

Data Path : Z:\HPCHEM1\BNA G\DATA\BG022716\  
 Data File : BG021105.D  
 Acq On : 27 Feb 2016 8:35  
 Operator : UM/SJ  
 Sample : H1558-14  
 Misc :  
 ALS Vial : 28 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 S-16

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:\HPCHEM1\BNA\_G\METHODS\8270-BG022616.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 3.093       | 37            | 43          | 55           | rBV      | 361160         | 548850        | 99.21%          | 18.132%       |
| 2         | 3.234       | 61            | 67          | 76           | rVB3     | 8554           | 18531         | 3.35%           | 0.612%        |
| 3         | 5.261       | 406           | 412         | 418          | rBV      | 22178          | 32595         | 5.89%           | 1.077%        |
| 4         | 5.720       | 484           | 490         | 498          | rBB      | 106569         | 166482        | 30.09%          | 5.500%        |
| 5         | 7.312       | 754           | 761         | 768          | rBB      | 113529         | 186047        | 33.63%          | 6.146%        |
| 6         | 7.700       | 820           | 827         | 834          | rBB      | 106892         | 182401        | 32.97%          | 6.026%        |
| 7         | 8.170       | 901           | 907         | 913          | rBB      | 20547          | 32928         | 5.95%           | 1.088%        |
| 8         | 8.493       | 955           | 962         | 969          | rBB      | 80088          | 130723        | 23.63%          | 4.319%        |
| 9         | 9.333       | 1098          | 1105        | 1112         | rBB      | 66608          | 115491        | 20.88%          | 3.815%        |
| 10        | 10.978      | 1379          | 1385        | 1392         | rBB      | 29024          | 49363         | 8.92%           | 1.631%        |
| 11        | 13.405      | 1791          | 1798        | 1804         | rBB      | 186147         | 272793        | 49.31%          | 9.012%        |
| 12        | 14.221      | 1932          | 1937        | 1942         | rBB      | 19126          | 26427         | 4.78%           | 0.873%        |
| 13        | 14.774      | 2026          | 2031        | 2037         | rBB2     | 46967          | 70893         | 12.81%          | 2.342%        |
| 14        | 15.602      | 2168          | 2172        | 2176         | rBB2     | 6103           | 7379          | 1.33%           | 0.244%        |
| 15        | 16.254      | 2275          | 2283        | 2289         | rBB3     | 160614         | 231471        | 41.84%          | 7.647%        |
| 16        | 16.460      | 2314          | 2318        | 2329         | rBB      | 6387           | 10475         | 1.89%           | 0.346%        |
| 17        | 17.512      | 2491          | 2497        | 2503         | rVB2     | 67614          | 97807         | 17.68%          | 3.231%        |
| 18        | 18.375      | 2640          | 2644        | 2650         | rBV3     | 6633           | 7841          | 1.42%           | 0.259%        |
| 19        | 20.103      | 2932          | 2938        | 2944         | rVB      | 432349         | 553207        | 100.00%         | 18.276%       |
| 20        | 21.401      | 3154          | 3159        | 3165         | rBV2     | 15331          | 19455         | 3.52%           | 0.643%        |
| 21        | 21.771      | 3216          | 3222        | 3230         | rVB      | 81581          | 132821        | 24.01%          | 4.388%        |
| 22        | 23.493      | 3510          | 3515        | 3521         | rVB3     | 9696           | 17365         | 3.14%           | 0.574%        |
| 23        | 25.056      | 3771          | 3781        | 3790         | rBV2     | 42629          | 115673        | 20.91%          | 3.821%        |

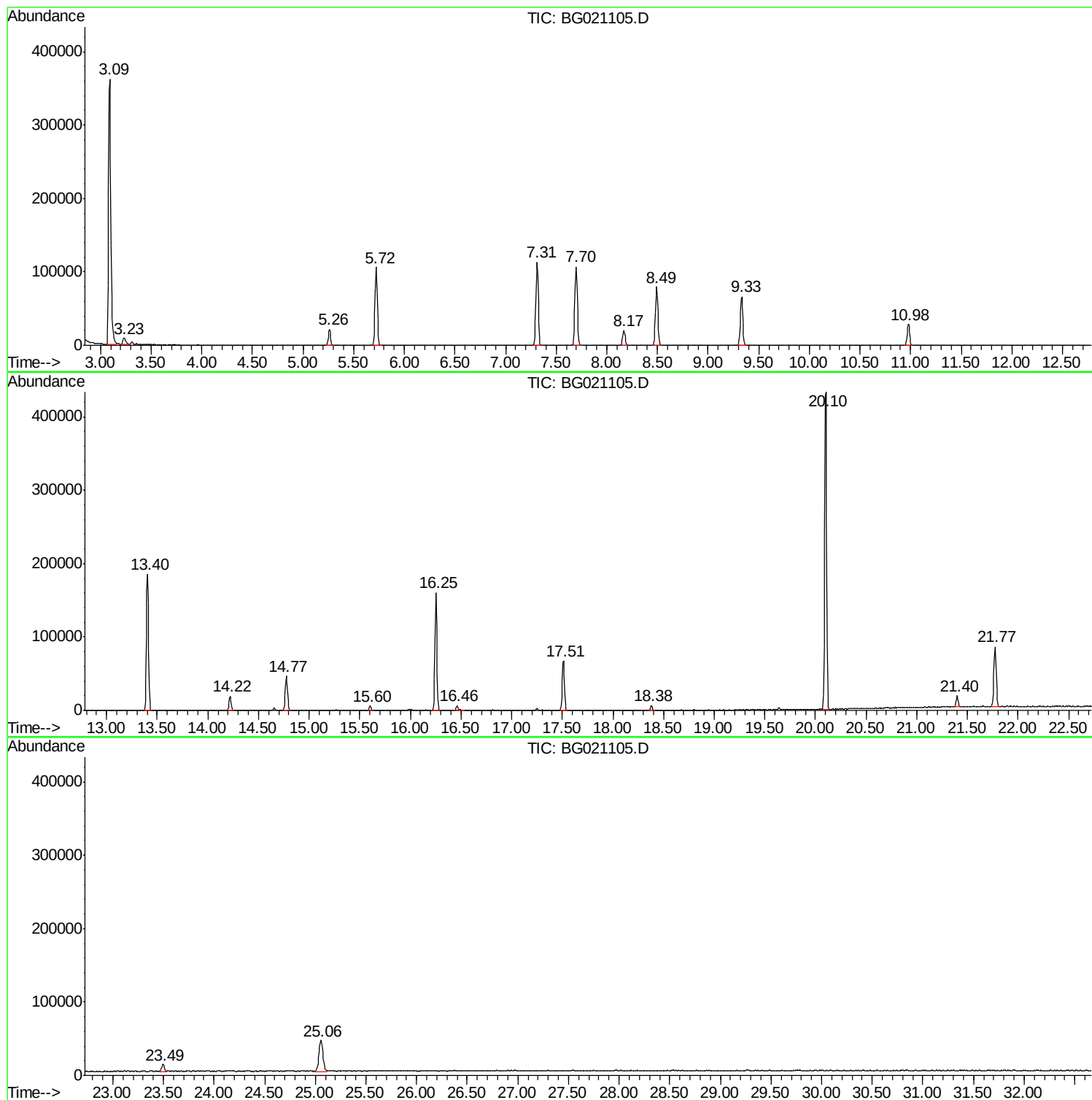
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Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
S-16

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022616.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P



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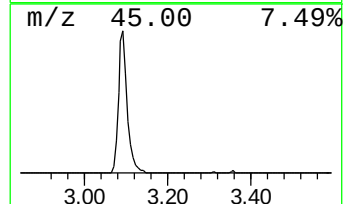
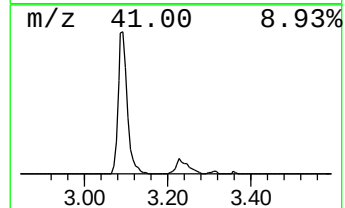
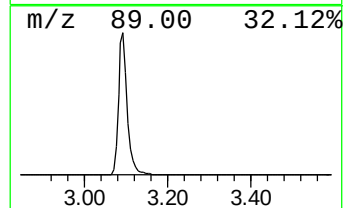
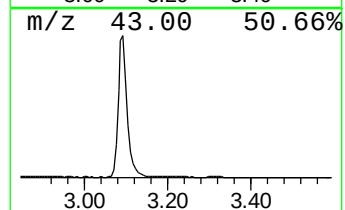
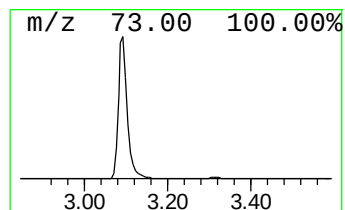
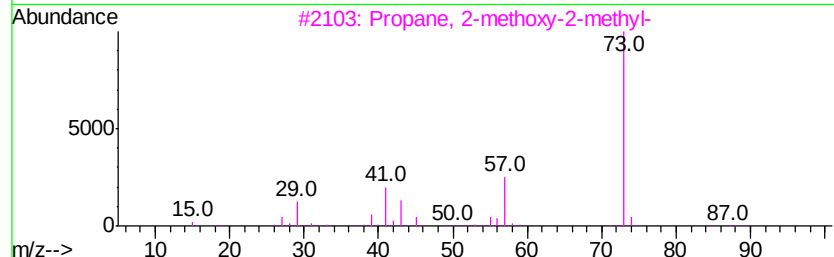
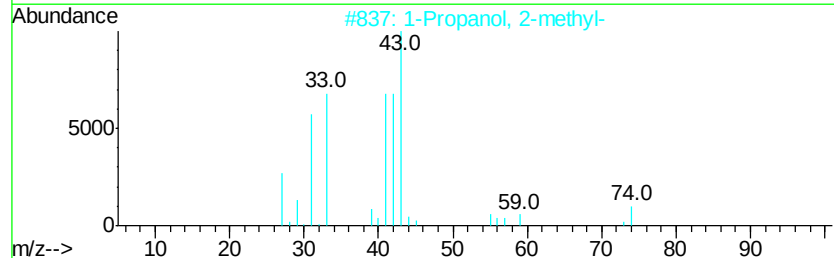
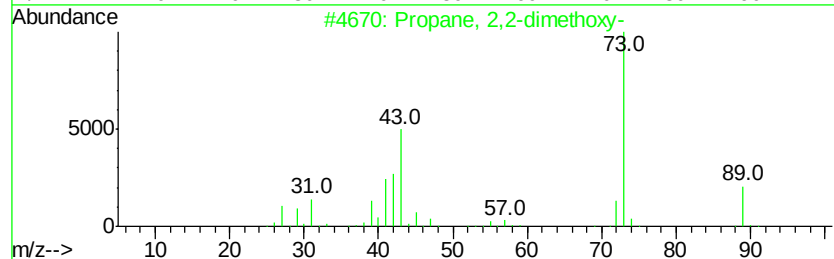
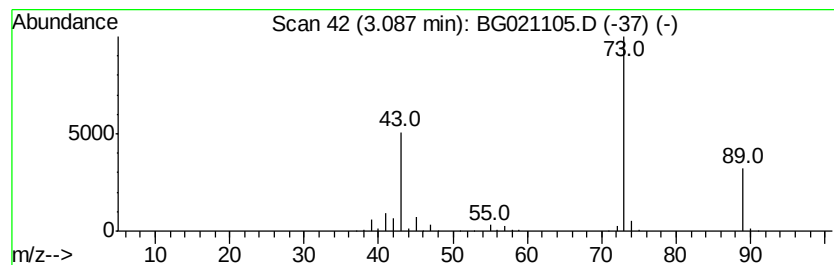
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 unknown3.09 Concentration Rank 1

| R.T. | EstConc   | Area   | Relative to ISTD       | R.T. |
|------|-----------|--------|------------------------|------|
| 3.09 | 333.36 ng | 548850 | 1,4-Dichlorobenzene-d4 | 8.17 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | Propane, 2,2-dimethoxy-        | 104 | C5H12O2 | 000077-76-9 | 38   |
| 2    |    | 1-Propanol, 2-methyl-          | 74  | C4H10O  | 000078-83-1 | 9    |
| 3    |    | Propane, 2-methoxy-2-methyl-   | 88  | C5H12O  | 001634-04-4 | 9    |
| 4    |    | Pentane, 3-methoxy-            | 102 | C6H14O  | 036839-67-5 | 9    |
| 5    |    | 2-Hydroxy-2-methylbutyric acid | 118 | C5H10O3 | 003739-30-8 | 9    |



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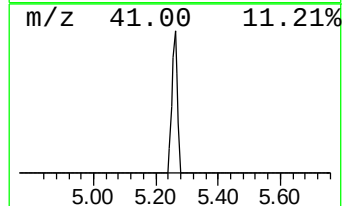
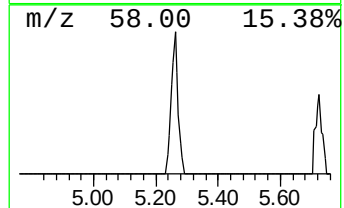
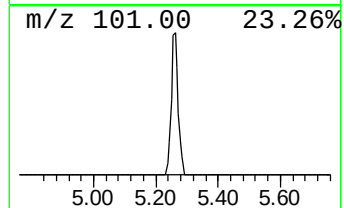
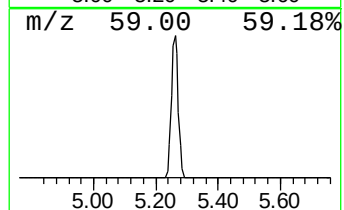
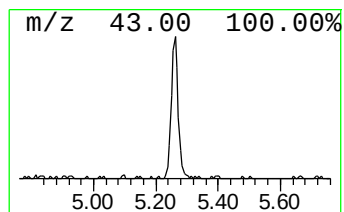
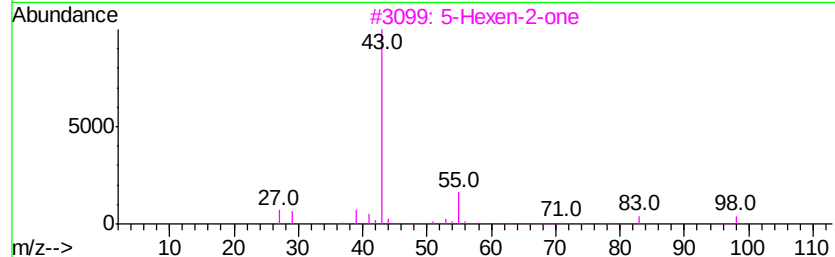
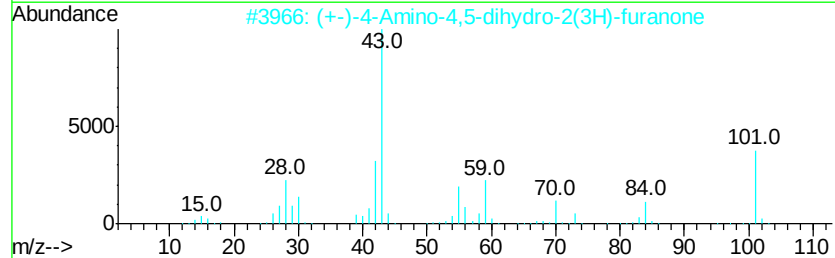
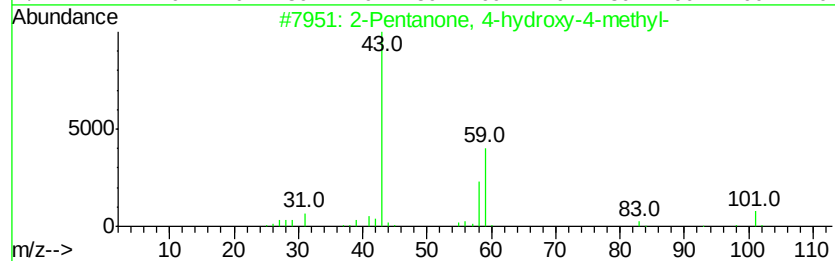
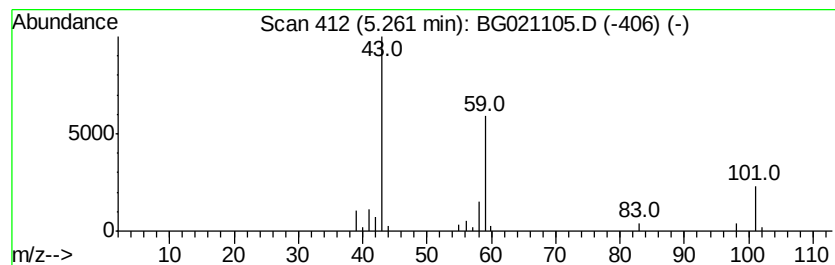
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TIC Integration Parameters: LSCINT.P

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Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

| R.T. | EstConc  | Area  | Relative to ISTD       | R.T. |
|------|----------|-------|------------------------|------|
| 5.26 | 19.80 ng | 32595 | 1,4-Dichlorobenzene-d4 | 8.17 |

| Hit# | of | Tentative ID                           | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 2-Pentanone, 4-hydroxy-4-methyl-       | 116 | C6H12O2 | 000123-42-2 | 64   |
| 2    |    | (+)-4-Amino-4,5-dihydro-2(3H)-furanone | 101 | C4H7NO2 | 016504-58-8 | 9    |
| 3    |    | 5-Hexen-2-one                          | 98  | C6H10O  | 000109-49-9 | 9    |
| 4    |    | Morpholine, 4-methyl-                  | 101 | C5H11NO | 000109-02-4 | 9    |
| 5    |    | Butane, 1-ethoxy-                      | 102 | C6H14O  | 000628-81-9 | 9    |



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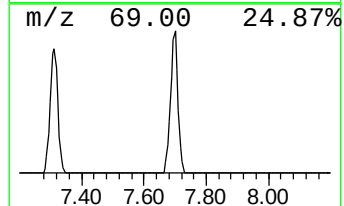
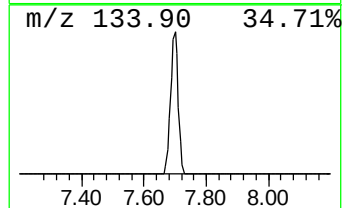
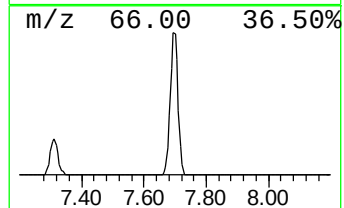
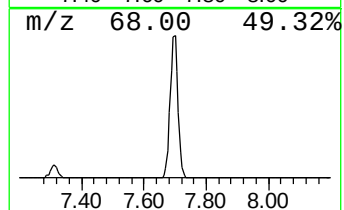
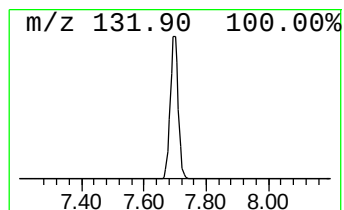
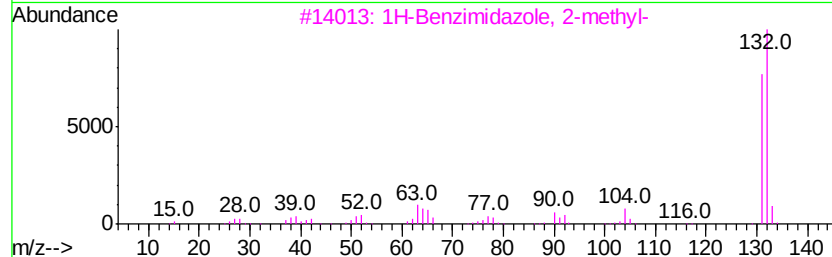
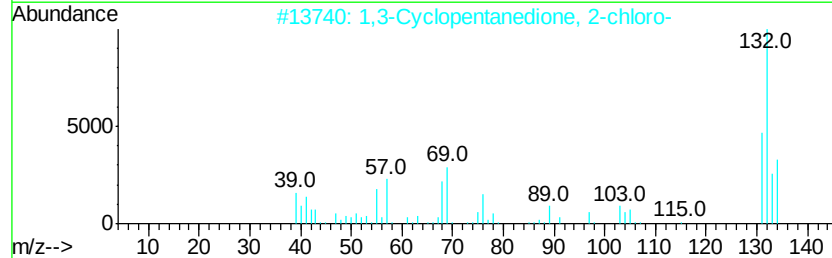
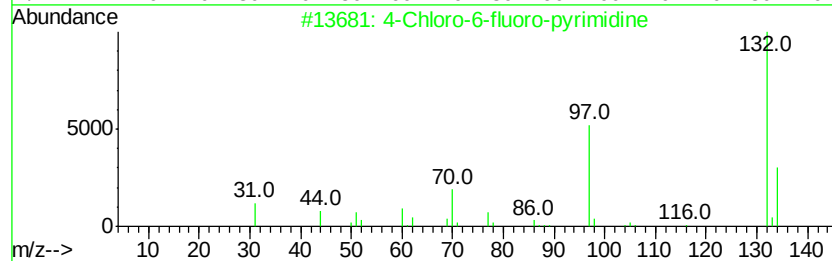
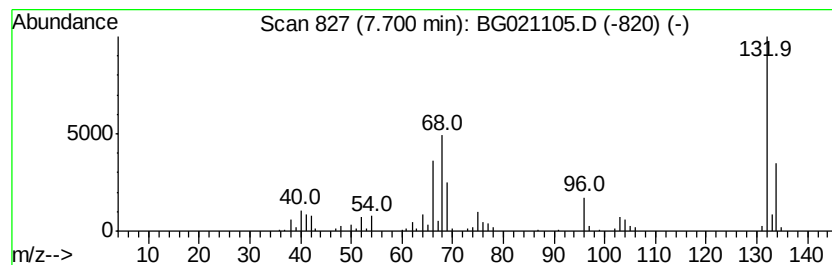
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 unknown7.70 Concentration Rank 2

| R.T. | EstConc   | Area   | Relative to ISTD       | R.T. |
|------|-----------|--------|------------------------|------|
| 7.70 | 110.79 ng | 182401 | 1,4-Dichlorobenzene-d4 | 8.17 |

| Hit# | of | Tentative ID                     | MW  | MolForm   | CAS#         | Qual |
|------|----|----------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 4-Chloro-6-fluoro-pyrimidine     | 132 | C4H2ClFN2 | 051422-01-6  | 27   |
| 2    |    | 1,3-Cyclopentanedione, 2-chloro- | 132 | C5H5ClO2  | 014203-19-1  | 25   |
| 3    |    | 1H-Benzimidazole, 2-methyl-      | 132 | C8H8N2    | 000615-15-6  | 22   |
| 4    |    | 5-Aminoindole                    | 132 | C8H8N2    | 005192-03-0  | 22   |
| 5    |    | (5-METHYL-2-PYRIDYL)ACETONITRILE | 132 | C8H8N2    | 1000241-93-9 | 22   |



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S-16

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022616.M  
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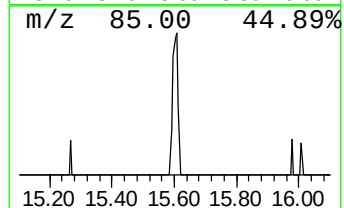
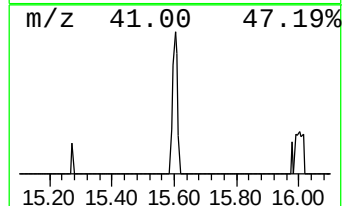
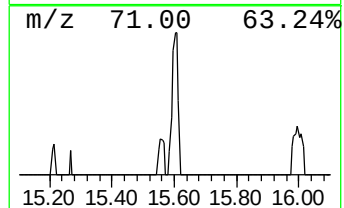
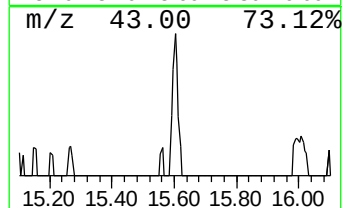
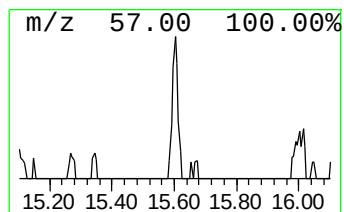
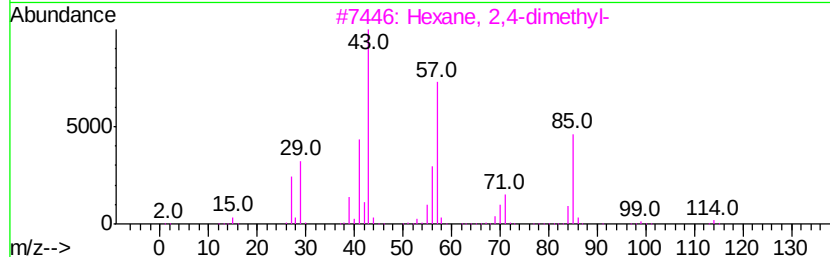
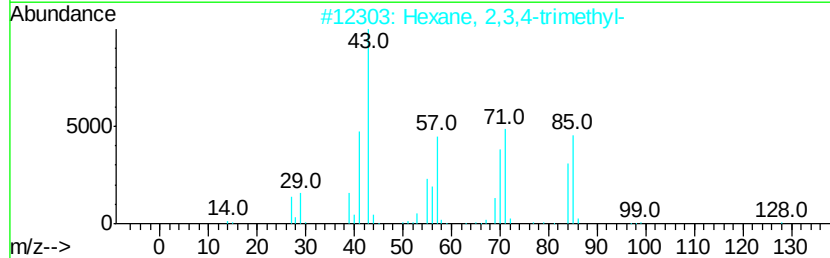
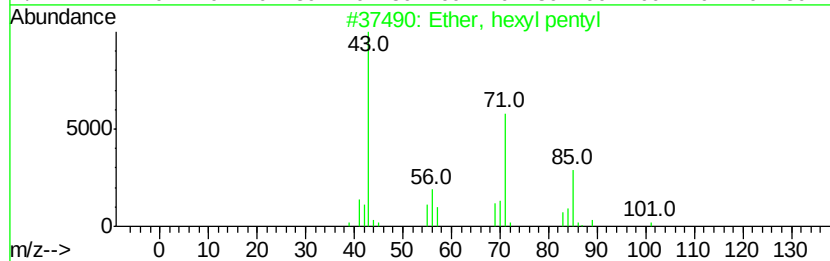
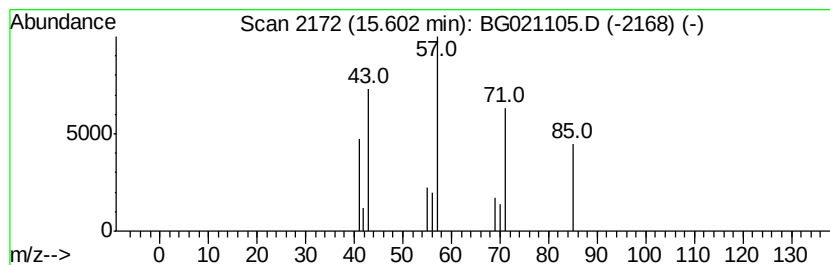
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 Ether, hexyl pentyl Concentration Rank 9

| R.T.  | EstConc | Area | Relative to ISTD | R.T.  |
|-------|---------|------|------------------|-------|
| 15.60 | 2.08 ng | 7379 | Acenaphthene-d10 | 14.77 |

| Hit# of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|---------|---|-------------------------------------|-----|----------|-------------|------|
| 1       |   | Ether, hexyl pentyl                 | 172 | C11H24O  | 032357-83-8 | 59   |
| 2       |   | Hexane, 2,3,4-trimethyl-            | 128 | C9H20    | 000921-47-1 | 42   |
| 3       |   | Hexane, 2,4-dimethyl-               | 114 | C8H18    | 000589-43-5 | 33   |
| 4       |   | Hexane, 3-ethyl-                    | 114 | C8H18    | 000619-99-8 | 28   |
| 5       |   | 2,2-Dimethylpropionic acid, 4-me... | 186 | C11H22O2 | 150588-62-8 | 25   |



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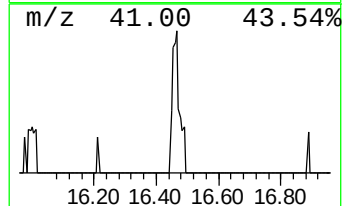
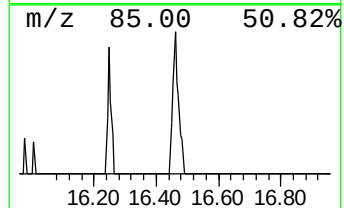
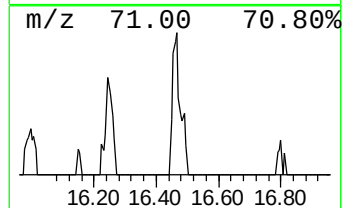
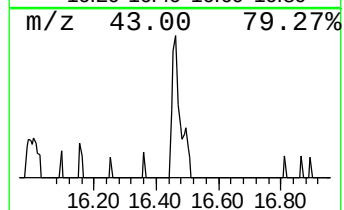
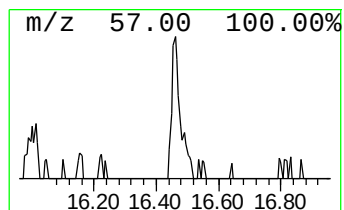
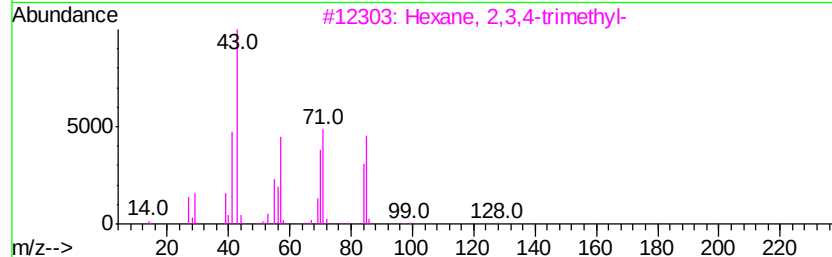
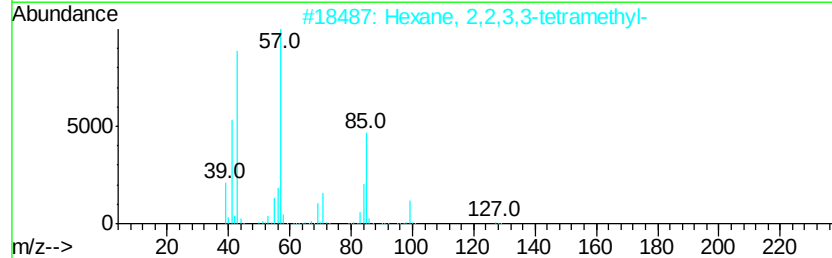
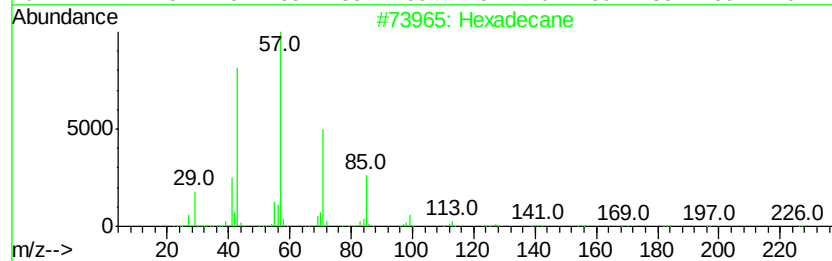
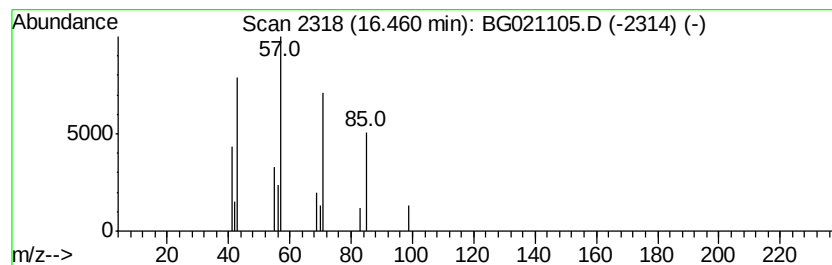
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 Hexadecane Concentration Rank 8

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 16.46 | 2.14 ng | 10475 | Phenanthrene-d10 | 17.51 |

| Hit# of | 5                            | Tentative ID | MW  | MolForm | CAS#        | Qual |
|---------|------------------------------|--------------|-----|---------|-------------|------|
| 1       | Hexadecane                   |              | 226 | C16H34  | 000544-76-3 | 72   |
| 2       | Hexane, 2,2,3,3-tetramethyl- |              | 142 | C10H22  | 013475-81-5 | 59   |
| 3       | Hexane, 2,3,4-trimethyl-     |              | 128 | C9H20   | 000921-47-1 | 45   |
| 4       | Heptane, 2,4,6-trimethyl-    |              | 142 | C10H22  | 002613-61-8 | 42   |
| 5       | Hexane, 2,4-dimethyl-        |              | 114 | C8H18   | 000589-43-5 | 40   |



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ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
S-16

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022616.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

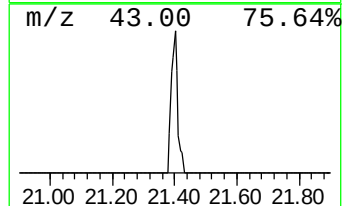
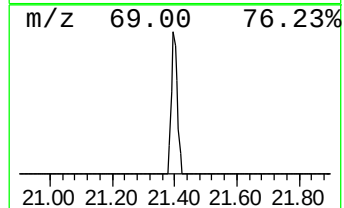
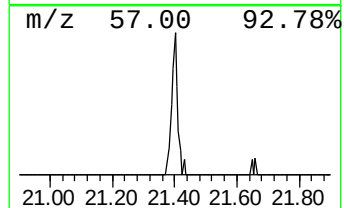
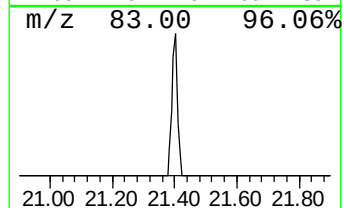
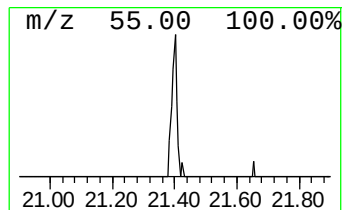
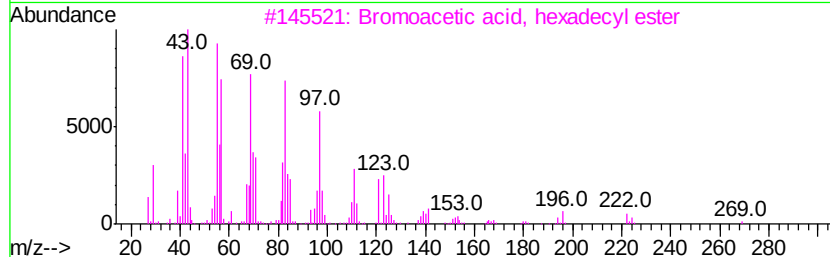
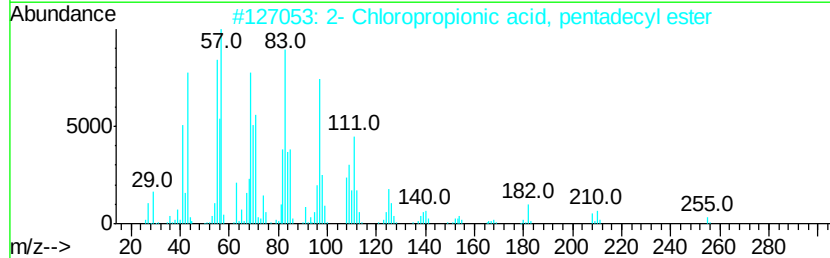
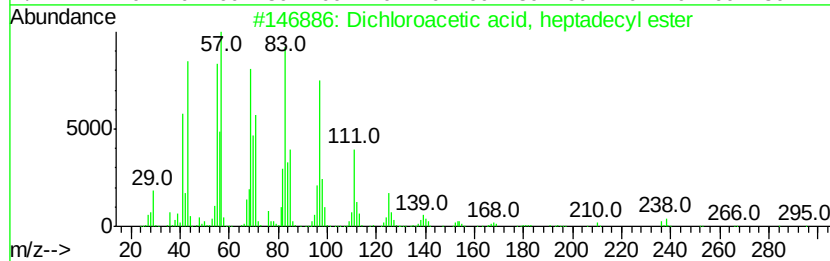
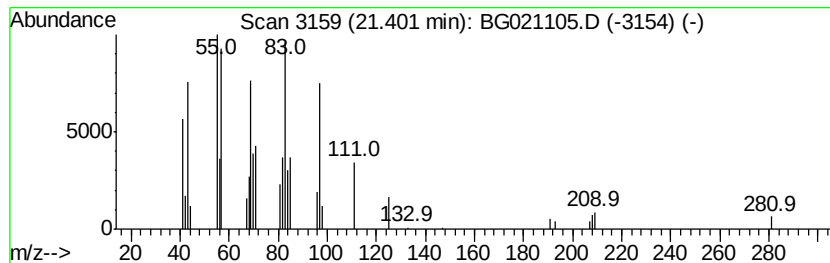
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 Dichloroacetic acid, heptad... Concentration Rank 7

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 21.40 | 2.93 ng | 19455 | Chrysene-d12     | 21.77 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-------------|--------------|------|
| 1    |    |   | Dichloroacetic acid, heptadecyl ... | 366 | C19H36Cl2O2 | 1000282-98-2 | 91   |
| 2    |    |   | 2- Chloropropionic acid, pentade... | 318 | C18H35ClO2  | 1000292-44-1 | 91   |
| 3    |    |   | Bromoacetic acid, hexadecyl ester   | 362 | C18H35BrO2  | 005454-48-8  | 90   |
| 4    |    |   | Acetic acid, chloro-, octadecyl ... | 346 | C20H39ClO2  | 005348-82-3  | 90   |
| 5    |    |   | 11-Tricosene                        | 322 | C23H46      | 052078-56-5  | 87   |





Data Path : Z:\HPCHEM1\BNA G\DATA\BG022716\  
Data File : BG021105.D  
Acq On : 27 Feb 2016 8:35  
Operator : UM/SJ  
Sample : H1558-14  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
S-16

Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022616.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

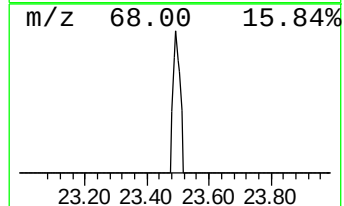
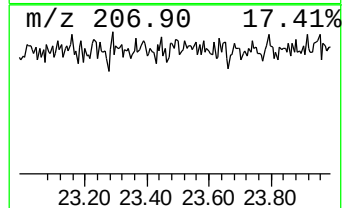
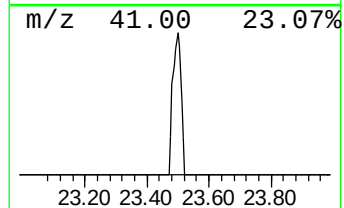
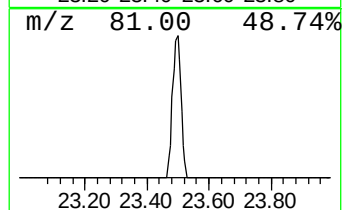
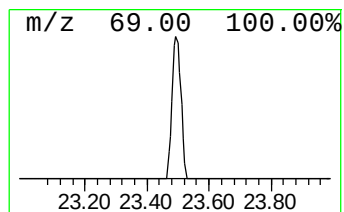
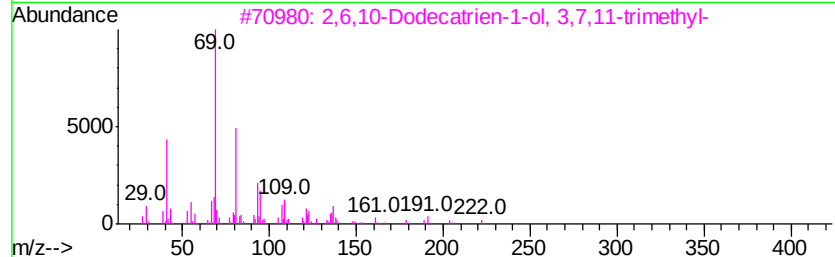
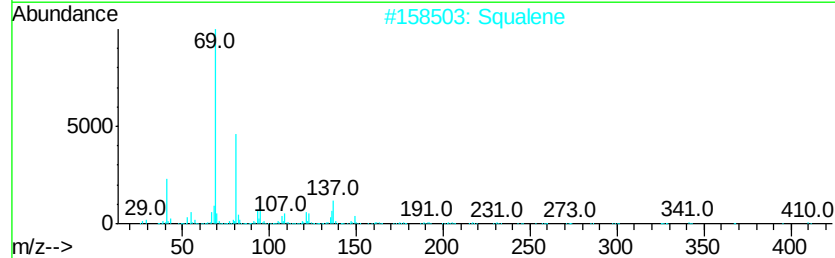
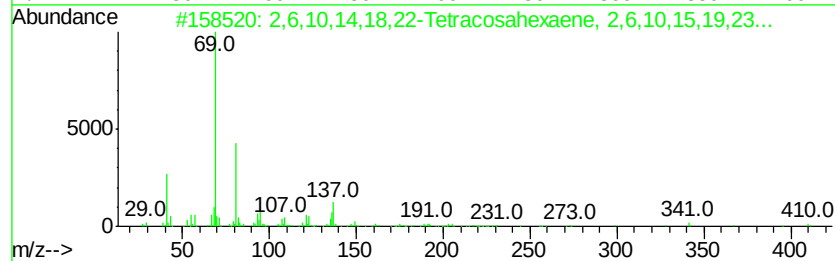
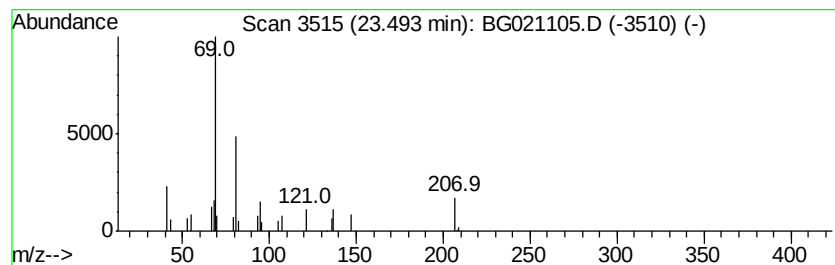
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 2,6,10,14,18,22-Tetracosahexaene... Concentration Rank 6

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 23.49 | 3.00 ng | 17365 | Perylene-d12     | 25.06 |

| Hit# | of                                  | 5   | Tentative ID | MW           | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|--------------|---------|------|------|
| 1    | 2,6,10,14,18,22-Tetracosahexaene... | 410 | C30H50       | 000111-02-4  | 72      |      |      |
| 2    | Squalene                            | 410 | C30H50       | 007683-64-9  | 56      |      |      |
| 3    | 2,6,10-Dodecatrien-1-ol, 3,7,11-... | 222 | C15H26O      | 004602-84-0  | 56      |      |      |
| 4    | 4,8,12-Tetradecatrienal, 5,9,13-... | 248 | C17H28O      | 066408-55-7  | 50      |      |      |
| 5    | Docosa-2,6,10,14,18-pentaen-22-a... | 384 | C27H44O      | 1000163-04-7 | 50      |      |      |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG022716\  
Data File : BG021105.D  
Acq On : 27 Feb 2016 8:35  
Operator : UM/SJ  
Sample : H1558-14  
Misc :  
ALS Vial : 28 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
S-16

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG022616.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |        |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|--------|------|
|                      |       |         |       |          | #                     | RT    | Resp   | Conc |
| unknown3.09          | 3.09  | 333.4   | ng    | 548850   | 1                     | 8.17  | 32928  | 20.0 |
| 2-Pentanone, 4-hy... | 5.26  | 19.8    | ng    | 32595    | 1                     | 8.17  | 32928  | 20.0 |
| unknown7.70          | 7.70  | 110.8   | ng    | 182401   | 1                     | 8.17  | 32928  | 20.0 |
| Ether, hexyl pentyl  | 15.60 | 2.1     | ng    | 7379     | 3                     | 14.77 | 70893  | 20.0 |
| Hexadecane           | 16.46 | 2.1     | ng    | 10475    | 4                     | 17.51 | 97807  | 20.0 |
| Dichloroacetic ac... | 21.40 | 2.9     | ng    | 19455    | 5                     | 21.77 | 132821 | 20.0 |
| 2,6,10,14,18,22-T... | 23.49 | 3.0     | ng    | 17365    | 6                     | 25.06 | 115673 | 20.0 |