

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071818\  
 Data File : BF107502.D  
 Acq On : 19 Jul 2018 8:28  
 Operator : JU/SJ  
 Sample : J4033-03  
 Misc :  
 ALS Vial : 38 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 071618-DBRI-F-D-B

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

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

Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF071618.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         | 2.902       | 91            | 96          | 102          | rBV      | 149153         | 218881        | 2.14%           | 0.401%        |
| 2         | 2.978       | 102           | 109         | 115          | rVB2     | 115512         | 204164        | 1.99%           | 0.374%        |
| 3         | 3.819       | 245           | 252         | 259          | rBV      | 565644         | 806387        | 7.87%           | 1.477%        |
| 4         | 4.496       | 360           | 367         | 371          | rBV4     | 73426          | 137593        | 1.34%           | 0.252%        |
| 5         | 4.660       | 384           | 395         | 398          | rVV      | 5211813        | 10241550      | 100.00%         | 18.757%       |
| 6         | 4.784       | 414           | 416         | 420          | rVB      | 129887         | 122335        | 1.19%           | 0.224%        |
| 7         | 4.872       | 426           | 431         | 440          | rBV2     | 2763146        | 3572418       | 34.88%          | 6.543%        |
| 8         | 5.607       | 549           | 556         | 560          | rBV      | 1447694        | 1400281       | 13.67%          | 2.564%        |
| 9         | 6.584       | 718           | 722         | 732          | rBV      | 1917668        | 2345542       | 22.90%          | 4.296%        |
| 10        | 6.754       | 746           | 751         | 755          | rBV2     | 2509446        | 2417255       | 23.60%          | 4.427%        |
| 11        | 6.984       | 786           | 790         | 793          | rBV      | 1018346        | 881563        | 8.61%           | 1.615%        |
| 12        | 7.031       | 795           | 798         | 805          | rVB      | 90013          | 112694        | 1.10%           | 0.206%        |
| 13        | 7.143       | 812           | 817         | 820          | rVB      | 3618431        | 3303916       | 32.26%          | 6.051%        |
| 14        | 7.213       | 826           | 829         | 833          | rBV      | 115715         | 158485        | 1.55%           | 0.290%        |
| 15        | 7.548       | 879           | 886         | 889          | rBV      | 2819871        | 3182291       | 31.07%          | 5.828%        |
| 16        | 7.707       | 908           | 913         | 918          | rBV3     | 717150         | 1013139       | 9.89%           | 1.855%        |
| 17        | 8.089       | 973           | 978         | 981          | rBV      | 548659         | 524951        | 5.13%           | 0.961%        |
| 18        | 8.266       | 1003          | 1008        | 1011         | rBV2     | 1020818        | 1057715       | 10.33%          | 1.937%        |
| 19        | 8.295       | 1011          | 1013        | 1016         | rVB      | 401259         | 352858        | 3.45%           | 0.646%        |
| 20        | 8.354       | 1019          | 1023        | 1027         | rBV      | 216630         | 215722        | 2.11%           | 0.395%        |
| 21        | 8.760       | 1089          | 1092        | 1094         | rBV      | 157881         | 171282        | 1.67%           | 0.314%        |
| 22        | 8.778       | 1094          | 1095        | 1098         | rVB      | 185896         | 106311        | 1.04%           | 0.195%        |
| 23        | 9.031       | 1134          | 1138        | 1140         | rBV      | 540749         | 474429        | 4.63%           | 0.869%        |
| 24        | 9.066       | 1140          | 1144        | 1147         | rVB      | 1000655        | 878597        | 8.58%           | 1.609%        |
| 25        | 9.278       | 1177          | 1180        | 1185         | rVB      | 211480         | 159511        | 1.56%           | 0.292%        |
| 26        | 9.342       | 1185          | 1191        | 1194         | rBV      | 4983545        | 5014226       | 48.96%          | 9.183%        |
| 27        | 10.025      | 1304          | 1307        | 1310         | rVB3     | 250781         | 218778        | 2.14%           | 0.401%        |
| 28        | 10.819      | 1437          | 1442        | 1445         | rBV      | 2582637        | 2489974       | 24.31%          | 4.560%        |
| 29        | 11.519      | 1556          | 1561        | 1572         | rBV4     | 478446         | 656030        | 6.41%           | 1.201%        |
| 30        | 11.895      | 1620          | 1625        | 1628         | rBV      | 578414         | 544576        | 5.32%           | 0.997%        |
| 31        | 12.036      | 1644          | 1649        | 1660         | rBV7     | 224089         | 552676        | 5.40%           | 1.012%        |
| 32        | 12.589      | 1740          | 1743        | 1745         | rVB      | 168304         | 137768        | 1.35%           | 0.252%        |
| 33        | 12.689      | 1756          | 1760        | 1763         | rBV      | 481789         | 424057        | 4.14%           | 0.777%        |
| 34        | 12.907      | 1793          | 1797        | 1800         | rBV2     | 190878         | 202416        | 1.98%           | 0.371%        |

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF071818\  
Data File : BF107502.D  
Acq On : 19 Jul 2018 8:28  
Operator : JU/SJ  
Sample : J4033-03  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
071618-DBRI-F-D-B

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:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF071618.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |       |         |         |        |        |
|----|--------|------|------|------|-------|---------|---------|--------|--------|
| 35 | 12.960 | 1804 | 1806 | 1809 | rVV   | 285296  | 204927  | 2.00%  | 0.375% |
| 36 | 13.107 | 1827 | 1831 | 1836 | rVB3  | 2081588 | 2450735 | 23.93% | 4.488% |
| 37 | 13.319 | 1863 | 1867 | 1870 | rVB   | 497545  | 426746  | 4.17%  | 0.782% |
| 38 | 13.660 | 1922 | 1925 | 1928 | rBV   | 764311  | 611802  | 5.97%  | 1.120% |
| 39 | 13.989 | 1978 | 1981 | 1984 | rBV   | 878366  | 721409  | 7.04%  | 1.321% |
| 40 | 14.166 | 2003 | 2011 | 2013 | rBV4  | 126787  | 190475  | 1.86%  | 0.349% |
| 41 | 14.307 | 2031 | 2035 | 2038 | rBV   | 1106140 | 924678  | 9.03%  | 1.693% |
| 42 | 14.501 | 2063 | 2068 | 2070 | rVB6  | 83341   | 108531  | 1.06%  | 0.199% |
| 43 | 14.613 | 2083 | 2087 | 2091 | rVV   | 1083046 | 981808  | 9.59%  | 1.798% |
| 44 | 14.936 | 2138 | 2142 | 2147 | rVB   | 860966  | 917982  | 8.96%  | 1.681% |
| 45 | 15.295 | 2199 | 2203 | 2208 | rBV   | 702165  | 807245  | 7.88%  | 1.478% |
| 46 | 15.701 | 2267 | 2272 | 2281 | rBV2  | 567332  | 795878  | 7.77%  | 1.458% |
| 47 | 16.171 | 2346 | 2352 | 2358 | rBV2  | 382053  | 598511  | 5.84%  | 1.096% |
| 48 | 16.712 | 2439 | 2444 | 2452 | rVB3  | 190423  | 372681  | 3.64%  | 0.683% |
| 49 | 17.359 | 2549 | 2554 | 2562 | rVB10 | 86534   | 188887  | 1.84%  | 0.346% |

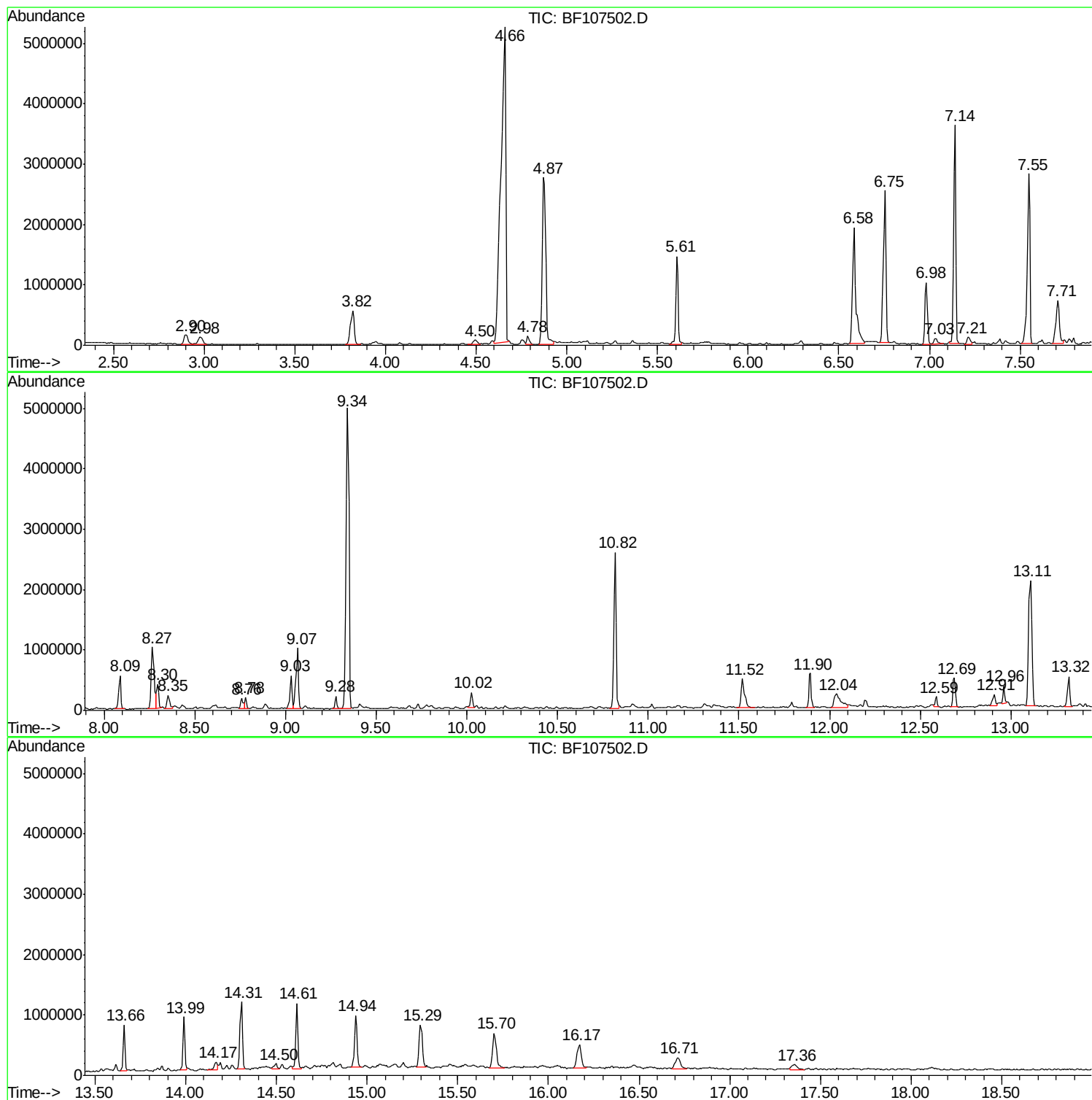
Sum of corrected areas: 54602666

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071818\  
Data File : BF107502.D  
Acq On : 19 Jul 2018 8:28  
Operator : JU/SJ  
Sample : J4033-03  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
071618-DBRI-F-D-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071618.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071818\  
Data File : BF107502.D  
Acq On : 19 Jul 2018 8:28  
Operator : JU/SJ  
Sample : J4033-03  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
071618-DBRI-F-D-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071618.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

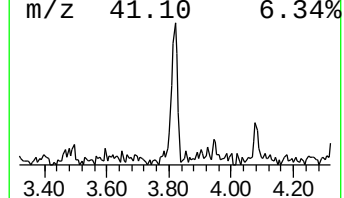
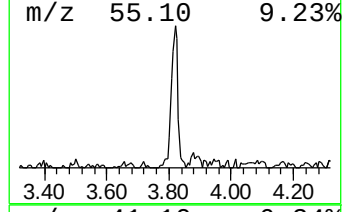
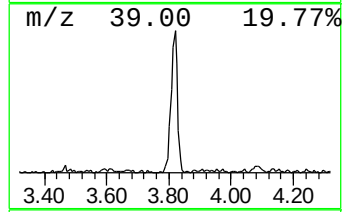
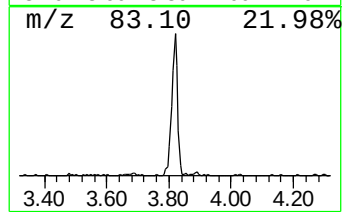
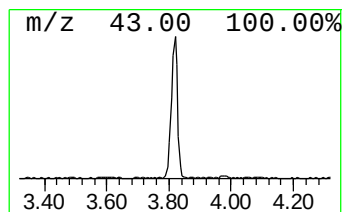
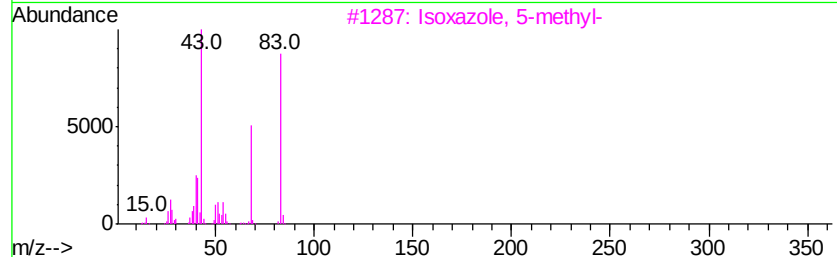
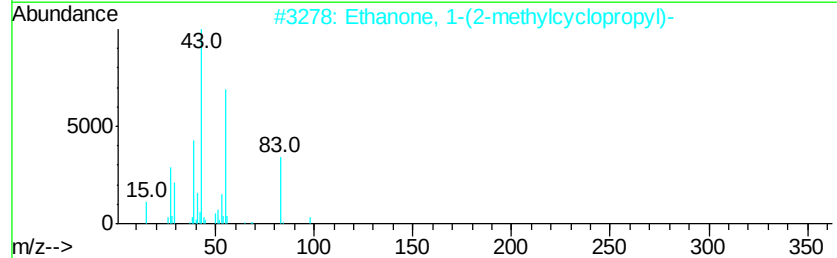
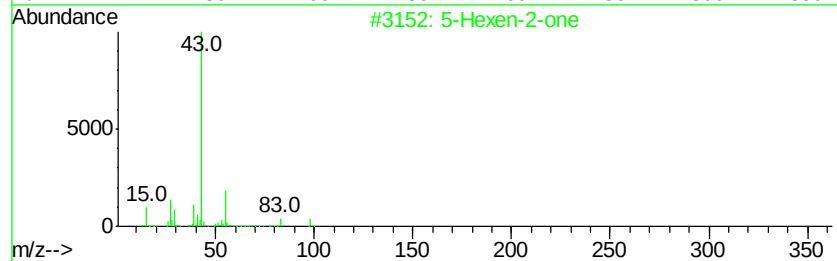
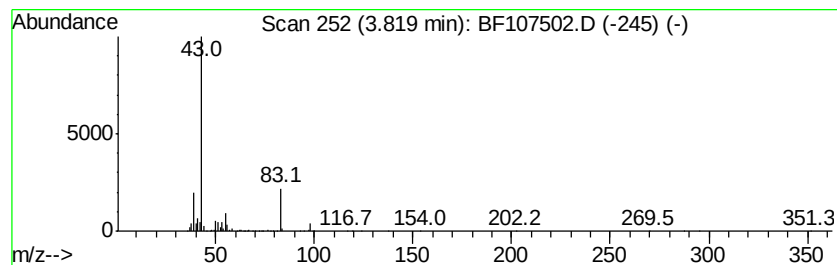
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 1 unknown3.82 Concentration Rank 15

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 3.82 | 18.29 ng | 806387 | 1,4-Dichlorobenzene-d4 | 6.98 |

| Hit# | of | Tentative ID                       | MW | MolForm | CAS#        | Qual |
|------|----|------------------------------------|----|---------|-------------|------|
| 1    | 5  | 5-Hexen-2-one                      | 98 | C6H10O  | 000109-49-9 | 47   |
| 2    |    | Ethanone, 1-(2-methylcyclopropyl)- | 98 | C6H10O  | 000930-56-3 | 40   |
| 3    |    | Isoxazole, 5-methyl-               | 83 | C4H5NO  | 005765-44-6 | 25   |
| 4    |    | 2-Methyl-3-oxobutyronitrile        | 97 | C5H7NO  | 004468-47-7 | 9    |
| 5    |    | 4-Penten-2-one, 4-methyl-          | 98 | C6H10O  | 003744-02-3 | 9    |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071818\  
Data File : BF107502.D  
Acq On : 19 Jul 2018 8:28  
Operator : JU/SJ  
Sample : J4033-03  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
071618-DBRI-F-D-B

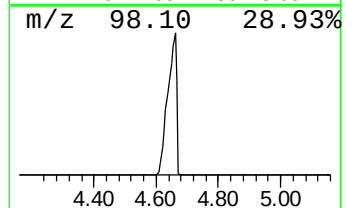
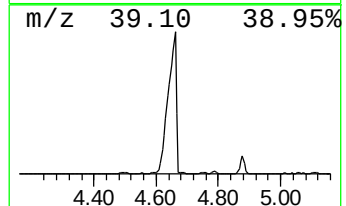
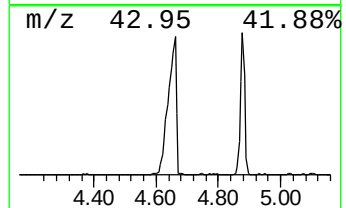
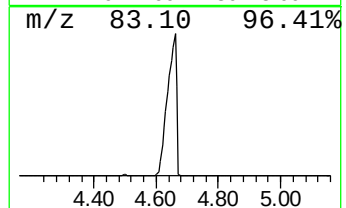
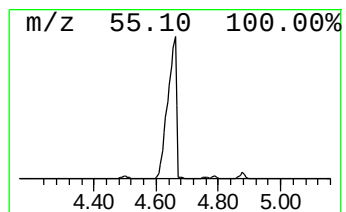
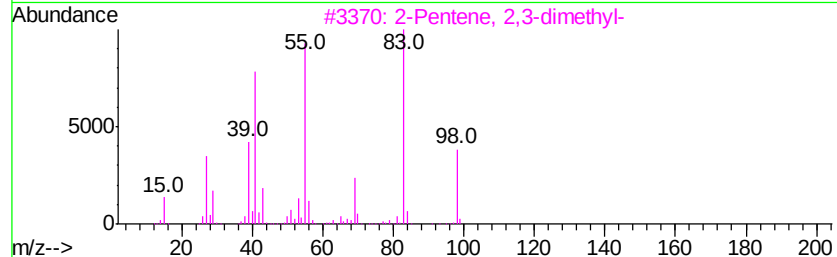
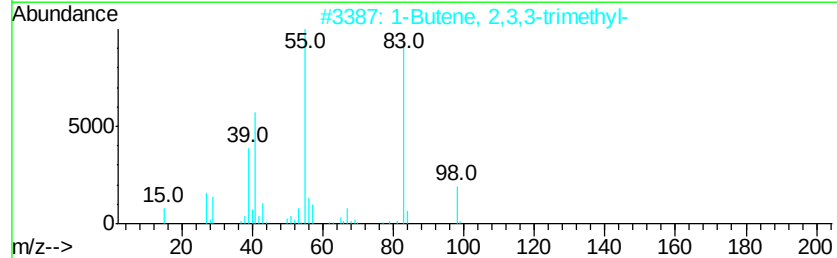
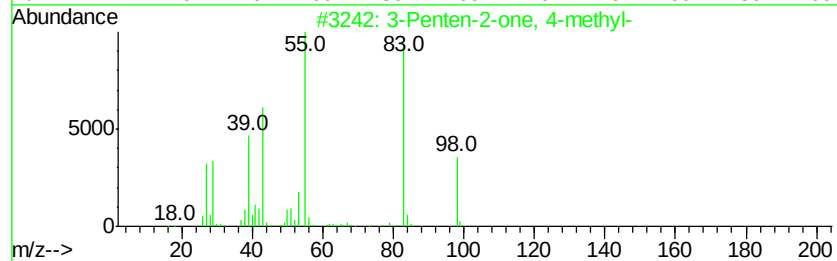
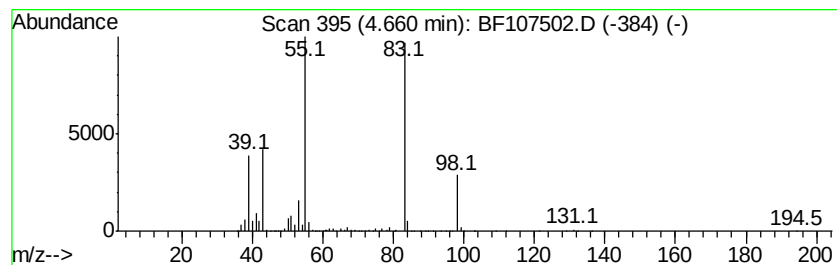
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071618.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 2 3-Penten-2-one, 4-methyl- Concentration Rank 1

| R.T. | EstConc   | Area     | Relative to ISTD       | R.T. |
|------|-----------|----------|------------------------|------|
| 4.66 | 232.35 ng | 10241600 | 1,4-Dichlorobenzene-d4 | 6.98 |

| Hit# | of | 5 | Tentative ID                   | MW | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|----|---------|-------------|------|
| 1    |    |   | 3-Penten-2-one, 4-methyl-      | 98 | C6H10O  | 000141-79-7 | 94   |
| 2    |    |   | 1-Butene, 2,3,3-trimethyl-     | 98 | C7H14   | 000594-56-9 | 86   |
| 3    |    |   | 2-Pentene, 2,3-dimethyl-       | 98 | C7H14   | 010574-37-5 | 83   |
| 4    |    |   | 2-Pentene, 2,4-dimethyl-       | 98 | C7H14   | 000625-65-0 | 64   |
| 5    |    |   | 2-Pentene, 3,4-dimethyl-, (E)- | 98 | C7H14   | 004914-92-5 | 64   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071818\  
Data File : BF107502.D  
Acq On : 19 Jul 2018 8:28  
Operator : JU/SJ  
Sample : J4033-03  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
071618-DBRI-F-D-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071618.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

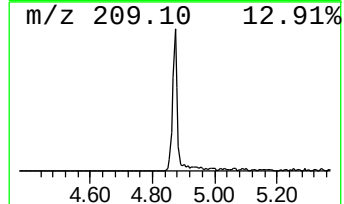
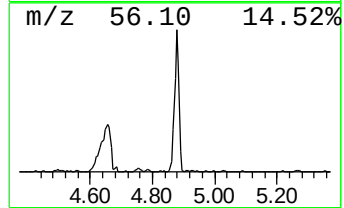
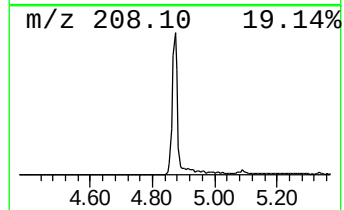
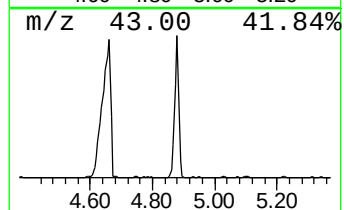
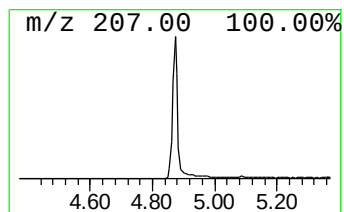
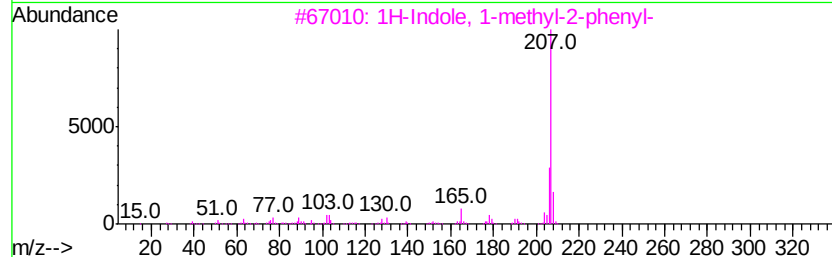
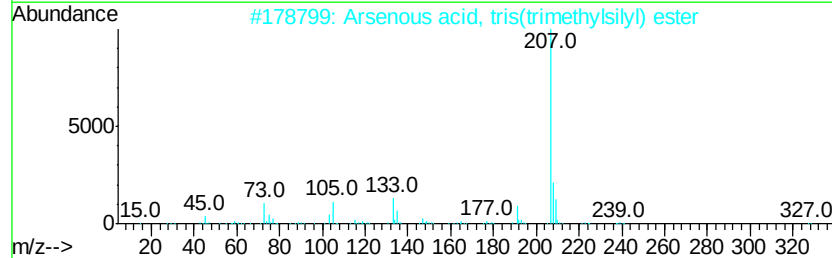
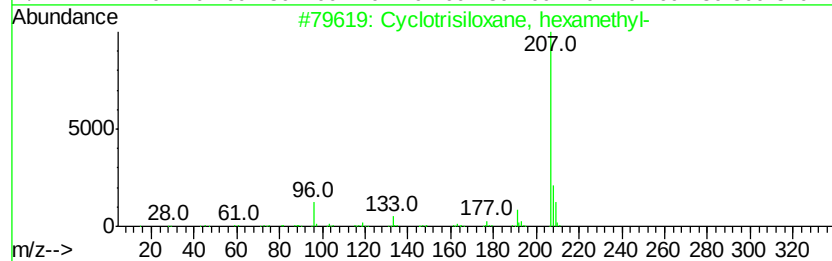
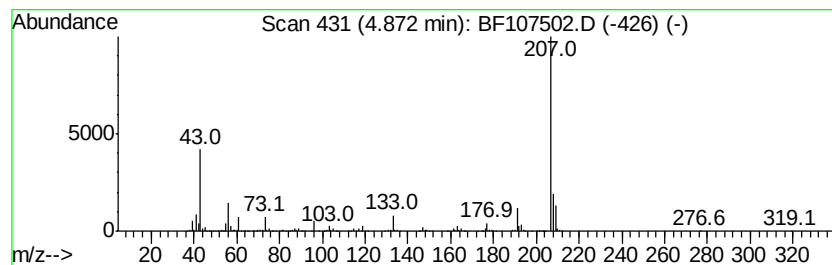
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 Cyclotrisiloxane, hexamethyl- Concentration Rank 4

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 4.87 | 81.05 ng | 3572420 | 1,4-Dichlorobenzene-d4 | 6.98 |

| Hit# | of | Tentative ID                        | MW  | MolForm      | CAS#        | Qual |
|------|----|-------------------------------------|-----|--------------|-------------|------|
| 1    | 5  | Cyclotrisiloxane, hexamethyl-       | 222 | C6H18O3Si3   | 000541-05-9 | 90   |
| 2    |    | Arsenous acid, tris(trimethylsil... | 342 | C9H27AsO3Si3 | 055429-29-3 | 64   |
| 3    |    | 1H-Indole, 1-methyl-2-phenyl-       | 207 | C15H13N      | 003558-24-5 | 42   |
| 4    |    | 4-Cyclohexene-1,2-dicarboximide,... | 207 | C12H17N02    | 028916-00-9 | 42   |
| 5    |    | Benzo[h]quinoline, 2,4-dimethyl-    | 207 | C15H13N      | 000605-67-4 | 39   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071818\  
Data File : BF107502.D  
Acq On : 19 Jul 2018 8:28  
Operator : JU/SJ  
Sample : J4033-03  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
071618-DBRI-F-D-B

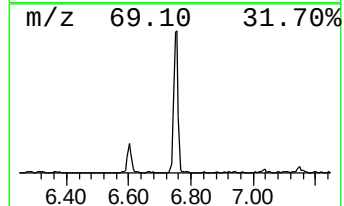
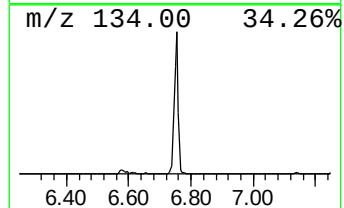
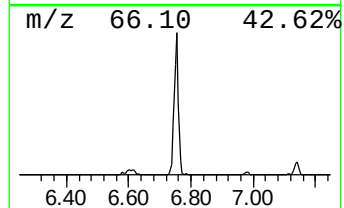
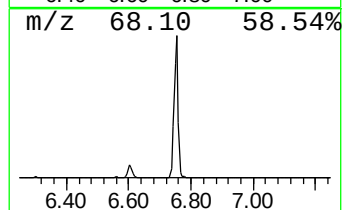
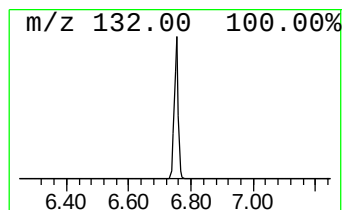
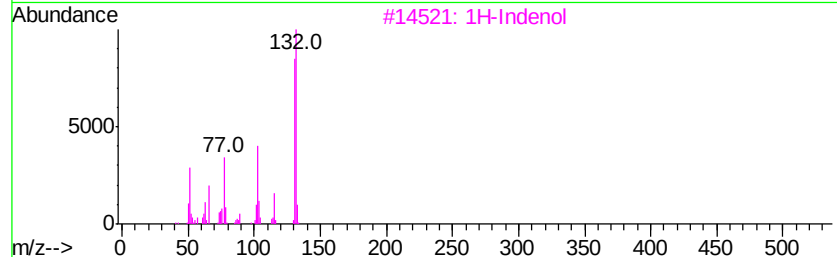
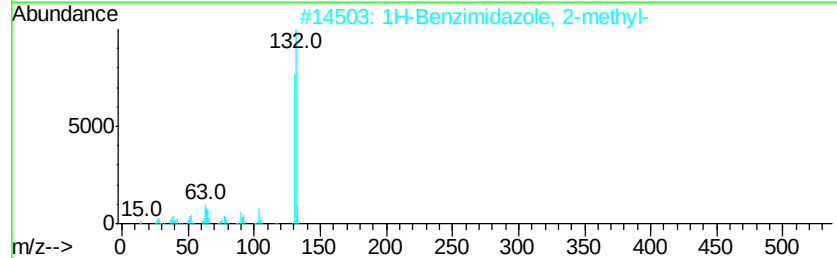
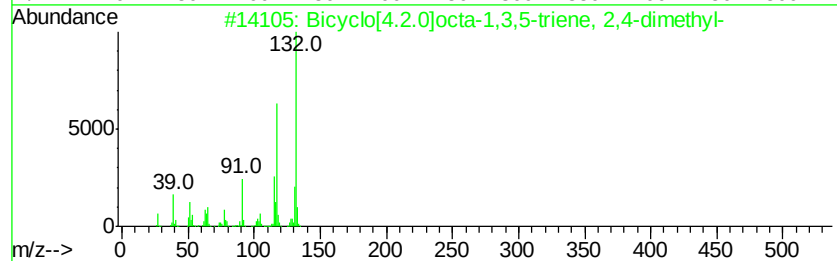
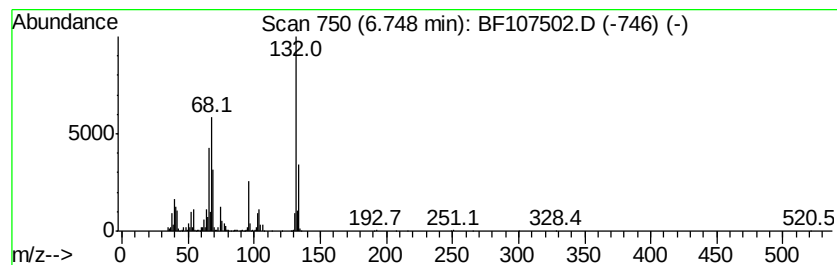
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071618.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 4 unknown6.75 Concentration Rank 8

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.75 | 54.84 ng | 2417260 | 1,4-Dichlorobenzene-d4 | 6.98 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | Bicyclo[4.2.0]octa-1,3,5-triene,... | 132 | C10H12   | 028749-81-7  | 14   |
| 2    |    | 1H-Benzimidazole, 2-methyl-         | 132 | C8H8N2   | 000615-15-6  | 14   |
| 3    |    | 1H-Indenol                          | 132 | C9H8O    | 056631-57-3  | 14   |
| 4    |    | 1H-Indazole, 7-methyl-              | 132 | C8H8N2   | 1000316-00-7 | 10   |
| 5    |    | 1,3,4-Thiadiazole-2(3H)-thione, ... | 132 | C3H4N2S2 | 029490-19-5  | 10   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071818\  
Data File : BF107502.D  
Acq On : 19 Jul 2018 8:28  
Operator : JU/SJ  
Sample : J4033-03  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
071618-DBRI-F-D-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071618.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

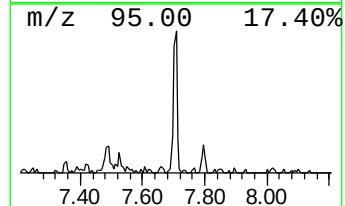
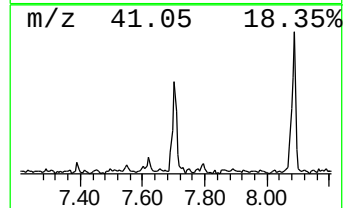
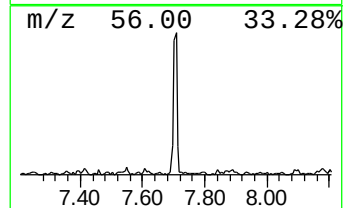
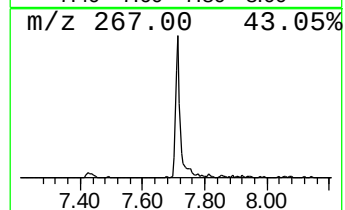
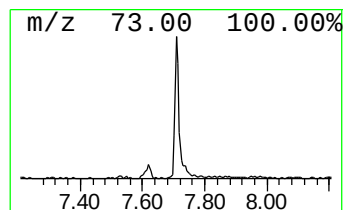
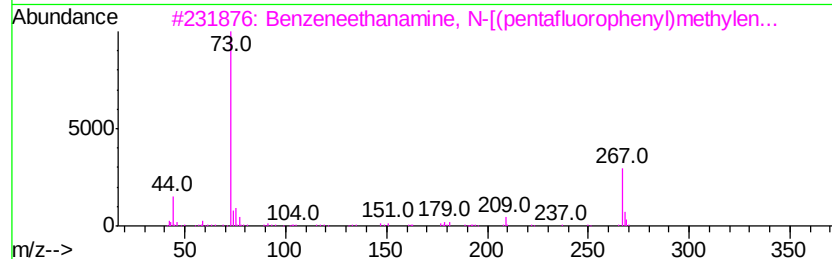
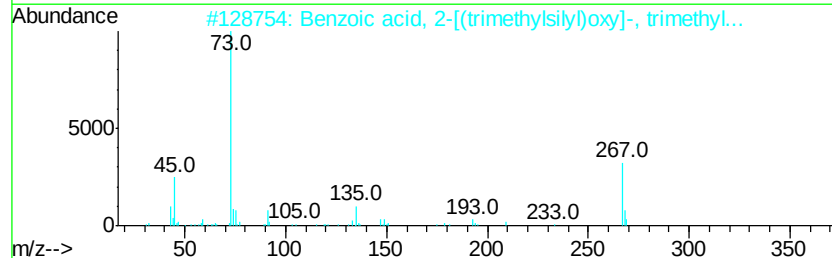
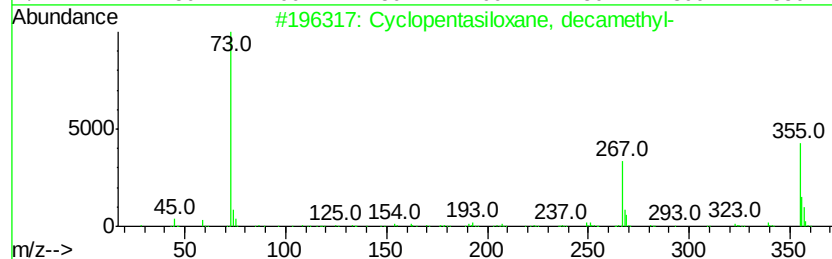
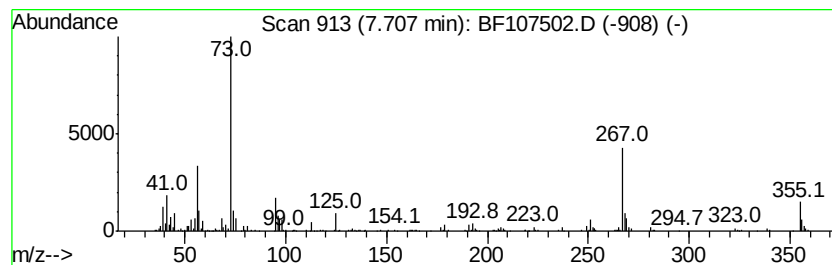
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 6 Cyclopentasiloxane, decamet... Concentration Rank 14

| R.T. | EstConc  | Area    | Relative to ISTD | R.T. |
|------|----------|---------|------------------|------|
| 7.71 | 19.16 ng | 1013140 | Naphthalene-d8   | 8.27 |

| Hit# | of | Tentative ID                                           | MW  | MolForm        | CAS#         | Qual |
|------|----|--------------------------------------------------------|-----|----------------|--------------|------|
| 1    | 5  | Cyclopentasiloxane, decamethyl-                        | 370 | C10H30O5Si5    | 000541-02-6  | 53   |
| 2    |    | Benzoic acid, 2-[(trimethylsilyl)oxy]-, trimethyl-     | 282 | C13H22O3Si2    | 003789-85-3  | 37   |
| 3    |    | Benzeneethanamine, N-[(pentafluorophenyl)methyl]-      | 475 | C21H26F5N02Si2 | 055429-85-1  | 37   |
| 4    |    | 3-Amino-2-phenazolin ditms                             | 355 | C18H25N3O3Si2  | 1000332-15-5 | 10   |
| 5    |    | Phosphonic acid, [3-[(trimethylsilyl)oxy]-, trimethyl- | 355 | C12H34N03PSi3  | 056051-92-4  | 10   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071818\  
Data File : BF107502.D  
Acq On : 19 Jul 2018 8:28  
Operator : JU/SJ  
Sample : J4033-03  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
071618-DBRI-F-D-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071618.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

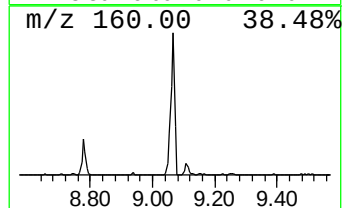
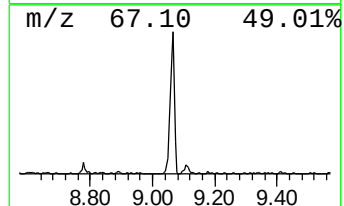
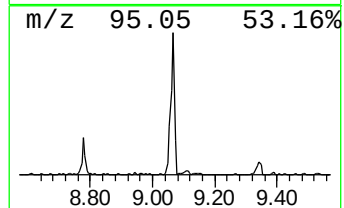
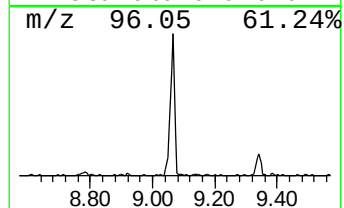
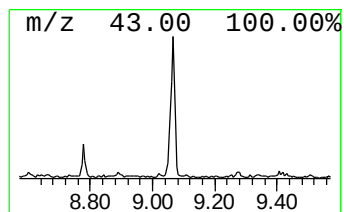
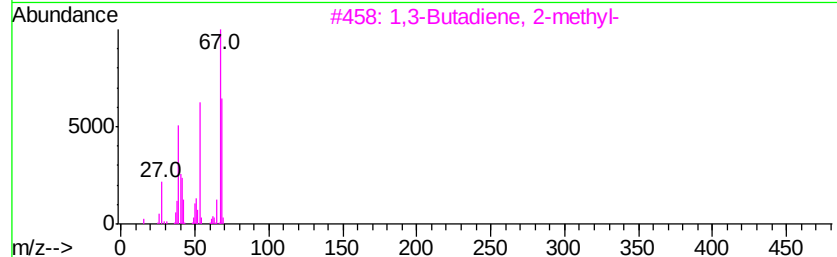
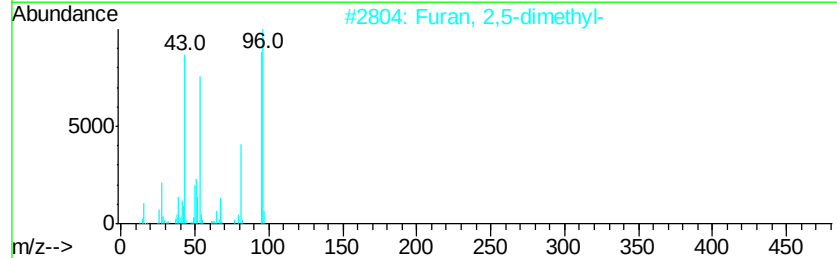
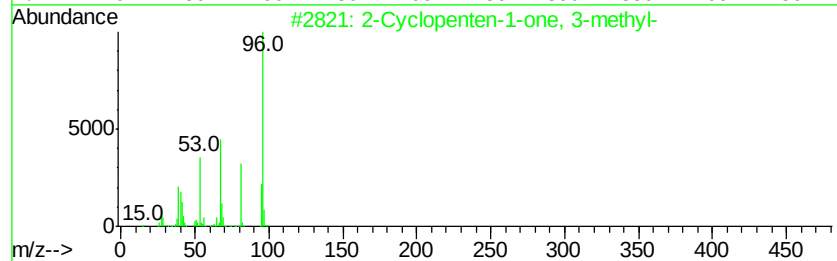
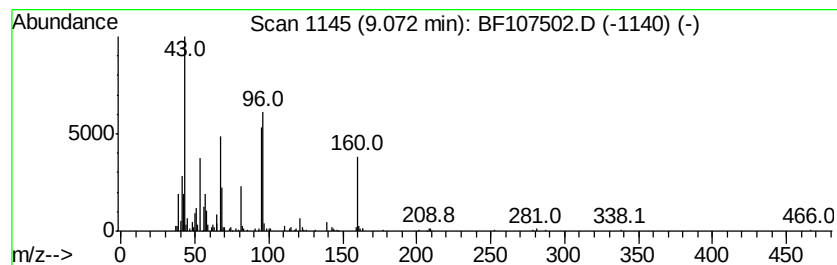
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 2-Cyclopenten-1-one, 3-methyl- Concentration Rank 17

| R.T. | EstConc  | Area   | Relative to ISTD | R.T. |
|------|----------|--------|------------------|------|
| 9.07 | 16.61 ng | 878597 | Napthalene-d8    | 8.27 |

| Hit# | of | Tentative ID                        | MW | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|----|---------|-------------|------|
| 1    | 5  | 2-Cyclopenten-1-one, 3-methyl-      | 96 | C6H8O   | 002758-18-1 | 93   |
| 2    |    | Furan, 2,5-dimethyl-                | 96 | C6H8O   | 000625-86-5 | 76   |
| 3    |    | 1,3-Butadiene, 2-methyl-            | 68 | C5H8    | 000078-79-5 | 50   |
| 4    |    | 1-Butyne, 3-methyl-                 | 68 | C5H8    | 000598-23-2 | 50   |
| 5    |    | Ethanone, 1-(methylenecyclopropyl)- | 96 | C6H8O   | 062266-35-7 | 50   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071818\  
Data File : BF107502.D  
Acq On : 19 Jul 2018 8:28  
Operator : JU/SJ  
Sample : J4033-03  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
071618-DBRI-F-D-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071618.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

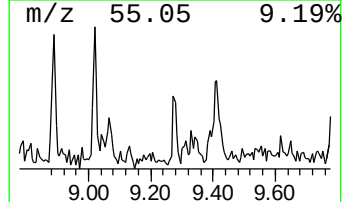
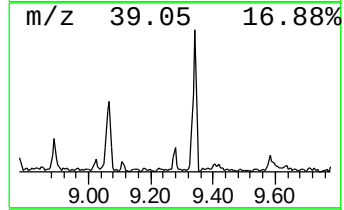
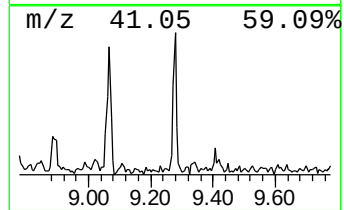
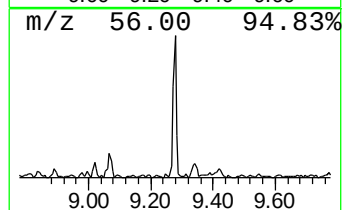
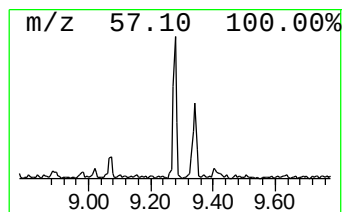
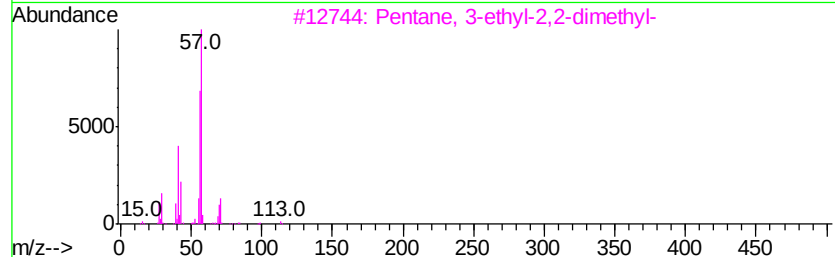
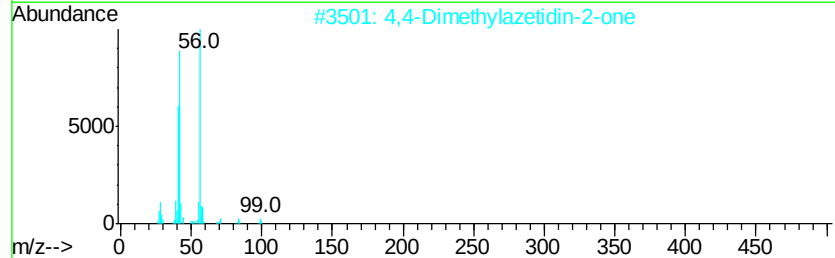
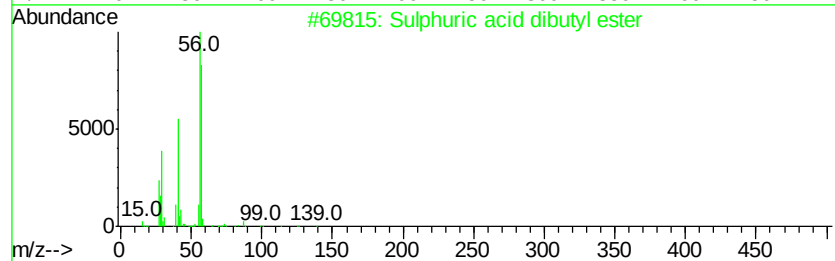
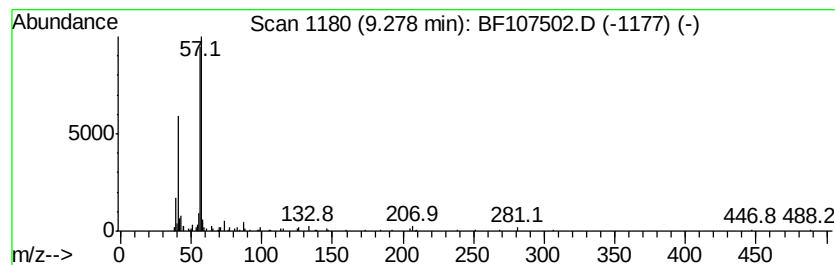
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 Sulphuric acid dibutyl ester Concentration Rank 20

| R.T. | EstConc  | Area   | Relative to ISTD | R.T.  |
|------|----------|--------|------------------|-------|
| 9.28 | 14.58 ng | 159511 | Acenaphthene-d10 | 10.02 |

| Hit# | of | Tentative ID                   | MW  | MolForm  | CAS#        | Qual |
|------|----|--------------------------------|-----|----------|-------------|------|
| 1    | 5  | Sulphuric acid dibutyl ester   | 210 | C8H18O4S | 000625-22-9 | 59   |
| 2    |    | 4,4-Dimethylazetidin-2-one     | 99  | C5H9NO   | 004879-95-2 | 50   |
| 3    |    | Pentane, 3-ethyl-2,2-dimethyl- | 128 | C9H20    | 016747-32-3 | 50   |
| 4    |    | Pentane, 3-methyl-             | 86  | C6H14    | 000096-14-0 | 50   |
| 5    |    | Pentane, 2,2,3-trimethyl-      | 114 | C8H18    | 000564-02-3 | 50   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071818\  
Data File : BF107502.D  
Acq On : 19 Jul 2018 8:28  
Operator : JU/SJ  
Sample : J4033-03  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
071618-DBRI-F-D-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071618.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

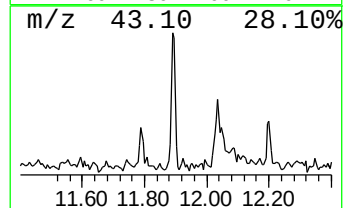
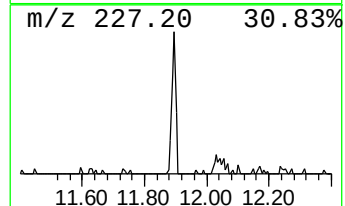
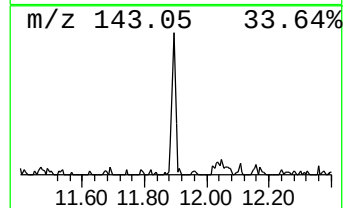
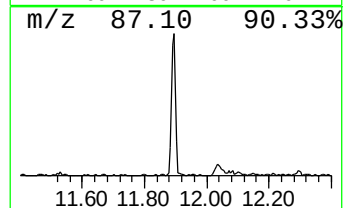
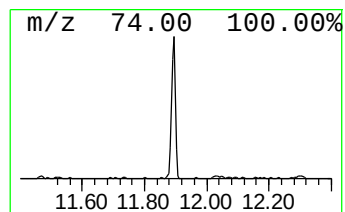
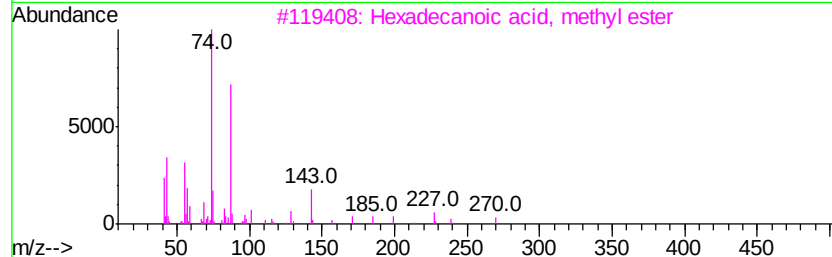
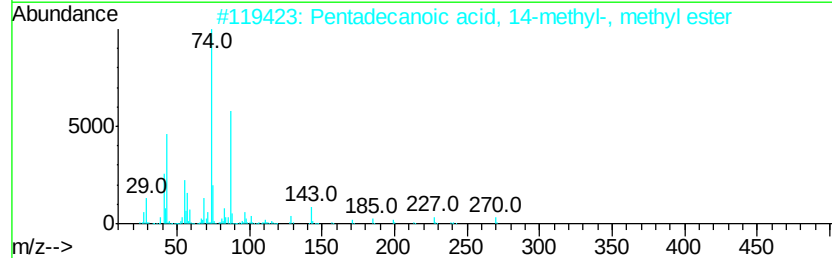
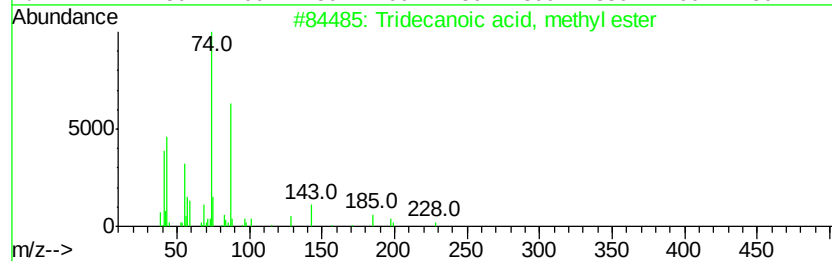
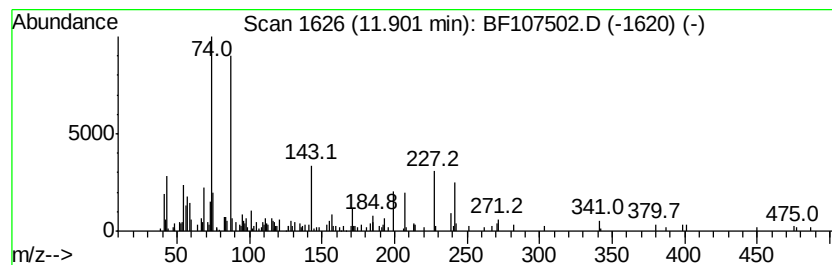
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 9 Tridecanoic acid, methyl ester Concentration Rank 18

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 11.90 | 16.60 ng | 544576 | Phenanthrene-d10 | 11.52 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Tridecanoic acid, methyl ester      | 228 | C14H28O2 | 001731-88-0 | 95   |
| 2    |    | Pentadecanoic acid, 14-methyl-, ... | 270 | C17H34O2 | 005129-60-2 | 90   |
| 3    |    | Hexadecanoic acid, methyl ester     | 270 | C17H34O2 | 000112-39-0 | 80   |
| 4    |    | Methyl stearate                     | 298 | C19H38O2 | 000112-61-8 | 80   |
| 5    |    | Decanoic acid, methyl ester         | 186 | C11H22O2 | 000110-42-9 | 72   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071818\  
Data File : BF107502.D  
Acq On : 19 Jul 2018 8:28  
Operator : JU/SJ  
Sample : J4033-03  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
071618-DBRI-F-D-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071618.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

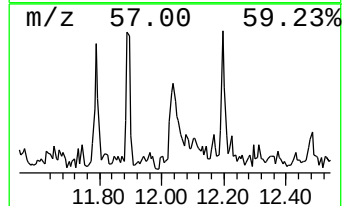
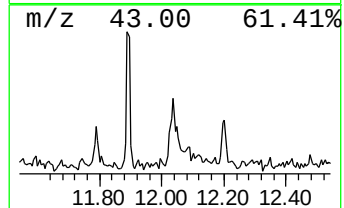
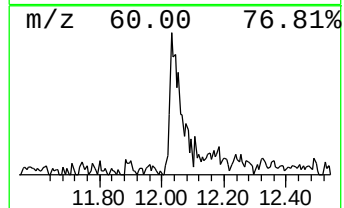
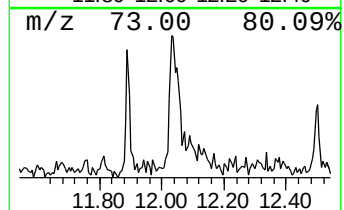
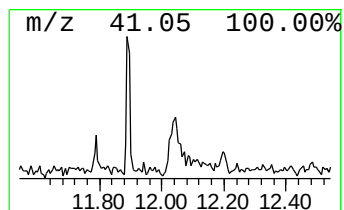
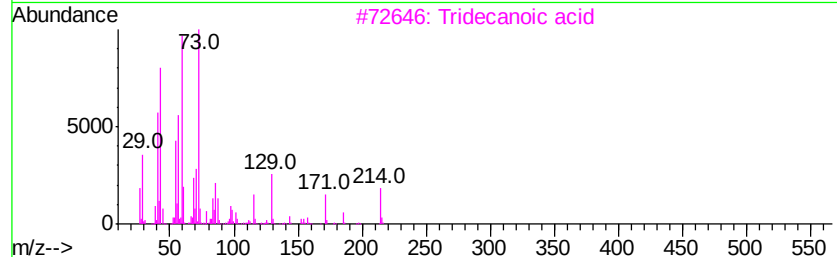
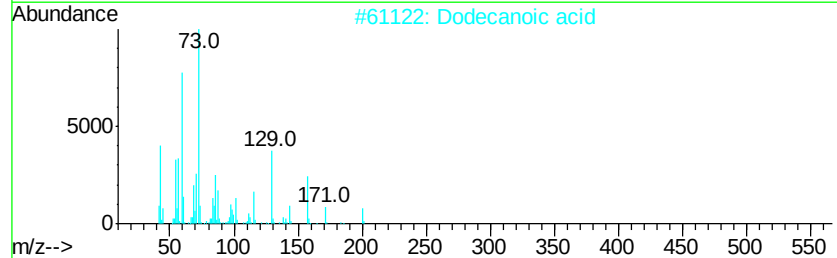
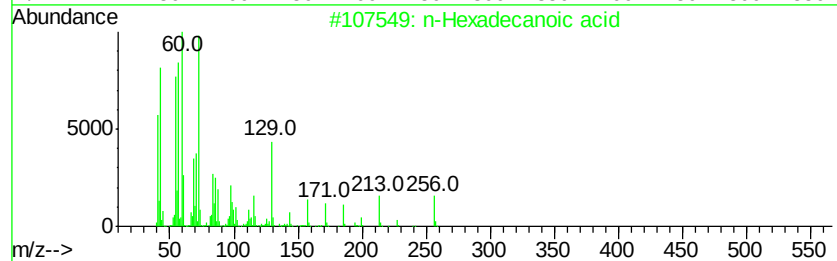
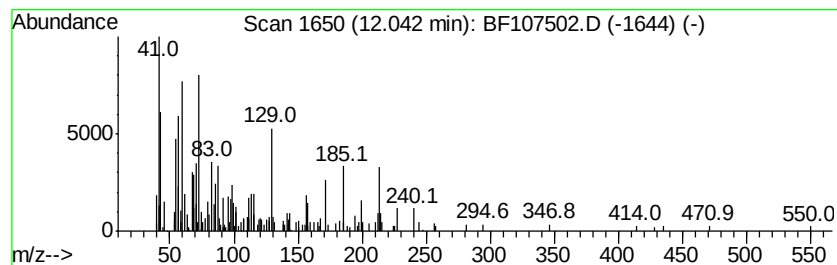
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 n-Hexadecanoic acid Concentration Rank 16

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 12.04 | 16.85 ng | 552676 | Phenanthrene-d10 | 11.52 |

| Hit# | of | Tentative ID        | MW  | MolForm  | CAS#        | Qual |
|------|----|---------------------|-----|----------|-------------|------|
| 1    | 5  | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 50   |
| 2    |    | Dodecanoic acid     | 200 | C12H24O2 | 000143-07-7 | 50   |
| 3    |    | Tridecanoic acid    | 214 | C13H26O2 | 000638-53-9 | 46   |
| 4    |    | Tetradecanoic acid  | 228 | C14H28O2 | 000544-63-8 | 45   |
| 5    |    | Pentadecanoic acid  | 242 | C15H30O2 | 001002-84-2 | 43   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071818\  
Data File : BF107502.D  
Acq On : 19 Jul 2018 8:28  
Operator : JU/SJ  
Sample : J4033-03  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
071618-DBRI-F-D-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071618.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

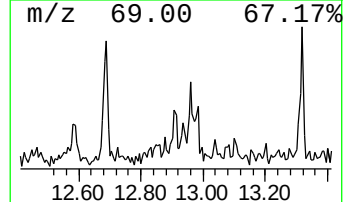
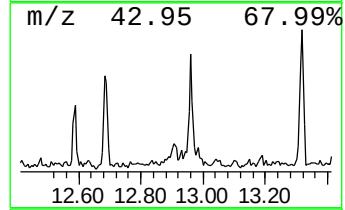
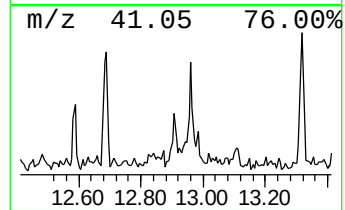
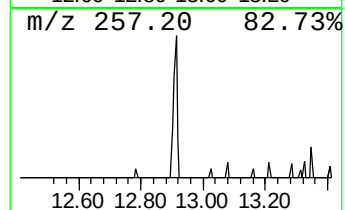
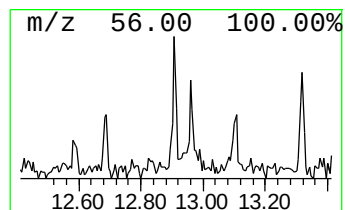
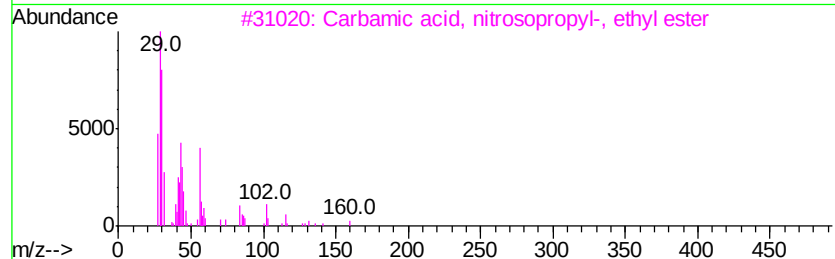
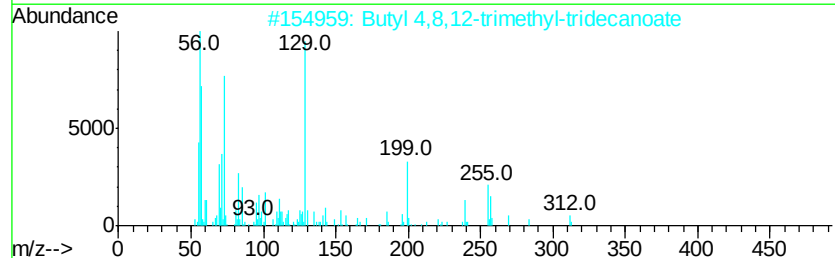
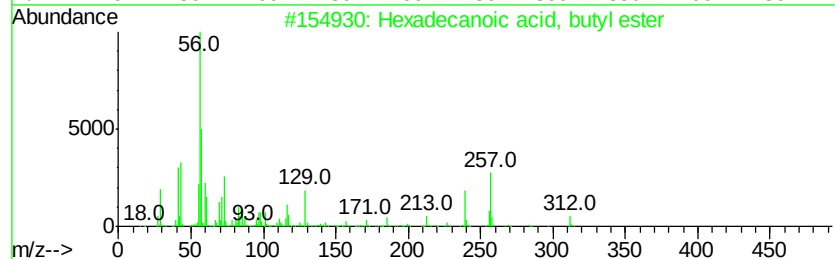
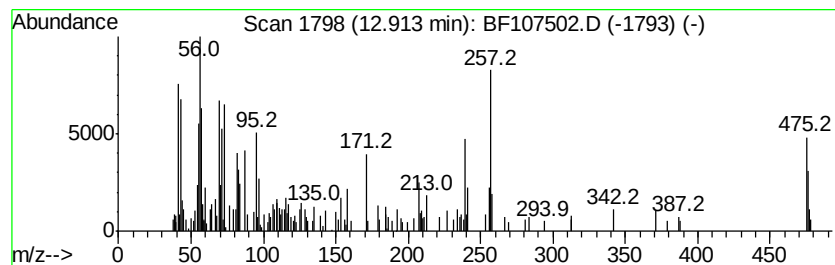
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 unknown12.91 Concentration Rank 12

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 12.91 | 21.25 ng | 202416 | Chrysene-d12     | 14.17 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Hexadecanoic acid, butyl ester      | 312 | C20H40O2  | 000111-06-8  | 45   |
| 2    |    | Butyl 4,8,12-trimethyl-tridecanoate | 312 | C20H40O2  | 1000336-63-2 | 38   |
| 3    |    | Carbamic acid, nitrosopropyl-, e... | 160 | C6H12N2O3 | 019935-86-5  | 12   |
| 4    |    | L-Homocitrulline                    | 189 | C7H15N3O3 | 001190-49-4  | 10   |
| 5    |    | 1,5-Pentanediol                     | 104 | C5H12O2   | 000111-29-5  | 10   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071818\  
Data File : BF107502.D  
Acq On : 19 Jul 2018 8:28  
Operator : JU/SJ  
Sample : J4033-03  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
071618-DBRI-F-D-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071618.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

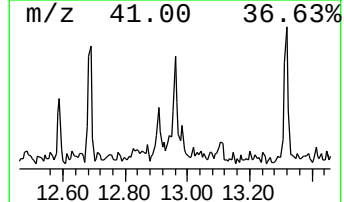
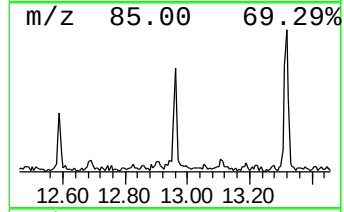
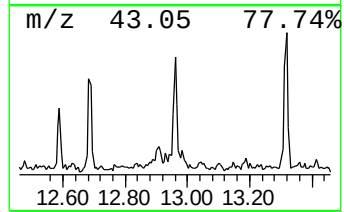
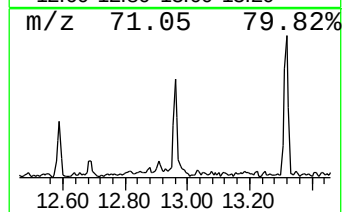
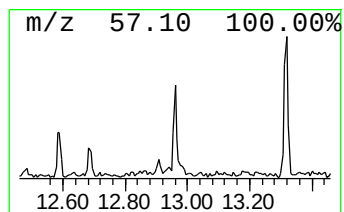
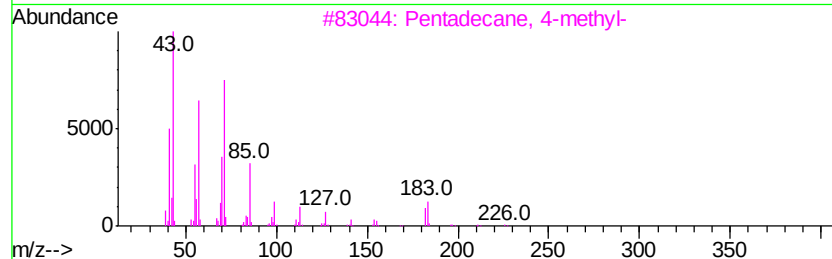
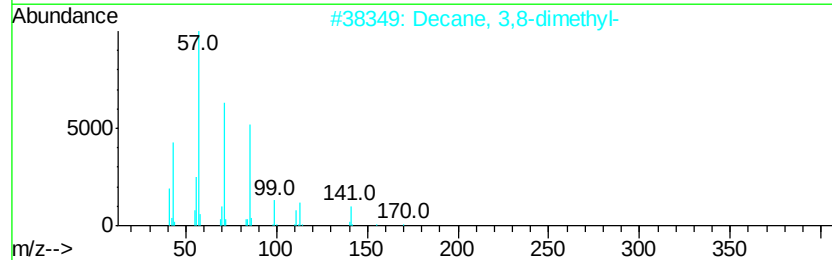
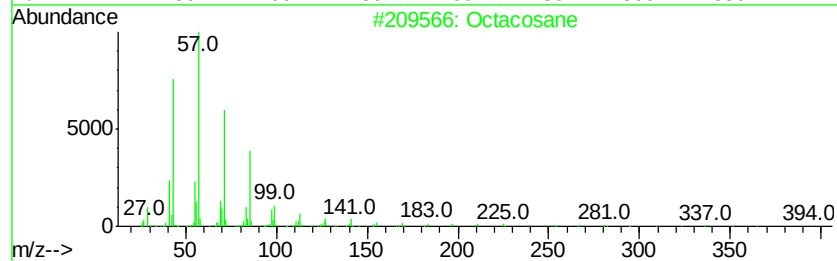
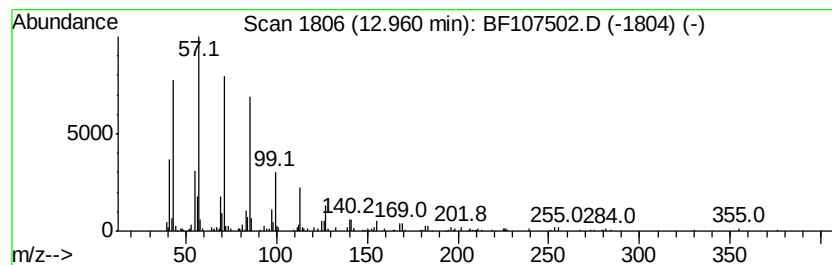
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 Octacosane Concentration Rank 11

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 12.96 | 21.52 ng | 204927 | Chrysene-d12     | 14.17 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Octacosane                          | 394 | C28H58    | 000630-02-4  | 72   |
| 2    |    | Decane, 3,8-dimethyl-               | 170 | C12H26    | 017312-55-9  | 64   |
| 3    |    | Pentadecane, 4-methyl-              | 226 | C16H34    | 002801-87-8  | 64   |
| 4    |    | Nonadecane                          | 268 | C19H40    | 000629-92-5  | 64   |
| 5    |    | Sulfurous acid, decyl 2-propyl e... | 264 | C13H28O3S | 1000309-12-1 | 64   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071818\  
Data File : BF107502.D  
Acq On : 19 Jul 2018 8:28  
Operator : JU/SJ  
Sample : J4033-03  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
071618-DBRI-F-D-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071618.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

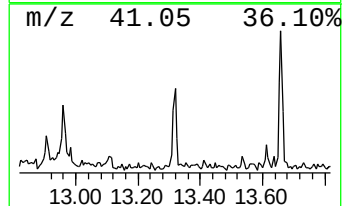
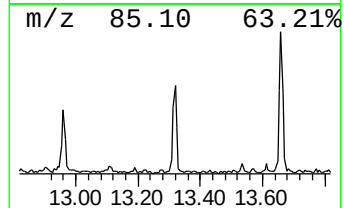
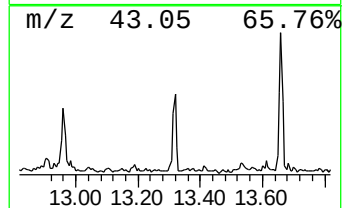
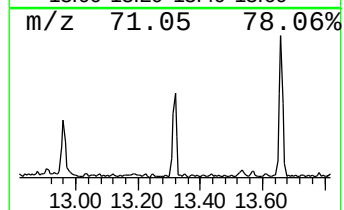
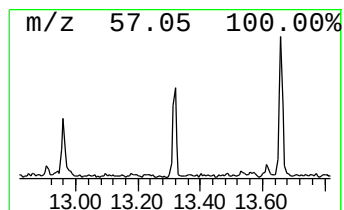
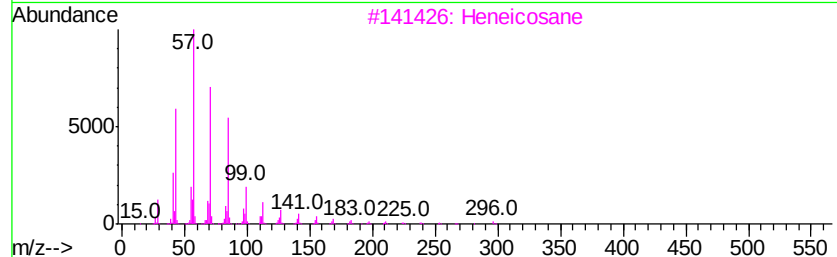
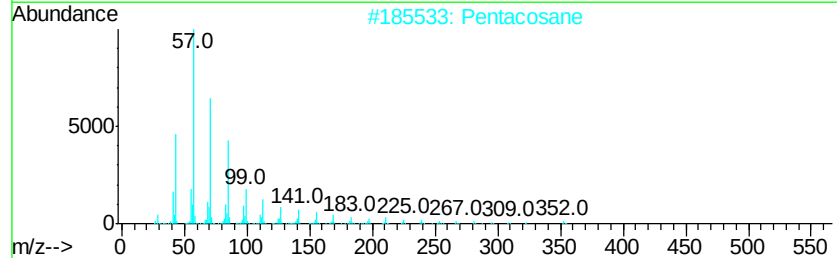
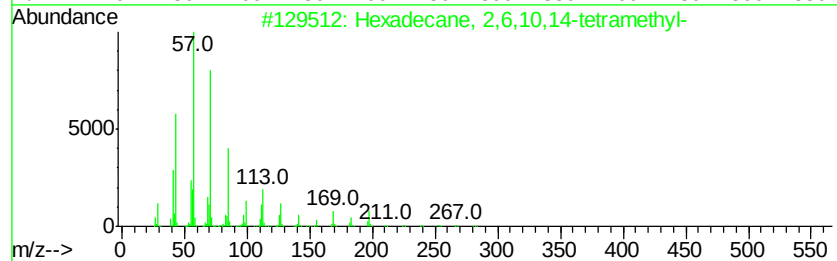
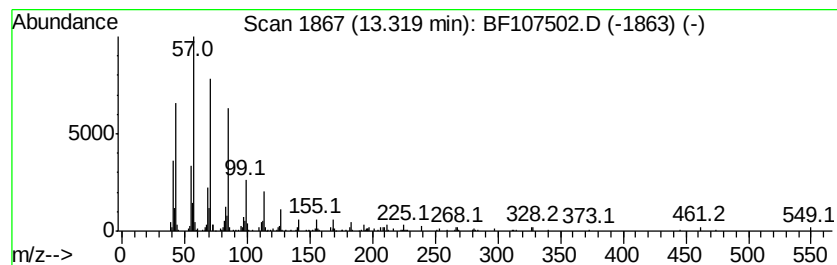
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 Hexadecane, 2,6,10,14-tetra... Concentration Rank 9

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 13.32 | 44.81 ng | 426746 | Chrysene-d12     | 14.17 |

| Hit# | of | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Hexadecane, 2,6,10,14-tetramethyl- | 282 | C20H42  | 000638-36-8 | 90   |
| 2    |    | Pentacosane                        | 352 | C25H52  | 000629-99-2 | 87   |
| 3    |    | Heneicosane                        | 296 | C21H44  | 000629-94-7 | 83   |
| 4    |    | triacontane                        | 422 | C30H62  | 000638-68-6 | 83   |
| 5    |    | Octadecane, 2-methyl-              | 268 | C19H40  | 001560-88-9 | 83   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071818\  
Data File : BF107502.D  
Acq On : 19 Jul 2018 8:28  
Operator : JU/SJ  
Sample : J4033-03  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
071618-DBRI-F-D-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071618.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

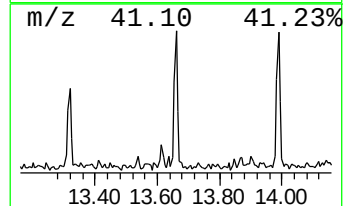
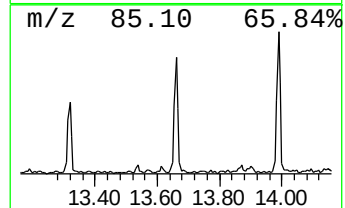
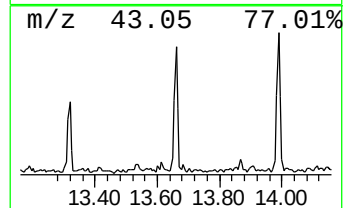
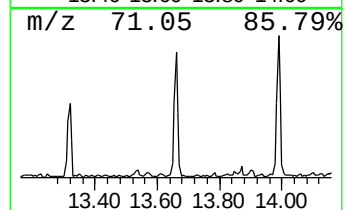
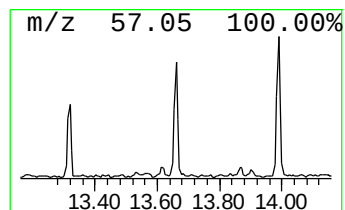
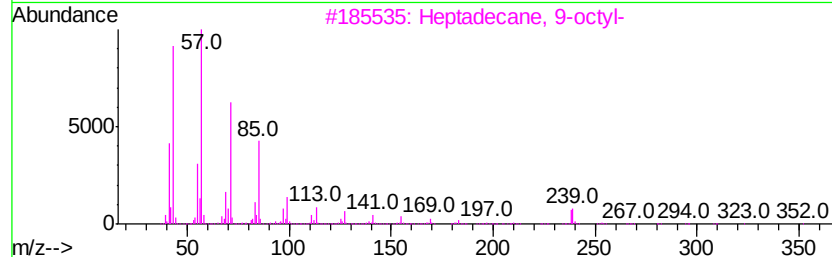
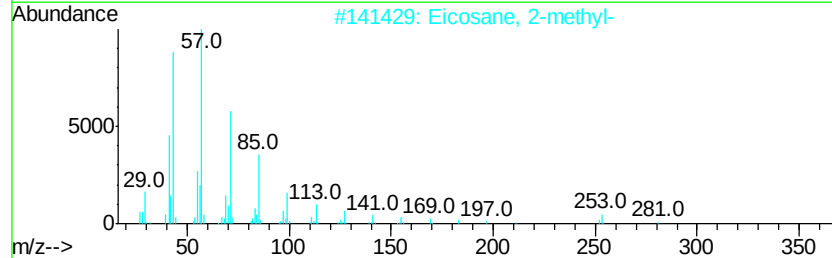
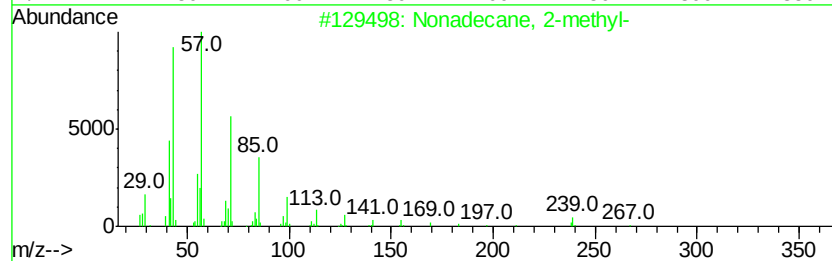
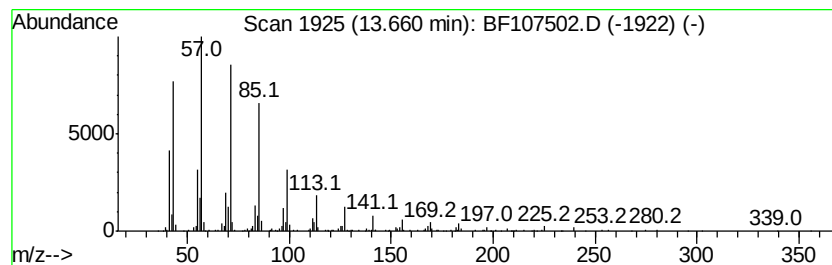
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 Nonadecane, 2-methyl- Concentration Rank 7

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 13.66 | 64.24 ng | 611802 | Chrysene-d12     | 14.17 |

| Hit# | of | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------|-----|---------|-------------|------|
| 1    | 5  | Nonadecane, 2-methyl- | 282 | C20H42  | 001560-86-7 | 90   |
| 2    |    | Eicosane, 2-methyl-   | 296 | C21H44  | 001560-84-5 | 90   |
| 3    |    | Heptadecane, 9-octyl- | 352 | C25H52  | 007225-64-1 | 87   |
| 4    |    | Octacosane            | 394 | C28H58  | 000630-02-4 | 87   |
| 5    |    | Heptacosane           | 380 | C27H56  | 000593-49-7 | 87   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071818\  
Data File : BF107502.D  
Acq On : 19 Jul 2018 8:28  
Operator : JU/SJ  
Sample : J4033-03  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
071618-DBRI-F-D-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071618.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

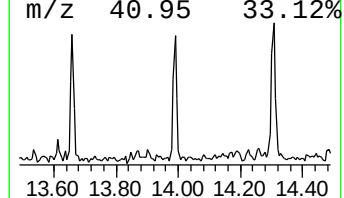
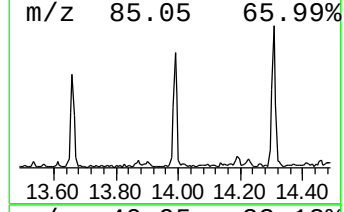
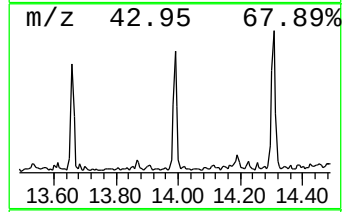
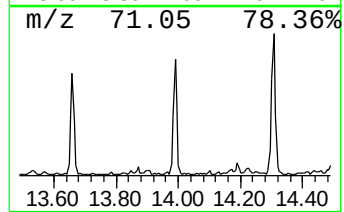
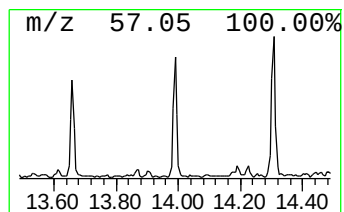
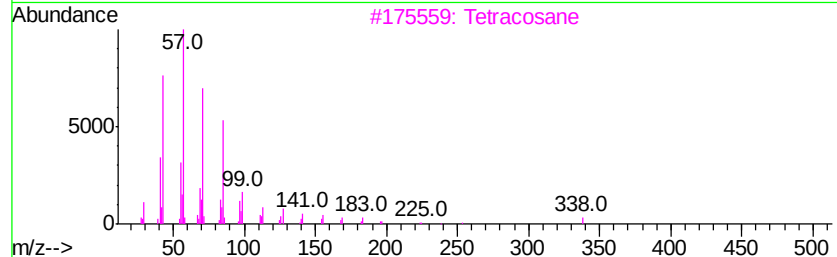
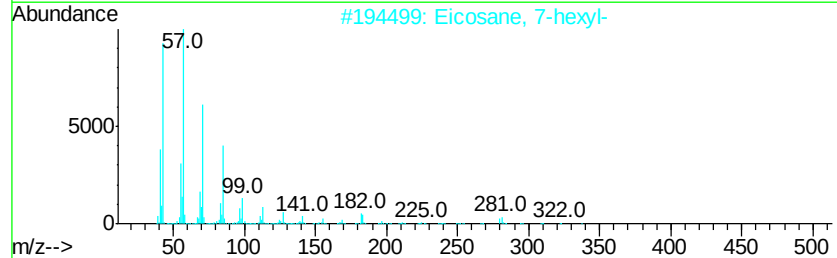
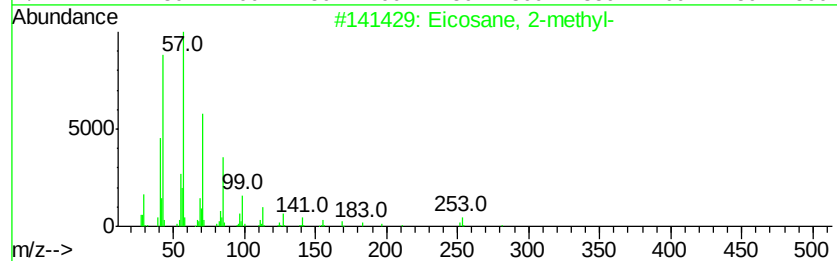
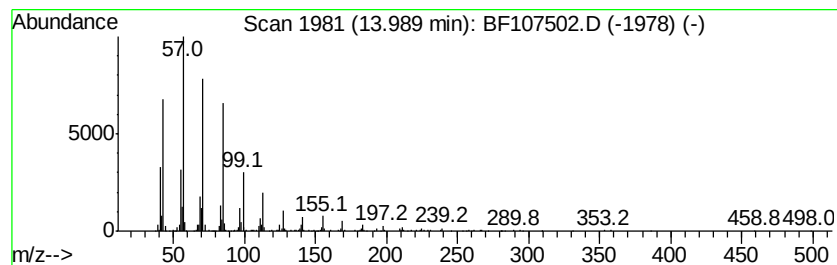
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 Eicosane, 2-methyl- Concentration Rank 5

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 13.99 | 75.75 ng | 721409 | Chrysene-d12     | 14.17 |

| Hit# | of | Tentative ID        | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------|-----|---------|-------------|------|
| 1    | 5  | Eicosane, 2-methyl- | 296 | C21H44  | 001560-84-5 | 90   |
| 2    |    | Eicosane, 7-hexyl-  | 366 | C26H54  | 055333-99-8 | 86   |
| 3    |    | Tetracosane         | 338 | C24H50  | 000646-31-1 | 83   |
| 4    |    | Hexacosane          | 366 | C26H54  | 000630-01-3 | 81   |
| 5    |    | triacontane         | 422 | C30H62  | 000638-68-6 | 80   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071818\  
Data File : BF107502.D  
Acq On : 19 Jul 2018 8:28  
Operator : JU/SJ  
Sample : J4033-03  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
071618-DBRI-F-D-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071618.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

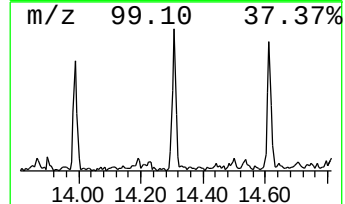
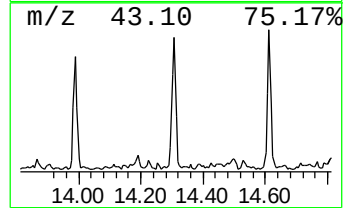
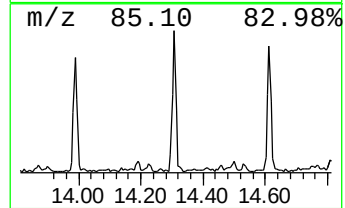
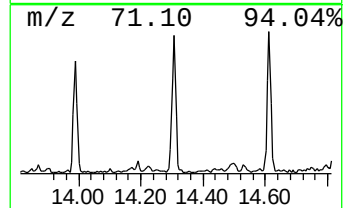
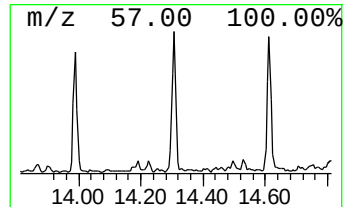
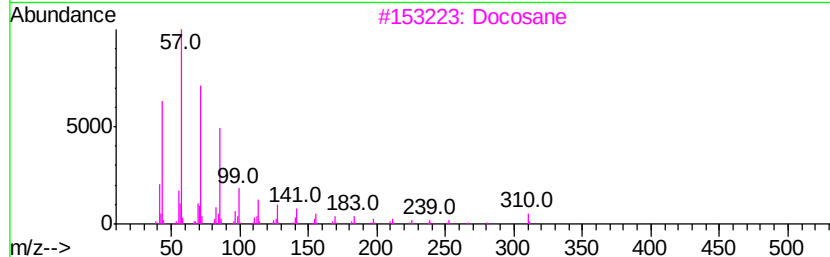
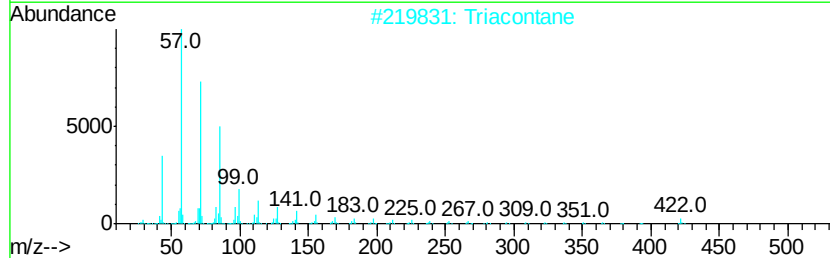
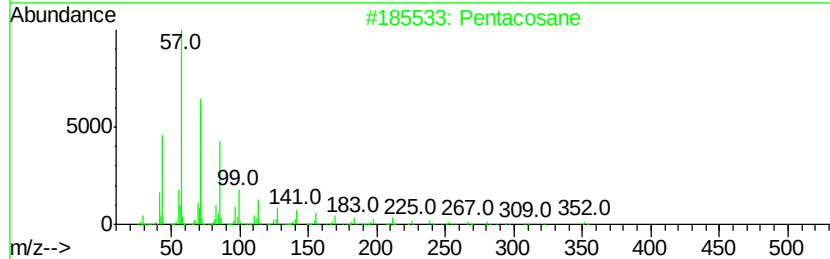
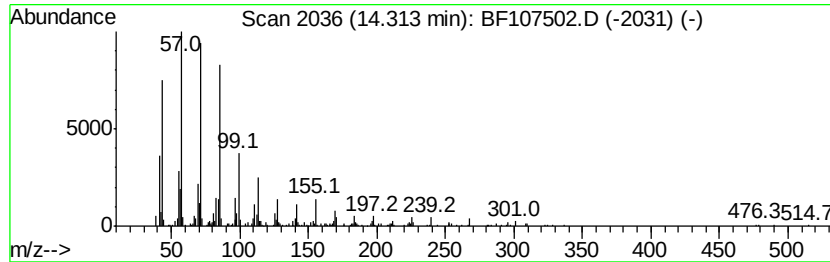
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 Pentacosane Concentration Rank 3

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 14.31 | 97.09 ng | 924678 | Chrysene-d12     | 14.17 |

| Hit# | of | Tentative ID        | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------|-----|---------|-------------|------|
| 1    | 5  | Pentacosane         | 352 | C25H52  | 000629-99-2 | 83   |
| 2    |    | Triacontane         | 422 | C30H62  | 000638-68-6 | 83   |
| 3    |    | Docosane            | 310 | C22H46  | 000629-97-0 | 83   |
| 4    |    | Hexadecane, 1-iodo- | 352 | C16H33I | 000544-77-4 | 80   |
| 5    |    | Heneicosane         | 296 | C21H44  | 000629-94-7 | 80   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071818\  
Data File : BF107502.D  
Acq On : 19 Jul 2018 8:28  
Operator : JU/SJ  
Sample : J4033-03  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
071618-DBRI-F-D-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071618.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

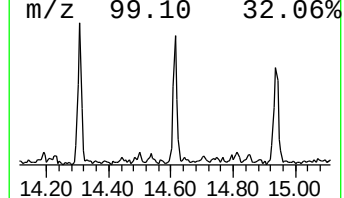
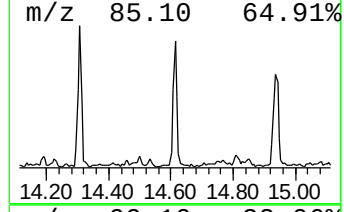
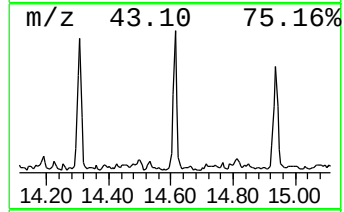
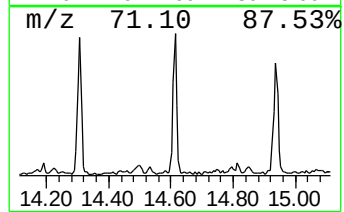
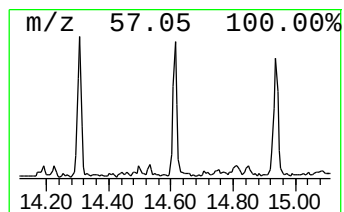
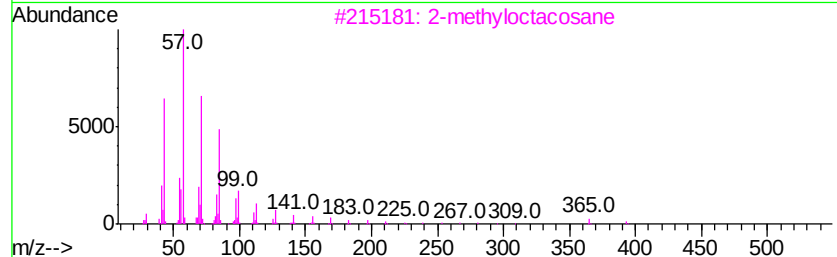
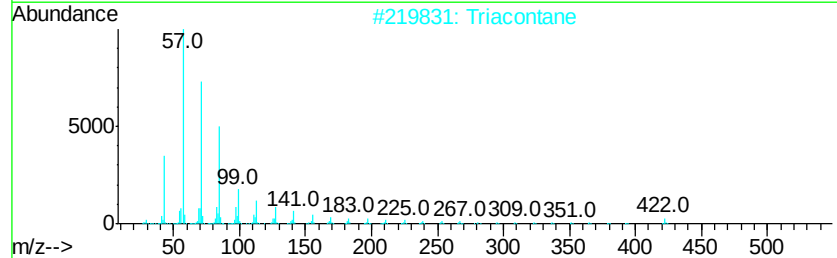
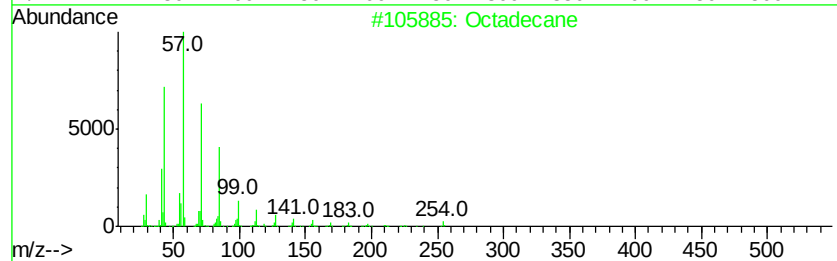
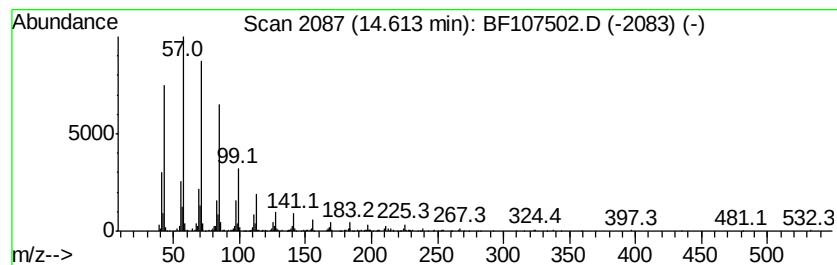
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 17 Octadecane Concentration Rank 2

| R.T.  | EstConc   | Area   | Relative to ISTD | R.T.  |
|-------|-----------|--------|------------------|-------|
| 14.61 | 103.09 ng | 981808 | Chrysene-d12     | 14.17 |

| Hit# | of | Tentative ID       | MW  | MolForm | CAS#         | Qual |
|------|----|--------------------|-----|---------|--------------|------|
| 1    | 5  | Octadecane         | 254 | C18H38  | 000593-45-3  | 87   |
| 2    |    | Triacontane        | 422 | C30H62  | 000638-68-6  | 83   |
| 3    |    | 2-methyloctacosane | 408 | C29H60  | 1000376-72-8 | 83   |
| 4    |    | Hentriacontane     | 437 | C31H64  | 000630-04-6  | 81   |
| 5    |    | Heneicosane        | 296 | C21H44  | 000629-94-7  | 80   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071818\  
Data File : BF107502.D  
Acq On : 19 Jul 2018 8:28  
Operator : JU/SJ  
Sample : J4033-03  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
071618-DBRI-F-D-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071618.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

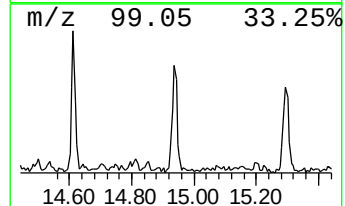
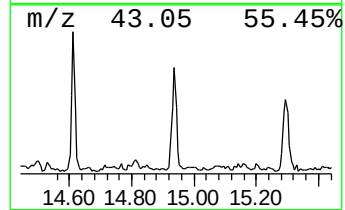
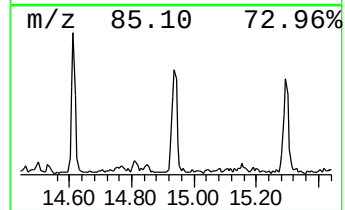
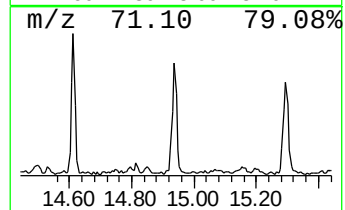
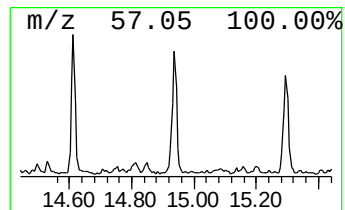
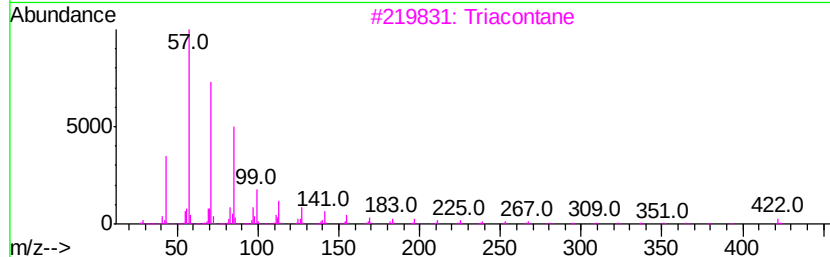
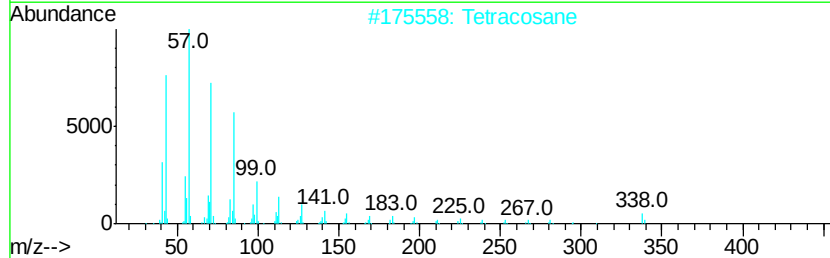
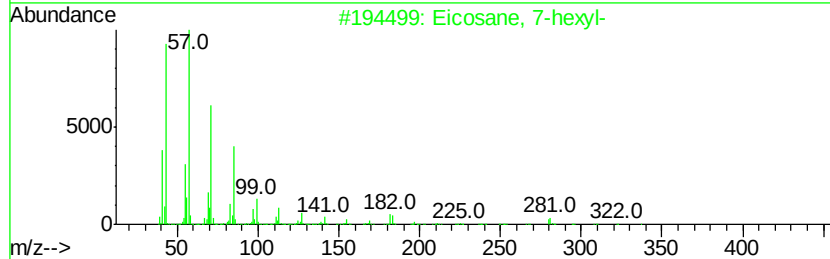
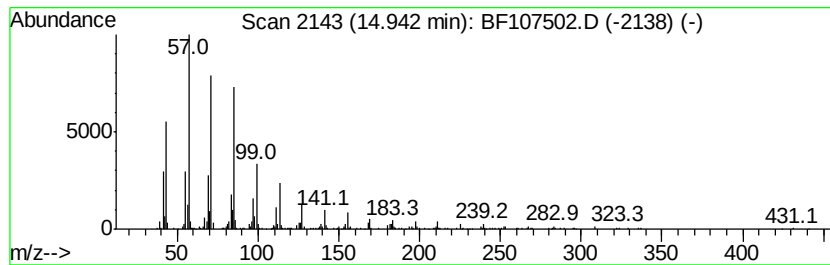
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 18 Eicosane, 7-hexyl- Concentration Rank 10

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 14.94 | 23.07 ng | 917982 | Perylene-d12     | 15.71 |

| Hit# | of | Tentative ID       | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------|-----|---------|-------------|------|
| 1    | 5  | Eicosane, 7-hexyl- | 366 | C26H54  | 055333-99-8 | 86   |
| 2    |    | Tetracosane        | 338 | C24H50  | 000646-31-1 | 86   |
| 3    |    | Triacontane        | 422 | C30H62  | 000638-68-6 | 80   |
| 4    |    | Docosane           | 310 | C22H46  | 000629-97-0 | 80   |
| 5    |    | Hentriacontane     | 437 | C31H64  | 000630-04-6 | 80   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071818\  
Data File : BF107502.D  
Acq On : 19 Jul 2018 8:28  
Operator : JU/SJ  
Sample : J4033-03  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
071618-DBRI-F-D-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071618.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

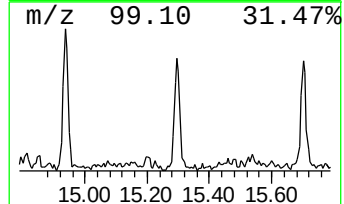
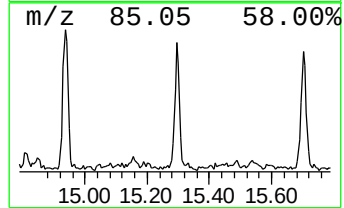
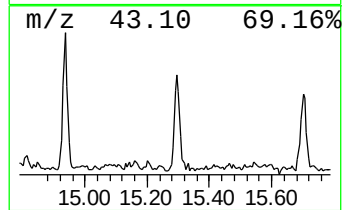
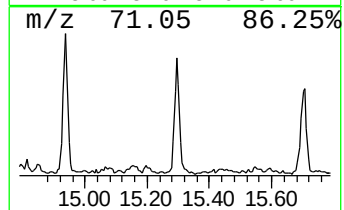
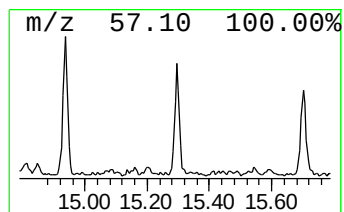
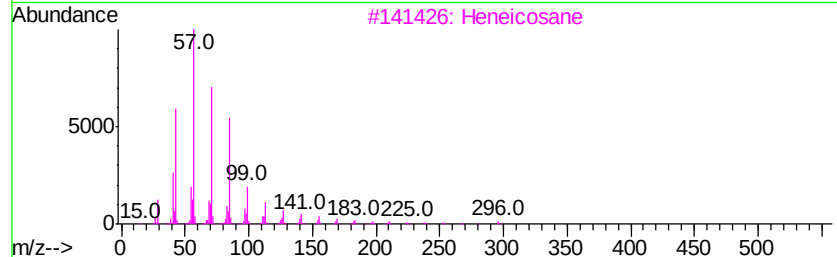
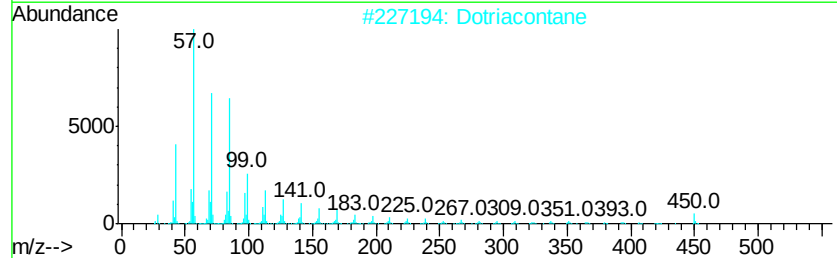
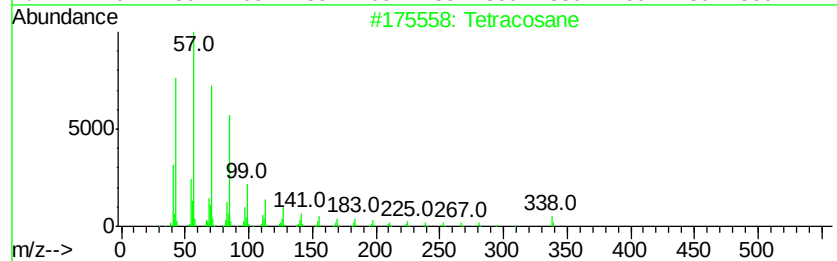
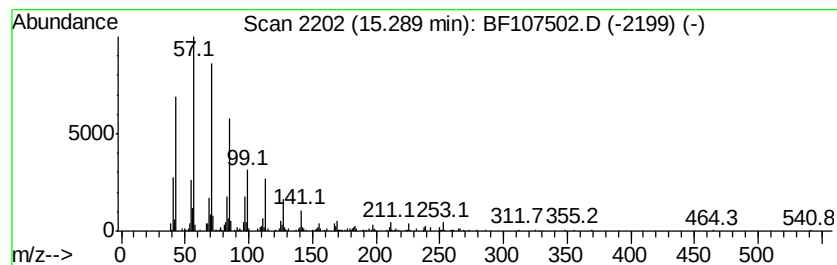
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 19 Tetracosane Concentration Rank 13

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 15.29 | 20.29 ng | 807245 | Perylene-d12     | 15.71 |

| Hit# | of | Tentative ID  | MW  | MolForm | CAS#        | Qual |
|------|----|---------------|-----|---------|-------------|------|
| 1    | 5  | Tetracosane   | 338 | C24H50  | 000646-31-1 | 93   |
| 2    |    | Dotriacontane | 451 | C32H66  | 000544-85-4 | 90   |
| 3    |    | Heneicosane   | 296 | C21H44  | 000629-94-7 | 87   |
| 4    |    | Docosane      | 310 | C22H46  | 000629-97-0 | 87   |
| 5    |    | Pentacosane   | 352 | C25H52  | 000629-99-2 | 83   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071818\  
Data File : BF107502.D  
Acq On : 19 Jul 2018 8:28  
Operator : JU/SJ  
Sample : J4033-03  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
071618-DBRI-F-D-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF071618.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

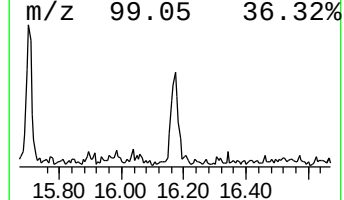
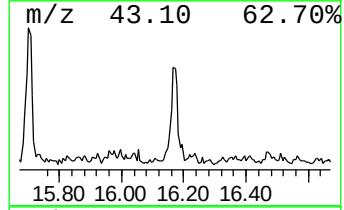
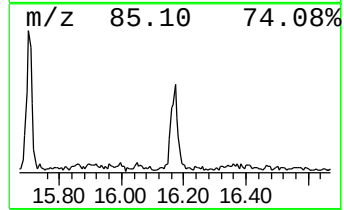
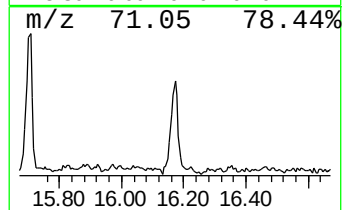
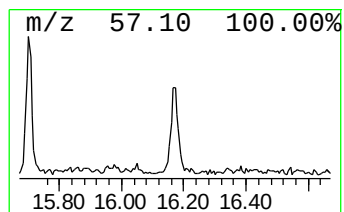
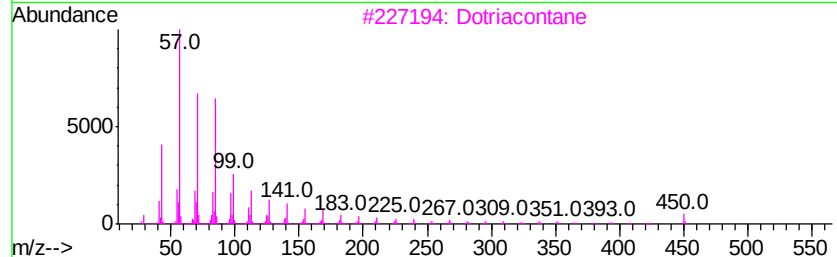
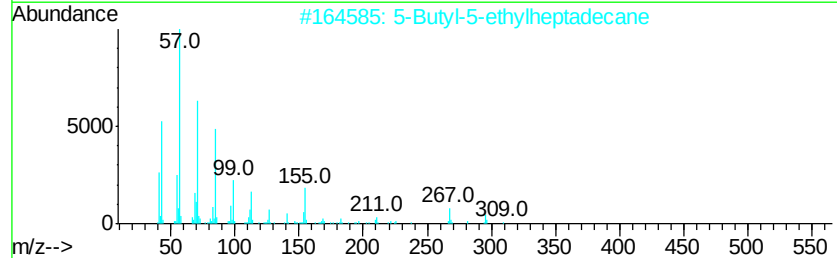
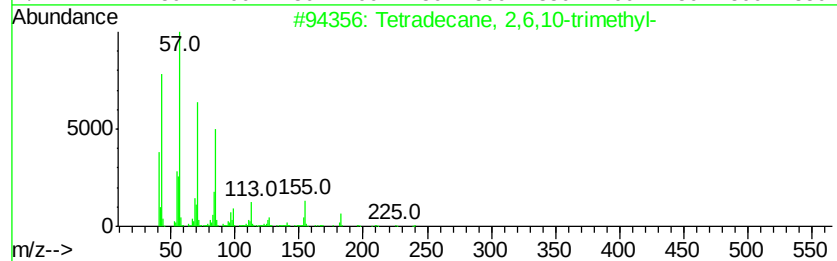
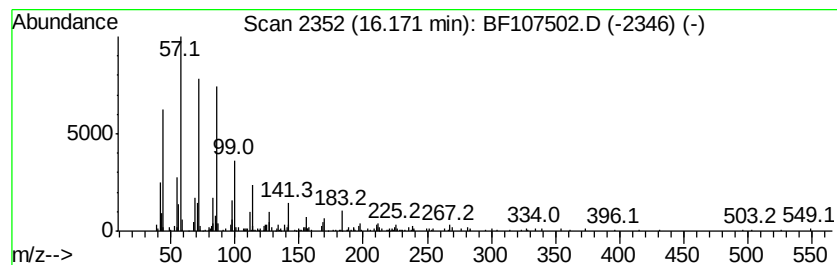
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 20 Tetradecane, 2,6,10-trimethyl- Concentration Rank 19

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 16.17 | 15.04 ng | 598511 | Perylene-d12     | 15.71 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#         | Qual |
|------|----|--------------------------------|-----|---------|--------------|------|
| 1    | 5  | Tetradecane, 2,6,10-trimethyl- | 240 | C17H36  | 014905-56-7  | 87   |
| 2    |    | 5-Butyl-5-ethylheptadecane     | 324 | C23H48  | 1000360-43-0 | 86   |
| 3    |    | Dotriacontane                  | 451 | C32H66  | 000544-85-4  | 86   |
| 4    |    | Tetracosane                    | 338 | C24H50  | 000646-31-1  | 80   |
| 5    |    | Heptadecane, 2-methyl-         | 254 | C18H38  | 001560-89-0  | 80   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF071818\  
 Data File : BF107502.D  
 Acq On : 19 Jul 2018 8:28  
 Operator : JU/SJ  
 Sample : J4033-03  
 Misc :  
 ALS Vial : 38 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 071618-DBRI-F-D-B

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF071618.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                      |       |         |       |          | #                     | RT    | Resp    | Conc |
| unknown3.82          | 3.82  | 18.3    | ng    | 806387   | 1                     | 6.98  | 881563  | 20.0 |
| 3-Penten-2-one, 4... | 4.66  | 232.3   | ng    | 10241600 | 1                     | 6.98  | 881563  | 20.0 |
| Cyclotrisiloxane,... | 4.87  | 81.0    | ng    | 3572420  | 1                     | 6.98  | 881563  | 20.0 |
| unknown6.75          | 6.75  | 54.8    | ng    | 2417260  | 1                     | 6.98  | 881563  | 20.0 |
| Cyclopentasiloxan... | 7.71  | 19.2    | ng    | 1013140  | 2                     | 8.27  | 1057720 | 20.0 |
| 2-Cyclopenten-1-o... | 9.07  | 16.6    | ng    | 878597   | 2                     | 8.27  | 1057720 | 20.0 |
| Sulphuric acid di... | 9.28  | 14.6    | ng    | 159511   | 3                     | 10.02 | 218778  | 20.0 |
| Tridecanoic acid,... | 11.90 | 16.6    | ng    | 544576   | 4                     | 11.52 | 656030  | 20.0 |
| n-Hexadecanoic acid  | 12.04 | 16.9    | ng    | 552676   | 4                     | 11.52 | 656030  | 20.0 |
| unknown12.91         | 12.91 | 21.3    | ng    | 202416   | 5                     | 14.17 | 190475  | 20.0 |
| Octacosane           | 12.96 | 21.5    | ng    | 204927   | 5                     | 14.17 | 190475  | 20.0 |
| Hexadecane, 2,6,1... | 13.32 | 44.8    | ng    | 426746   | 5                     | 14.17 | 190475  | 20.0 |
| Nonadecane, 2-met... | 13.66 | 64.2    | ng    | 611802   | 5                     | 14.17 | 190475  | 20.0 |
| Eicosane, 2-methyl-  | 13.99 | 75.8    | ng    | 721409   | 5                     | 14.17 | 190475  | 20.0 |
| Pentacosane          | 14.31 | 97.1    | ng    | 924678   | 5                     | 14.17 | 190475  | 20.0 |
| Octadecane           | 14.61 | 103.1   | ng    | 981808   | 5                     | 14.17 | 190475  | 20.0 |
| Eicosane, 7-hexyl-   | 14.94 | 23.1    | ng    | 917982   | 6                     | 15.71 | 795878  | 20.0 |
| Tetracosane          | 15.29 | 20.3    | ng    | 807245   | 6                     | 15.71 | 795878  | 20.0 |
| Tetradecane, 2,6,... | 16.17 | 15.0    | ng    | 598511   | 6                     | 15.71 | 795878  | 20.0 |