LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN072618\
Data File : VN050085.D
Acq On : 26 Jul 2018 16:26
Operator : MD\SY
Sample : J4142-02
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 14 Sample Multiplier: 1

**Instrument :** MSVOA_N
**ClientSampleId :** MW-5

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N072418W.M
Title : SW846 8260

Signal : TIC

| peak # | R.T. min | first scan | max scan | last scan | PK TY | peak height | corr. area | corr. % max. | % of total |
|---|---|---|---|---|---|---|---|---|---|
| 1 | 1.661 | 3 | 22 | 49 | rBV | 2797729 | 4365565 | 100.00% | 12.914% |
| 2 | 7.596 | 1847 | 1868 | 1879 | rBV | 549334 | 1382228 | 31.66% | 4.089% |
| 3 | 7.670 | 1879 | 1891 | 1917 | rVB | 779373 | 1872792 | 42.90% | 5.540% |
| 4 | 8.030 | 1985 | 2003 | 2029 | rBV | 557433 | 1308186 | 29.97% | 3.870% |
| 5 | 8.252 | 2040 | 2072 | 2075 | rBV2 | 133543 | 512691 | 11.74% | 1.517% |
| 6 | 8.590 | 2160 | 2177 | 2209 | rBV | 1176558 | 2489049 | 57.02% | 7.363% |
| 7 | 10.095 | 2630 | 2645 | 2686 | rBV | 2086994 | 3884333 | 88.98% | 11.491% |
| 8 | 11.410 | 3039 | 3054 | 3075 | rBV | 1545528 | 2752217 | 63.04% | 8.142% |
| 9 | 12.252 | 3307 | 3316 | 3326 | rVB | 47362 | 82106 | 1.88% | 0.243% |
| 10 | 12.403 | 3350 | 3363 | 3377 | rBV | 1361093 | 2332776 | 53.44% | 6.901% |
| 11 | 12.725 | 3455 | 3463 | 3475 | rVB | 22361 | 47122 | 1.08% | 0.139% |
| 12 | 13.175 | 3592 | 3603 | 3610 | rBV2 | 48296 | 87728 | 2.01% | 0.260% |
| 13 | 13.345 | 3644 | 3656 | 3671 | rBV | 1450617 | 2401267 | 55.00% | 7.104% |
| 14 | 13.445 | 3680 | 3687 | 3698 | rVB | 69926 | 104963 | 2.40% | 0.311% |
| 15 | 13.512 | 3700 | 3708 | 3714 | rBV | 171115 | 270026 | 6.19% | 0.799% |
| 16 | 13.596 | 3728 | 3734 | 3746 | rVB2 | 44346 | 80197 | 1.84% | 0.237% |
| 17 | 13.673 | 3747 | 3758 | 3770 | rVB2 | 266383 | 463405 | 10.62% | 1.371% |
| 18 | 13.892 | 3816 | 3826 | 3835 | rVB | 352119 | 534791 | 12.25% | 1.582% |
| 19 | 13.950 | 3836 | 3844 | 3847 | rBV | 60043 | 86036 | 1.97% | 0.255% |
| 20 | 13.995 | 3848 | 3858 | 3868 | rVV3 | 519318 | 888243 | 20.35% | 2.628% |
| 21 | 14.049 | 3869 | 3875 | 3890 | rVB8 | 79907 | 193806 | 4.44% | 0.573% |
| 22 | 14.133 | 3894 | 3901 | 3908 | rBV2 | 79369 | 111348 | 2.55% | 0.329% |
| 23 | 14.181 | 3909 | 3916 | 3918 | rBV4 | 70768 | 88678 | 2.03% | 0.262% |
| 24 | 14.213 | 3919 | 3926 | 3934 | rVB2 | 222198 | 329531 | 7.55% | 0.975% |
| 25 | 14.297 | 3944 | 3952 | 3959 | rBV2 | 150525 | 239257 | 5.48% | 0.708% |
| 26 | 14.368 | 3968 | 3974 | 3981 | rVB | 103347 | 132839 | 3.04% | 0.393% |
| 27 | 14.419 | 3981 | 3990 | 4002 | rVB3 | 215556 | 445062 | 10.19% | 1.317% |
| 28 | 14.487 | 4002 | 4011 | 4017 | rBV4 | 102072 | 183878 | 4.21% | 0.544% |
| 29 | 14.528 | 4017 | 4024 | 4032 | rVB2 | 195363 | 284424 | 6.52% | 0.841% |
| 30 | 14.583 | 4032 | 4041 | 4055 | rBV2 | 295057 | 511601 | 11.72% | 1.513% |
| 31 | 14.657 | 4056 | 4064 | 4073 | rBV6 | 123803 | 237432 | 5.44% | 0.702% |
| 32 | 14.786 | 4095 | 4104 | 4110 | rBV3 | 87698 | 169742 | 3.89% | 0.502% |
| 33 | 14.853 | 4117 | 4125 | 4131 | rBV5 | 270042 | 447215 | 10.24% | 1.323% |
| 34 | 14.959 | 4149 | 4158 | 4177 | rVB4 | 471065 | 1031178 | 23.62% | 3.050% |

LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN072618\
Data File : VN050085.D
Acq On : 26 Jul 2018 16:26
Operator : MD\SY
Sample : J4142-02
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 14 Sample Multiplier: 1

**Instrument :** MSVOA_N
**ClientSampleId :** MW-5

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N072418W.M
Title : SW846 8260

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 35 | 15.049 | 4181 | 4186 | 4193 | rVV2 | 63125 | 94077 | 2.15% | 0.278% |
| 36 | 15.094 | 4195 | 4200 | 4213 | rVB5 | 81660 | 147094 | 3.37% | 0.435% |
| 37 | 15.194 | 4222 | 4231 | 4239 | rBV | 269784 | 452743 | 10.37% | 1.339% |
| 38 | 15.265 | 4241 | 4253 | 4270 | rVB3 | 301977 | 808208 | 18.51% | 2.391% |
| 39 | 15.416 | 4292 | 4300 | 4311 | rBV4 | 60897 | 117263 | 2.69% | 0.347% |
| 40 | 15.705 | 4382 | 4390 | 4399 | rVB3 | 113672 | 189853 | 4.35% | 0.562% |
| 41 | 15.917 | 4442 | 4456 | 4468 | rBV2 | 370285 | 798127 | 18.28% | 2.361% |
| 42 | 16.107 | 4506 | 4515 | 4525 | rVB2 | 159189 | 307543 | 7.04% | 0.910% |
| 43 | 16.168 | 4527 | 4534 | 4545 | rBV4 | 93963 | 162388 | 3.72% | 0.480% |
| 44 | 16.358 | 4580 | 4593 | 4603 | rBV5 | 139253 | 374974 | 8.59% | 1.109% |

Sum of corrected areas: 33803982

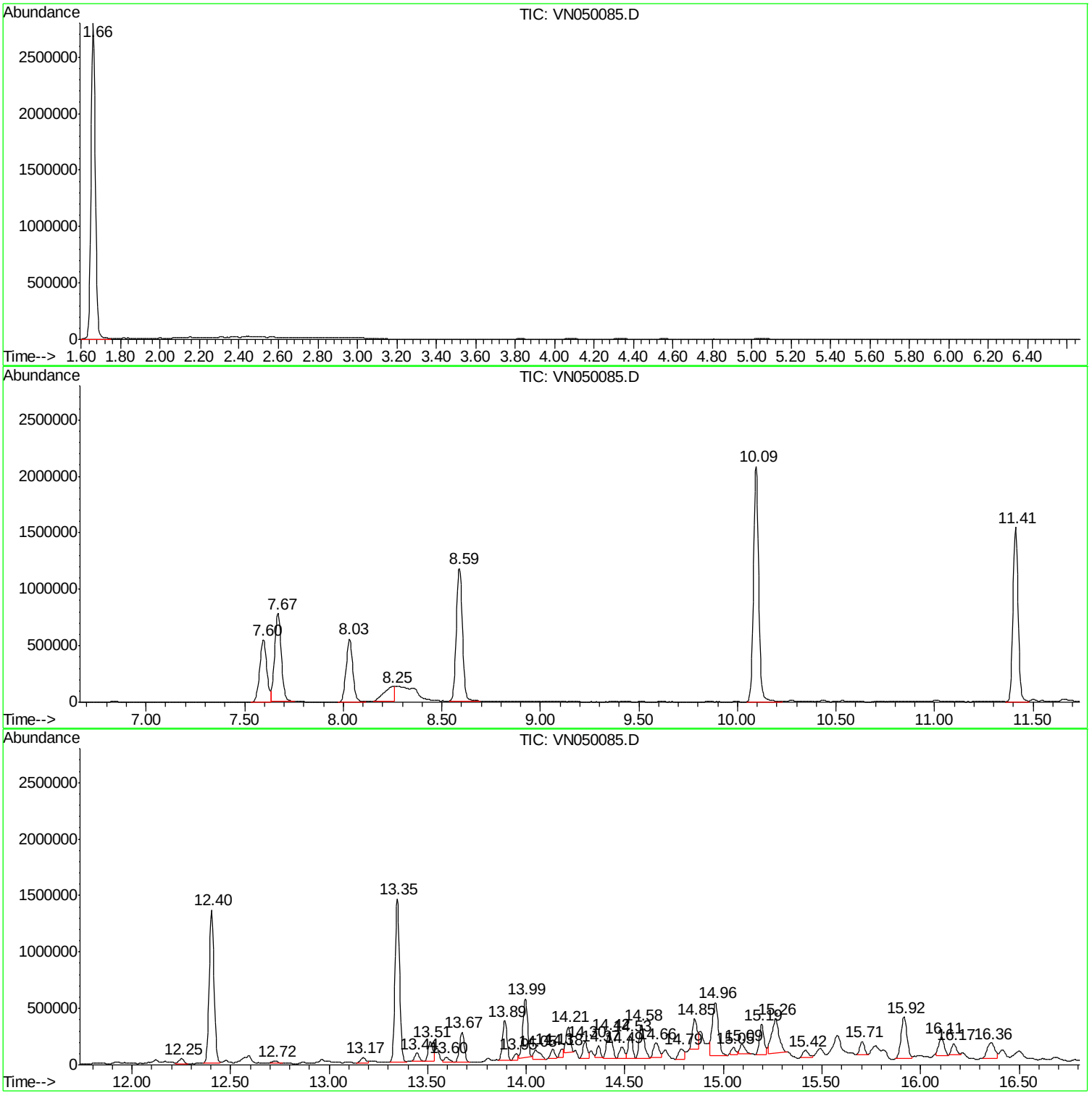
LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN072618\
Data File : VN050085.D
Acq On : 26 Jul 2018 16:26
Operator : MD\SY
Sample : J4142-02
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 14 Sample Multiplier: 1

**Instrument :** MSVOA_N
**ClientSampleId :** MW-5

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N072418W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



# Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN072618\
Data File : VN050085.D
Acq On : 26 Jul 2018 16:26
Operator : MD\SY
Sample : J4142-02
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 14 Sample Multiplier: 1

**Instrument :** MSVOA_N
**ClientSampleId :** MW-5

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N072418W.M
Quant Title : SW846 8260

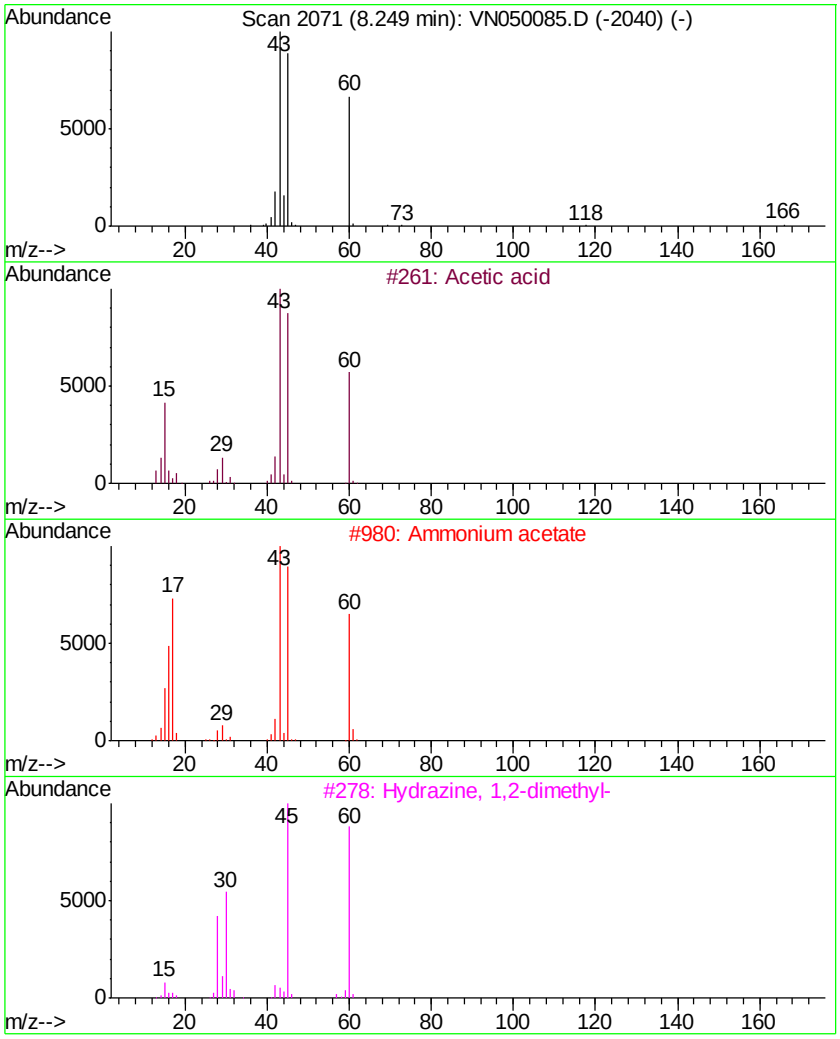
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

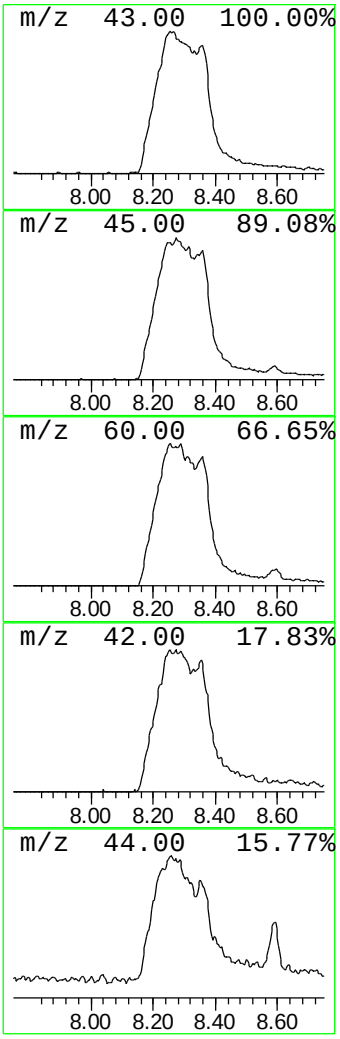
**********

Peak Number 2 Acetic acid Concentration Rank 8

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 8.25 | 10.30 ug/l | 512691 | 1,4-Difluorobenzene | 8.59 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Acetic acid | 60 | C2H4O2 | 000064-19-7 | 90 |
| 2 | Ammonium acetate | 77 | C2H7NO2 | 000631-61-8 | 64 |
| 3 | Hydrazine, 1,2-dimethyl- | 60 | C2H8N2 | 000540-73-8 | 50 |
| 4 | Cyclohexan-1,4,5-triol-3-one-1-c... | 190 | C7H10O6 | 1000128-45-5 | 43 |
| 5 | Hydrazine, 1,1-dimethyl- | 60 | C2H8N2 | 000057-14-7 | 9 |



# Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN072618\
Data File : VN050085.D
Acq On : 26 Jul 2018 16:26
Operator : MD\SY
Sample : J4142-02
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 14 Sample Multiplier: 1

**Instrument :** MSVOA_N
**ClientSampleId :** MW-5

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N072418W.M
Quant Title : SW846 8260

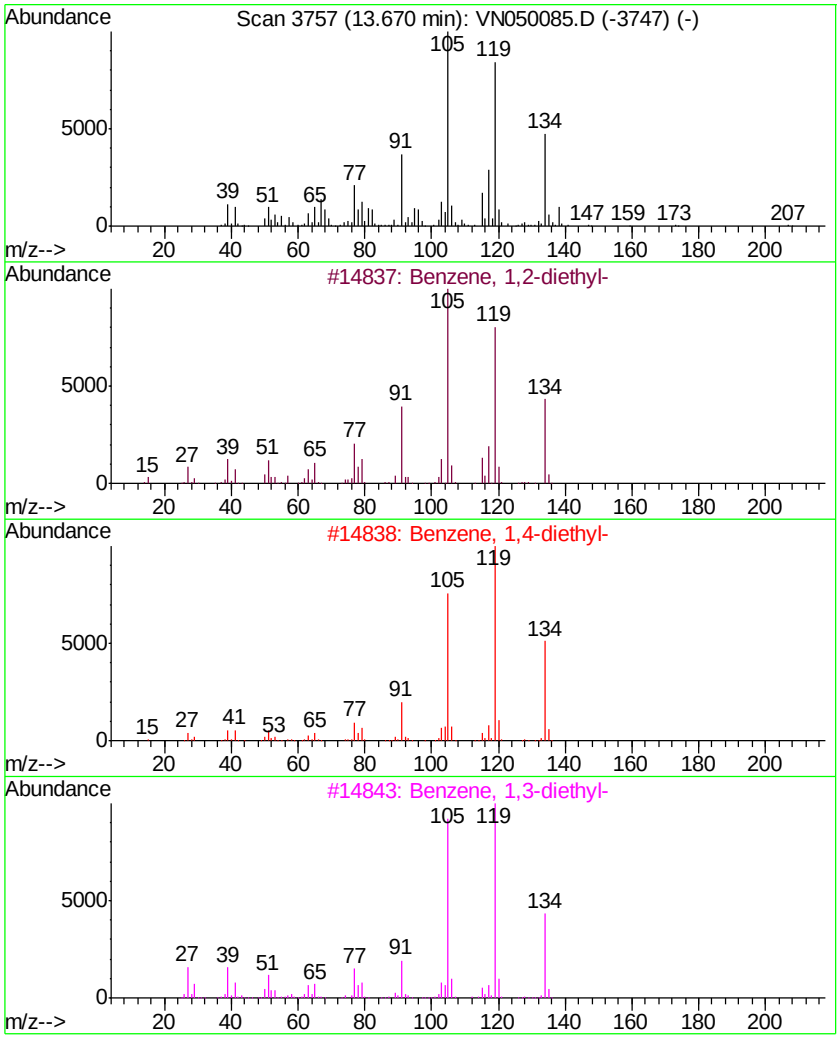
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

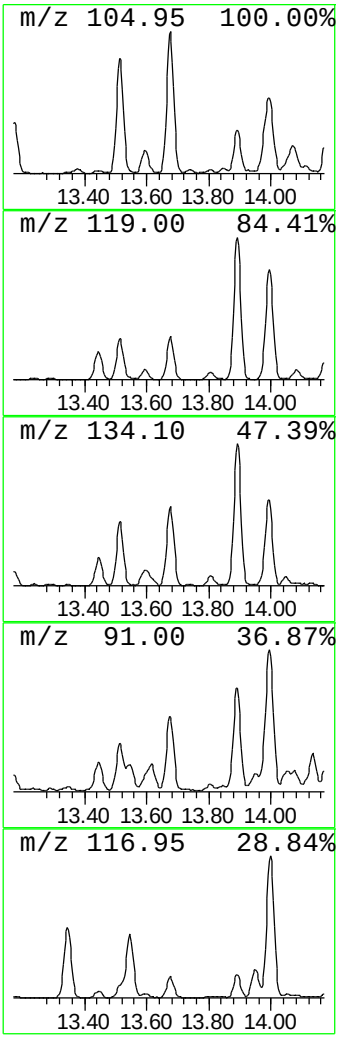
**********

Peak Number 3 Benzene, 1,2-diethyl- Concentration Rank 9

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.67 | 9.65 ug/l | 463405 | 1,4-Dichlorobenzene-d4 | 13.35 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, 1,2-diethyl- | 134 | C10H14 | 000135-01-3 | 96 |
| 2 | Benzene, 1,4-diethyl- | 134 | C10H14 | 000105-05-5 | 90 |
| 3 | Benzene, 1,3-diethyl- | 134 | C10H14 | 000141-93-5 | 87 |
| 4 | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14 | 000933-98-2 | 72 |
| 5 | Benzene, 4-ethyl-1,2-dimethyl- | 134 | C10H14 | 000934-80-5 | 62 |



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN072618\
Data File : VN050085.D
Acq On : 26 Jul 2018 16:26
Operator : MD\SY
Sample : J4142-02
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 14 Sample Multiplier: 1

**Instrument :** MSVOA_N
**ClientSampleId :** MW-5

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N072418W.M
Quant Title : SW846 8260

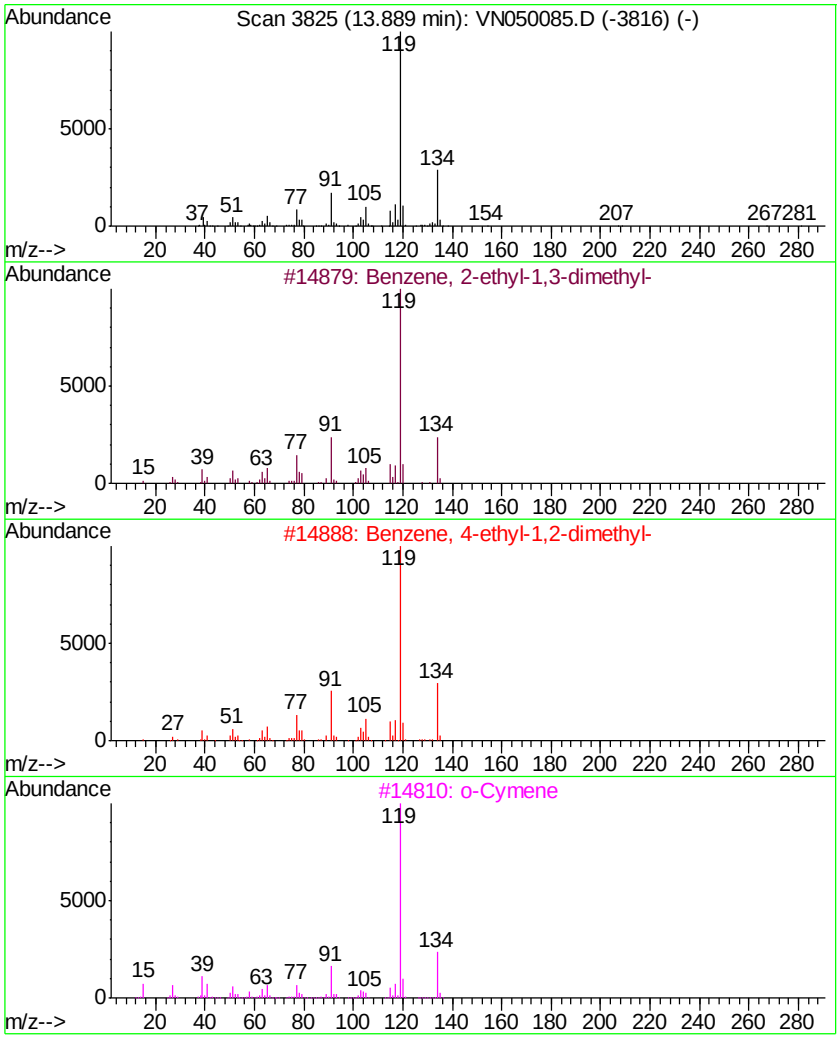
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

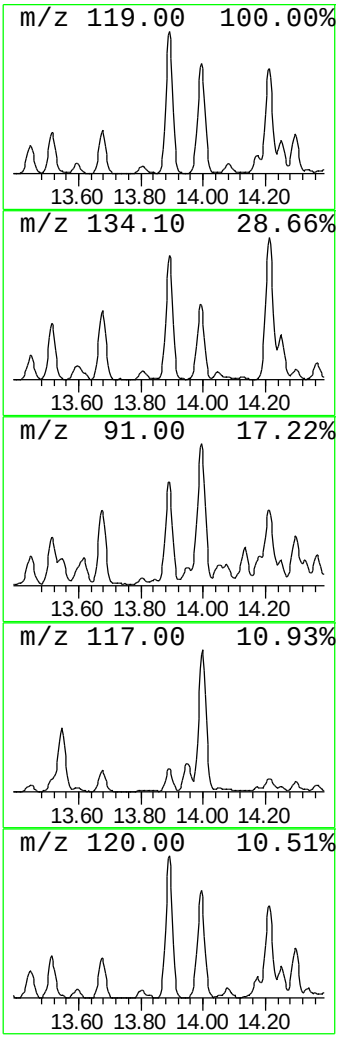
**********

Peak Number 4 Benzene, 2-ethyl-1,3-dimethyl- Concentration Rank 6

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.89 | 11.14 ug/l | 534791 | 1,4-Dichlorobenzene-d4 | 13.35 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, 2-ethyl-1,3-dimethyl- | 134 | C10H14 | 002870-04-4 | 95 |
| 2 | Benzene, 4-ethyl-1,2-dimethyl- | 134 | C10H14 | 000934-80-5 | 95 |
| 3 | o-Cymene | 134 | C10H14 | 000527-84-4 | 94 |
| 4 | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14 | 000933-98-2 | 94 |
| 5 | Benzene, 1-methyl-3-(1-methyleth... | 134 | C10H14 | 000535-77-3 | 93 |



Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN072618\
Data File : VN050085.D
Acq On : 26 Jul 2018 16:26
Operator : MD\SY
Sample : J4142-02
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 14 Sample Multiplier: 1

**Instrument :** MSVOA_N
**ClientSampleId :** MW-5

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N072418W.M
Quant Title : SW846 8260

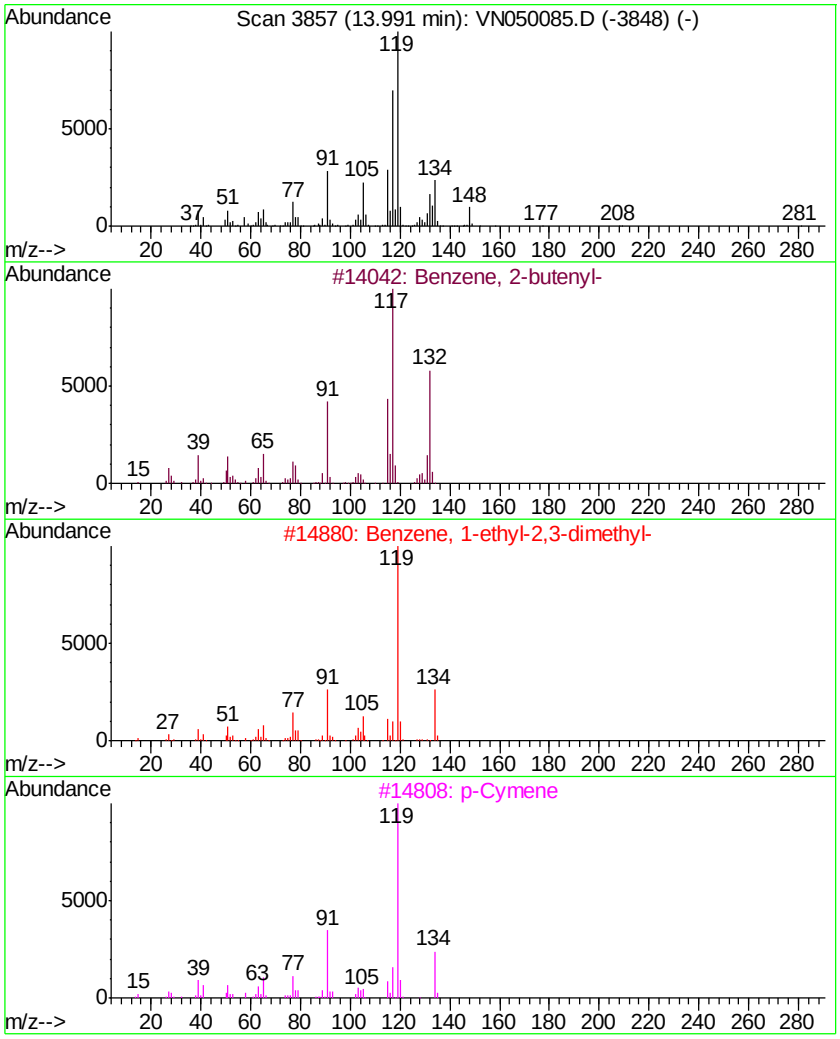
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

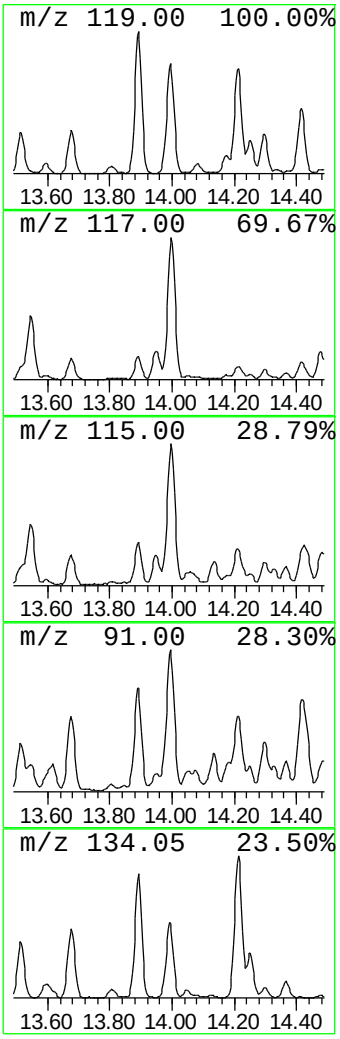
**********

Peak Number 5 Benzene, 2-butenyl- Concentration Rank 3

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.99 | 18.50 ug/l | 888243 | 1,4-Dichlorobenzene-d4 | 13.35 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, 2-butenyl- | 132 | C10H12 | 001560-06-1 | 83 |
| 2 | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14 | 000933-98-2 | 50 |
| 3 | p-Cymene | 134 | C10H14 | 000099-87-6 | 50 |
| 4 | Benzene, 1-methyl-3-(1-methyleth... | 134 | C10H14 | 000535-77-3 | 50 |
| 5 | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14 | 001758-88-9 | 46 |



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN072618\
Data File : VN050085.D
Acq On : 26 Jul 2018 16:26
Operator : MD\SY
Sample : J4142-02
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 14 Sample Multiplier: 1

**Instrument :** MSVOA_N
**ClientSampleId :** MW-5

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N072418W.M
Quant Title : SW846 8260

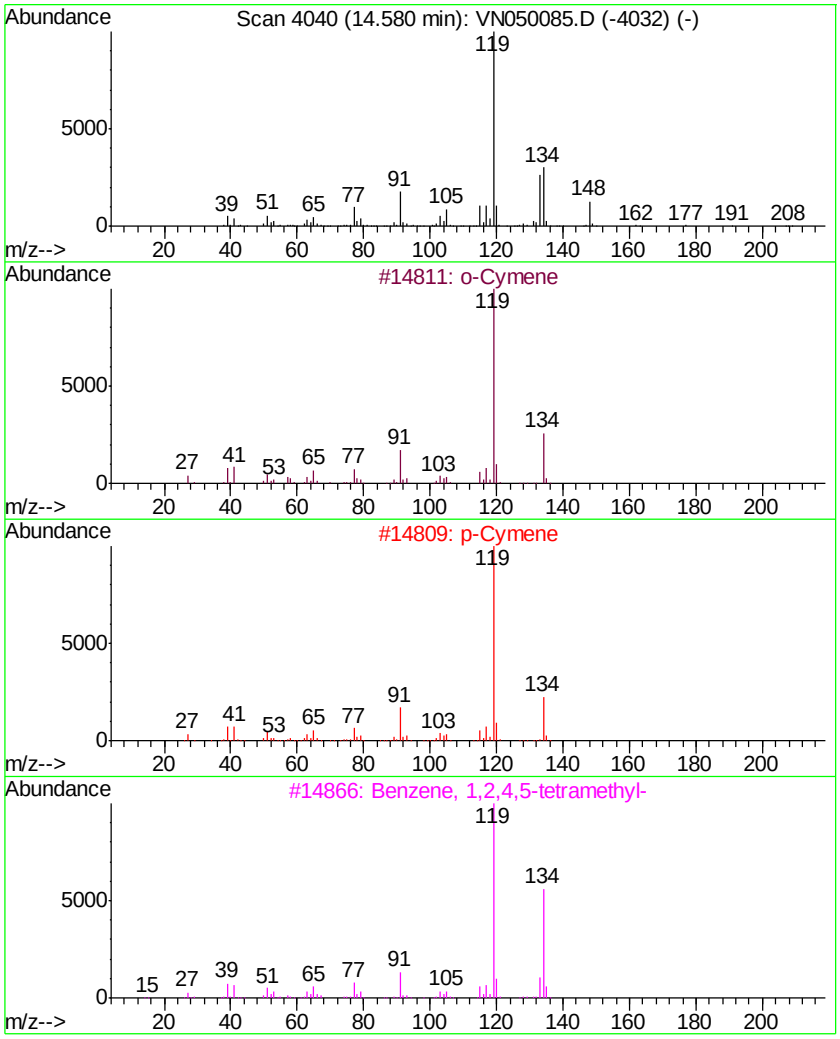
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

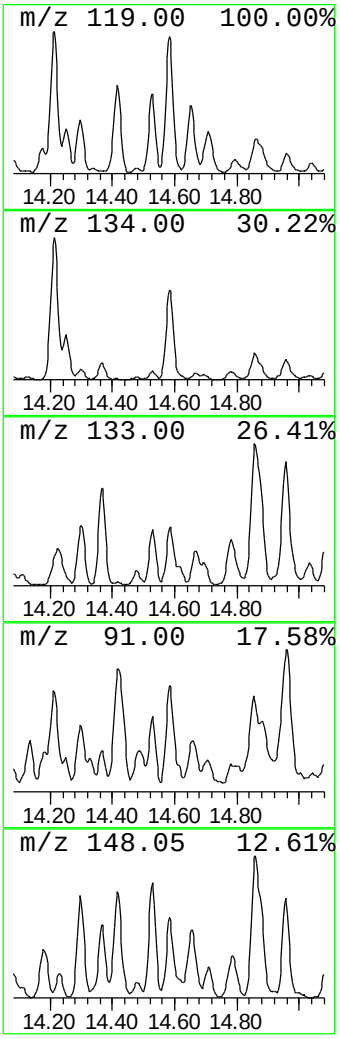
**********

Peak Number 6 o-Cymene Concentration Rank 7

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.58 | 10.65 ug/l | 511601 | 1,4-Dichlorobenzene-d4 | 13.35 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | o-Cymene | 134 | C10H14 | 000527-84-4 | 94 |
| 2 | p-Cymene | 134 | C10H14 | 000099-87-6 | 93 |
| 3 | Benzene, 1,2,4,5-tetramethyl- | 134 | C10H14 | 000095-93-2 | 87 |
| 4 | Benzene, 4-ethyl-1,2-dimethyl- | 134 | C10H14 | 000934-80-5 | 87 |
| 5 | Benzene, 1,2,3,4-tetramethyl- | 134 | C10H14 | 000488-23-3 | 87 |



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN072618\
Data File : VN050085.D
Acq On : 26 Jul 2018 16:26
Operator : MD\SY
Sample : J4142-02
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 14 Sample Multiplier: 1

**Instrument :** MSVOA_N
**ClientSampleId :** MW-5

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N072418W.M
Quant Title : SW846 8260

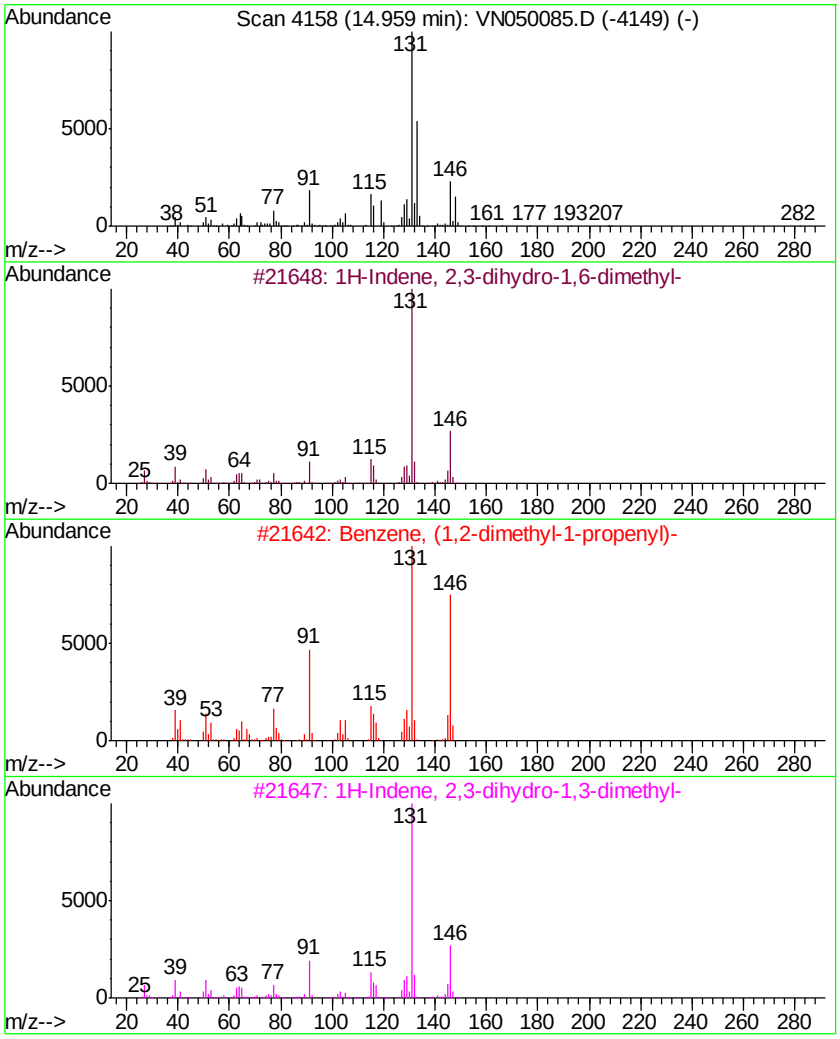
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

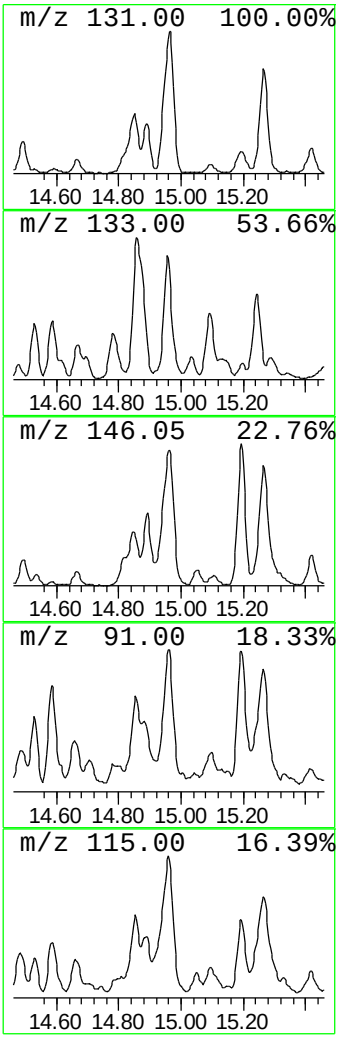
**********

Peak Number 7 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 2

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.96 | 21.47 ug/l | 1031180 | 1,4-Dichlorobenzene-d4 | 13.35 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1H-Indene, 2,3-dihydro-1,6-dimet... | 146 | C11H14 | 017059-48-2 | 83 |
| 2 | Benzene, (1,2-dimethyl-1-propenyl)- | 146 | C11H14 | 000769-57-3 | 64 |
| 3 | 1H-Indene, 2,3-dihydro-1,3-dimet... | 146 | C11H14 | 004175-53-5 | 64 |
| 4 | 1H-Indene, 2,3-dihydro-1,1-dimet... | 146 | C11H14 | 004912-92-9 | 64 |
| 5 | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 | C11H14 | 006682-71-9 | 64 |



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN072618\
Data File : VN050085.D
Acq On : 26 Jul 2018 16:26
Operator : MD\SY
Sample : J4142-02
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 14 Sample Multiplier: 1

**Instrument :** MSVOA_N
**ClientSampleId :** MW-5

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N072418W.M
Quant Title : SW846 8260

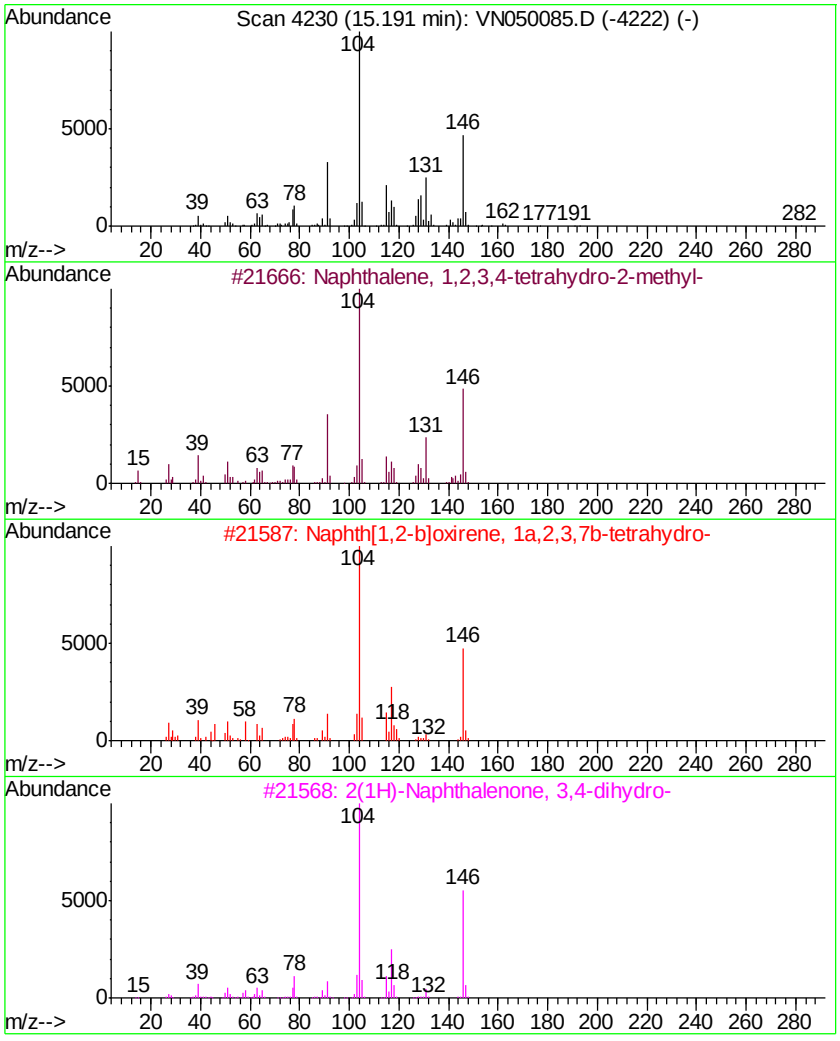
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

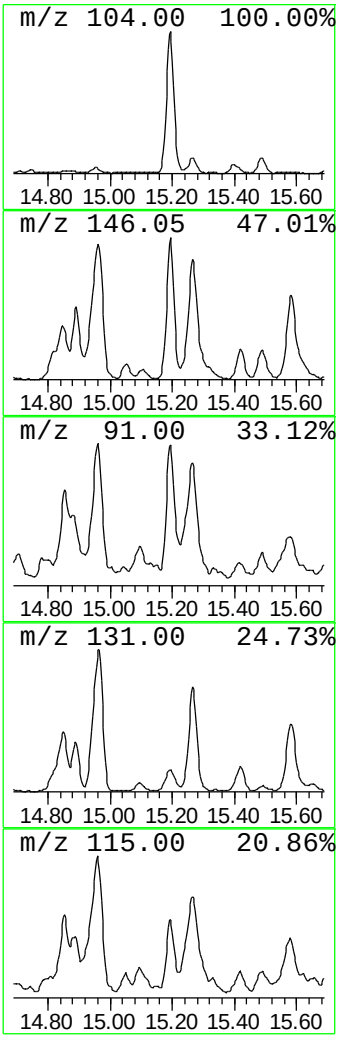
**********

Peak Number 8 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 10

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.19 | 9.43 ug/l | 452743 | 1,4-Dichlorobenzene-d4 | 13.35 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14 | 003877-19-8 | 97 |
| 2 | Naphth[1,2-b]oxirene, 1a,2,3,7b-... | 146 | C10H10O | 002461-34-9 | 70 |
| 3 | 2(1H)-Naphthalenone, 3,4-dihydro- | 146 | C10H10O | 000530-93-8 | 68 |
| 4 | 2-Methylbenzyl cyanide | 131 | C9H9N | 022364-68-7 | 50 |
| 5 | Naphthalene, 2-ethyl-1,2,3,4-tet... | 160 | C12H16 | 032367-54-7 | 47 |



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN072618\
Data File : VN050085.D
Acq On : 26 Jul 2018 16:26
Operator : MD\SY
Sample : J4142-02
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 14 Sample Multiplier: 1

**Instrument :** MSVOA_N
**ClientSampleId :** MW-5

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N072418W.M
Quant Title : SW846 8260

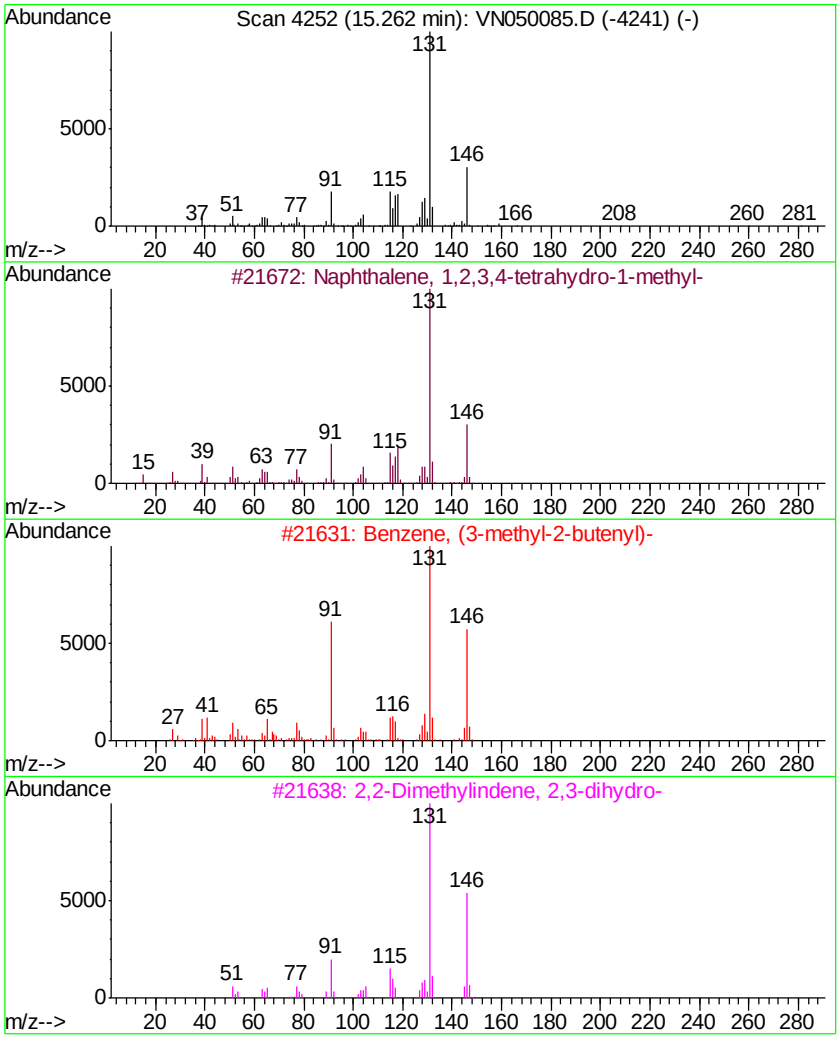
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

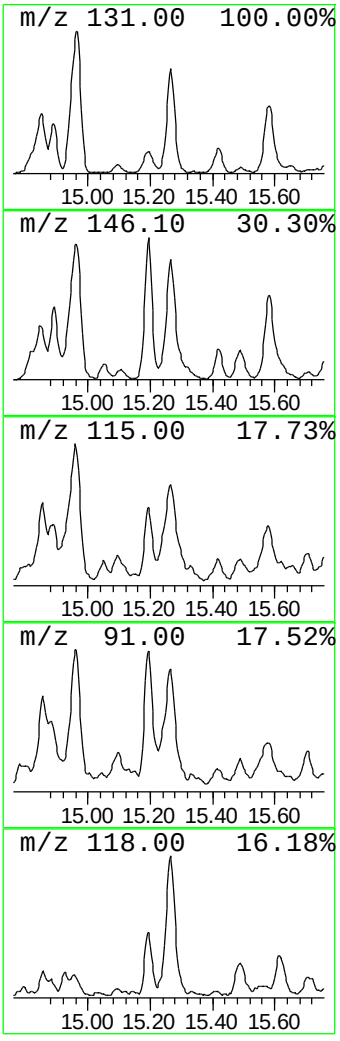
**********

Peak Number 9 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 4

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.26 | 16.83 ug/l | 808208 | 1,4-Dichlorobenzene-d4 | 13.35 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14 | 001559-81-5 | 94 |
| 2 | Benzene, (3-methyl-2-butenyl)- | 146 | C11H14 | 004489-84-3 | 93 |
| 3 | 2,2-Dimethylindene, 2,3-dihydro- | 146 | C11H14 | 020836-11-7 | 90 |
| 4 | 1H-Indene, 2,3-dihydro-1,6-dimet... | 146 | C11H14 | 017059-48-2 | 90 |
| 5 | 1H-Indene, 2,3-dihydro-1,2-dimet... | 146 | C11H14 | 017057-82-8 | 87 |



# Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN072618\
Data File : VN050085.D
Acq On : 26 Jul 2018 16:26
Operator : MD\SY
Sample : J4142-02
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 14 Sample Multiplier: 1

**Instrument :** MSVOA_N
**ClientSampleId :** MW-5

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N072418W.M
Quant Title : SW846 8260

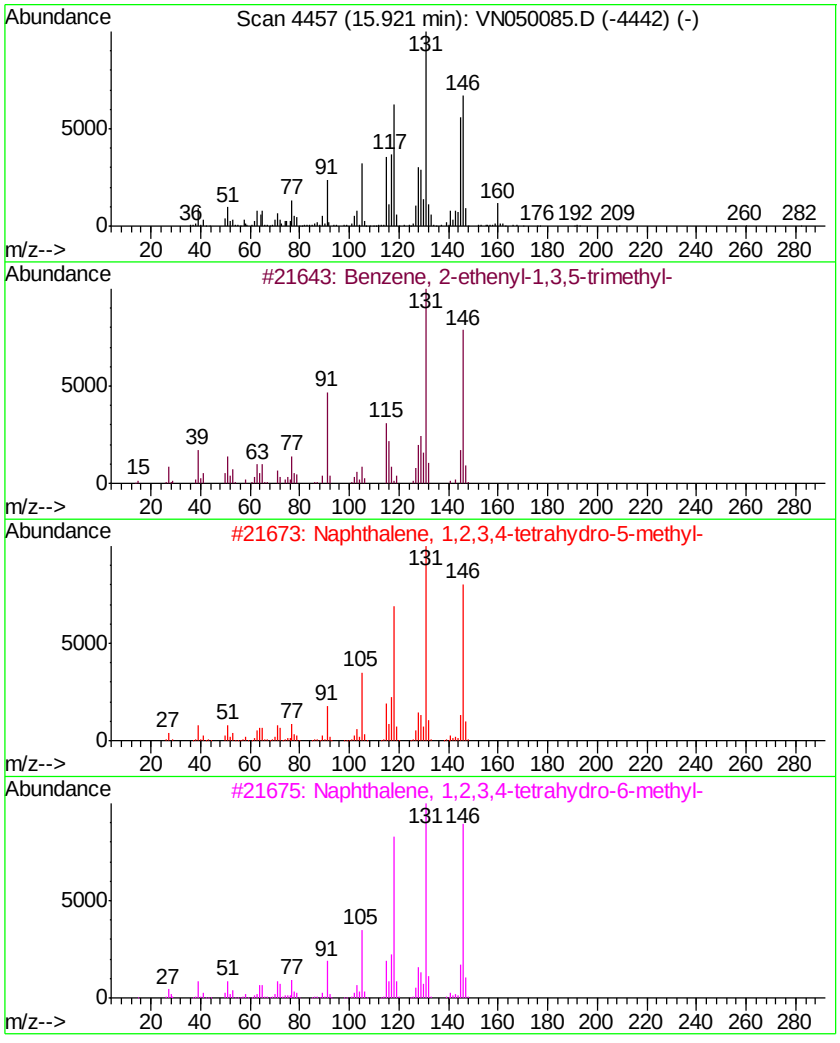
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

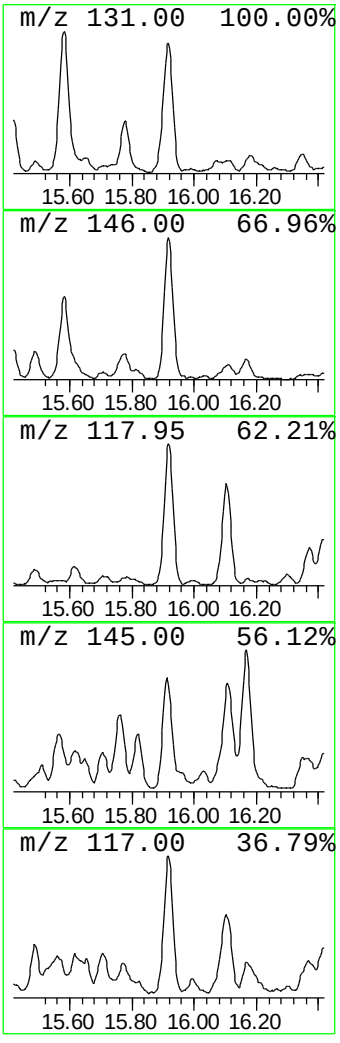
**********

Peak Number 10 Benzene, 2-ethenyl-1,3,5-tr... Concentration Rank 5

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.92 | 16.62 ug/l | 798127 | 1,4-Dichlorobenzene-d4 | 13.35 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, 2-ethenyl-1,3,5-trimethyl- | 146 | C11H14 | 000769-25-5 | 78 |
| 2 | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14 | 002809-64-5 | 76 |
| 3 | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14 | 001680-51-9 | 76 |
| 4 | Naphthalene, 1,2,3,4-tetrahydro-... | 160 | C12H16 | 004175-54-6 | 50 |
| 5 | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 | C11H14 | 006682-71-9 | 49 |



Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN072618\
Data File : VN050085.D
Acq On : 26 Jul 2018 16:26
Operator : MD\SY
Sample : J4142-02
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 14 Sample Multiplier: 1

**Instrument :** MSVOA_N
**ClientSampleId :** MW-5

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N072418W.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

| | | | | | Internal Standard | | | |
|---|---|---|---|---|---|---|---|---|
| TIC Top Hit name | RT | EstConc | Units | Response | # | RT | Resp | Conc |
| Acetic acid | 8.25 | 10.3 | ug/l | 512691 | 2 | 8.59 | 2489050 | 50.0 |
| Benzene, 1,2-diet... | 13.67 | 9.7 | ug/l | 463405 | 4 | 13.35 | 2401270 | 50.0 |
| Benzene, 2-ethyl-... | 13.89 | 11.1 | ug/l | 534791 | 4 | 13.35 | 2401270 | 50.0 |
| Benzene, 2-butenyl- | 13.99 | 18.5 | ug/l | 888243 | 4 | 13.35 | 2401270 | 50.0 |
| o-Cymene | 14.58 | 10.7 | ug/l | 511601 | 4 | 13.35 | 2401270 | 50.0 |
| 1H-Indene, 2,3-di... | 14.96 | 21.5 | ug/l | 1031180 | 4 | 13.35 | 2401270 | 50.0 |
| Naphthalene, 1,2,... | 15.19 | 9.4 | ug/l | 452743 | 4 | 13.35 | 2401270 | 50.0 |
| Naphthalene, 1,2,... | 15.26 | 16.8 | ug/l | 808208 | 4 | 13.35 | 2401270 | 50.0 |
| Benzene, 2-etheny... | 15.92 | 16.6 | ug/l | 798127 | 4 | 13.35 | 2401270 | 50.0 |