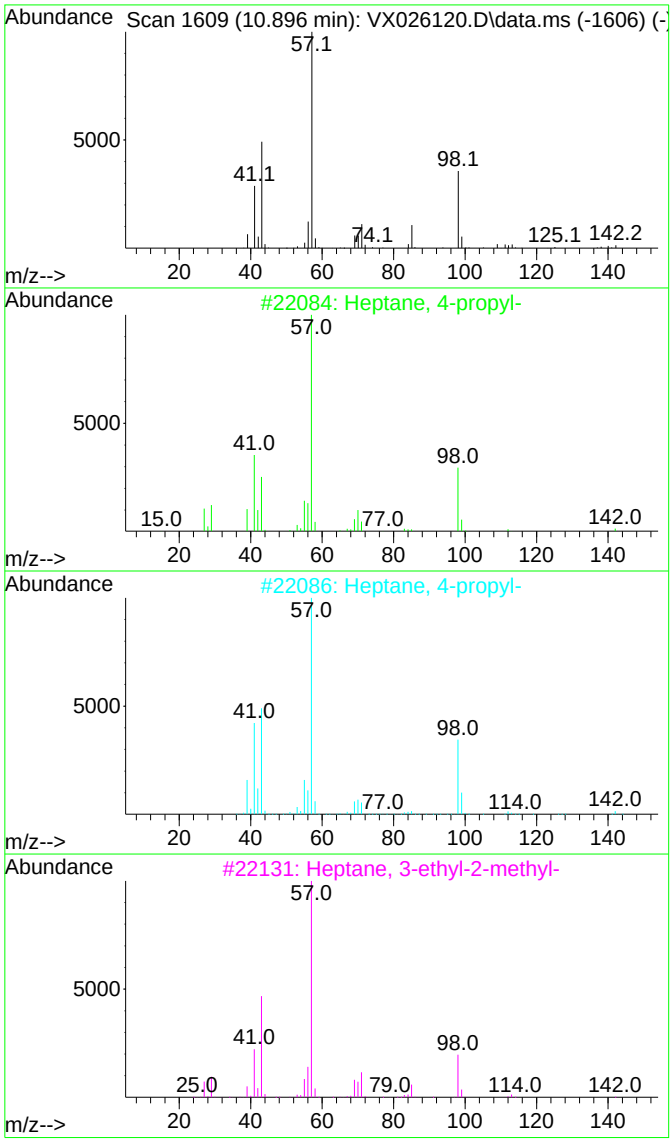
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

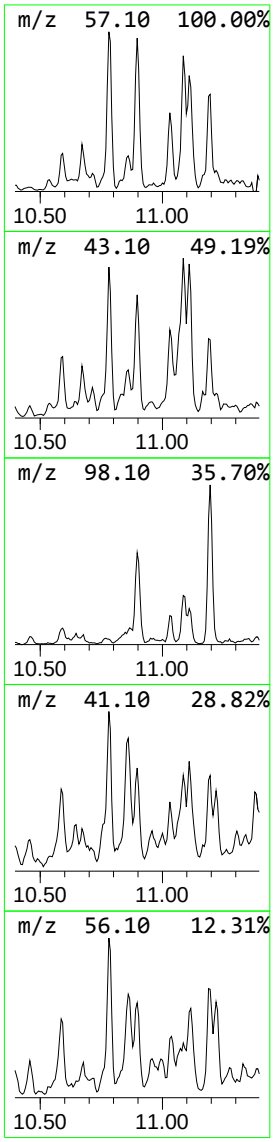
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 28 (DEL) Alkane: Straight-Chai... Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.896	15.77 ug/L	174378	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 4-propyl-	142	C10H22	003178-29-8	64
2	Heptane, 4-propyl-	142	C10H22	003178-29-8	58
3	Heptane, 3-ethyl-2-methyl-	142	C10H22	014676-29-0	53
4	Hexane, 3-ethyl-2,5-dimethyl-	142	C10H22	052897-04-8	50
5	Heptane, 3-ethyl-2-methyl-	142	C10H22	014676-29-0	50





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

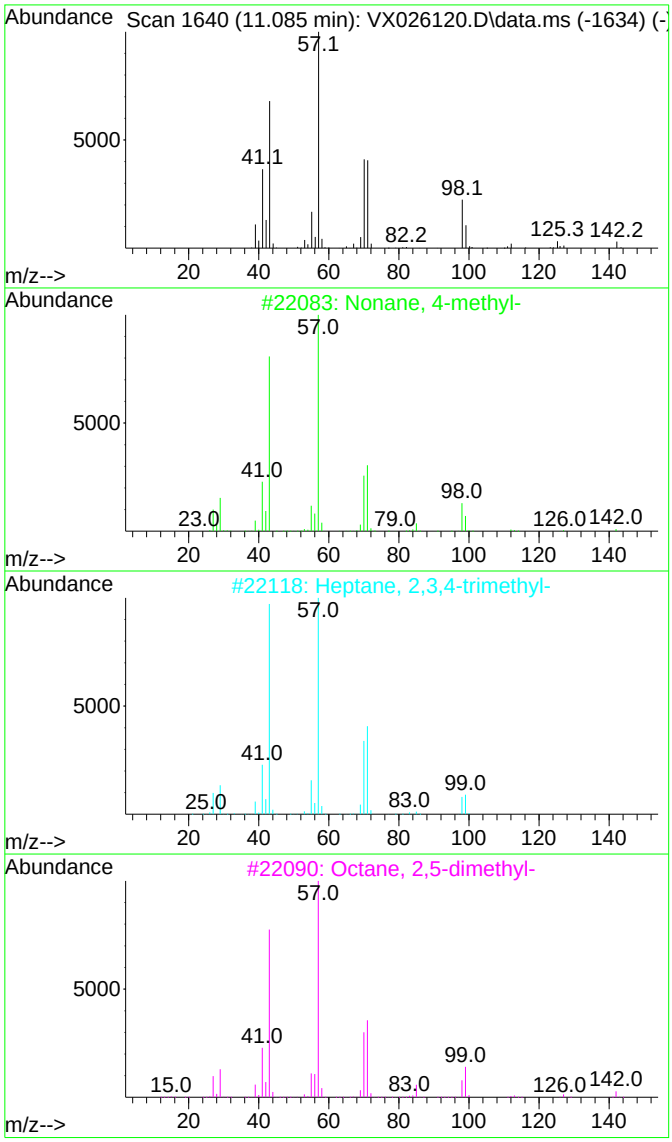
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

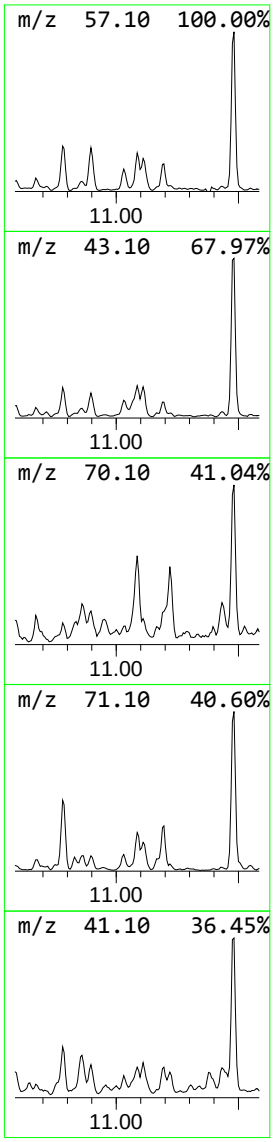
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 29 (DEL) Alkane: Straight-Chai... Concentration Rank 54

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.085	9.58 ug/L	146710	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane, 4-methyl-	142	C10H22	017301-94-9	80
2	Heptane, 2,3,4-trimethyl-	142	C10H22	052896-95-4	64
3	Octane, 2,5-dimethyl-	142	C10H22	015869-89-3	59
4	Nonane, 4-methyl-	142	C10H22	017301-94-9	53
5	Octane, 4-ethyl-	142	C10H22	015869-86-0	53





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

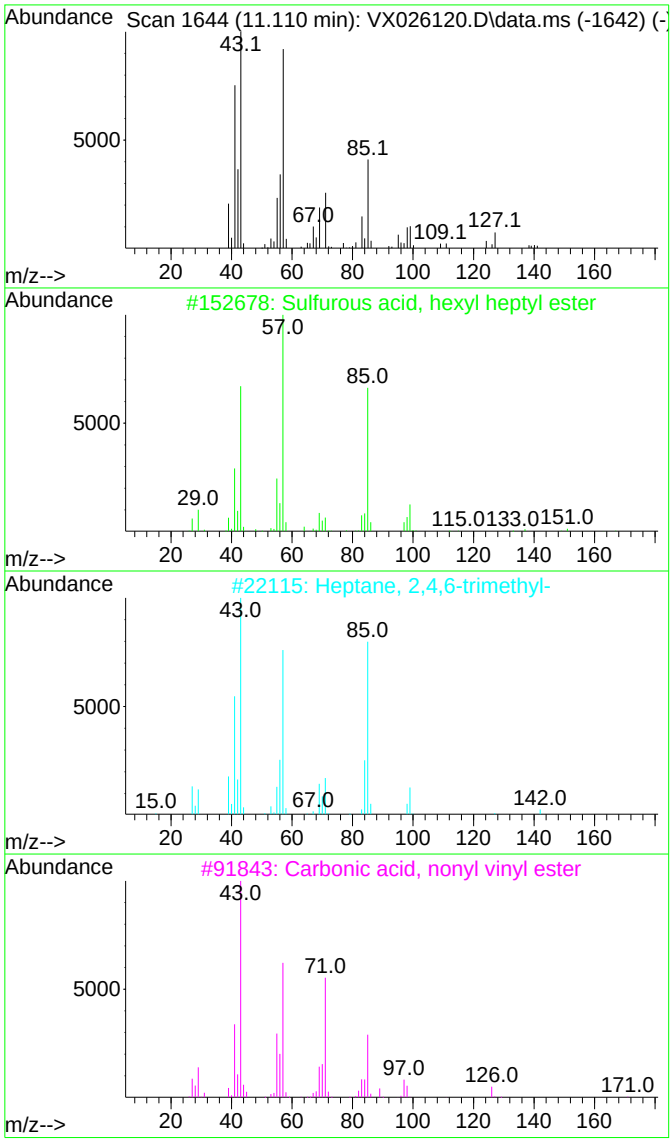
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

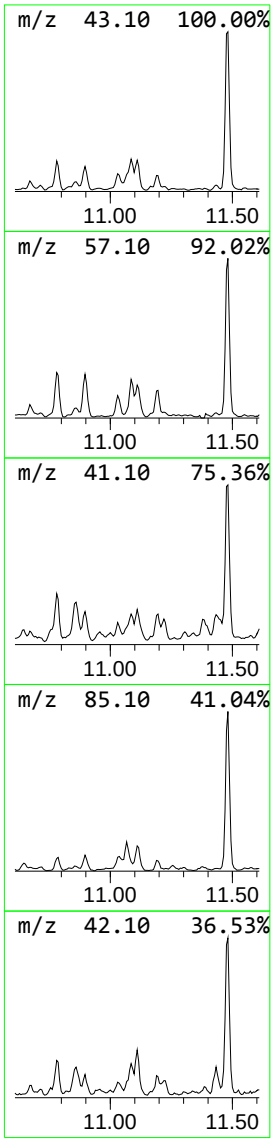
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 30 Sulfurous acid, hexyl hepty... Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.110	13.08 ug/L	200319	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Sulfurous acid, hexyl heptyl ester	264	C13H28O3S	1000309-12-9	50
2	Heptane, 2,4,6-trimethyl-	142	C10H22	002613-61-8	50
3	Carbonic acid, nonyl vinyl ester	214	C12H22O3	1000383-25-6	50
4	Ether, heptyl hexyl	200	C13H28O	007289-40-9	50
5	Nonane, 2-methyl-	142	C10H22	000871-83-0	50





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

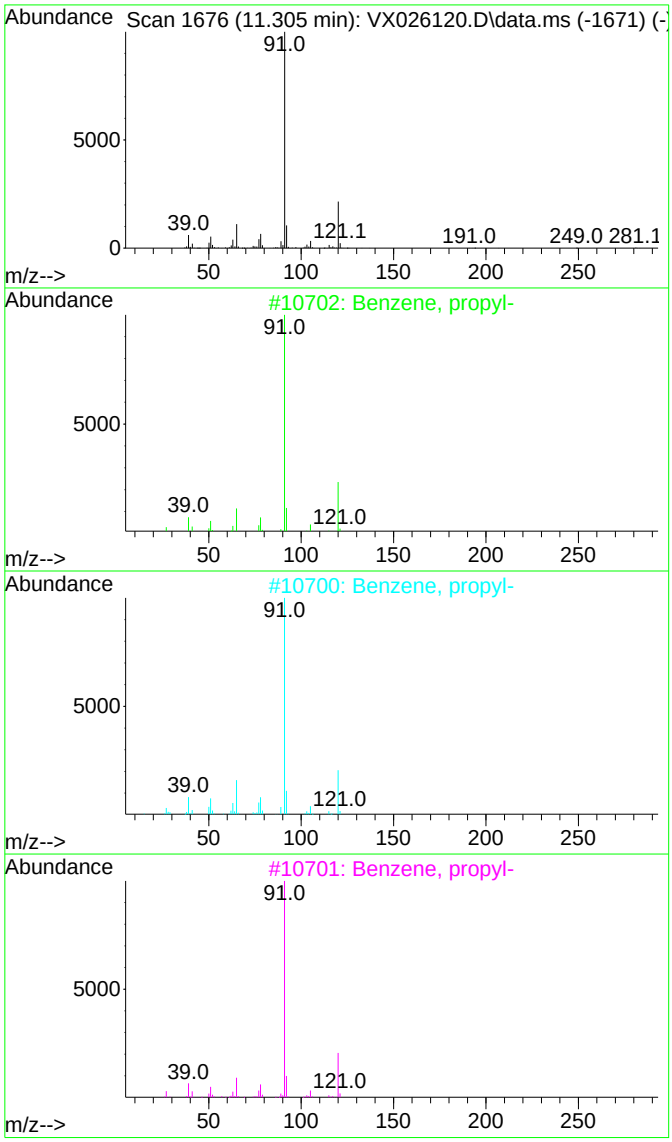
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

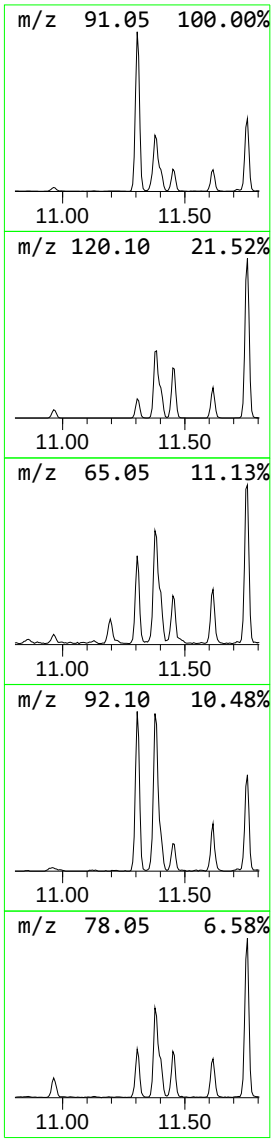
****

Peak Number 31 Benzene, propyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.305	27.77 ug/L	425259	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, propyl-	120	C9H12	000103-65-1	94
2	Benzene, propyl-	120	C9H12	000103-65-1	91
3	Benzene, propyl-	120	C9H12	000103-65-1	90
4	Benzene, propyl-	120	C9H12	000103-65-1	86
5	Ethanol, 2-[(phenylmethyl)amino]-	151	C9H13NO	000104-63-2	78





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

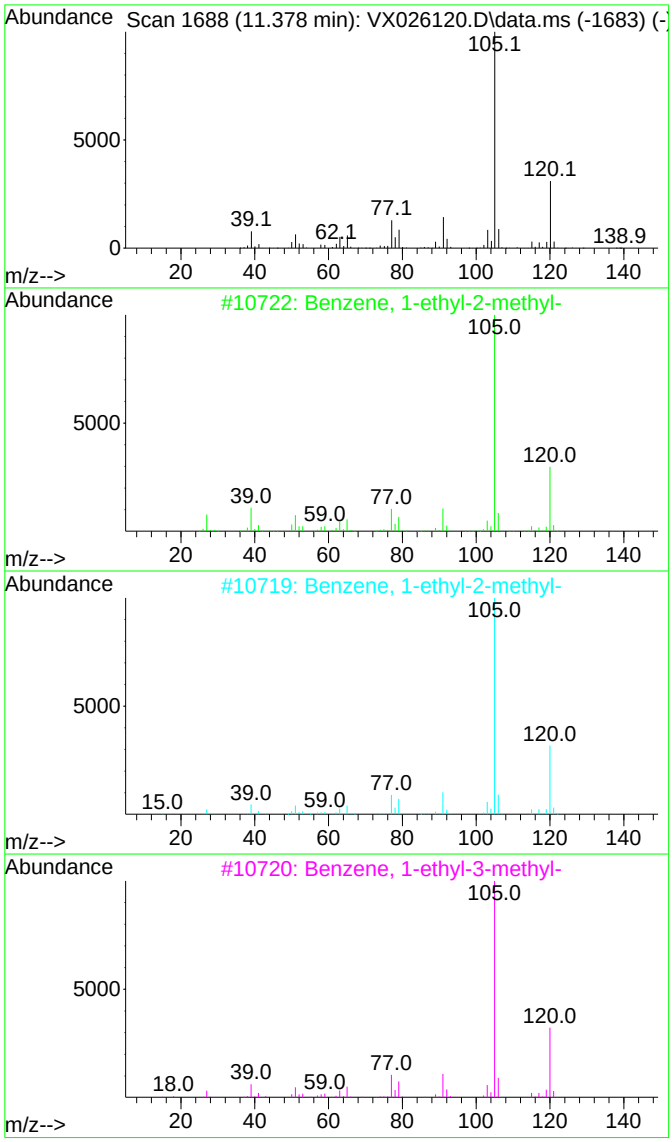
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

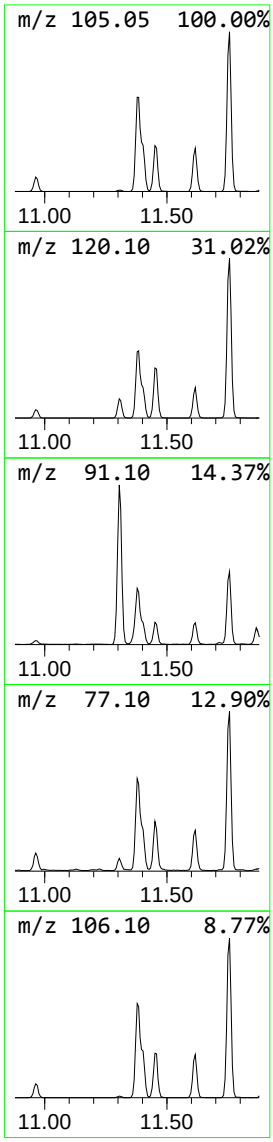
****

Peak Number 32 Benzene, 1-ethyl-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.378	128.14 ug/L	1962120	1,4-Dichlorobenzene-d4	12.024

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-2-methyl-	120	C9H12	000611-14-3	94
3	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

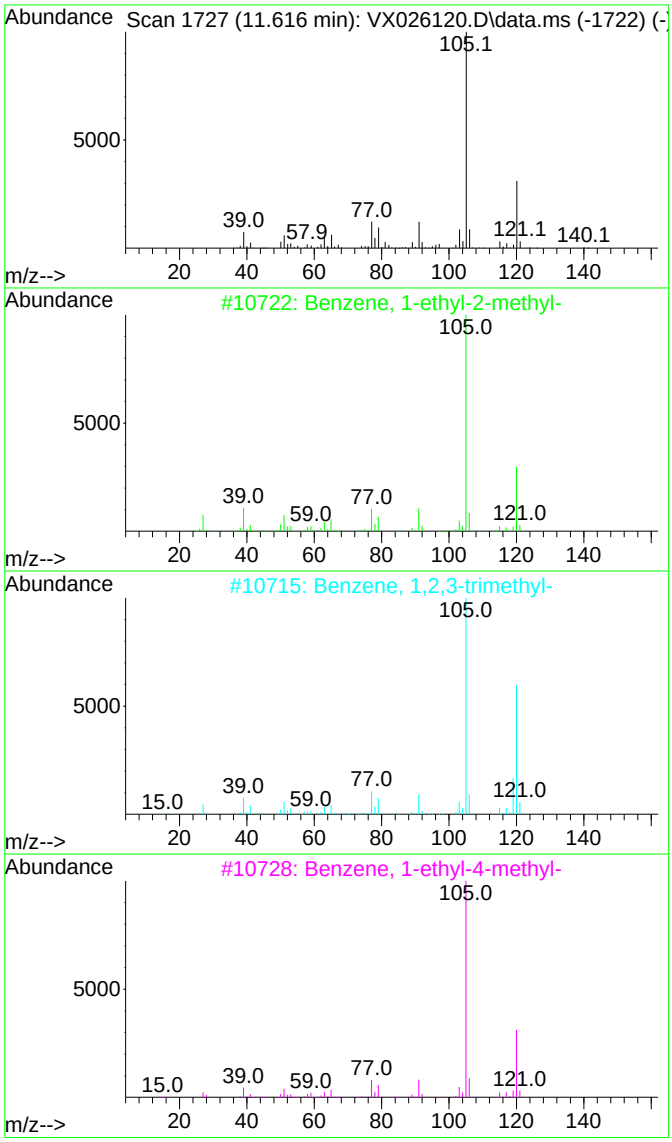
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

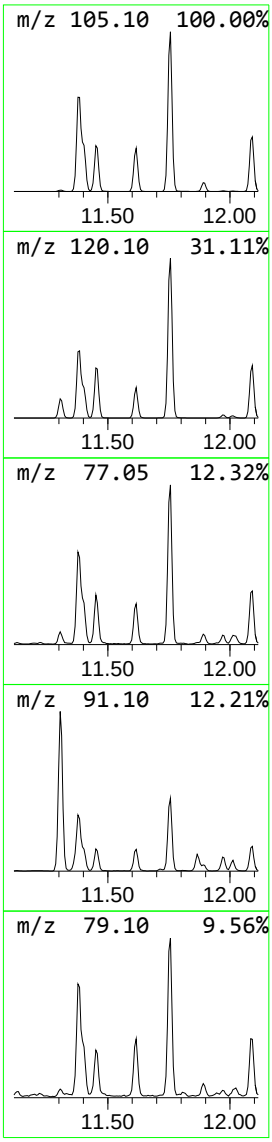
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 33 Benzene, 1,2,3-trimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.616	43.66 ug/L	668562	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
3	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

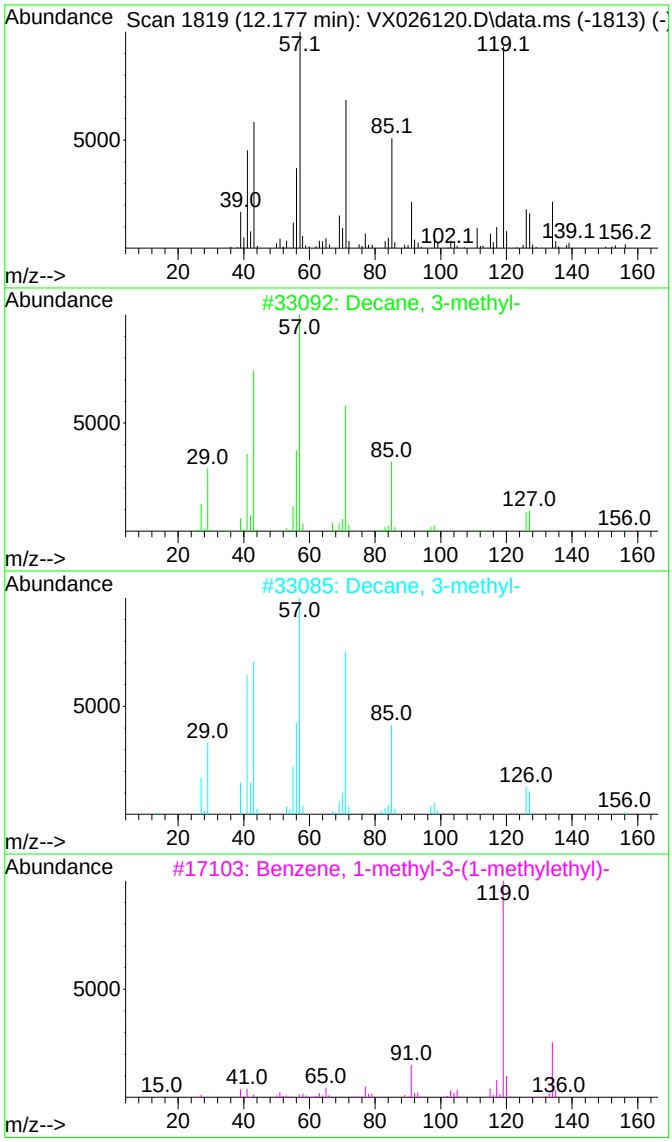
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

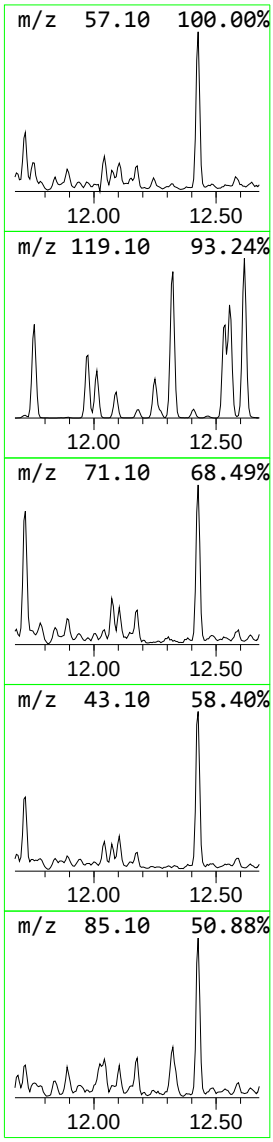
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 36 (DEL) Alkane: Straight-Chai... Concentration Rank 57

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.177	9.29 ug/L	142297	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane, 3-methyl-	156	C11H24	013151-34-3	50
2	Decane, 3-methyl-	156	C11H24	013151-34-3	50
3	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	42
4	p-Cymene	134	C10H14	000099-87-6	42
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	42





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

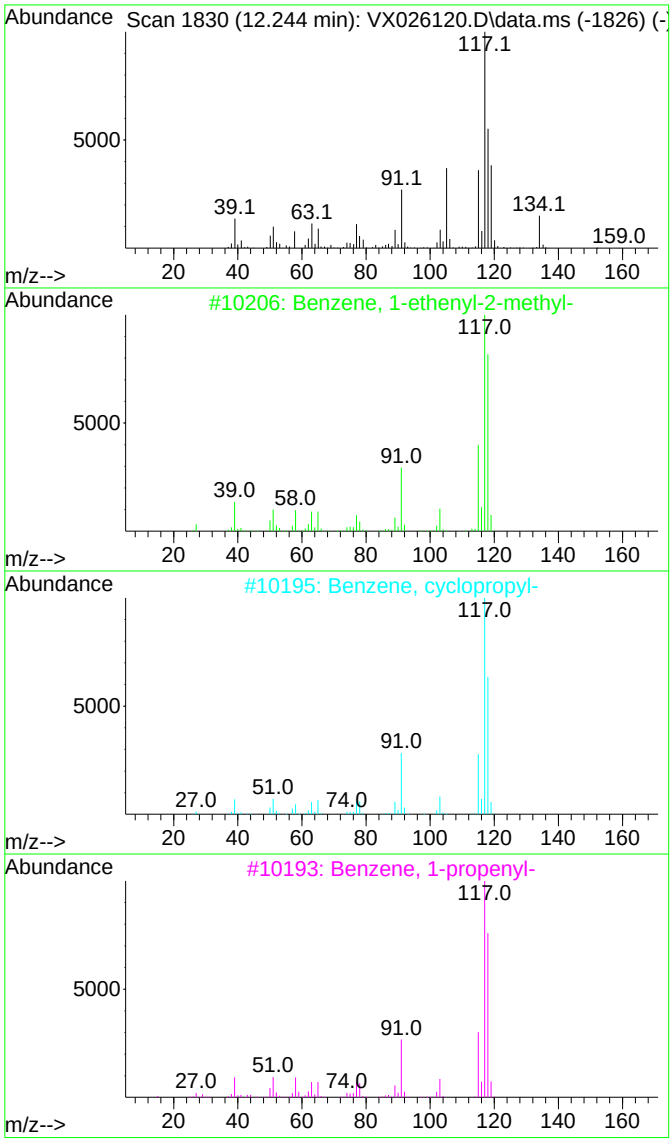
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

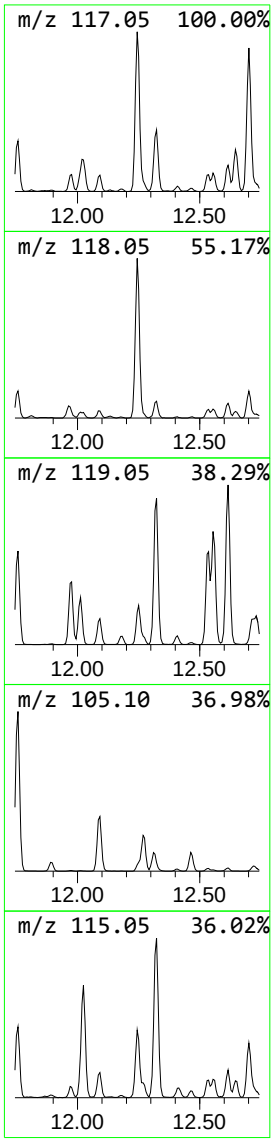
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 37 Benzene, 1-ethenyl-2-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	29.60 ug/L	453220	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	90
2	Benzene, cyclopropyl-	118	C9H10	000873-49-4	87
3	Benzene, 1-propenyl-	118	C9H10	000637-50-3	87
4	Benzene, 2-propenyl-	118	C9H10	000300-57-2	55
5	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	55





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

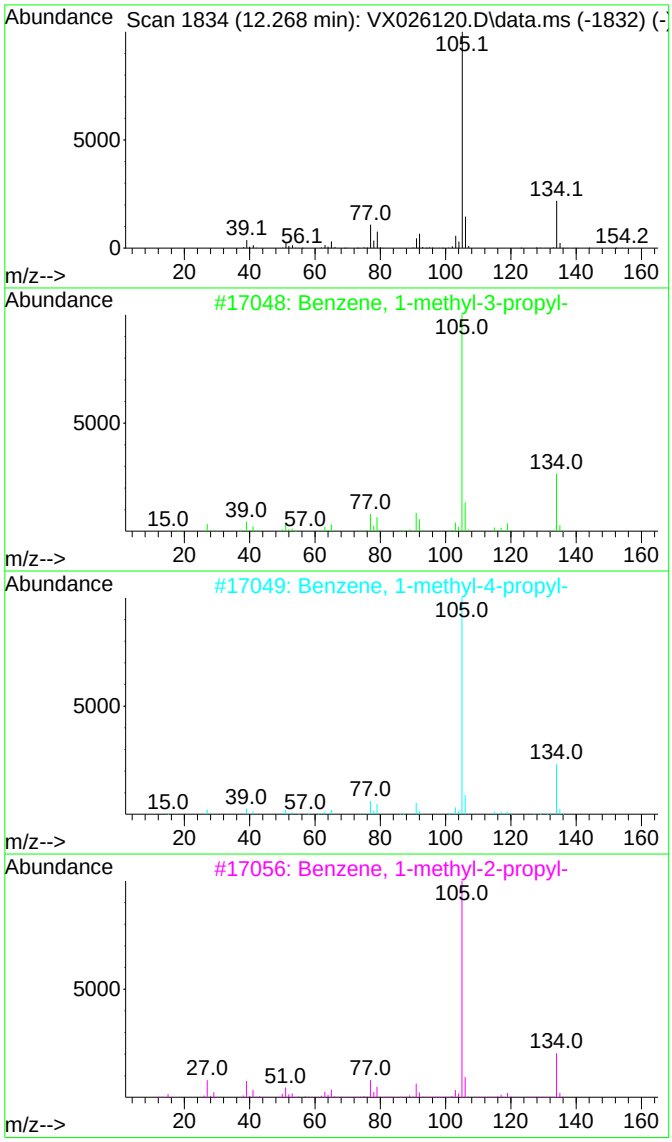
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

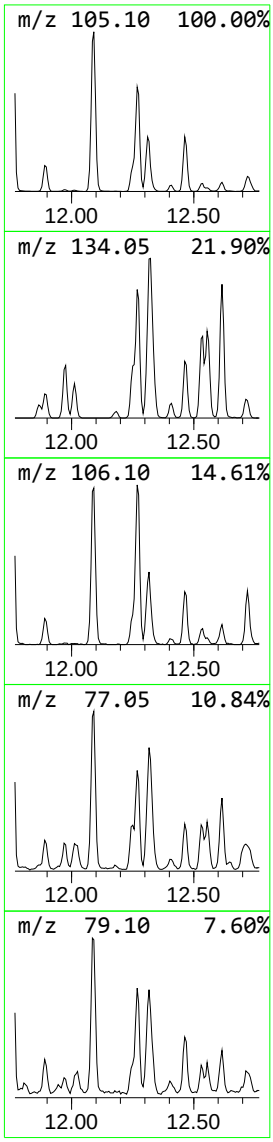
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 38 Benzene, 1-methyl-3-propyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.268	27.27 ug/L	417534	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	91
2	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	90
3	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	86
4	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	86
5	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	86





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

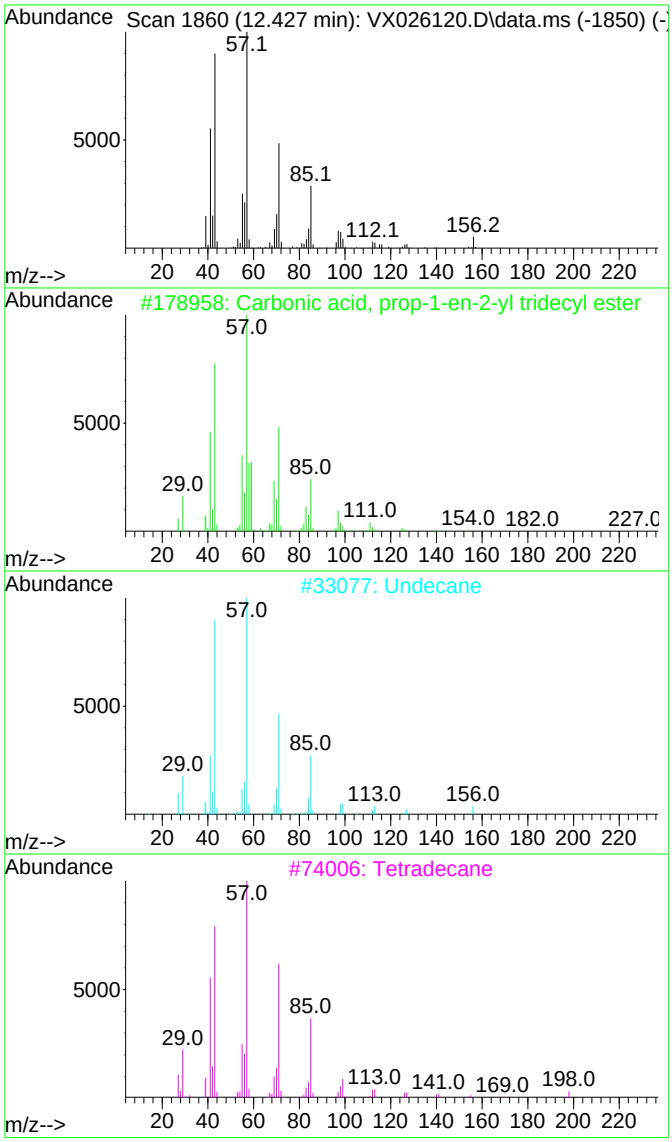
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

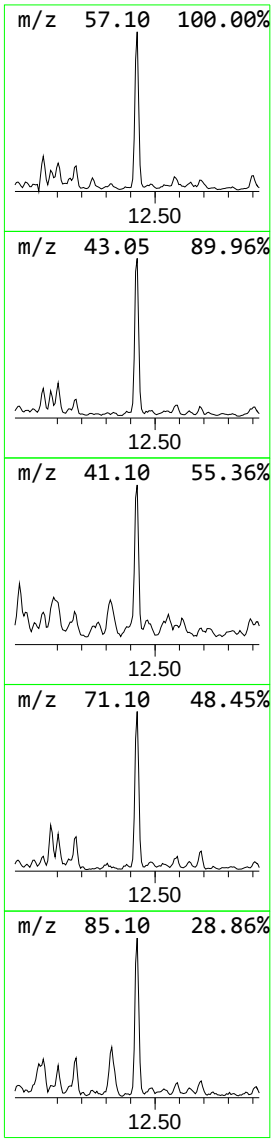
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 39 Carbonic acid, prop-1-en-2-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.427	32.09 ug/L	491362	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Carbonic acid, prop-1-en-2-yl tr...	284	C17H32O3	1000382-90-8	90
2	Undecane	156	C11H24	001120-21-4	87
3	Tetradecane	198	C14H30	000629-59-4	86
4	Undecane	156	C11H24	001120-21-4	81
5	Undecane	156	C11H24	001120-21-4	74





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

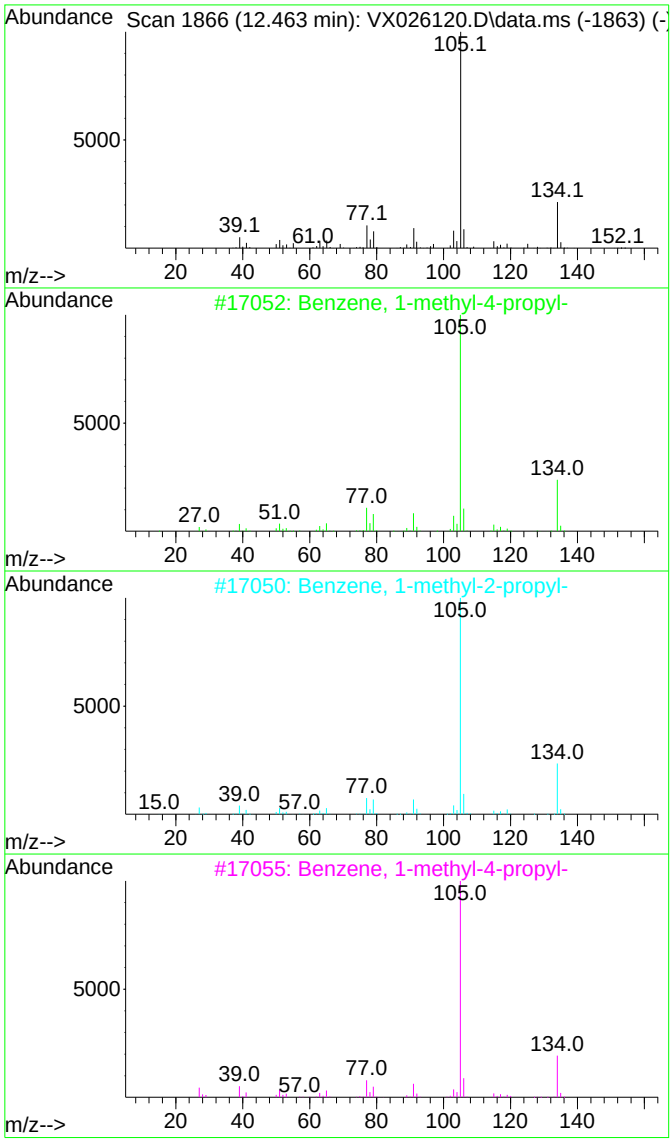
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

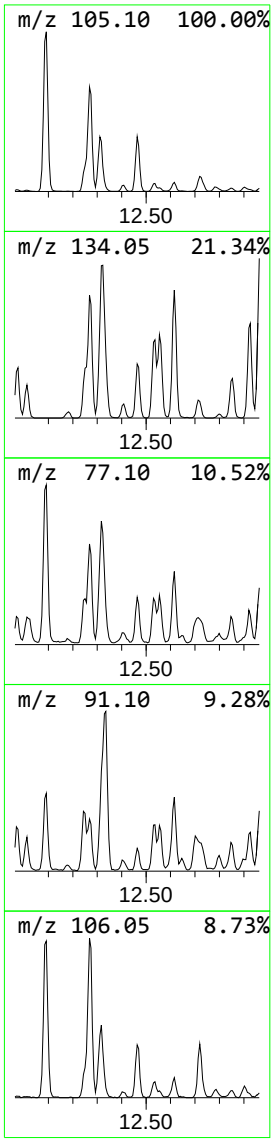
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 40 Benzene, 1-methyl-4-propyl- Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.463	17.21 ug/L	263510	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	95
2	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	94
3	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
4	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	91
5	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	90





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

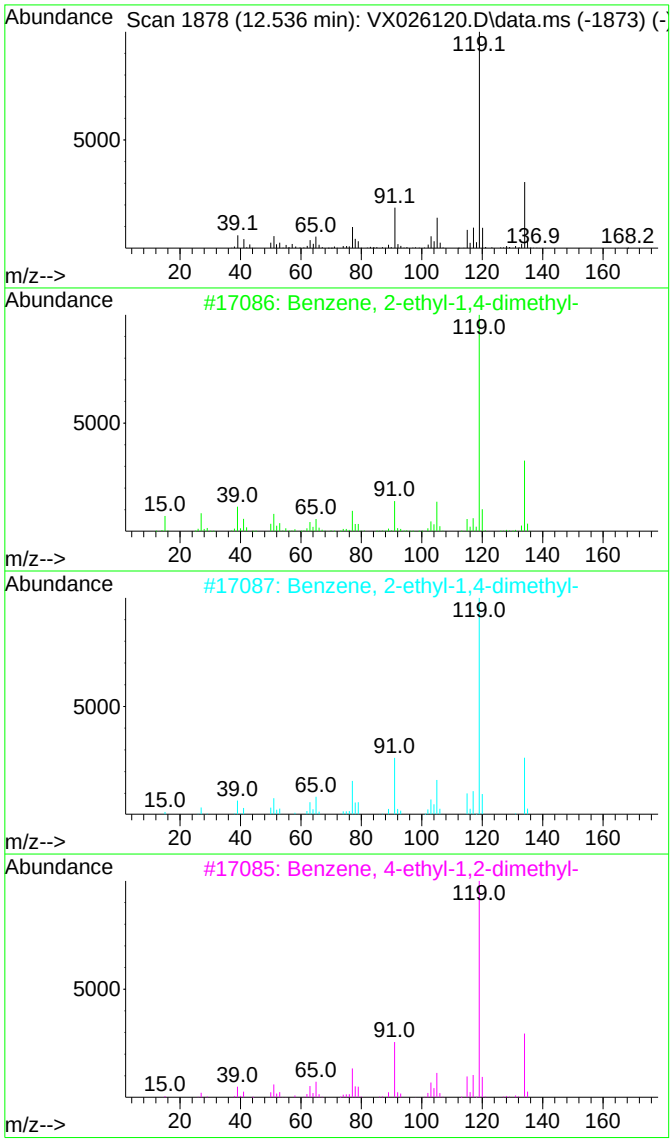
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

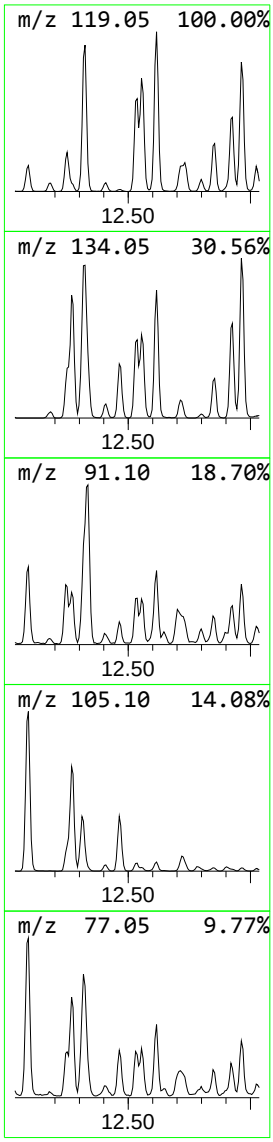
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 41 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.536	18.94 ug/L	289994	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	97
2	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
3	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
4	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
5	o-Cymene	134	C10H14	000527-84-4	95





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

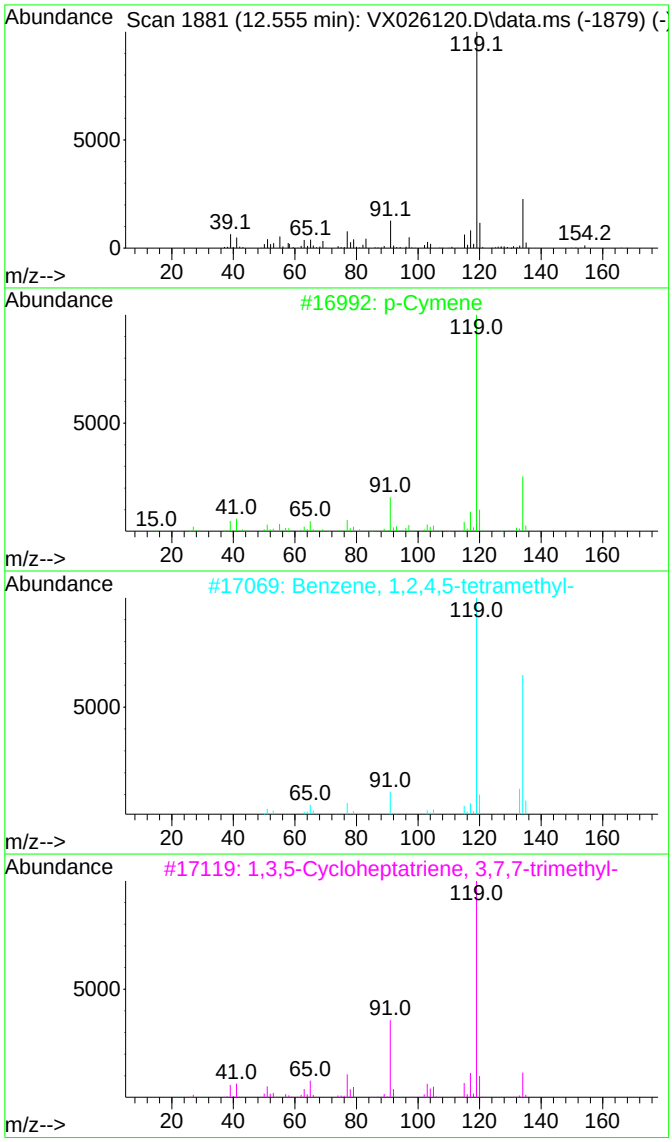
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

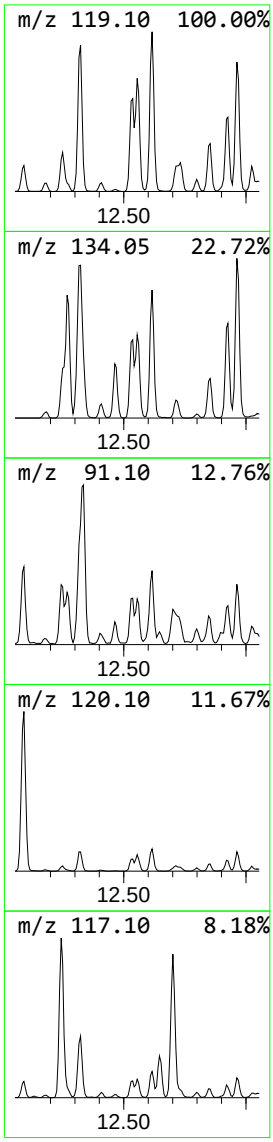
****

Peak Number 42 p-Cymene Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.555	18.25 ug/L	279516	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Cymene	134	C10H14	000099-87-6	94
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	93
3	1,3,5-Cycloheptatriene, 3,7,7-tr...	134	C10H14	003479-89-8	90
4	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	90
5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	90





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

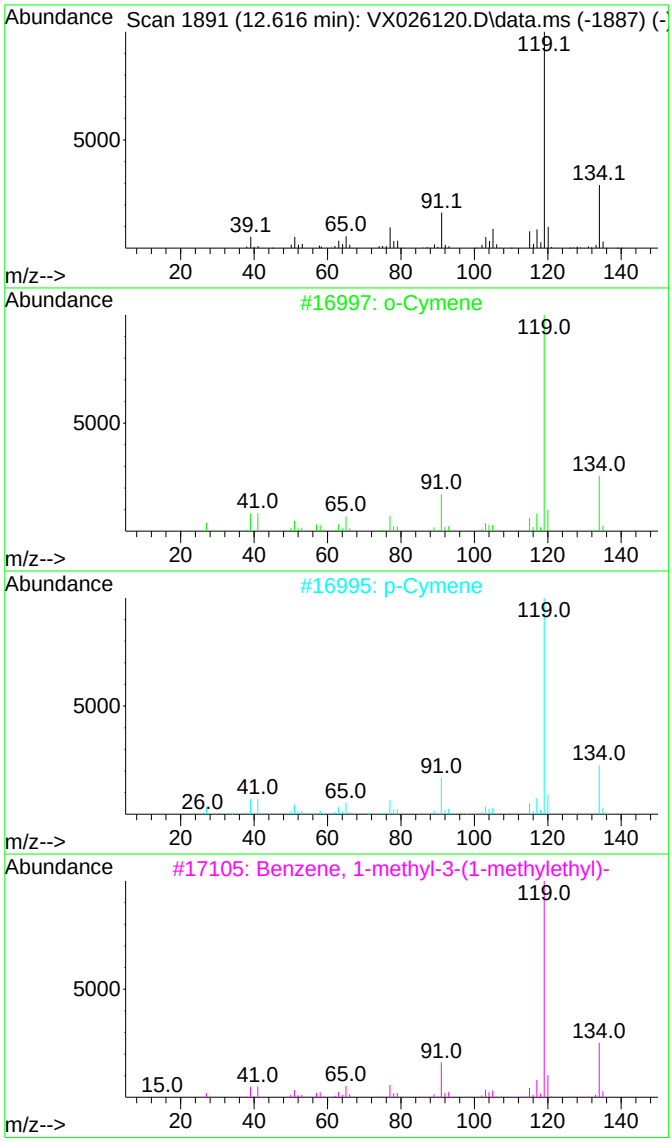
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

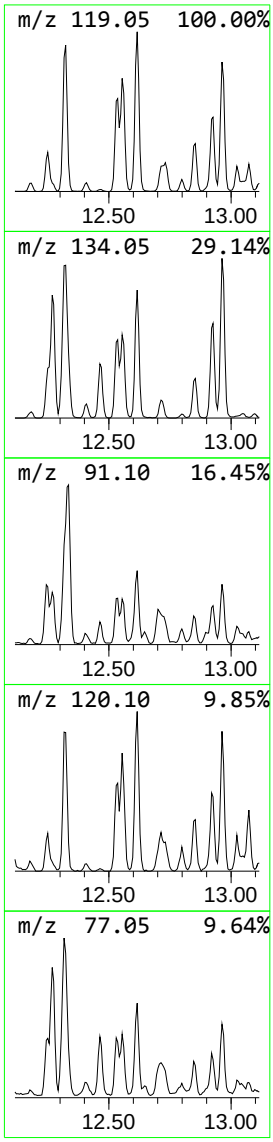
****

Peak Number 43 o-Cymene Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.616	26.27 ug/L	402190	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	p-Cymene	134	C10H14	000099-87-6	95
3	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
4	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

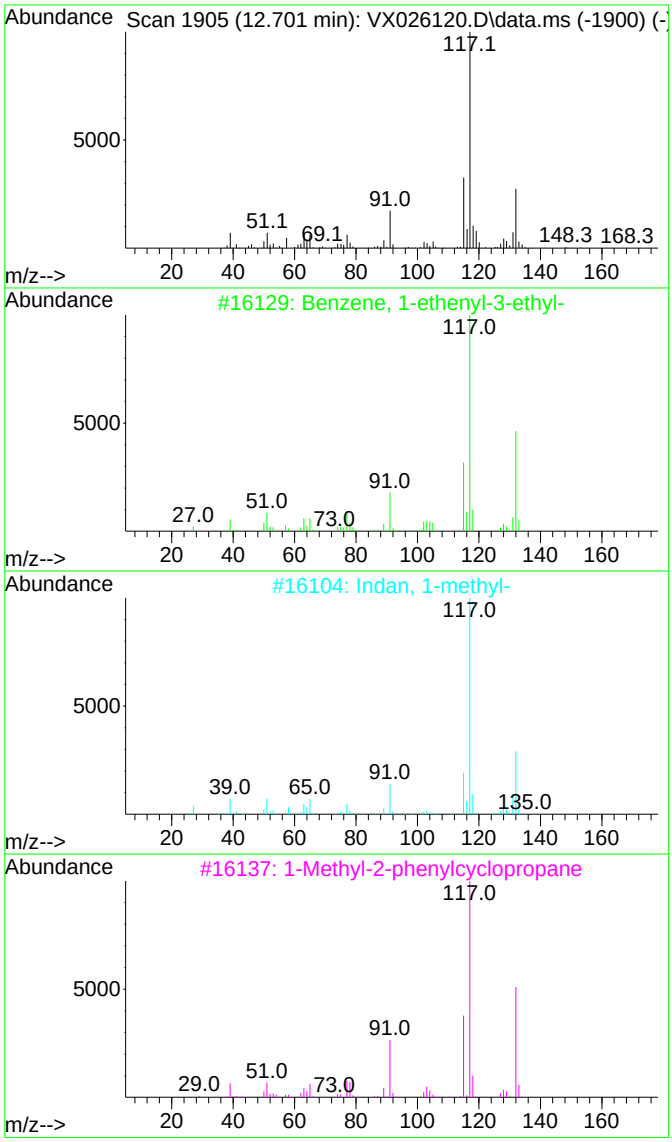
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

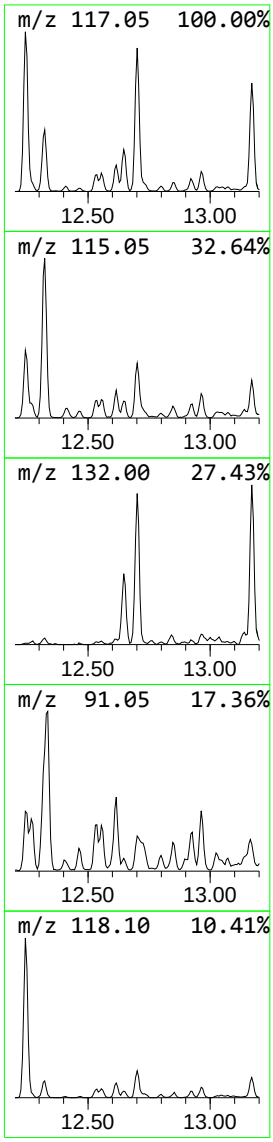
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 44 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	30.83 ug/L	472125	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	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	87
4	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	83
5	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	83





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

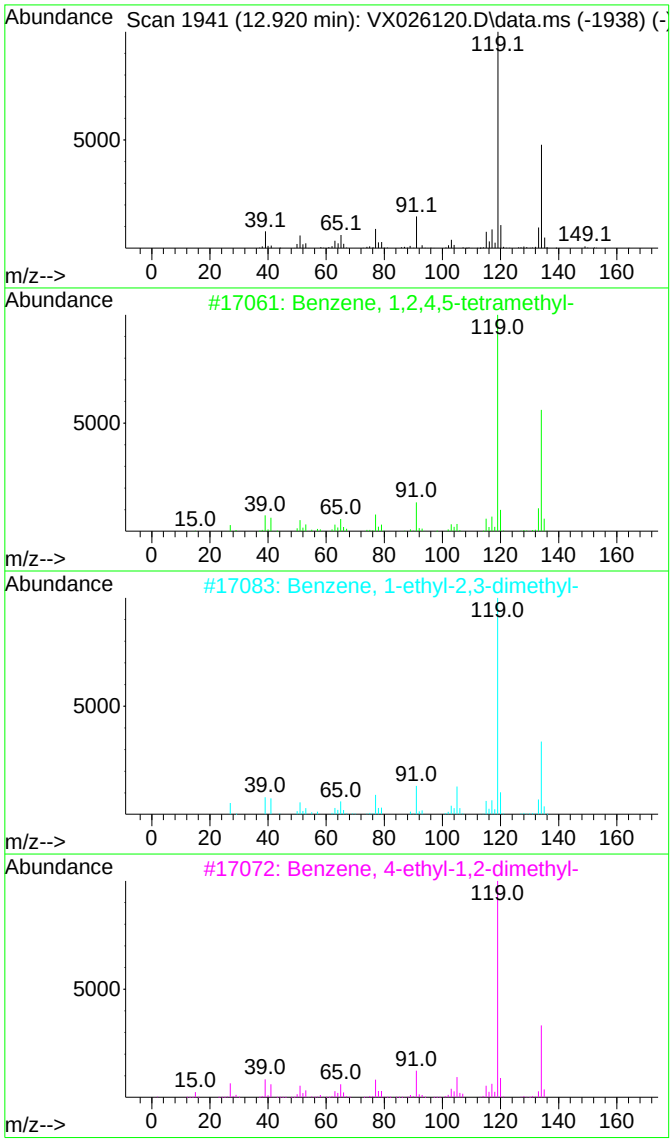
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

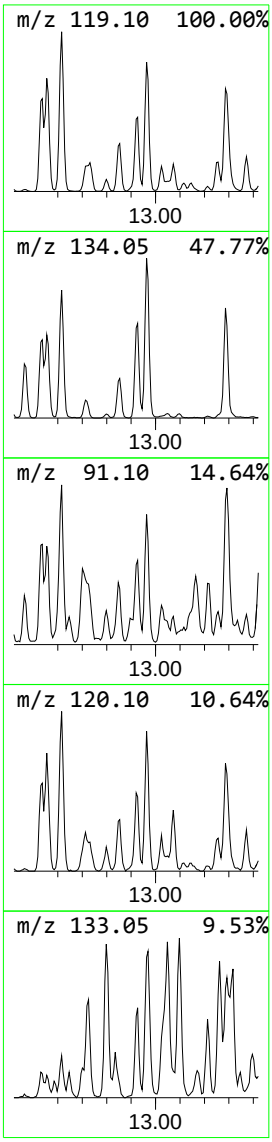
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 46 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 39

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.920	12.91 ug/L	197656	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	94
2	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94
3	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	91
4	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91
5	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

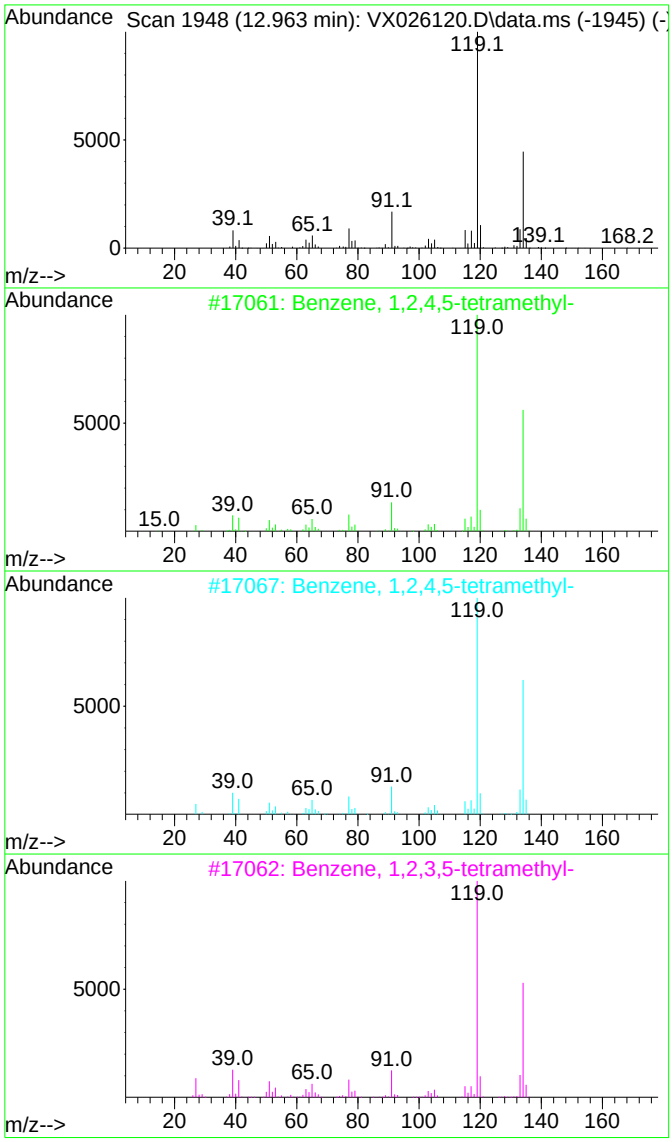
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

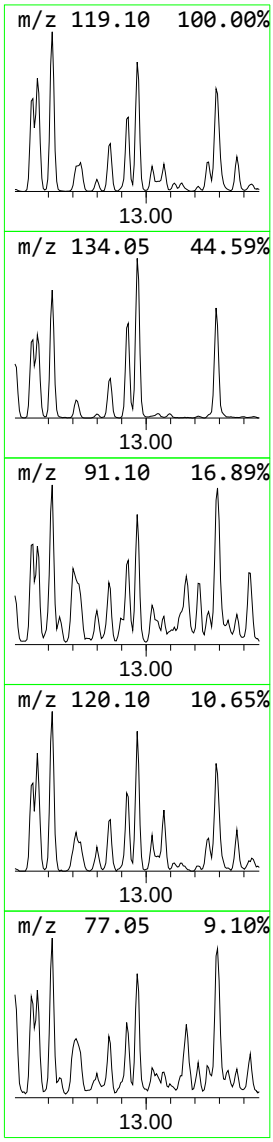
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 47 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.963	28.30 ug/L	433410	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
3	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
4	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

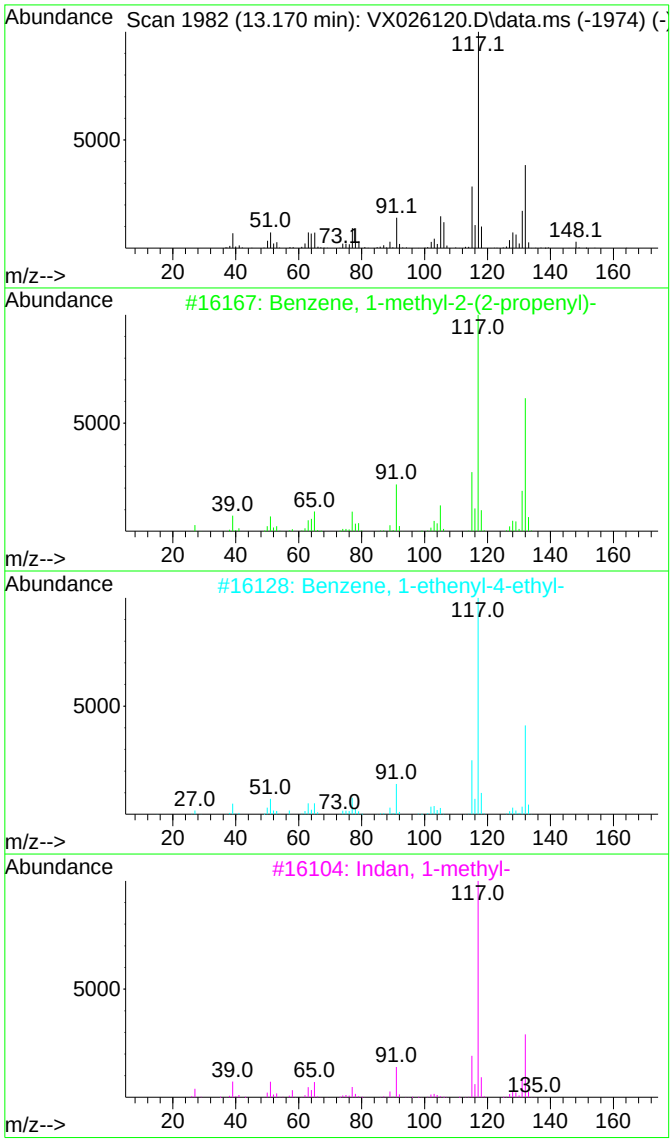
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

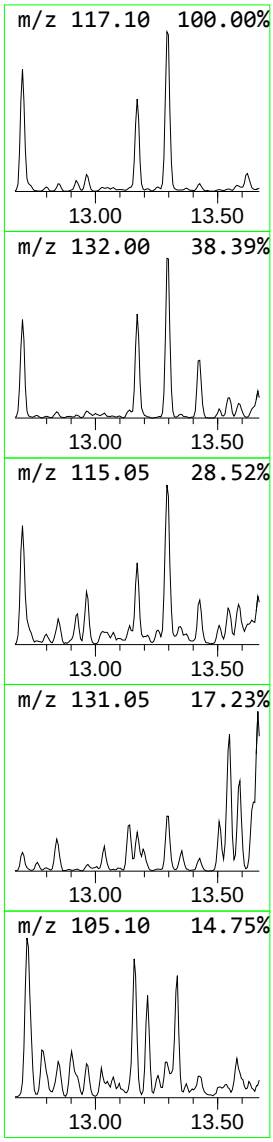
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 48 Benzene, 1-methyl-2-(2-prop... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.170	18.33 ug/L	280627	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	90
2	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
3	Indan, 1-methyl-	132	C10H12	000767-58-8	87
4	Benzene, 2-butenyl-	132	C10H12	001560-06-1	87
5	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	87





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

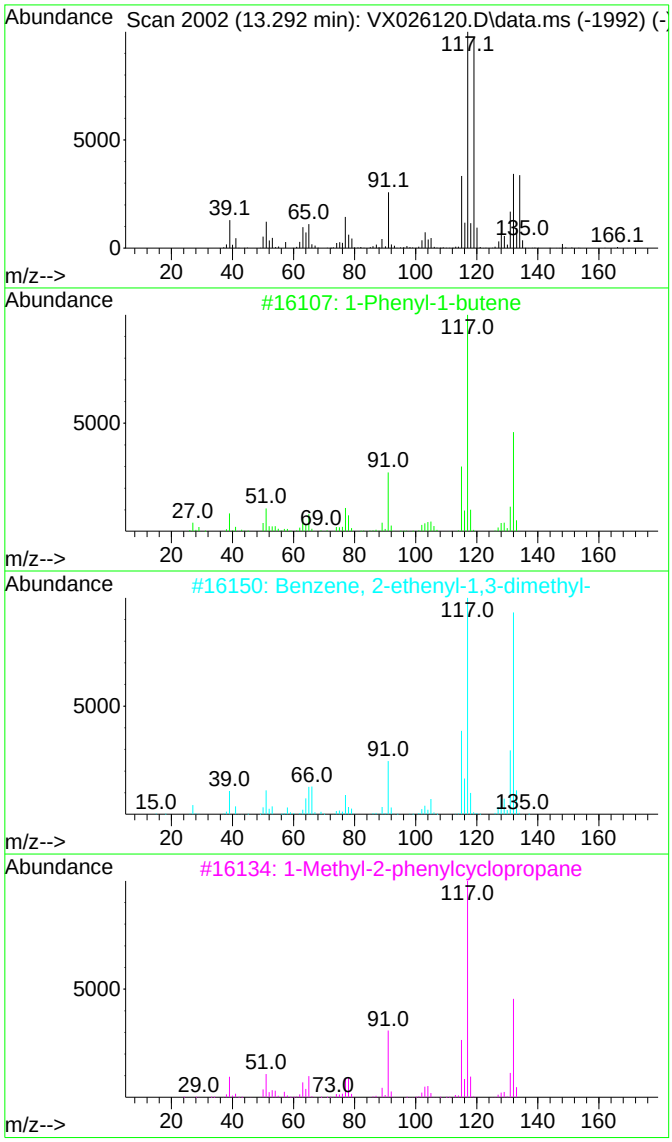
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

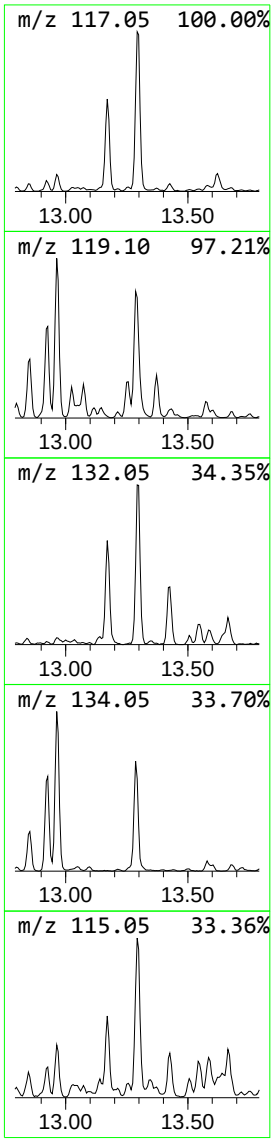
****

Peak Number 49 1-Phenyl-1-butene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	56.88 ug/L	870945	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	87
2	Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	70
3	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	55
4	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	55
5	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	55





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

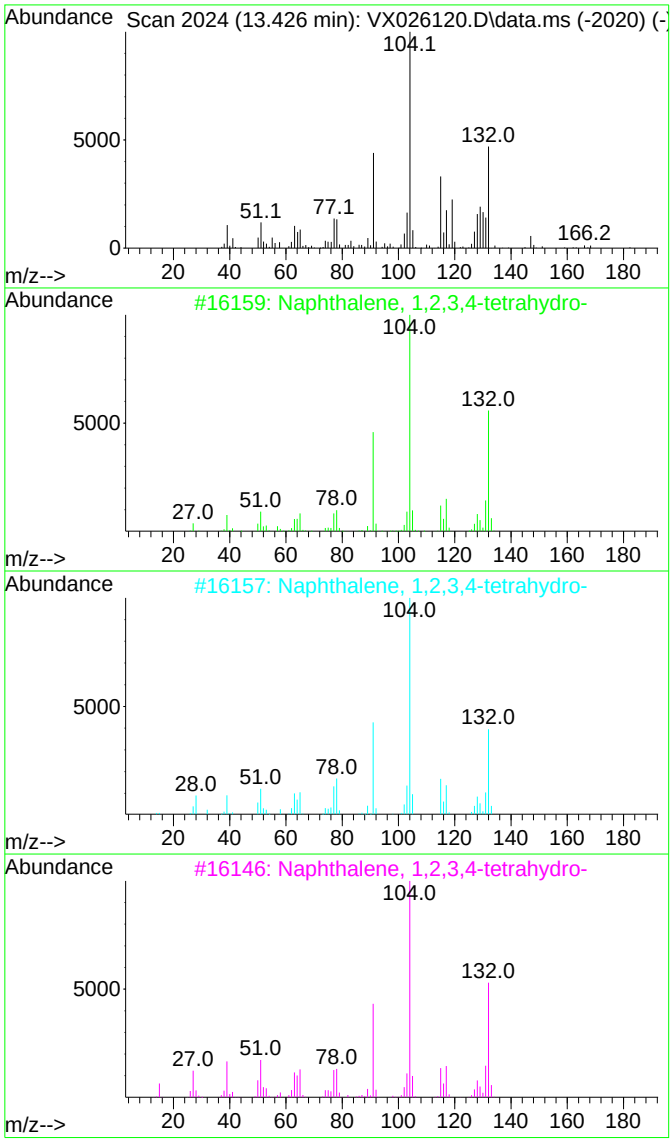
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

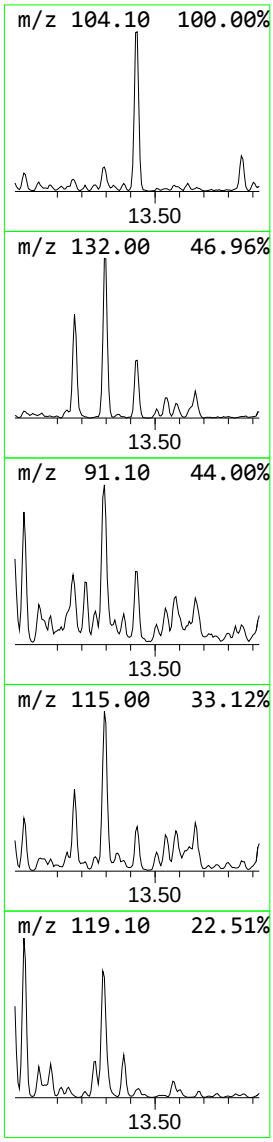
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 50 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.427	14.06 ug/L	215279	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	93
2	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	93
3	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	70
4	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	70
5	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	70





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

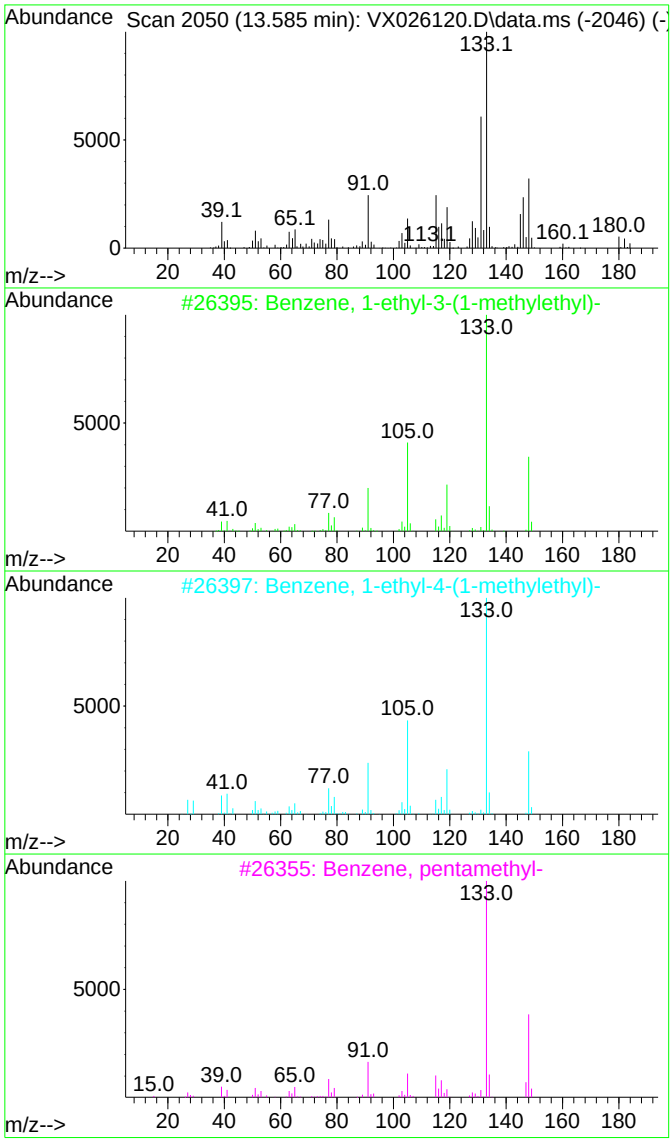
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

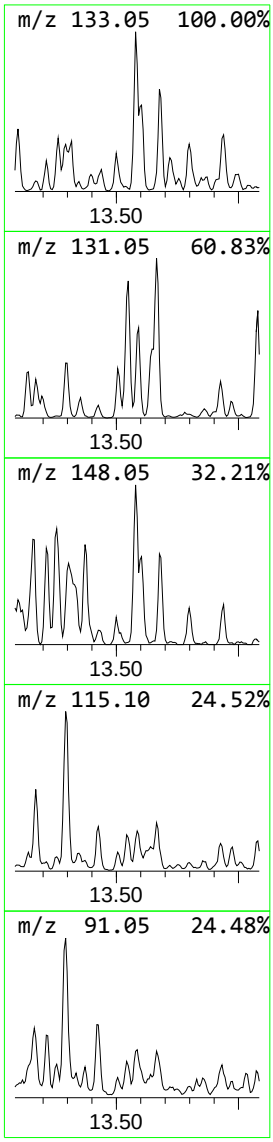
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 52 Benzene, 1-ethyl-3-(1-methy... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.585	16.48 ug/L	252302	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3-(1-methylethyl)-	148	C11H16	004920-99-4	53
2	Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	53
3	Benzene, pentamethyl-	148	C11H16	000700-12-9	53
4	Ethanone, 1-(4-ethylphenyl)-	148	C10H12O	000937-30-4	50
5	Ethanone, 1-(4-ethylphenyl)-	148	C10H12O	000937-30-4	50





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

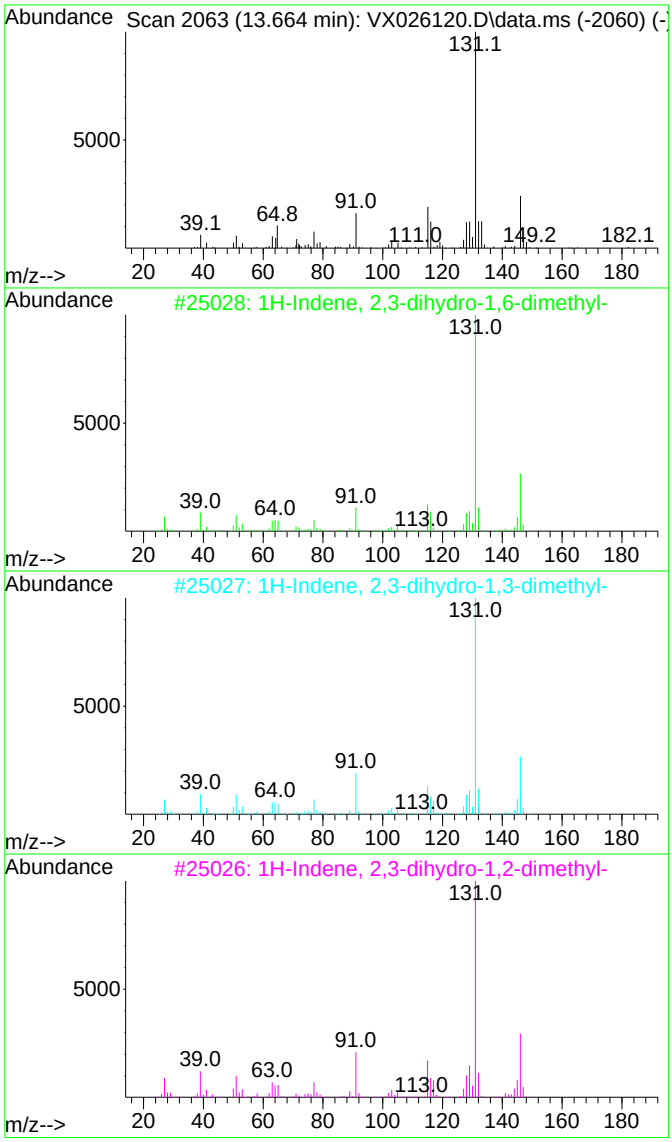
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

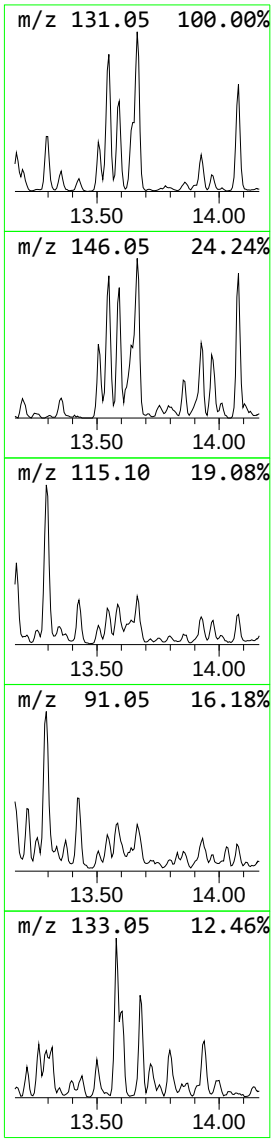
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 53 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 40

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.664	12.27 ug/L	187933	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	91
2	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	91
3	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	91
4	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	91
5	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

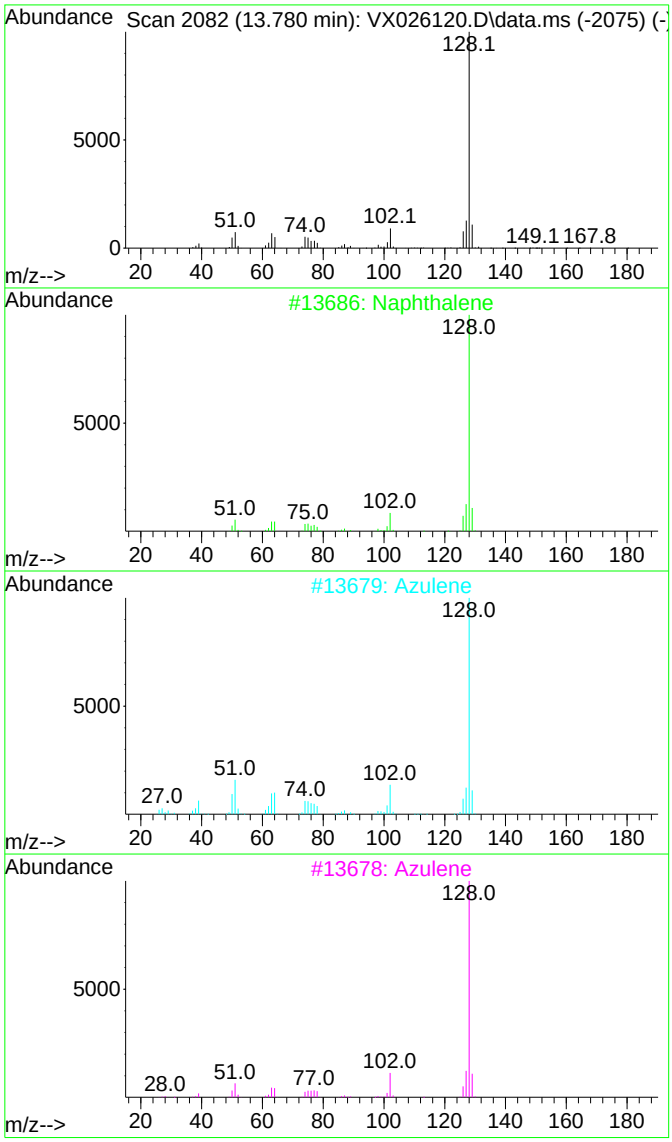
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

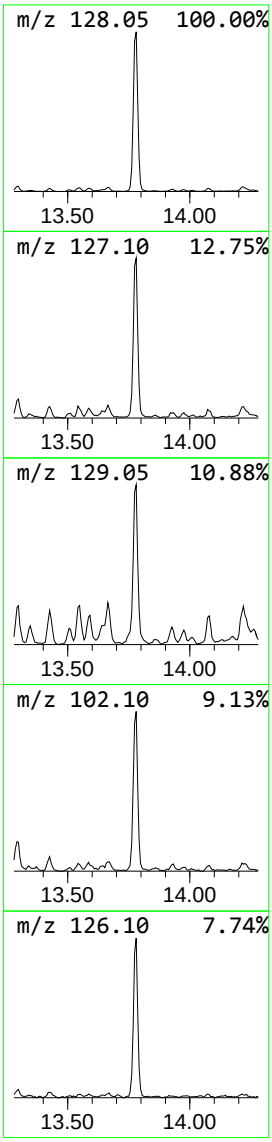
****

Peak Number 54 Azulene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.780	34.39 ug/L	526566	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene	128	C10H8	000091-20-3	95
2	Azulene	128	C10H8	000275-51-4	91
3	Azulene	128	C10H8	000275-51-4	91
4	Azulene	128	C10H8	000275-51-4	91
5	Naphthalene	128	C10H8	000091-20-3	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

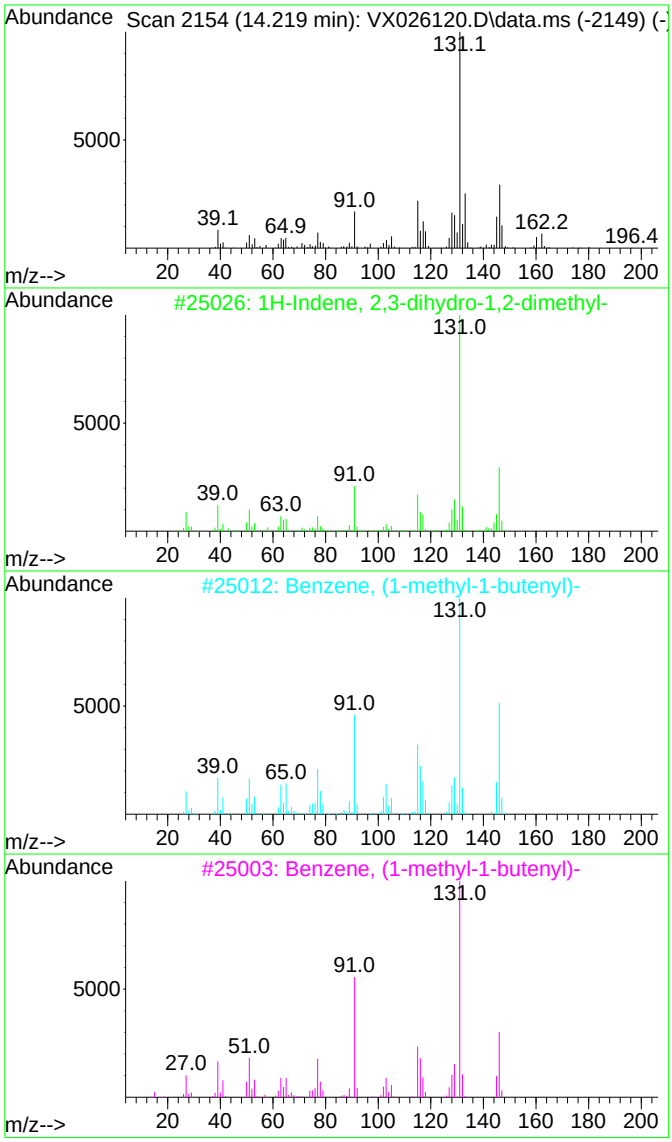
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

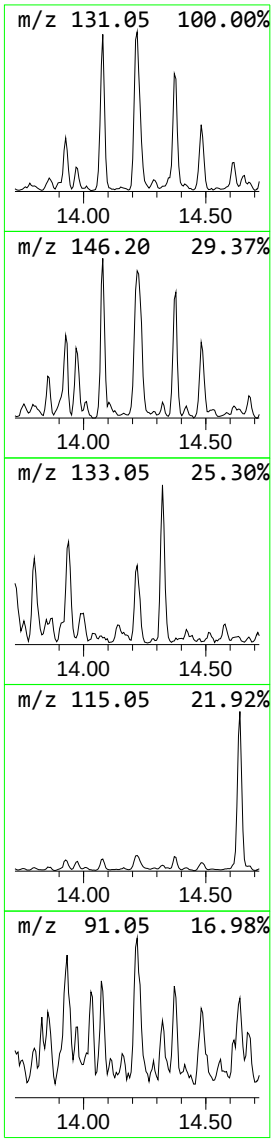
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 57 1H-Indene, 2,3-dihydro-1,2-... Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.219	14.83 ug/L	227082	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	87
2	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	87
3	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	83
4	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	81
5	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	81





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

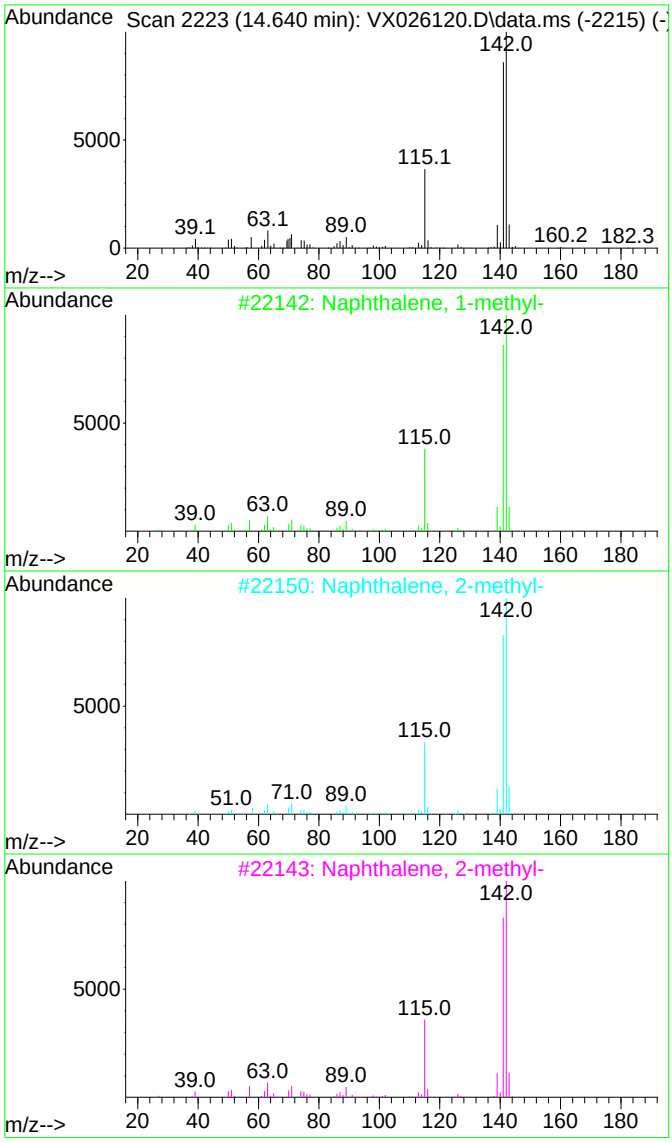
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

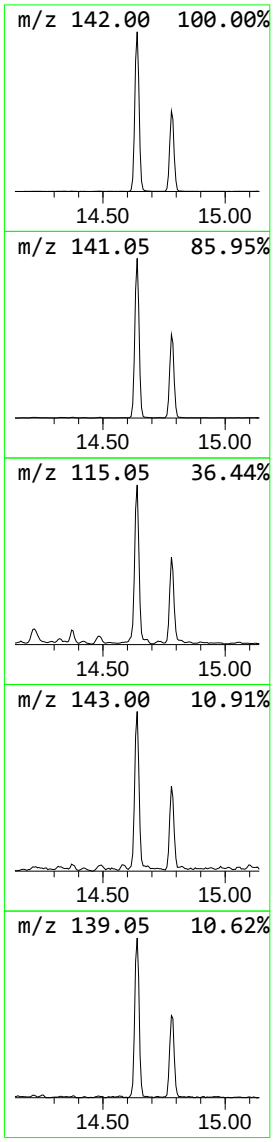
****

Peak Number 60 Naphthalene, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.640	59.09 ug/L	904869	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	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
4	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	95
5	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

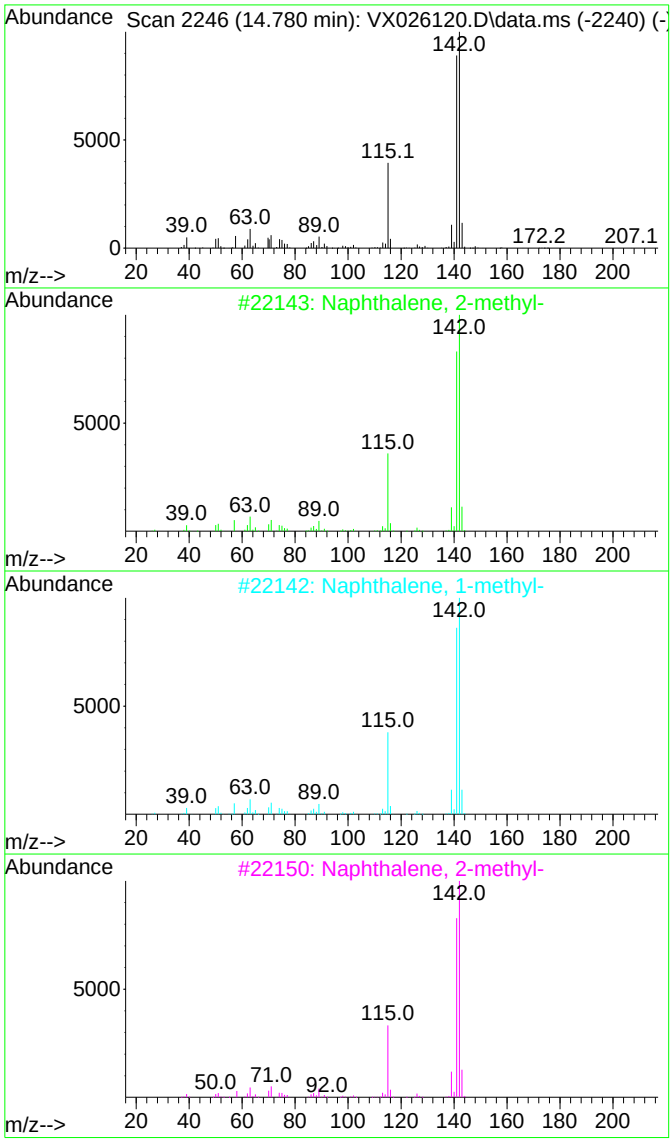
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

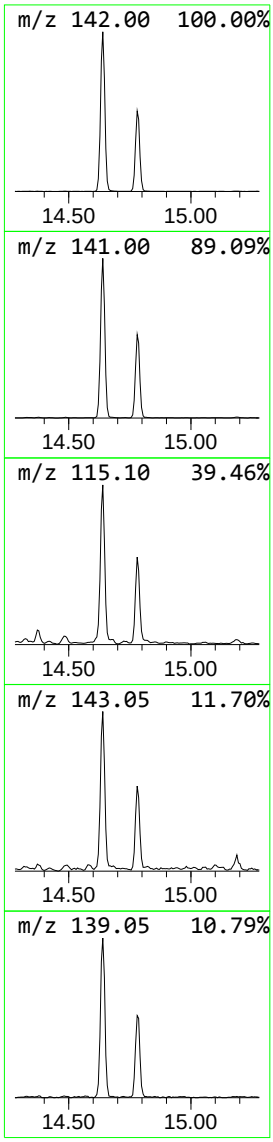
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 61 Naphthalene, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.780	32.87 ug/L	503316	1,4-Dichlorobenzene-d4	12.024

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	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
4	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	95
5	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

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

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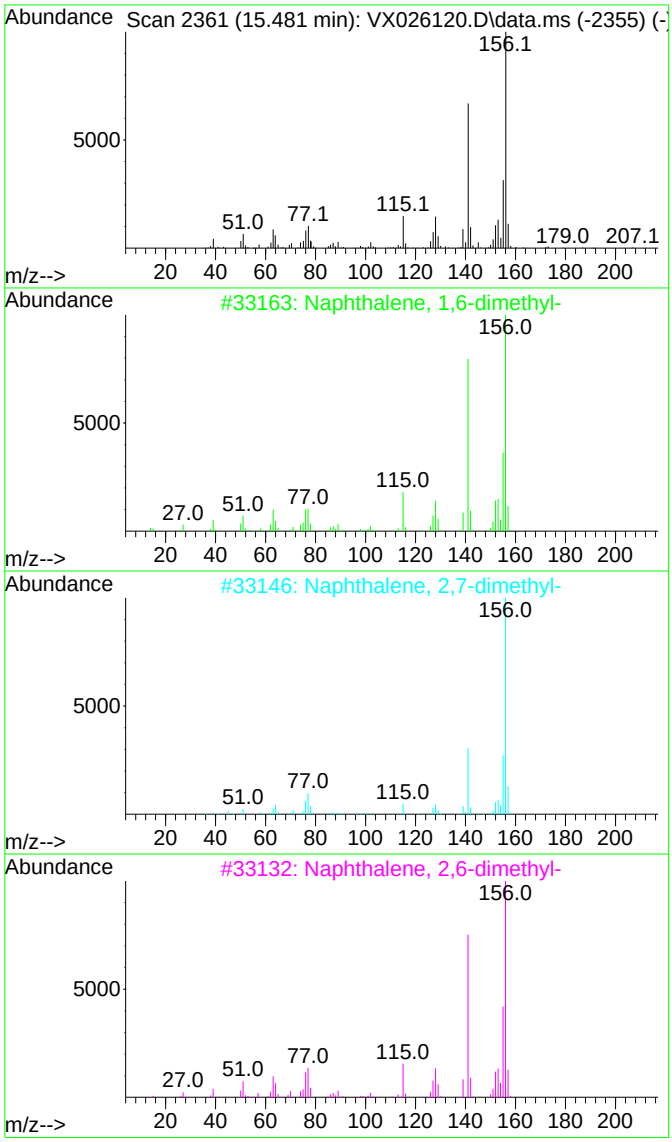
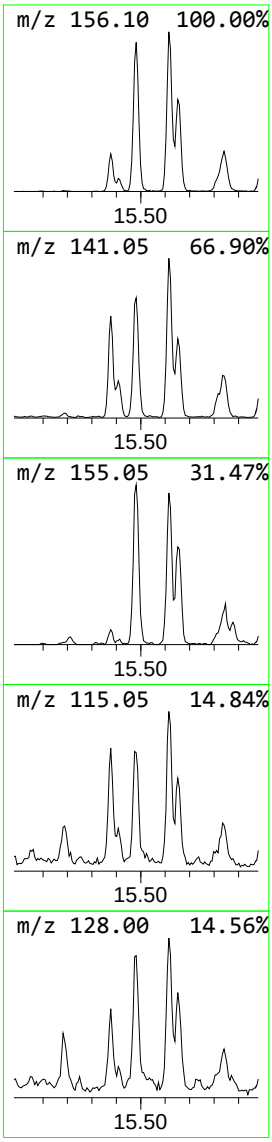
Peak Number 63 Naphthalene, 1,6-dimethyl- Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.481	18.28 ug/L	279942	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	98
2	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
3	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
4	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
5	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	97





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

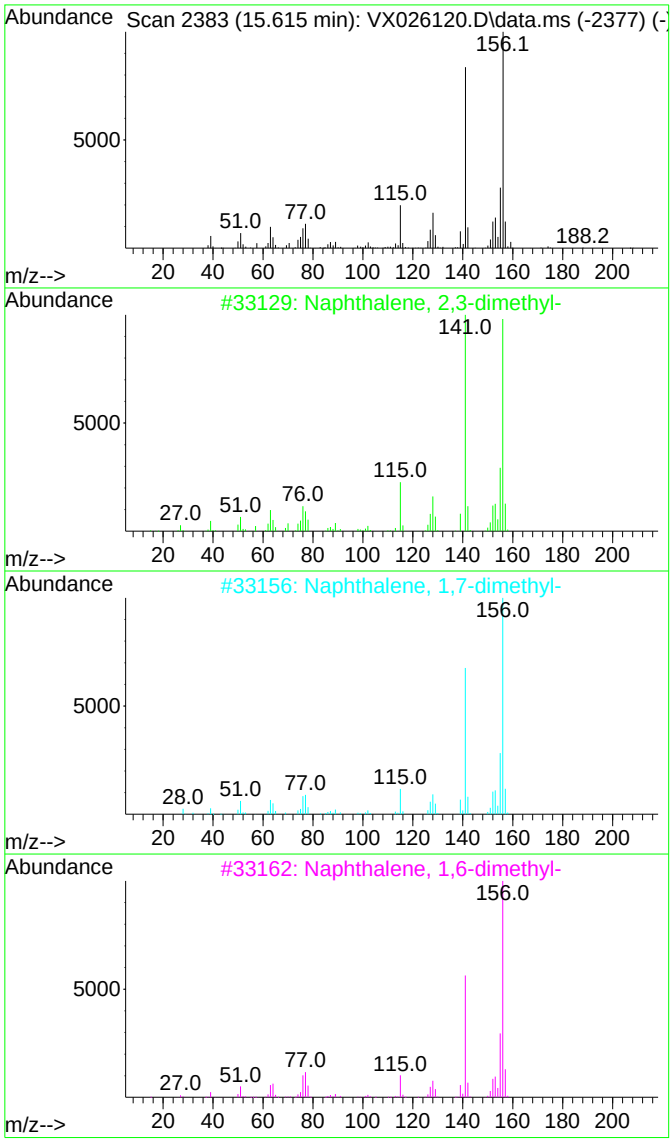
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

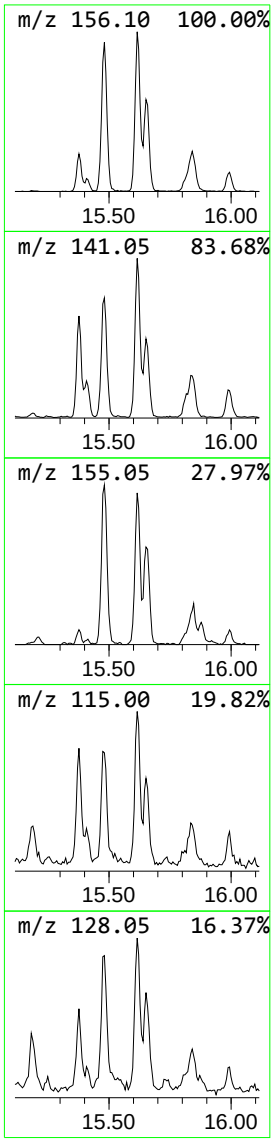
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 64 Naphthalene, 2,3-dimethyl- Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.615	16.47 ug/L	252131	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, 1,7-dimethyl-	156	C12H12	000575-37-1	97
3	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
4	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
5	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

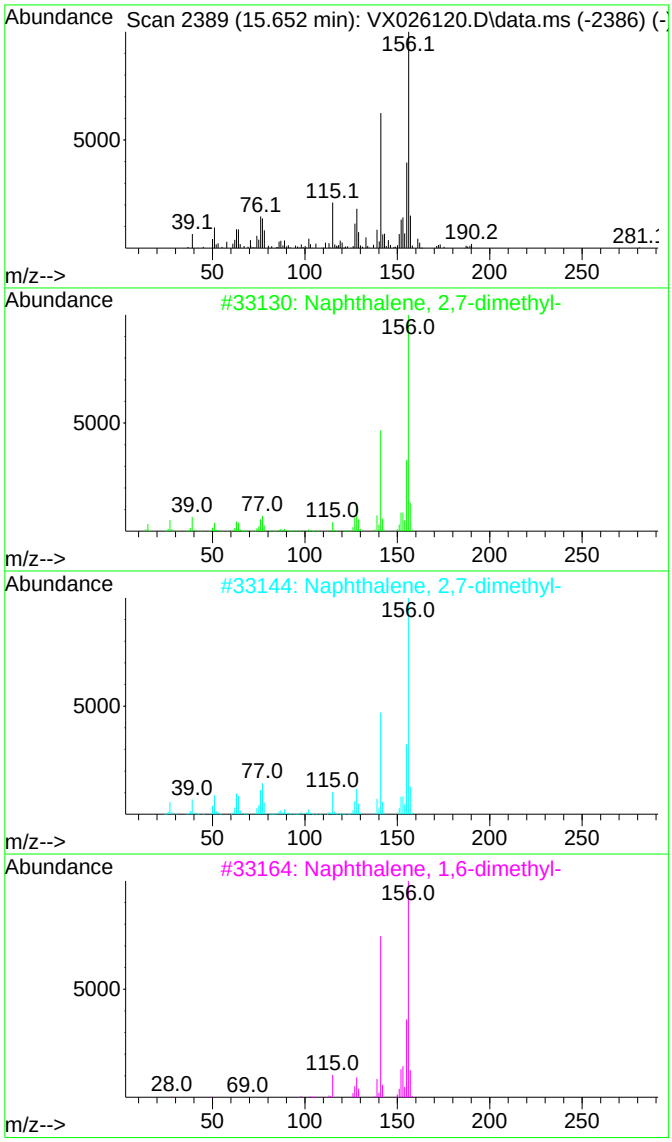
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

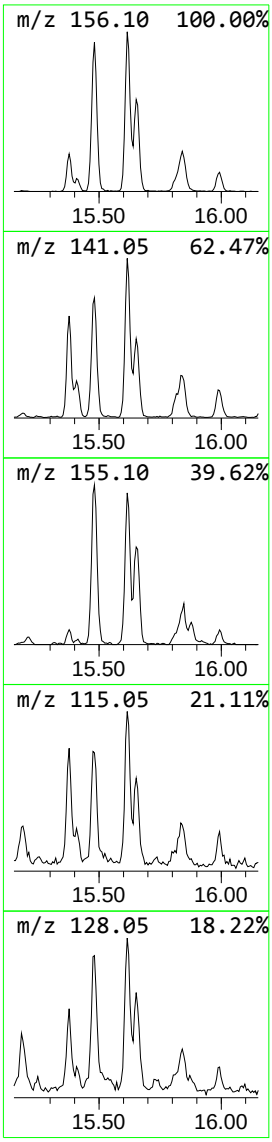
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 65 Naphthalene, 2,7-dimethyl- Concentration Rank 45

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.652	11.29 ug/L	172870	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96
2	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96
3	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	93
4	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	93
5	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	93





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122821\
Data File : VX026120.D
Acq On : 29 Dec 2021 00:14
Operator : JC/MD
Sample : M5233-01 10X
Misc : 10.51g/5mL/100uL/5.00mL/MSVOA_X/MEOH
ALS Vial : 27 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
GBCC1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\SFAMXLM122821WMA.M
Quant Title : VOC Analysis

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
(DEL) Alkane: S...	3.447	10.4	ug/L	99174	1	6.763	475762	50.0
Cyclopentene, 1...	5.190	16.2	ug/L	153880	1	6.763	475762	50.0
(DEL) Alkane: S...	5.623	14.2	ug/L	135416	1	6.763	475762	50.0
(DEL) Alkane: S...	5.830	14.2	ug/L	135520	1	6.763	475762	50.0
(DEL) Alkane: S...	6.257	18.3	ug/L	174173	1	6.763	475762	50.0
(DEL) Alkane: S...	6.367	11.5	ug/L	109086	1	6.763	475762	50.0
(DEL) Alkane: S...	6.617	25.6	ug/L	243654	1	6.763	475762	50.0
(DEL) Alkane: S...	6.848	14.2	ug/L	135346	1	6.763	475762	50.0
(DEL) Alkane: S...	7.178	11.1	ug/L	105919	1	6.763	475762	50.0
(DEL) Alkane: S...	7.988	15.1	ug/L	144035	1	6.763	475762	50.0
(DEL) Alkane: S...	8.104	11.4	ug/L	108718	1	6.763	475762	50.0
Cyclopentene, 4...	8.147	10.6	ug/L	100716	1	6.763	475762	50.0
(DEL) Alkane: S...	8.281	18.6	ug/L	176557	1	6.763	475762	50.0
(DEL) Alkane: S...	8.440	11.4	ug/L	125845	2	10.055	552959	50.0
Cyclohexene, 1-...	8.507	10.8	ug/L	119421	2	10.055	552959	50.0
unknown-01	9.500	12.1	ug/L	133899	2	10.055	552959	50.0
(DEL) Alkane: C...	9.567	9.2	ug/L	101260	2	10.055	552959	50.0
(DEL) Alkane: C...	9.616	10.0	ug/L	110890	2	10.055	552959	50.0
(DEL) Alkane: C...	9.836	9.6	ug/L	105807	2	10.055	552959	50.0
(DEL) Alkane: S...	9.909	14.4	ug/L	159650	2	10.055	552959	50.0
(DEL) Alkane: S...	10.360	63.4	ug/L	701595	2	10.055	552959	50.0
(DEL) Alkane: C...	10.458	8.8	ug/L	97389	2	10.055	552959	50.0
(DEL) Alkane: C...	10.586	19.7	ug/L	217543	2	10.055	552959	50.0
(DEL) Alkane: S...	10.781	35.7	ug/L	394353	2	10.055	552959	50.0
(DEL) Alkane: C...	10.860	27.6	ug/L	305461	2	10.055	552959	50.0
(DEL) Alkane: S...	10.896	15.8	ug/L	174378	2	10.055	552959	50.0
(DEL) Alkane: S...	11.085	9.6	ug/L	146710	3	12.024	765615	50.0
Sulfurous acid,...	11.110	13.1	ug/L	200319	3	12.024	765615	50.0
Benzene, propyl-	11.305	27.8	ug/L	425259	3	12.024	765615	50.0
Benzene, 1-ethy...	11.378	128.1	ug/L	1962120	3	12.024	765615	50.0
Benzene, 1,2,3-...	11.616	43.7	ug/L	668562	3	12.024	765615	50.0
(DEL) Alkane: S...	12.177	9.3	ug/L	142297	3	12.024	765615	50.0
Benzene, 1-ethe...	12.244	29.6	ug/L	453220	3	12.024	765615	50.0
Benzene, 1-meth...	12.268	27.3	ug/L	417534	3	12.024	765615	50.0
Carbonic acid, ...	12.427	32.1	ug/L	491362	3	12.024	765615	50.0
Benzene, 1-meth...	12.463	17.2	ug/L	263510	3	12.024	765615	50.0
Benzene, 2-ethy...	12.536	18.9	ug/L	289994	3	12.024	765615	50.0
p-Cymene	12.555	18.3	ug/L	279516	3	12.024	765615	50.0
o-Cymene	12.616	26.3	ug/L	402190	3	12.024	765615	50.0
Benzene, 1-ethe...	12.701	30.8	ug/L	472125	3	12.024	765615	50.0
Benzene, 1,2,4,...	12.920	12.9	ug/L	197656	3	12.024	765615	50.0
Benzene, 1,2,3,...	12.963	28.3	ug/L	433410	3	12.024	765615	50.0
Benzene, 1-meth...	13.170	18.3	ug/L	280627	3	12.024	765615	50.0
1-Phenyl-1-butene	13.292	56.9	ug/L	870945	3	12.024	765615	50.0
Naphthalene, 1,...	13.427	14.1	ug/L	215279	3	12.024	765615	50.0
Benzene, 1-ethy...	13.585	16.5	ug/L	252302	3	12.024	765615	50.0
1H-Indene, 2,3-...	13.664	12.3	ug/L	187933	3	12.024	765615	50.0
Azulene	13.780	34.4	ug/L	526566	3	12.024	765615	50.0
1H-Indene, 2,3-...	14.219	14.8	ug/L	227082	3	12.024	765615	50.0
Naphthalene, 1-...	14.640	59.1	ug/L	904869	3	12.024	765615	50.0
Naphthalene, 2-...	14.780	32.9	ug/L	503316	3	12.024	765615	50.0
Naphthalene, 1,...	15.481	18.3	ug/L	279942	3	12.024	765615	50.0

Naphthalene, 2,...	15.615	16.5	ug/L	252131	3	12.024	765615	50.0
Naphthalene, 2,...	15.652	11.3	ug/L	172870	3	12.024	765615	50.0

Instrument :
MSVOA_X
ClientSampleId :
GBCC1