LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092922\
Data File : VX031792.D
Acq On : 29 Sep 2022 23:39
Operator : JC/MD
Sample : N4864-10 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 35 Sample Multiplier: 1

**Instrument :**
MSVOA_X
**ClientSampleId :**
MW-44

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\82X092622W.M
Title : SW846 8260

Signal : TIC: VX031792.D\data.ms

| peak # | R.T. min | first scan | max scan | last scan | PK TY | peak height | corr. area | corr. % max. | % of total |
|---|---|---|---|---|---|---|---|---|---|
| 1 | 5.385 | 690 | 705 | 719 | rBV | 78430 | 228767 | 6.97% | 1.331% |
| 2 | 5.556 | 722 | 733 | 753 | rVB | 78153 | 231314 | 7.05% | 1.346% |
| 3 | 5.958 | 788 | 799 | 812 | rBV | 78183 | 212927 | 6.49% | 1.239% |
| 4 | 6.763 | 922 | 931 | 940 | rBV | 108459 | 261770 | 7.97% | 1.523% |
| 5 | 8.110 | 1145 | 1152 | 1161 | rBV2 | 14636 | 34583 | 1.05% | 0.201% |
| 6 | 8.647 | 1226 | 1240 | 1249 | rBV | 388631 | 627642 | 19.12% | 3.652% |
| 7 | 10.055 | 1466 | 1471 | 1481 | rBV | 282471 | 375658 | 11.44% | 2.186% |
| 8 | 10.195 | 1488 | 1494 | 1502 | rVB | 1208819 | 1486066 | 45.27% | 8.647% |
| 9 | 10.299 | 1502 | 1511 | 1519 | rBV | 855002 | 1168732 | 35.60% | 6.801% |
| 10 | 10.640 | 1562 | 1567 | 1580 | rVB | 47755 | 65576 | 2.00% | 0.382% |
| 11 | 10.963 | 1615 | 1620 | 1627 | rBV | 76422 | 94538 | 2.88% | 0.550% |
| 12 | 11.079 | 1634 | 1639 | 1651 | rBV | 355815 | 460374 | 14.02% | 2.679% |
| 13 | 11.305 | 1671 | 1676 | 1683 | rBV | 312952 | 365295 | 11.13% | 2.126% |
| 14 | 11.378 | 1683 | 1688 | 1690 | rVV | 280976 | 393148 | 11.98% | 2.288% |
| 15 | 11.396 | 1690 | 1691 | 1696 | rVV | 322371 | 314751 | 9.59% | 1.831% |
| 16 | 11.451 | 1696 | 1700 | 1712 | rVB | 305545 | 360705 | 10.99% | 2.099% |
| 17 | 11.610 | 1721 | 1726 | 1735 | rBV | 200515 | 264889 | 8.07% | 1.541% |
| 18 | 11.750 | 1744 | 1749 | 1756 | rBV | 2765159 | 3282769 | 100.00% | 19.102% |
| 19 | 12.024 | 1788 | 1794 | 1799 | rBV | 481724 | 606038 | 18.46% | 3.526% |
| 20 | 12.085 | 1799 | 1804 | 1809 | rVV | 663909 | 792054 | 24.13% | 4.609% |
| 21 | 12.244 | 1824 | 1830 | 1837 | rBV2 | 727873 | 969100 | 29.52% | 5.639% |
| 22 | 12.317 | 1837 | 1842 | 1852 | rVB4 | 219967 | 337313 | 10.28% | 1.963% |
| 23 | 12.408 | 1852 | 1857 | 1862 | rBV2 | 53363 | 73059 | 2.23% | 0.425% |
| 24 | 12.463 | 1862 | 1866 | 1872 | rVB | 76515 | 90372 | 2.75% | 0.526% |
| 25 | 12.530 | 1872 | 1877 | 1879 | rBV | 177738 | 217229 | 6.62% | 1.264% |
| 26 | 12.555 | 1879 | 1881 | 1886 | rVB | 147634 | 151785 | 4.62% | 0.883% |
| 27 | 12.616 | 1886 | 1891 | 1894 | rBV | 320637 | 395160 | 12.04% | 2.299% |
| 28 | 12.646 | 1894 | 1896 | 1900 | rVB | 45817 | 46554 | 1.42% | 0.271% |
| 29 | 12.701 | 1900 | 1905 | 1913 | rBV2 | 140627 | 212380 | 6.47% | 1.236% |
| 30 | 12.847 | 1924 | 1929 | 1934 | rVB | 61765 | 77020 | 2.35% | 0.448% |
| 31 | 12.920 | 1934 | 1941 | 1944 | rBV | 184506 | 214826 | 6.54% | 1.250% |
| 32 | 12.963 | 1944 | 1948 | 1953 | rVB | 243807 | 284929 | 8.68% | 1.658% |
| 33 | 13.024 | 1953 | 1958 | 1960 | rBV | 26488 | 34979 | 1.07% | 0.204% |
| 34 | 13.170 | 1977 | 1982 | 1986 | rVB | 167820 | 200315 | 6.10% | 1.166% |
| 35 | 13.256 | 1992 | 1996 | 1998 | rBV2 | 25382 | 37650 | 1.15% | 0.219% |
| 36 | 13.292 | 1998 | 2002 | 2007 | rVB2 | 356281 | 464125 | 14.14% | 2.701% |

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092922\
Data File : VX031792.D
Acq On : 29 Sep 2022 23:39
Operator : JC/MD
Sample : N4864-10 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 35 Sample Multiplier: 1

**Instrument :**
MSVOA_X
**ClientSampleId :**
MW-44

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\82X092622W.M
Title : SW846 8260

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 37 | 13.426 | 2019 | 2024 | 2029 | rVB | 55261 | 71028 | 2.16% | 0.413% |
| 38 | 13.542 | 2039 | 2043 | 2046 | rVV | 44334 | 62605 | 1.91% | 0.364% |
| 39 | 13.579 | 2046 | 2049 | 2056 | rVV3 | 53743 | 104569 | 3.19% | 0.608% |
| 40 | 13.664 | 2060 | 2063 | 2069 | rVB2 | 39763 | 64701 | 1.97% | 0.376% |
| 41 | 13.774 | 2076 | 2081 | 2091 | rBV | 769332 | 932123 | 28.39% | 5.424% |
| 42 | 13.865 | 2091 | 2096 | 2101 | rVB | 41436 | 54873 | 1.67% | 0.319% |
| 43 | 14.073 | 2125 | 2130 | 2136 | rBV | 36106 | 47008 | 1.43% | 0.274% |
| 44 | 14.213 | 2149 | 2153 | 2159 | rBV | 31331 | 48678 | 1.48% | 0.283% |
| 45 | 14.640 | 2218 | 2223 | 2233 | rVB | 172736 | 229508 | 6.99% | 1.335% |
| 46 | 14.780 | 2241 | 2246 | 2255 | rBV | 100196 | 136413 | 4.16% | 0.794% |

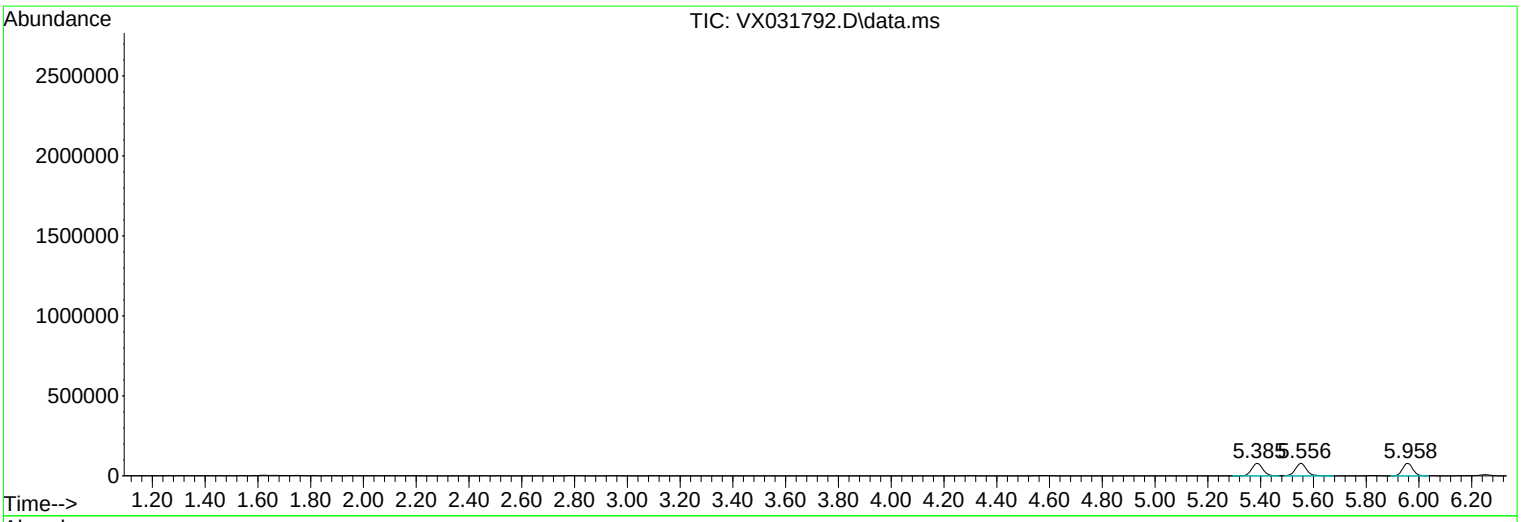
Sum of corrected areas: 17185898

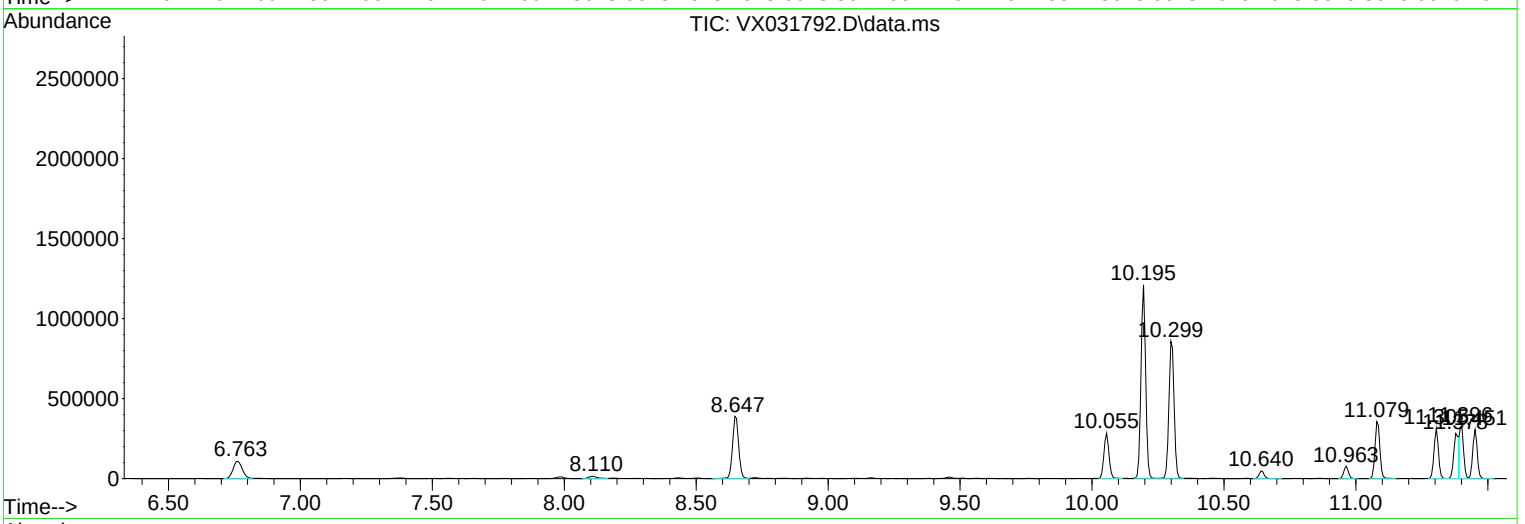
Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092922\
Data File : VX031792.D
Acq On : 29 Sep 2022 23:39
Operator : JC/MD
Sample : N4864-10 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 35 Sample Multiplier: 1

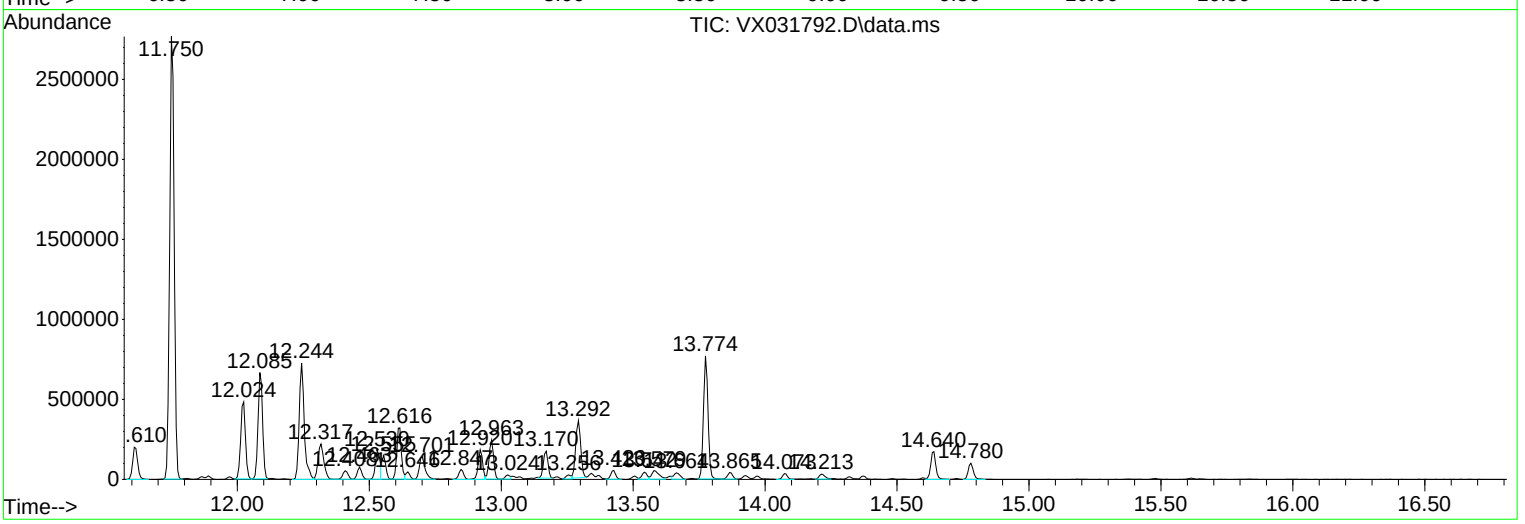
**Instrument :**
MSVOA_X
**ClientSampleId :**
MW-44

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X092622W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092922\
Data File : VX031792.D
Acq On : 29 Sep 2022 23:39
Operator : JC/MD
Sample : N4864-10 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 35 Sample Multiplier: 1

**Instrument :**
MSVOA_X
**ClientSampleId :**
MW-44

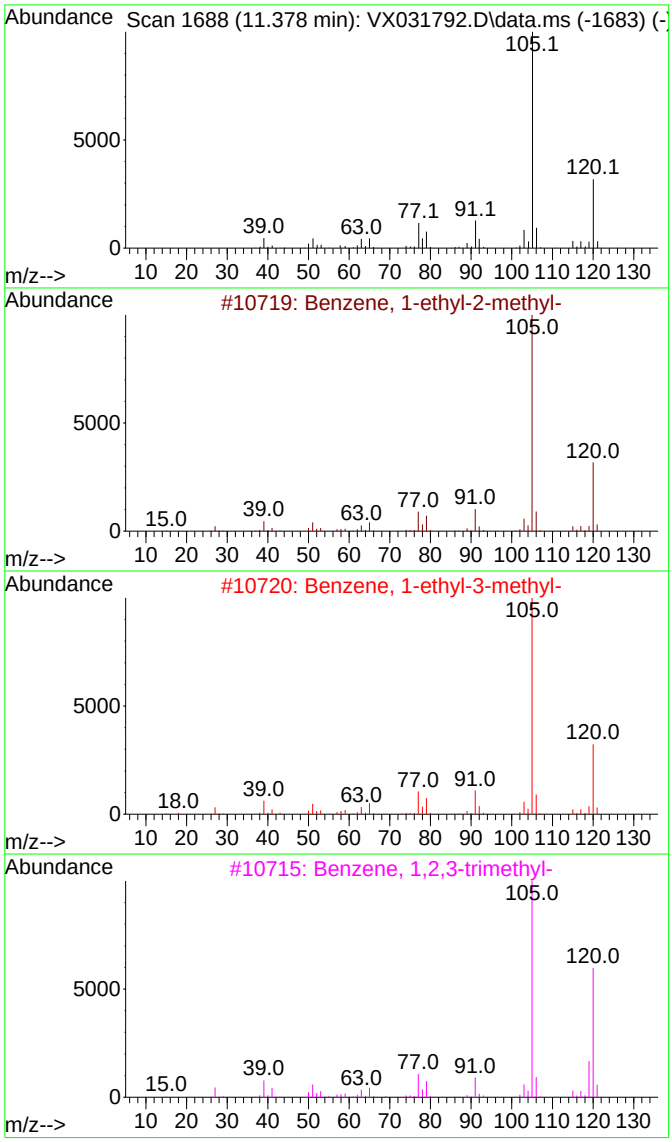
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X092622W.M
Quant Title : SW846 8260

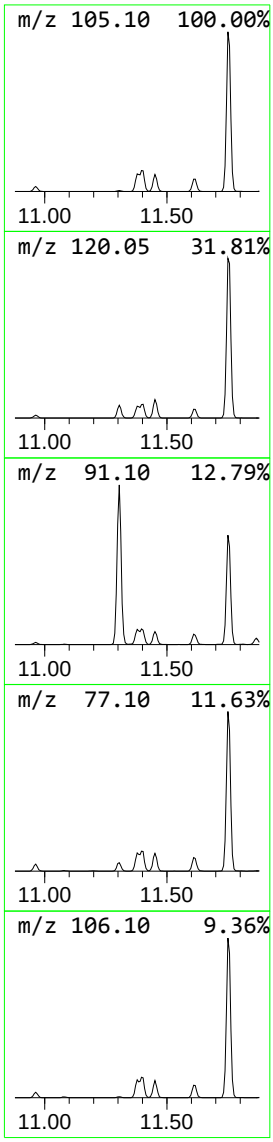
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 1 Benzene, 1-ethyl-2-methyl- Concentration Rank 5

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.378 | 32.44 ug/l | 393148 | 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-3-methyl- | 120 | C9H12 | 000620-14-4 | 95 |
| 3 | Benzene, 1,2,3-trimethyl- | 120 | C9H12 | 000526-73-8 | 91 |
| 4 | Benzene, 1-ethyl-4-methyl- | 120 | C9H12 | 000622-96-8 | 91 |
| 5 | Mesitylene | 120 | C9H12 | 000108-67-8 | 91 |



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092922\
Data File : VX031792.D
Acq On : 29 Sep 2022 23:39
Operator : JC/MD
Sample : N4864-10 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 35 Sample Multiplier: 1

**Instrument :**
MSVOA_X
**ClientSampleId :**
MW-44

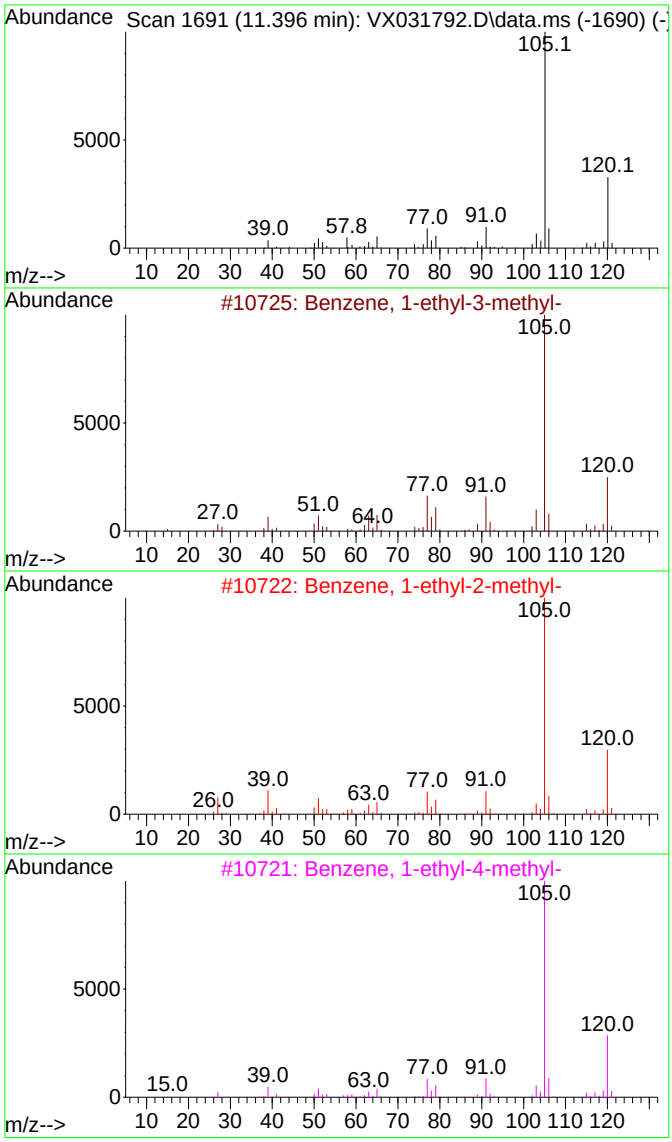
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X092622W.M
Quant Title : SW846 8260

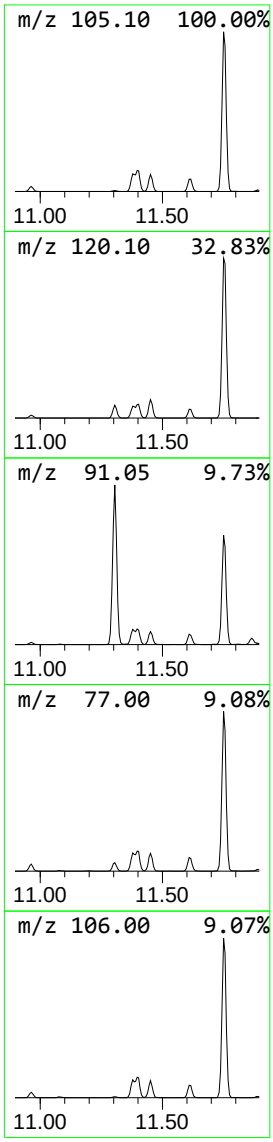
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 2 Benzene, 1-ethyl-3-methyl- Concentration Rank 6

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.396 | 25.97 ug/l | 314751 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, 1-ethyl-3-methyl- | 120 | C9H12 | 000620-14-4 | 91 |
| 2 | Benzene, 1-ethyl-2-methyl- | 120 | C9H12 | 000611-14-3 | 91 |
| 3 | Benzene, 1-ethyl-4-methyl- | 120 | C9H12 | 000622-96-8 | 91 |
| 4 | Mesitylene | 120 | C9H12 | 000108-67-8 | 91 |
| 5 | Benzene, 1,2,4-trimethyl- | 120 | C9H12 | 000095-63-6 | 91 |



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092922\
Data File : VX031792.D
Acq On : 29 Sep 2022 23:39
Operator : JC/MD
Sample : N4864-10 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 35 Sample Multiplier: 1

**Instrument :**
MSVOA_X
**ClientSampleId :**
MW-44

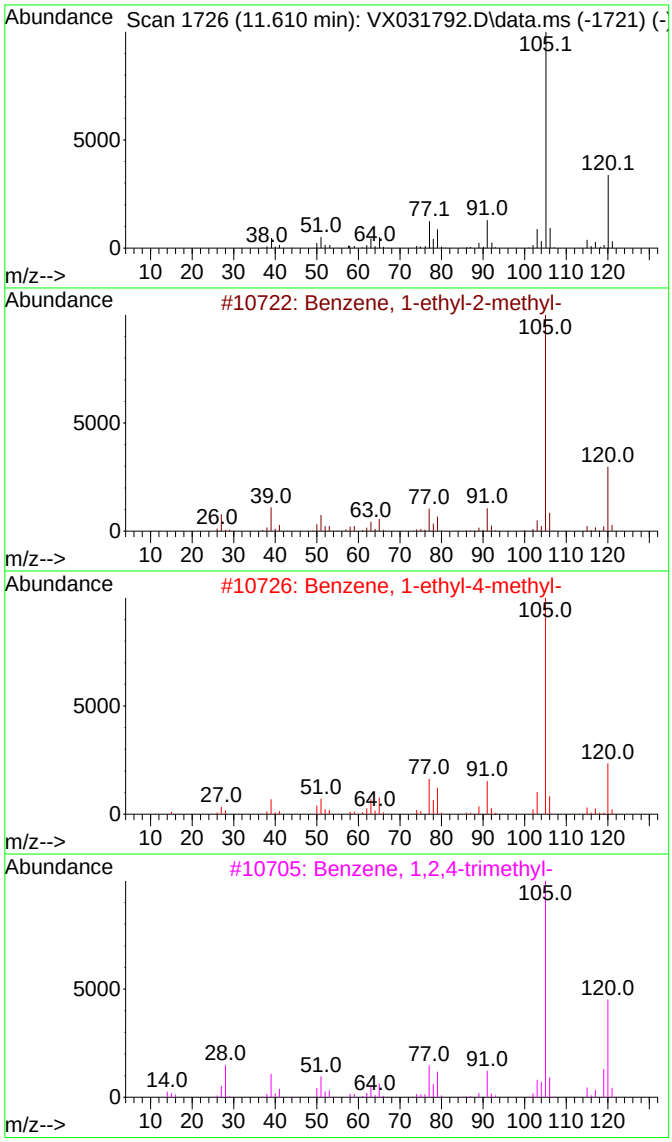
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X092622W.M
Quant Title : SW846 8260

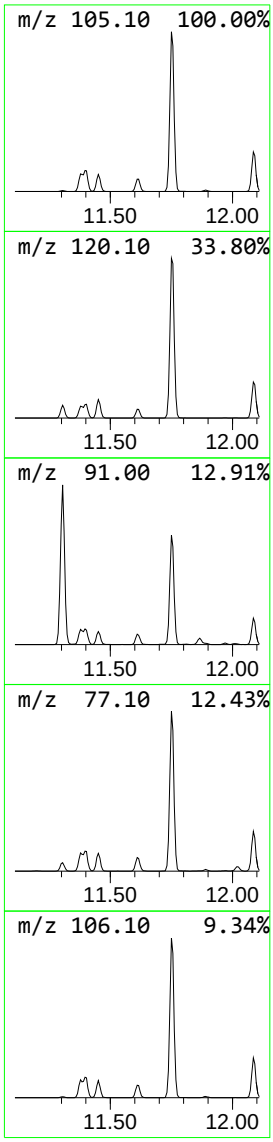
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 3 Benzene, 1-ethyl-4-methyl- Concentration Rank 8

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.610 | 21.85 ug/l | 264889 | 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 | 93 |
| 2 | Benzene, 1-ethyl-4-methyl- | 120 | C9H12 | 000622-96-8 | 93 |
| 3 | Benzene, 1,2,4-trimethyl- | 120 | C9H12 | 000095-63-6 | 90 |
| 4 | Benzene, 1-ethyl-3-methyl- | 120 | C9H12 | 000620-14-4 | 90 |
| 5 | Mesitylene | 120 | C9H12 | 000108-67-8 | 90 |



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092922\
Data File : VX031792.D
Acq On : 29 Sep 2022 23:39
Operator : JC/MD
Sample : N4864-10 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 35 Sample Multiplier: 1

**Instrument :**
MSVOA_X
**ClientSampleId :**
MW-44

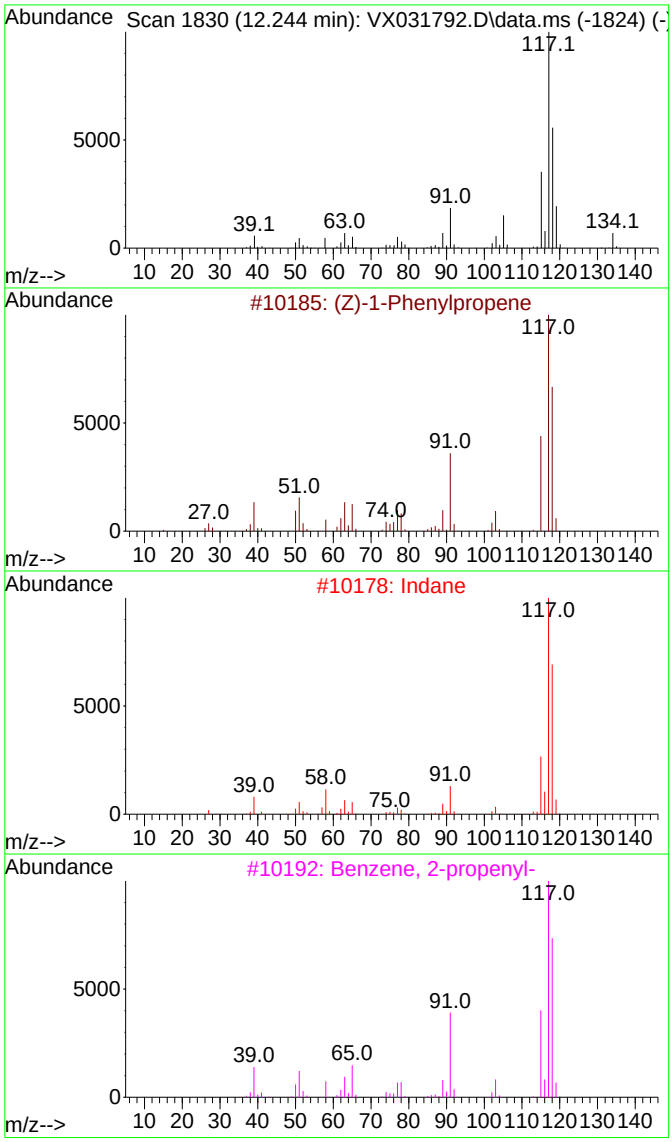
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X092622W.M
Quant Title : SW846 8260

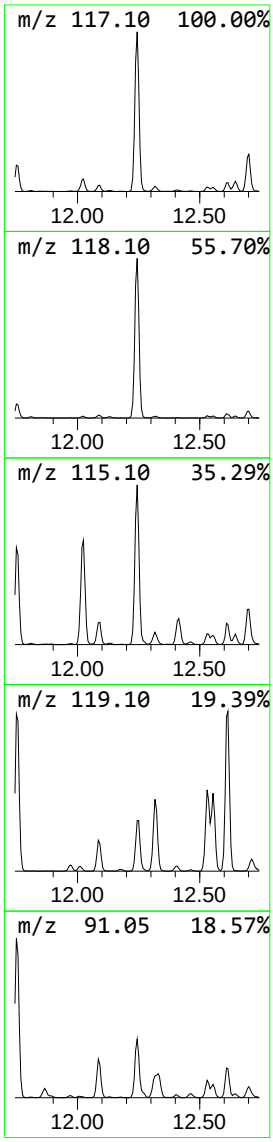
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 5 (Z)-1-Phenylpropene Concentration Rank 1

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.244 | 79.95 ug/l | 969100 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | (Z)-1-Phenylpropene | 118 | C9H10 | 000766-90-5 | 70 |
| 2 | Indane | 118 | C9H10 | 000496-11-7 | 70 |
| 3 | Benzene, 2-propenyl- | 118 | C9H10 | 000300-57-2 | 68 |
| 4 | Benzene, cyclopropyl- | 118 | C9H10 | 000873-49-4 | 62 |
| 5 | Tetracyclo[3.3.1.0(2,8).0(4,6)]-... | 118 | C9H10 | 1000191-13-7 | 53 |



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092922\
Data File : VX031792.D
Acq On : 29 Sep 2022 23:39
Operator : JC/MD
Sample : N4864-10 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 35 Sample Multiplier: 1

**Instrument :**
MSVOA_X
**ClientSampleId :**
MW-44

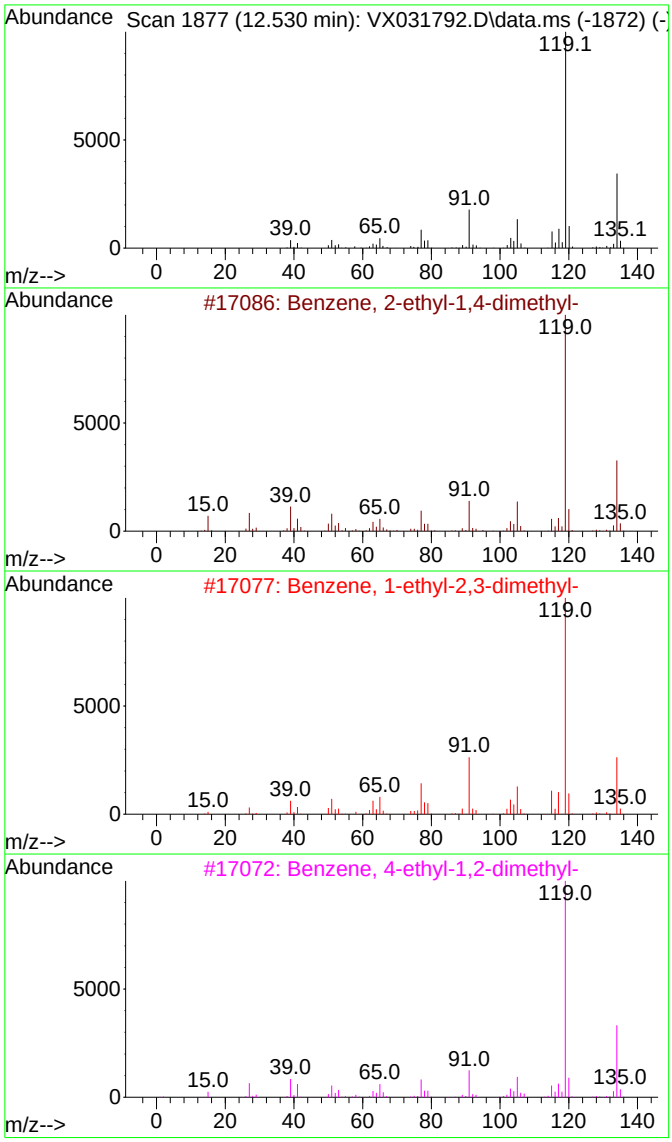
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X092622W.M
Quant Title : SW846 8260

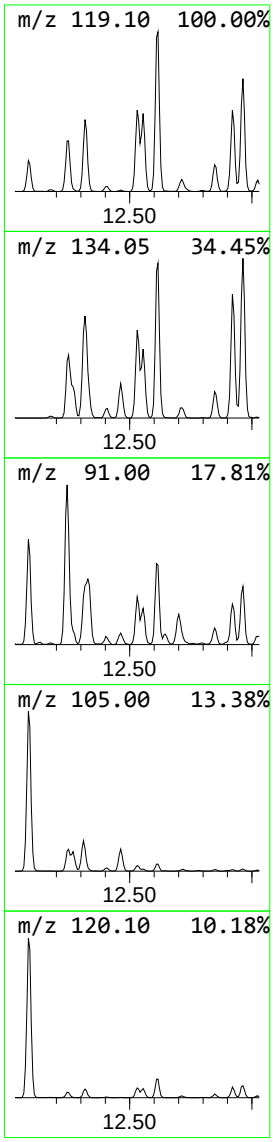
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 6 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 10

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.530 | 17.92 ug/l | 217229 | 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, 1-ethyl-2,3-dimethyl- | 134 | C10H14 | 000933-98-2 | 96 |
| 3 | Benzene, 4-ethyl-1,2-dimethyl- | 134 | C10H14 | 000934-80-5 | 95 |
| 4 | Benzene, 2-ethyl-1,3-dimethyl- | 134 | C10H14 | 002870-04-4 | 94 |
| 5 | Benzene, 1-methyl-3-(1-methyleth... | 134 | C10H14 | 000535-77-3 | 93 |



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092922\
Data File : VX031792.D
Acq On : 29 Sep 2022 23:39
Operator : JC/MD
Sample : N4864-10 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 35 Sample Multiplier: 1

**Instrument :**
MSVOA_X
**ClientSampleId :**
MW-44

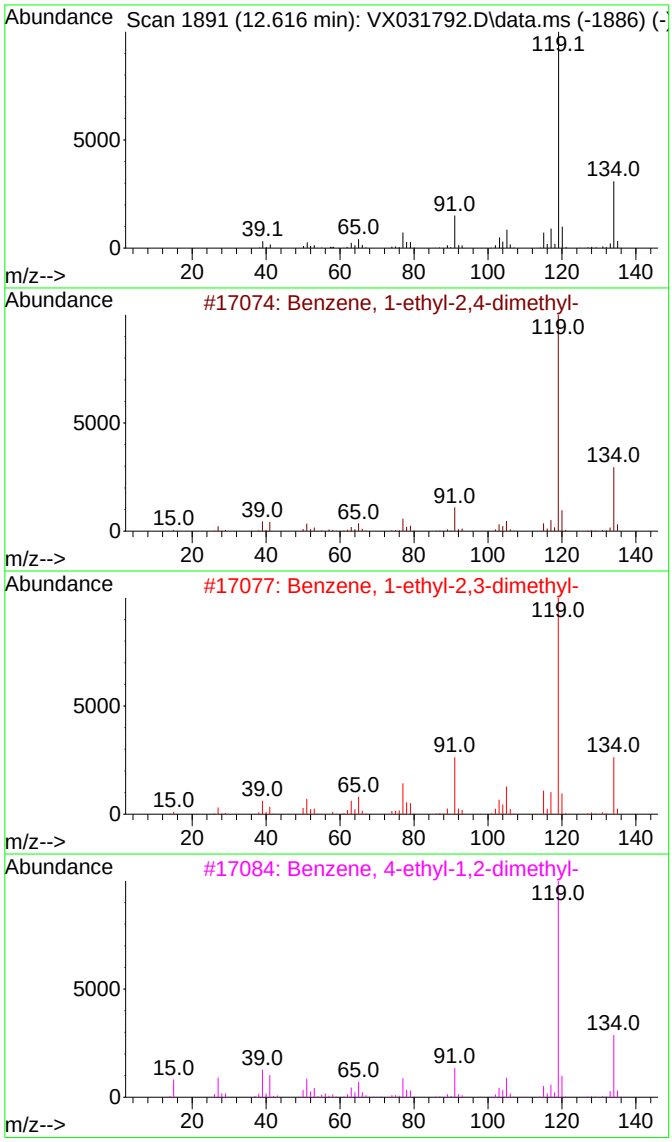
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X092622W.M
Quant Title : SW846 8260

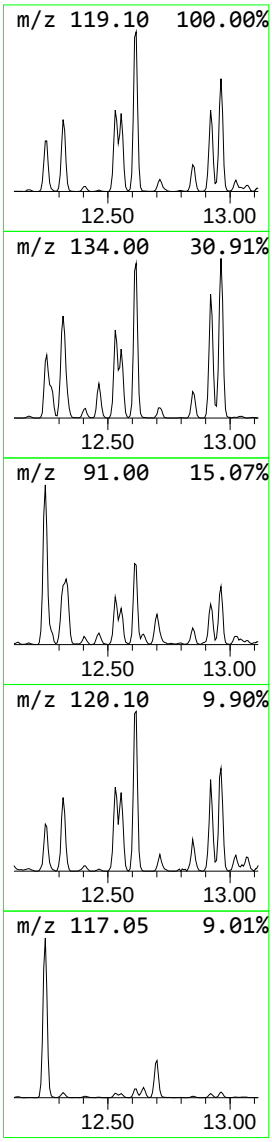
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 7 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 4

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.616 | 32.60 ug/l | 395160 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, 1-ethyl-2,4-dimethyl- | 134 | C10H14 | 000874-41-9 | 97 |
| 2 | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14 | 000933-98-2 | 96 |
| 3 | Benzene, 4-ethyl-1,2-dimethyl- | 134 | C10H14 | 000934-80-5 | 95 |
| 4 | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14 | 001758-88-9 | 95 |
| 5 | Benzene, 1-methyl-3-(1-methyleth... | 134 | C10H14 | 000535-77-3 | 95 |



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092922\
Data File : VX031792.D
Acq On : 29 Sep 2022 23:39
Operator : JC/MD
Sample : N4864-10 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 35 Sample Multiplier: 1

**Instrument :**
MSVOA_X
**ClientSampleId :**
MW-44

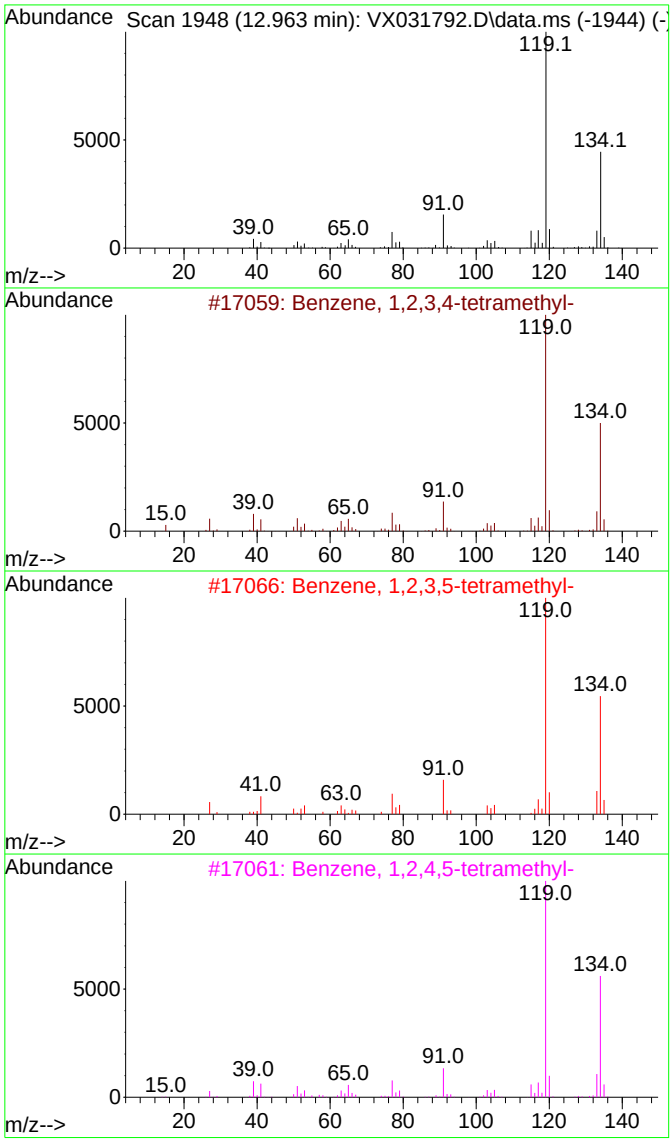
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X092622W.M
Quant Title : SW846 8260

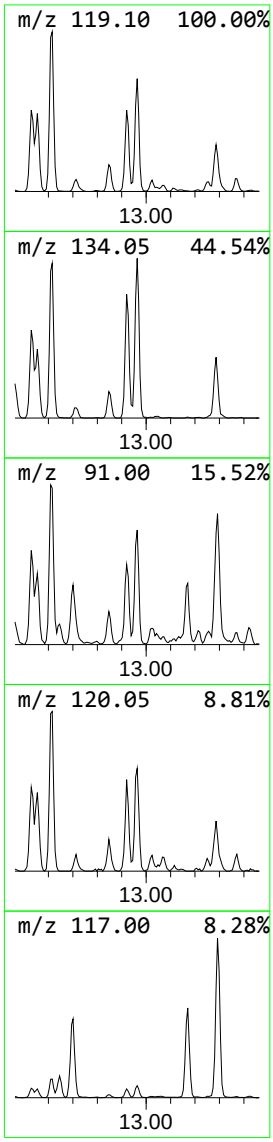
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 8 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 7

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.963 | 23.51 ug/l | 284929 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, 1,2,3,4-tetramethyl- | 134 | C10H14 | 000488-23-3 | 97 |
| 2 | Benzene, 1,2,3,5-tetramethyl- | 134 | C10H14 | 000527-53-7 | 96 |
| 3 | Benzene, 1,2,4,5-tetramethyl- | 134 | C10H14 | 000095-93-2 | 95 |
| 4 | o-Cymene | 134 | C10H14 | 000527-84-4 | 95 |
| 5 | Benzene, 1-ethyl-2,4-dimethyl- | 134 | C10H14 | 000874-41-9 | 94 |



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092922\
Data File : VX031792.D
Acq On : 29 Sep 2022 23:39
Operator : JC/MD
Sample : N4864-10 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 35 Sample Multiplier: 1

**Instrument :**
MSVOA_X
**ClientSampleId :**
MW-44

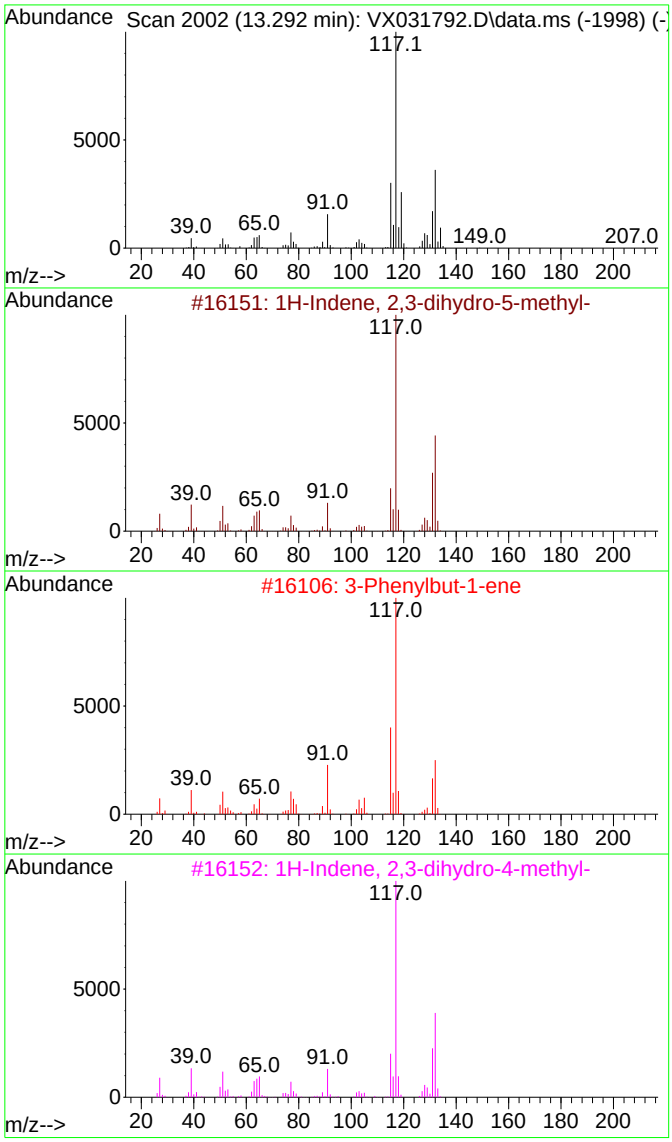
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X092622W.M
Quant Title : SW846 8260

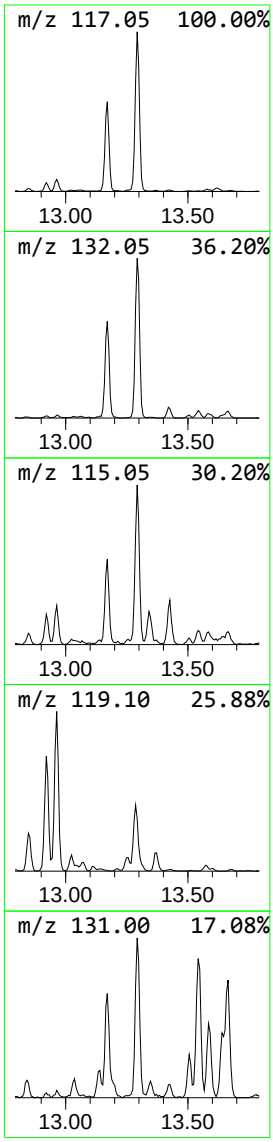
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 9 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 3

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.292 | 38.29 ug/l | 464125 | 1,4-Dichlorobenzene-d4 | 12.024 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1H-Indene, 2,3-dihydro-5-methyl- | 132 | C10H12 | 000874-35-1 | 89 |
| 2 | 3-Phenylbut-1-ene | 132 | C10H12 | 000934-10-1 | 81 |
| 3 | 1H-Indene, 2,3-dihydro-4-methyl- | 132 | C10H12 | 000824-22-6 | 76 |
| 4 | Benzene, 1-ethenyl-4-ethyl- | 132 | C10H12 | 003454-07-7 | 76 |
| 5 | Indan, 1-methyl- | 132 | C10H12 | 000767-58-8 | 76 |



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092922\
Data File : VX031792.D
Acq On : 29 Sep 2022 23:39
Operator : JC/MD
Sample : N4864-10 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 35 Sample Multiplier: 1

**Instrument :**
MSVOA_X
**ClientSampleId :**
MW-44

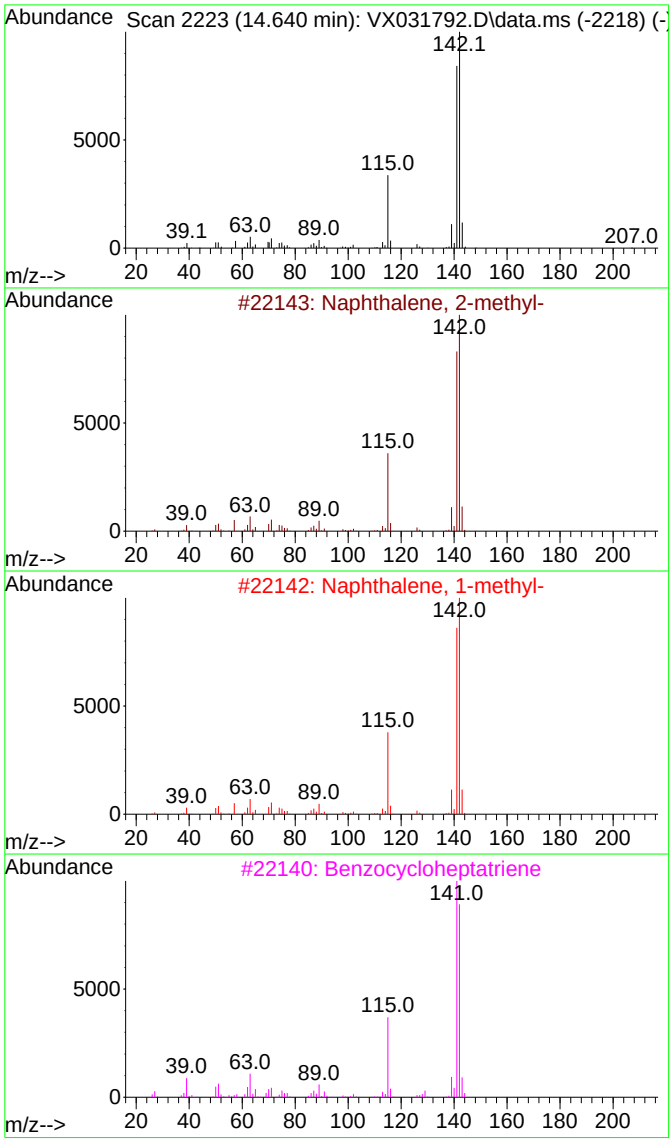
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X092622W.M
Quant Title : SW846 8260

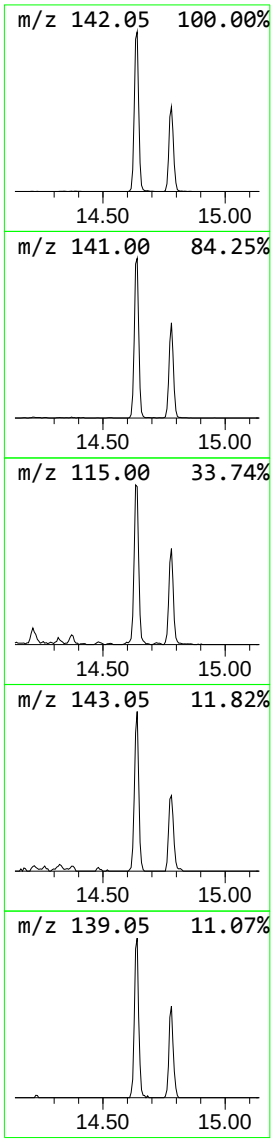
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.640 | 18.94 ug/l | 229508 | 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 | Benzocycloheptatriene | 142 | C11H10 | 000264-09-5 | 91 |
| 4 | 1,4-Methanonaphthalene, 1,4-dihy... | 142 | C11H10 | 004453-90-1 | 91 |
| 5 | 1H-Indene, 1-ethylidene- | 142 | C11H10 | 002471-83-2 | 64 |



Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX092922\
Data File : VX031792.D
Acq On : 29 Sep 2022 23:39
Operator : JC/MD
Sample : N4864-10 10X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 35 Sample Multiplier: 1

**Instrument :**
MSVOA_X
**ClientSampleId :**
MW-44

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X092622W.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 |
| Benzene, 1-ethy... | 11.378 | 32.4 | ug/l | 393148 | 4 | 12.024 | 606038 | 50.0 |
| Benzene, 1-ethy... | 11.396 | 26.0 | ug/l | 314751 | 4 | 12.024 | 606038 | 50.0 |
| Benzene, 1-ethy... | 11.610 | 21.9 | ug/l | 264889 | 4 | 12.024 | 606038 | 50.0 |
| (Z)-1-Phenylpro... | 12.244 | 80.0 | ug/l | 969100 | 4 | 12.024 | 606038 | 50.0 |
| Benzene, 2-ethy... | 12.530 | 17.9 | ug/l | 217229 | 4 | 12.024 | 606038 | 50.0 |
| Benzene, 1-ethy... | 12.616 | 32.6 | ug/l | 395160 | 4 | 12.024 | 606038 | 50.0 |
| Benzene, 1,2,3,... | 12.963 | 23.5 | ug/l | 284929 | 4 | 12.024 | 606038 | 50.0 |
| 1H-Indene, 2,3-... | 13.292 | 38.3 | ug/l | 464125 | 4 | 12.024 | 606038 | 50.0 |
| Naphthalene, 1-... | 14.640 | 18.9 | ug/l | 229508 | 4 | 12.024 | 606038 | 50.0 |