

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060916\  
 Data File : BF087854.D  
 Acq On : 9 Jun 2016 16:02  
 Operator : UM/SJ  
 Sample : H3399-16 5X  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SS-35

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:\HPCHEM1\BNA\_F\METHODS\8270-BF060716.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         | 1.747       | 13            | 14          | 24           | rVB      | 106772         | 273047        | 4.75%           | 1.195%        |
| 2         | 1.884       | 24            | 26          | 90           | rVB      | 2159105        | 5753470       | 100.00%         | 25.181%       |
| 3         | 4.970       | 294           | 296         | 300          | rBV      | 69525          | 77447         | 1.35%           | 0.339%        |
| 4         | 5.404       | 331           | 334         | 336          | rBV      | 699592         | 700235        | 12.17%          | 3.065%        |
| 5         | 6.444       | 422           | 425         | 427          | rBV      | 830701         | 743075        | 12.92%          | 3.252%        |
| 6         | 6.582       | 434           | 437         | 439          | rBV      | 876746         | 759272        | 13.20%          | 3.323%        |
| 7         | 6.822       | 455           | 458         | 460          | rVB      | 643157         | 714265        | 12.41%          | 3.126%        |
| 8         | 6.970       | 469           | 471         | 473          | rVB      | 670196         | 569393        | 9.90%           | 2.492%        |
| 9         | 7.382       | 504           | 507         | 509          | rBV      | 409879         | 428881        | 7.45%           | 1.877%        |
| 10        | 8.102       | 568           | 570         | 573          | rBV      | 1072878        | 940355        | 16.34%          | 4.116%        |
| 11        | 9.176       | 662           | 664         | 666          | rBV      | 1117997        | 861142        | 14.97%          | 3.769%        |
| 12        | 9.565       | 697           | 698         | 701          | rVB2     | 57876          | 66064         | 1.15%           | 0.289%        |
| 13        | 9.850       | 721           | 723         | 725          | rBV      | 1149562        | 985378        | 17.13%          | 4.313%        |
| 14        | 10.628      | 789           | 791         | 794          | rBV2     | 339927         | 440330        | 7.65%           | 1.927%        |
| 15        | 11.325      | 850           | 852         | 857          | rBV2     | 743751         | 1096492       | 19.06%          | 4.799%        |
| 16        | 11.405      | 857           | 859         | 860          | rVB2     | 57768          | 76407         | 1.33%           | 0.334%        |
| 17        | 11.942      | 904           | 906         | 909          | rVB      | 37713          | 58551         | 1.02%           | 0.256%        |
| 18        | 12.479      | 950           | 953         | 956          | rBV2     | 39821          | 58720         | 1.02%           | 0.257%        |
| 19        | 12.536      | 956           | 958         | 960          | rVB      | 265179         | 277605        | 4.83%           | 1.215%        |
| 20        | 12.754      | 974           | 977         | 979          | rBV2     | 173861         | 285802        | 4.97%           | 1.251%        |
| 21        | 12.902      | 989           | 990         | 993          | rVB      | 420347         | 549422        | 9.55%           | 2.405%        |
| 22        | 13.085      | 1001          | 1006        | 1008         | rBV      | 41401          | 68373         | 1.19%           | 0.299%        |
| 23        | 13.336      | 1024          | 1028        | 1034         | rVB6     | 30938          | 95683         | 1.66%           | 0.419%        |
| 24        | 13.702      | 1057          | 1060        | 1062         | rBV2     | 30406          | 60532         | 1.05%           | 0.265%        |
| 25        | 13.817      | 1067          | 1070        | 1073         | rVB3     | 50439          | 85508         | 1.49%           | 0.374%        |
| 26        | 13.954      | 1079          | 1082        | 1083         | rVV2     | 794862         | 889894        | 15.47%          | 3.895%        |
| 27        | 14.331      | 1111          | 1115        | 1117         | rBV      | 38340          | 98855         | 1.72%           | 0.433%        |
| 28        | 14.434      | 1122          | 1124        | 1127         | rBV3     | 65249          | 113190        | 1.97%           | 0.495%        |
| 29        | 14.799      | 1154          | 1156        | 1160         | rBV2     | 43285          | 70445         | 1.22%           | 0.308%        |
| 30        | 14.959      | 1167          | 1170        | 1175         | rBV      | 152848         | 305399        | 5.31%           | 1.337%        |
| 31        | 15.051      | 1175          | 1178        | 1184         | rVB2     | 244830         | 372931        | 6.48%           | 1.632%        |
| 32        | 15.245      | 1192          | 1195        | 1197         | rVB      | 84637          | 139865        | 2.43%           | 0.612%        |
| 33        | 15.302      | 1197          | 1200        | 1202         | rBV      | 94076          | 136276        | 2.37%           | 0.596%        |
| 34        | 15.360      | 1202          | 1205        | 1207         | rBV      | 544365         | 677238        | 11.77%          | 2.964%        |

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 BNA\_F  
 ClientSampleId :  
 SS-35

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\_F\METHODS\8270-BF060716.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |        |        |        |        |
|----|--------|------|------|------|------|--------|--------|--------|--------|
| 35 | 15.691 | 1232 | 1234 | 1237 | rVB3 | 37432  | 75141  | 1.31%  | 0.329% |
| 36 | 15.817 | 1242 | 1245 | 1247 | rBV  | 201393 | 304415 | 5.29%  | 1.332% |
| 37 | 15.862 | 1247 | 1249 | 1255 | rVB6 | 33237  | 71587  | 1.24%  | 0.313% |
| 38 | 16.057 | 1261 | 1266 | 1274 | rVB4 | 45779  | 150866 | 2.62%  | 0.660% |
| 39 | 16.720 | 1320 | 1324 | 1326 | rBV2 | 50937  | 120509 | 2.09%  | 0.527% |
| 40 | 16.834 | 1331 | 1334 | 1336 | rVV  | 40169  | 74575  | 1.30%  | 0.326% |
| 41 | 16.891 | 1336 | 1339 | 1345 | rVB4 | 46595  | 130591 | 2.27%  | 0.572% |
| 42 | 17.051 | 1350 | 1353 | 1356 | rVV4 | 32581  | 88895  | 1.55%  | 0.389% |
| 43 | 17.120 | 1357 | 1359 | 1365 | rVB  | 43696  | 107146 | 1.86%  | 0.469% |
| 44 | 17.280 | 1369 | 1373 | 1385 | rVV3 | 256485 | 959231 | 16.67% | 4.198% |
| 45 | 17.474 | 1387 | 1390 | 1395 | rVV  | 58449  | 153979 | 2.68%  | 0.674% |
| 46 | 17.588 | 1396 | 1400 | 1402 | rVV4 | 46478  | 135550 | 2.36%  | 0.593% |
| 47 | 17.703 | 1406 | 1410 | 1415 | rVV2 | 109164 | 375299 | 6.52%  | 1.643% |
| 48 | 17.794 | 1415 | 1418 | 1424 | rVB4 | 109431 | 301963 | 5.25%  | 1.322% |
| 49 | 18.023 | 1434 | 1438 | 1439 | rBV  | 103964 | 228445 | 3.97%  | 1.000% |
| 50 | 18.160 | 1447 | 1450 | 1453 | rBV4 | 21840  | 59066  | 1.03%  | 0.259% |
| 51 | 18.228 | 1453 | 1456 | 1462 | rVB5 | 38345  | 104609 | 1.82%  | 0.458% |
| 52 | 19.028 | 1522 | 1526 | 1533 | rVB4 | 49946  | 162357 | 2.82%  | 0.711% |
| 53 | 19.188 | 1535 | 1540 | 1550 | rVB2 | 123898 | 414633 | 7.21%  | 1.815% |

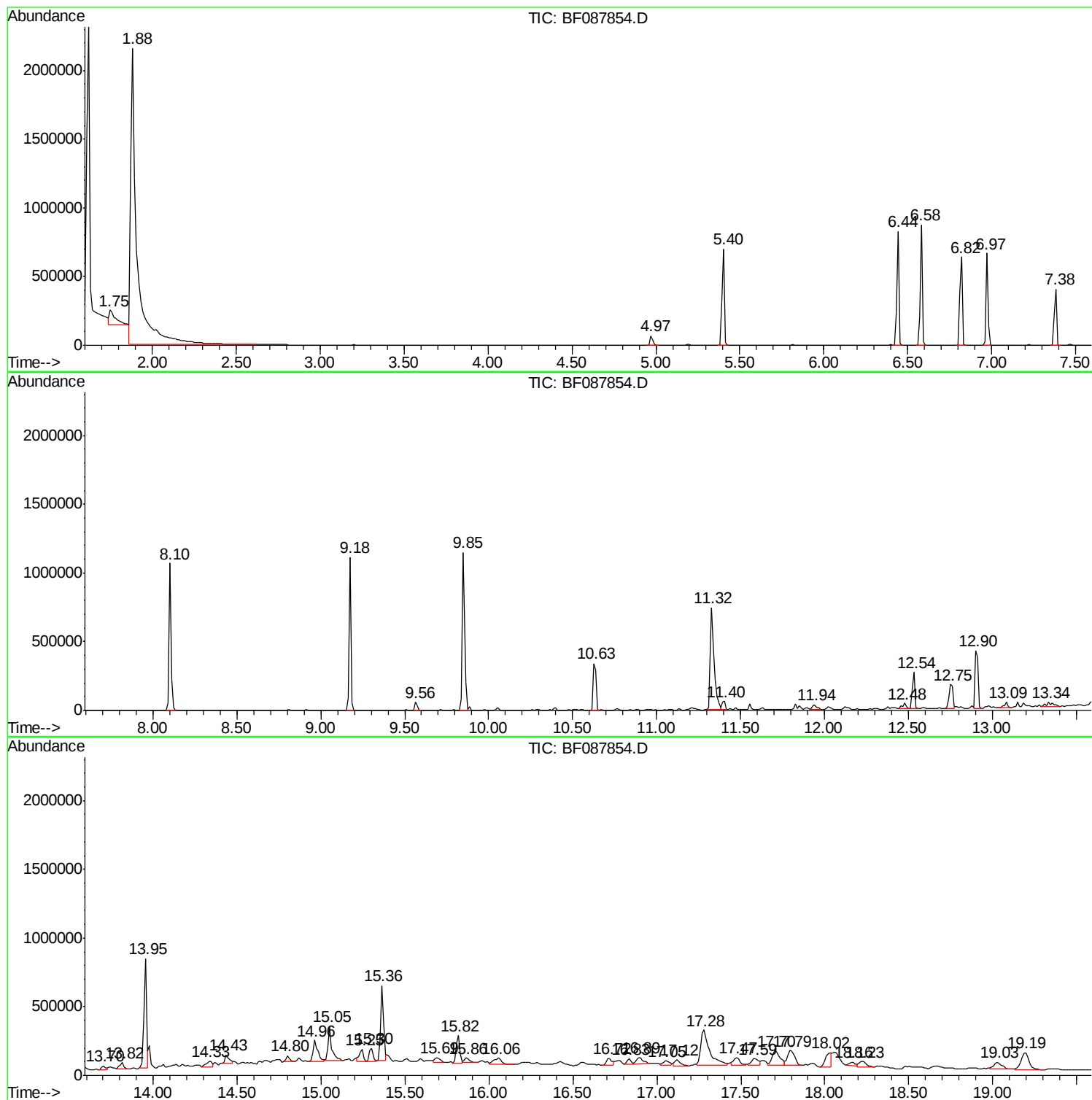
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Misc :  
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ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060916\  
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Operator : UM/SJ  
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Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
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ClientSampleId :  
SS-35

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

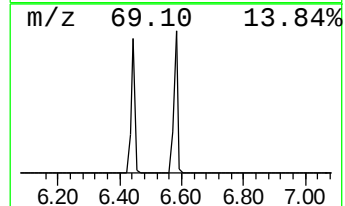
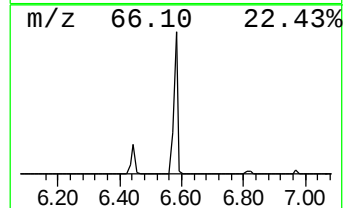
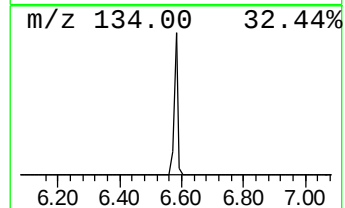
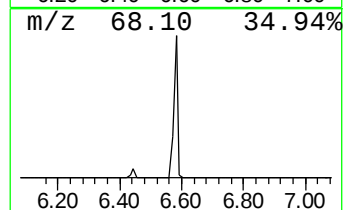
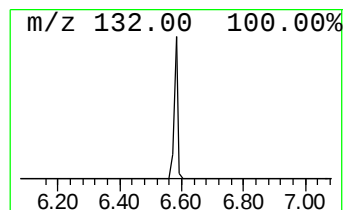
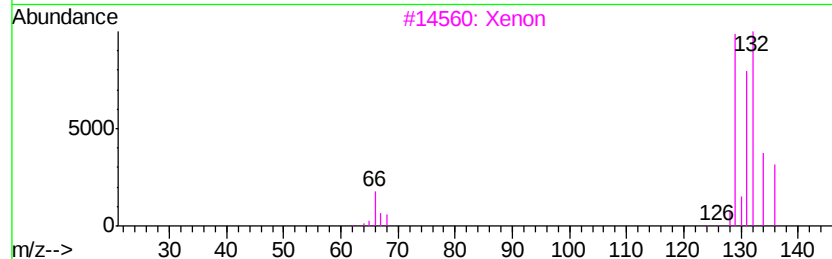
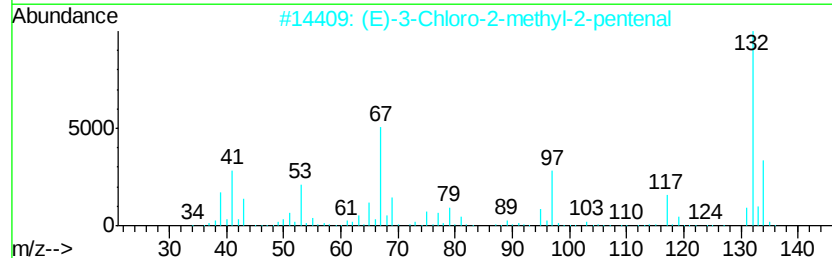
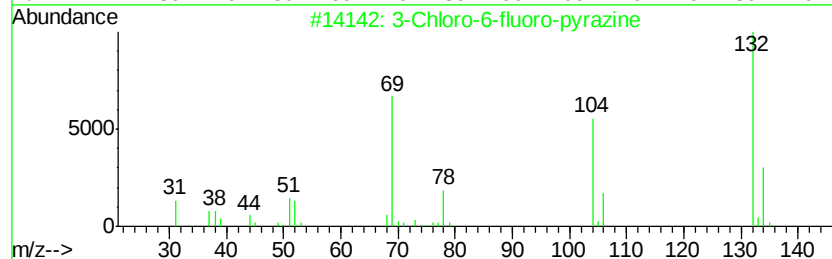
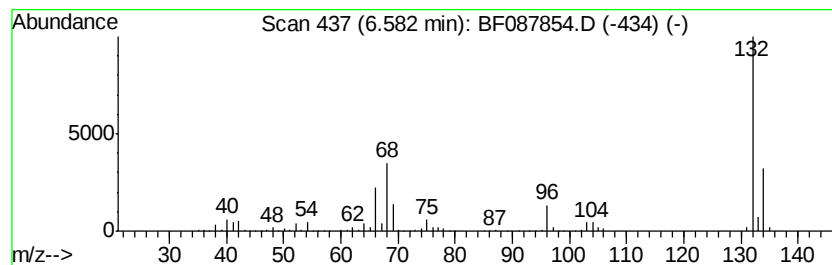
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 unknown6.58 Concentration Rank 3

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 6.58 | 21.26 ng | 759272 | 1,4-Dichlorobenzene-d4 | 6.82 |

| Hit# | of                               | 5 | Tentative ID | MW  | MolForm   | CAS#         | Qual |
|------|----------------------------------|---|--------------|-----|-----------|--------------|------|
| 1    | 3-Chloro-6-fluoro-pyrazine       |   |              | 132 | C4H2ClFN2 | 1000146-10-7 | 35   |
| 2    | (E)-3-Chloro-2-methyl-2-pentenal |   |              | 132 | C6H9ClO   | 031357-76-3  | 17   |
| 3    | Xenon                            |   |              | 132 | Xe        | 007440-63-3  | 9    |
| 4    | 4-Chloro-2-fluoro-pyrimidine     |   |              | 132 | C4H2ClFN2 | 051422-00-5  | 9    |
| 5    | 4-Methylpyrrolo[1,2-a]pyrazine   |   |              | 132 | C8H8N2    | 064608-60-2  | 9    |



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Instrument :  
BNA\_F  
ClientSampleId :  
SS-35

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

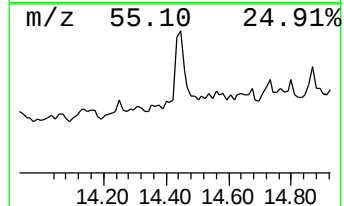
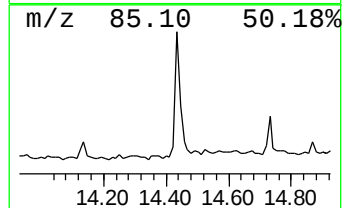
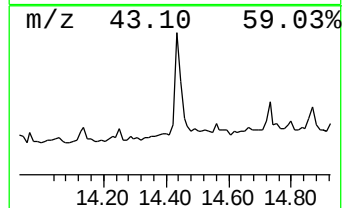
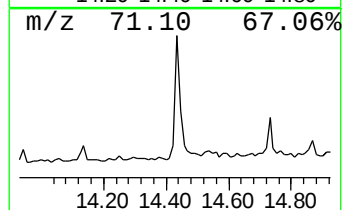
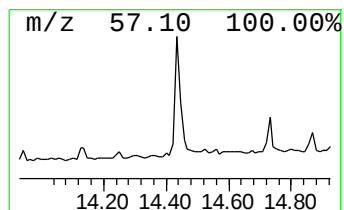
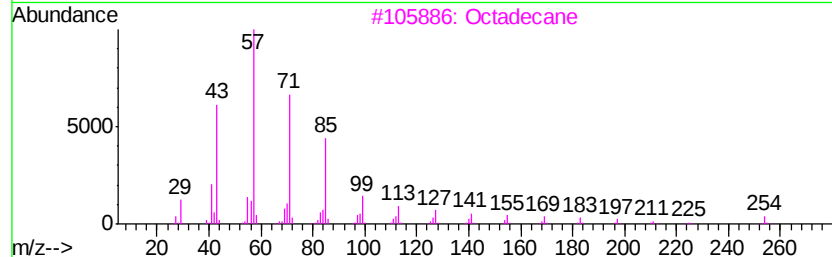
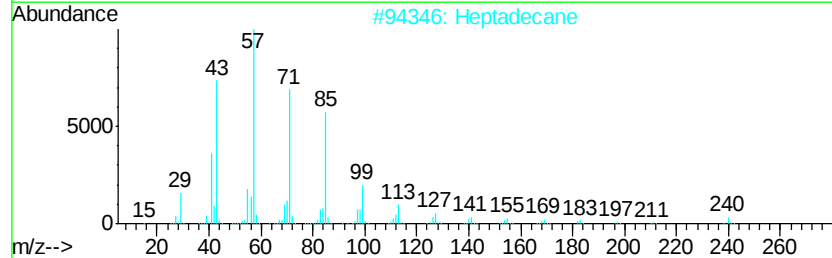
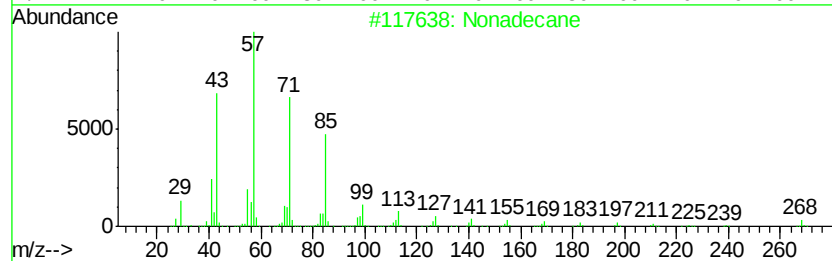
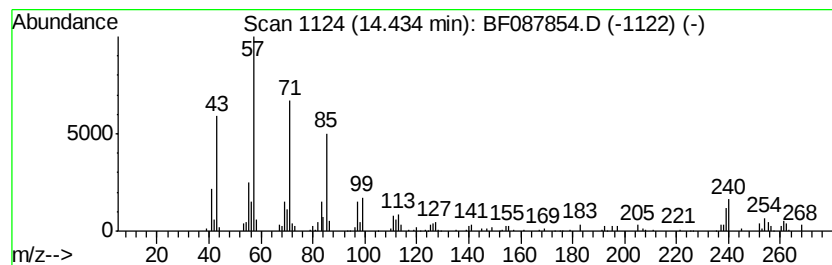
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 Nonadecane Concentration Rank 20

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 14.43 | 2.54 ng | 113190 | Chrysene-d12     | 13.95 |

| Hit# of | 5           | Tentative ID | MW  | MolForm | CAS#        | Qual |
|---------|-------------|--------------|-----|---------|-------------|------|
| 1       | Nonadecane  |              | 268 | C19H40  | 000629-92-5 | 92   |
| 2       | Heptadecane |              | 240 | C17H36  | 000629-78-7 | 86   |
| 3       | Octadecane  |              | 254 | C18H38  | 000593-45-3 | 83   |
| 4       | Hexacosane  |              | 366 | C26H54  | 000630-01-3 | 60   |
| 5       | Pentadecane |              | 212 | C15H32  | 000629-62-9 | 53   |



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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

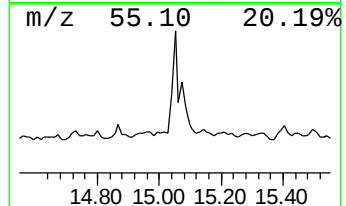
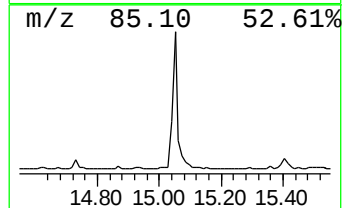
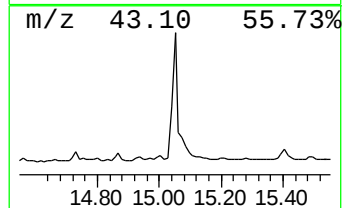
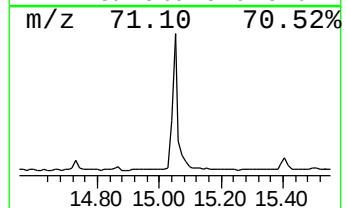
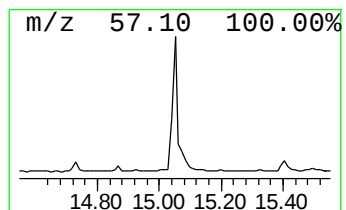
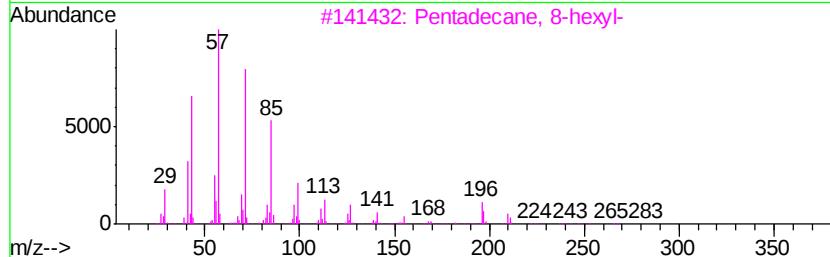
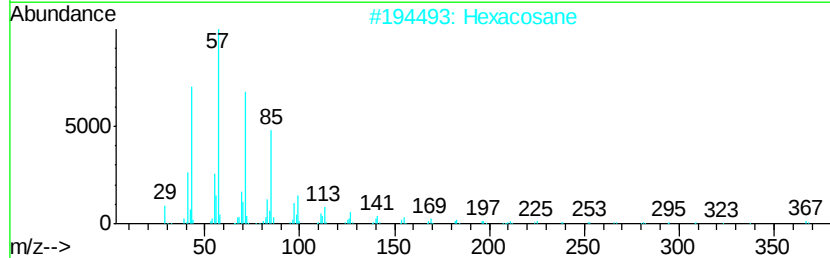
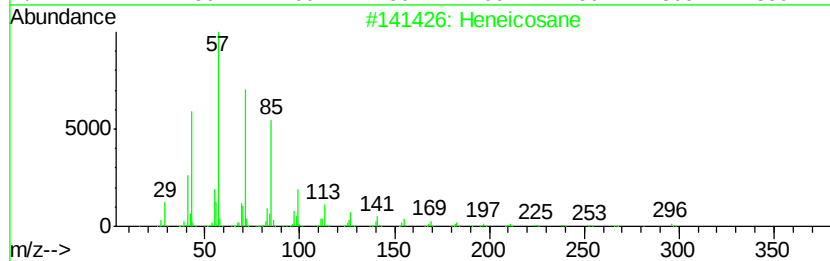
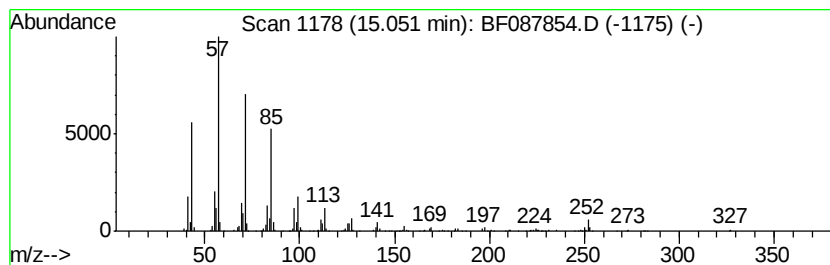
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 Heneicosane Concentration Rank 7

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 15.05 | 11.01 ng | 372931 | Perylene-d12     | 15.36 |

| Hit# of | 5                     | Tentative ID | MW  | MolForm | CAS#        | Qual |
|---------|-----------------------|--------------|-----|---------|-------------|------|
| 1       | Heneicosane           |              | 296 | C21H44  | 000629-94-7 | 90   |
| 2       | Hexacosane            |              | 366 | C26H54  | 000630-01-3 | 90   |
| 3       | Pentadecane, 8-hexyl- |              | 296 | C21H44  | 013475-75-7 | 87   |
| 4       | Pentacosane           |              | 352 | C25H52  | 000629-99-2 | 87   |
| 5       | Hexadecane            |              | 226 | C16H34  | 000544-76-3 | 83   |



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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

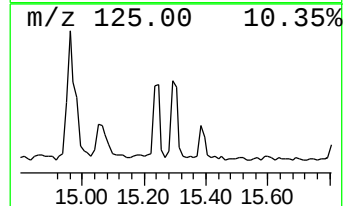
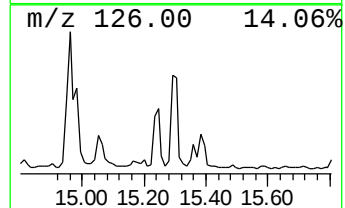
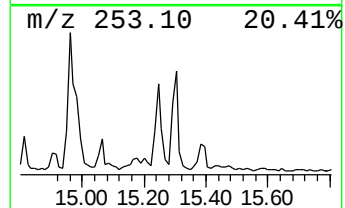
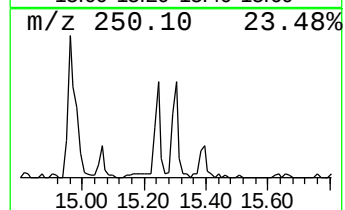
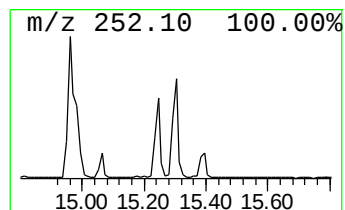
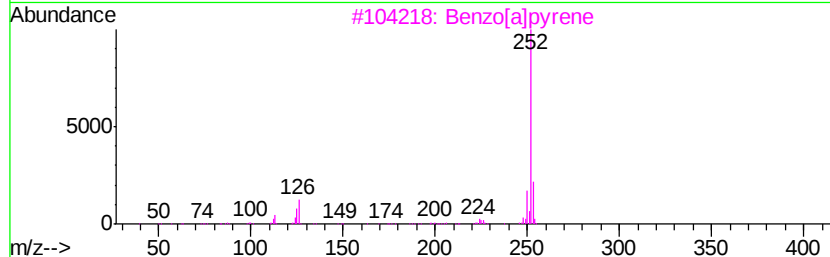
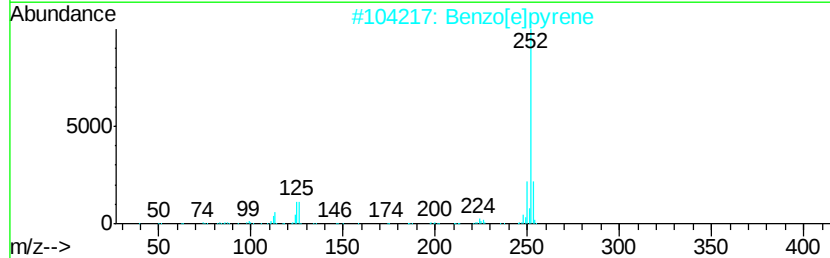
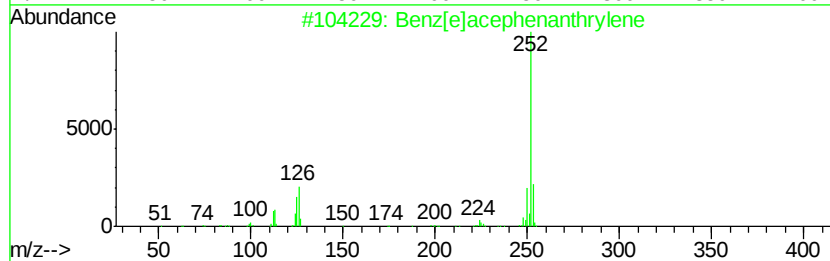
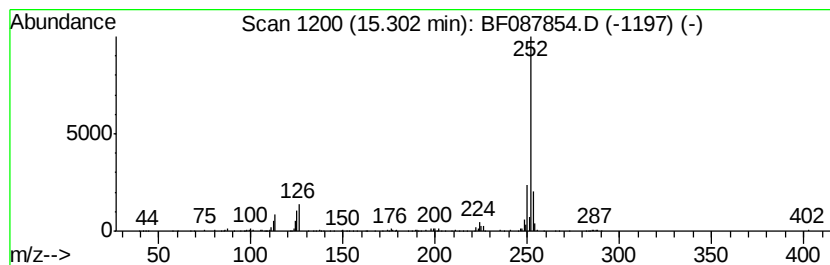
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 Benz[e]acephenanthrylene Concentration Rank 15

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 15.30 | 4.02 ng | 136276 | Perylene-d12     | 15.36 |

| Hit# of | 5 | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|---------|---|--------------------------|-----|---------|-------------|------|
| 1       |   | Benz[e]acephenanthrylene | 252 | C20H12  | 000205-99-2 | 96   |
| 2       |   | Benzo[e]pyrene           | 252 | C20H12  | 000192-97-2 | 95   |
| 3       |   | Benzo[a]pyrene           | 252 | C20H12  | 000050-32-8 | 94   |
| 4       |   | Benzo[k]fluoranthene     | 252 | C20H12  | 000207-08-9 | 93   |
| 5       |   | Perylene                 | 252 | C20H12  | 000198-55-0 | 93   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060916\  
Data File : BF087854.D  
Acq On : 9 Jun 2016 16:02  
Operator : UM/SJ  
Sample : H3399-16 5X  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SS-35

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

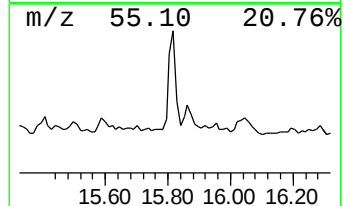
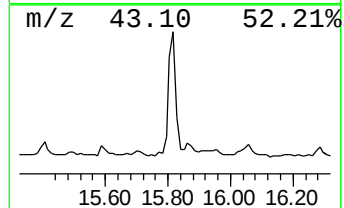
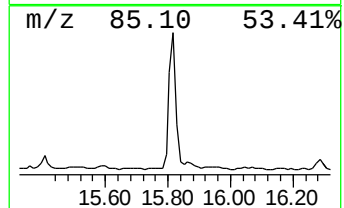
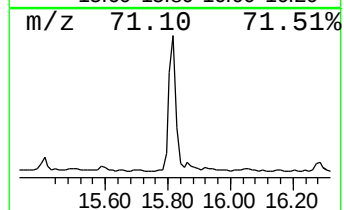
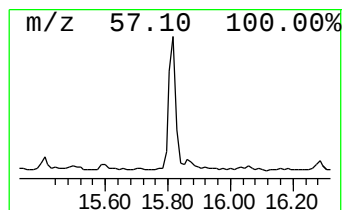
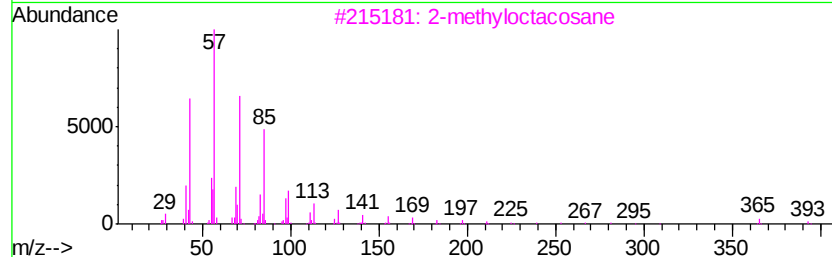
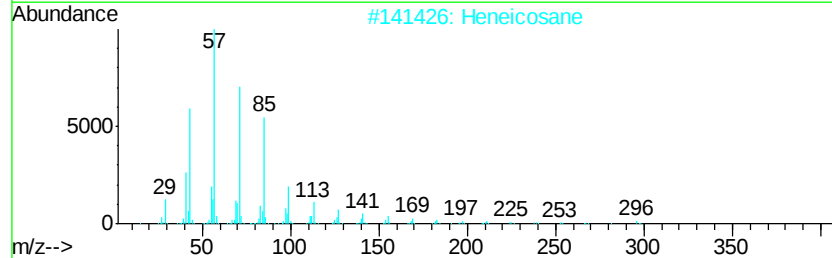
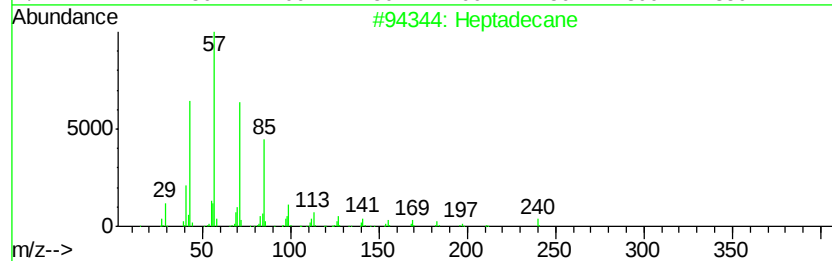
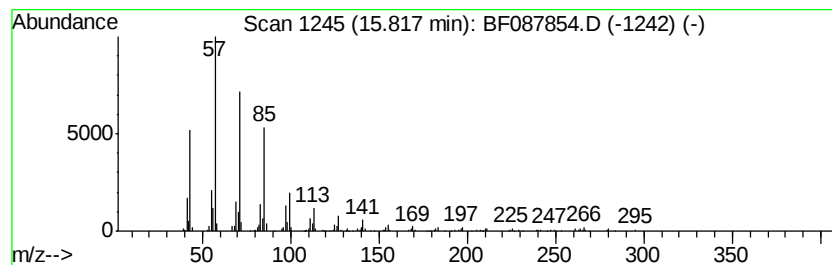
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TIC Integration Parameters: LSCINT.P

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Peak Number 8 Heptadecane Concentration Rank 8

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 15.82 | 8.99 ng | 304415 | Perylene-d12     | 15.36 |

| Hit# of | 5                  | Tentative ID | MW  | MolForm | CAS#         | Qual |
|---------|--------------------|--------------|-----|---------|--------------|------|
| 1       | Heptadecane        |              | 240 | C17H36  | 000629-78-7  | 94   |
| 2       | Heneicosane        |              | 296 | C21H44  | 000629-94-7  | 90   |
| 3       | 2-methyloctacosane |              | 408 | C29H60  | 1000376-72-8 | 90   |
| 4       | Octacosane         |              | 394 | C28H58  | 000630-02-4  | 87   |
| 5       | Hexacosane         |              | 366 | C26H54  | 000630-01-3  | 87   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060916\  
Data File : BF087854.D  
Acq On : 9 Jun 2016 16:02  
Operator : UM/SJ  
Sample : H3399-16 5X  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060716.M  
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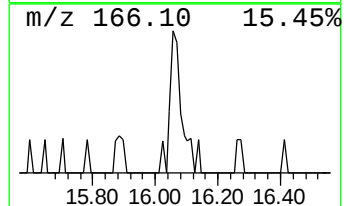
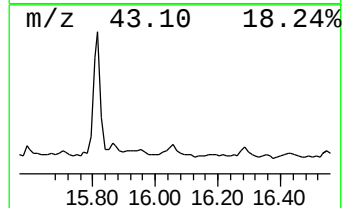
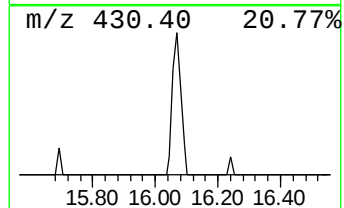
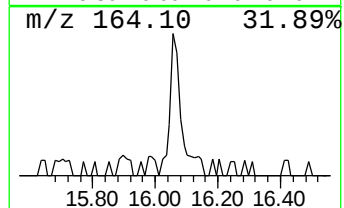
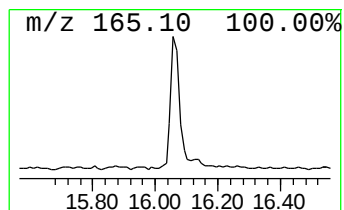
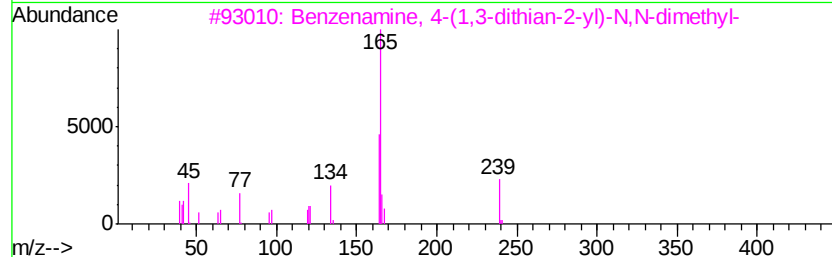
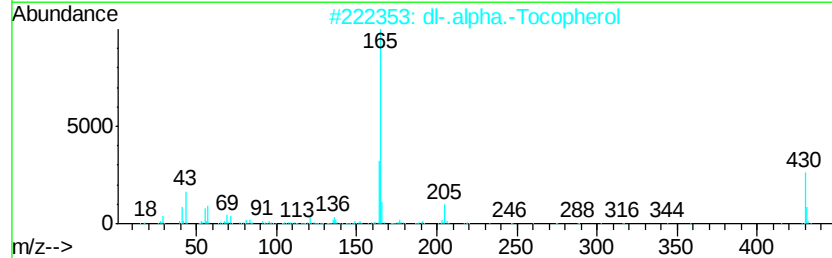
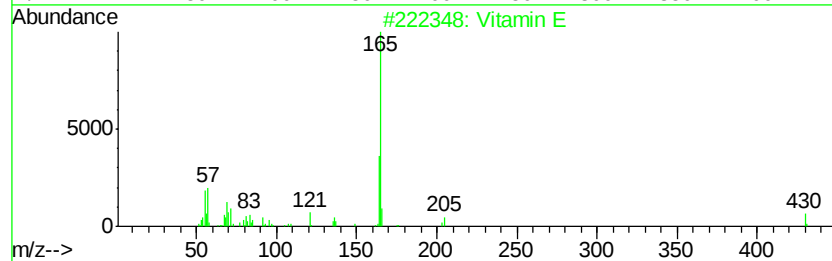
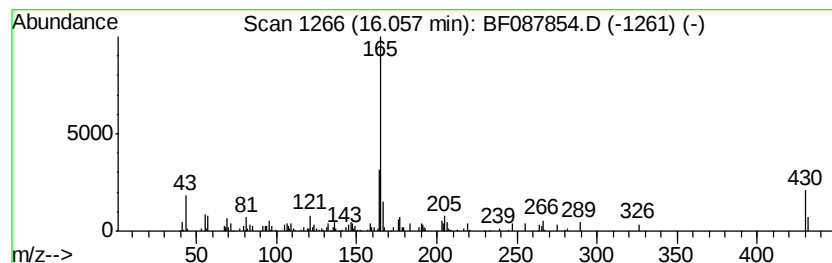
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TIC Integration Parameters: LSCINT.P

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Peak Number 9 Vitamin E Concentration Rank 14

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 16.06 | 4.46 ng | 150866 | Perylene-d12     | 15.36 |

| Hit# | of | Tentative ID                         | MW  | MolForm   | CAS#         | Qual |
|------|----|--------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Vitamin E                            | 430 | C29H50O2  | 000059-02-9  | 93   |
| 2    |    | dl-.alpha.-Tocopherol                | 430 | C29H50O2  | 010191-41-0  | 55   |
| 3    |    | Benzenamine, 4-(1,3-dithian-2-yl...) | 239 | C12H17NS2 | 024588-75-8  | 40   |
| 4    |    | 2,6-Pyridinedicarboxylic acid, p...  | 321 | C18H27N04 | 1000368-78-8 | 40   |
| 5    |    | 2-Amino-6-hydroxy-9-methylpurine     | 165 | C6H7N5O   | 005502-78-3  | 38   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060916\  
Data File : BF087854.D  
Acq On : 9 Jun 2016 16:02  
Operator : UM/SJ  
Sample : H3399-16 5X  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
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ClientSampleId :  
SS-35

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

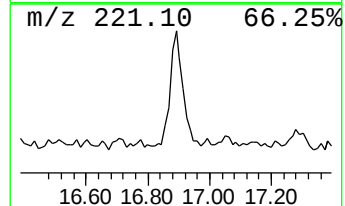
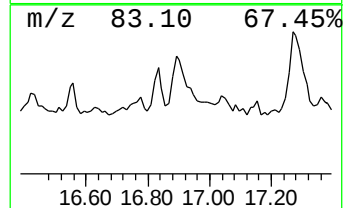
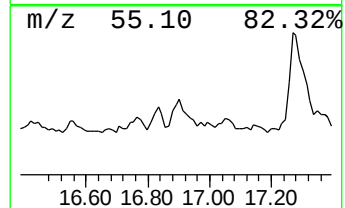
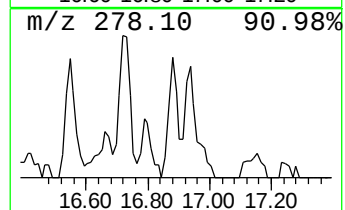
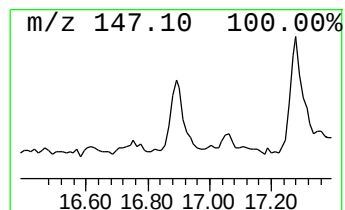
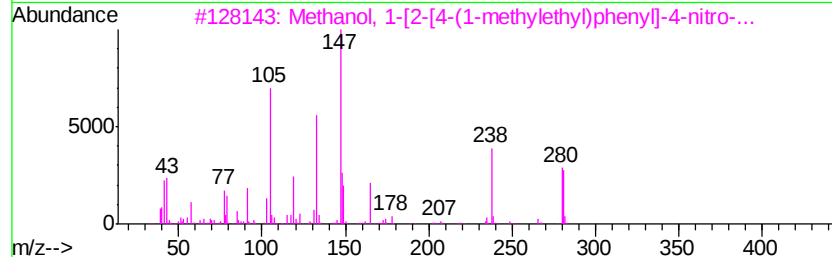
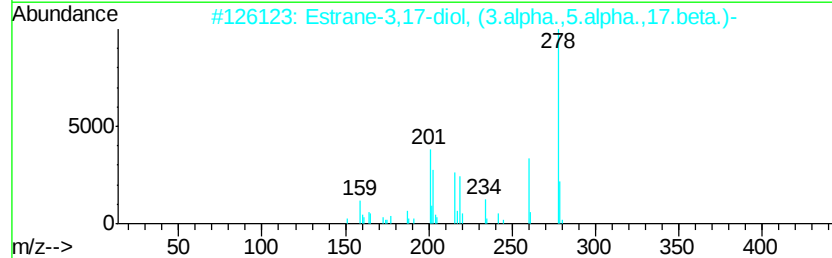
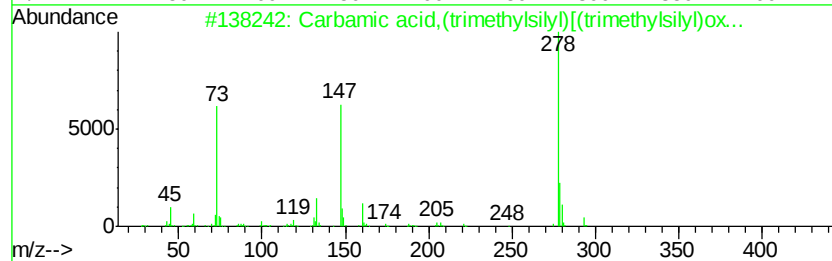
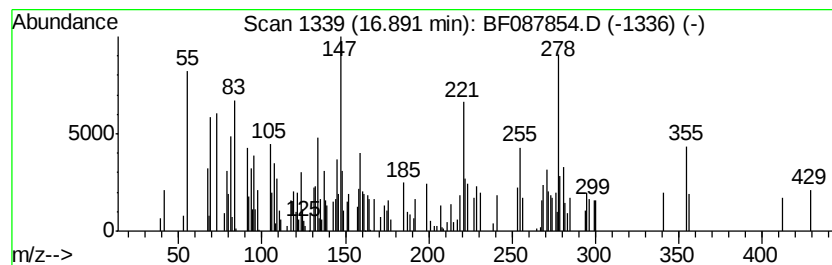
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TIC Integration Parameters: LSCINT.P

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Peak Number 10 unknown16.89 Concentration Rank 17

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 16.89 | 3.86 ng | 130591 | Perylene-d12     | 15.36 |

| Hit# | of | Tentative ID                         | MW  | MolForm      | CAS#         | Qual |
|------|----|--------------------------------------|-----|--------------|--------------|------|
| 1    | 5  | Carbamic acid,(trimethylsilyl)[(...  | 293 | C10H27N03Si3 | 070726-27-1  | 43   |
| 2    |    | Estrane-3,17-diol, (3.alpha.,5.a...  | 278 | C18H30O2     | 010002-95-6  | 22   |
| 3    |    | Methanol, 1-[2-[4-(1-methylethyl)... | 281 | C14H19NO5    | 1000271-97-1 | 20   |
| 4    |    | Picene                               | 278 | C22H14       | 000213-46-7  | 18   |
| 5    |    | Dibenz[a,h]anthracene                | 278 | C22H14       | 000053-70-3  | 18   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060916\  
Data File : BF087854.D  
Acq On : 9 Jun 2016 16:02  
Operator : UM/SJ  
Sample : H3399-16 5X  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SS-35

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

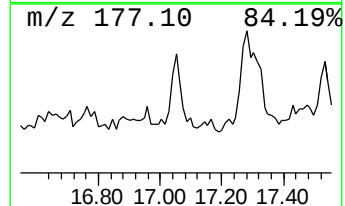
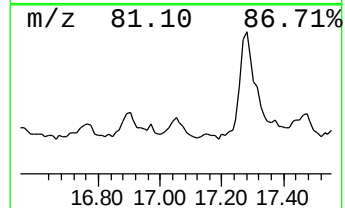
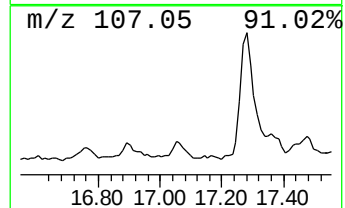
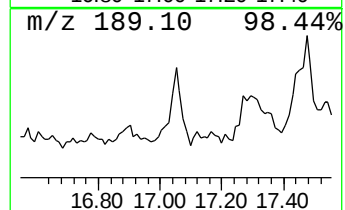
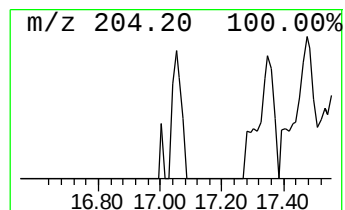
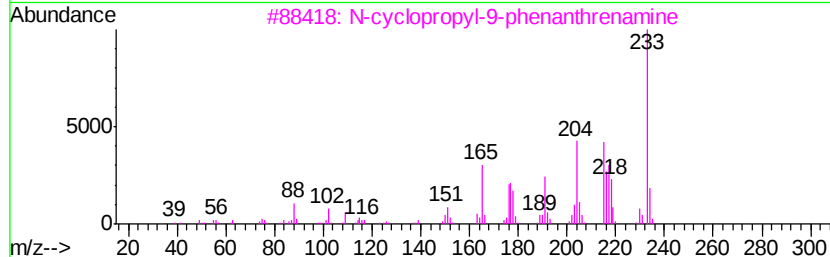
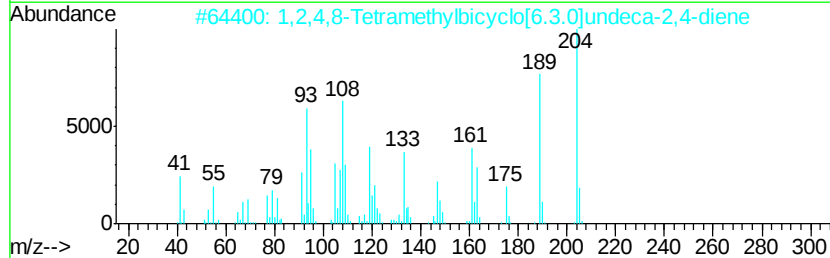
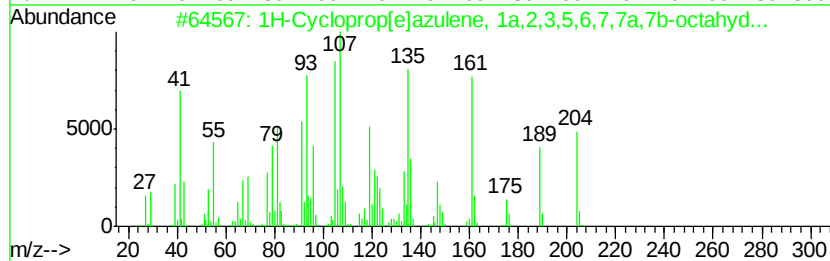
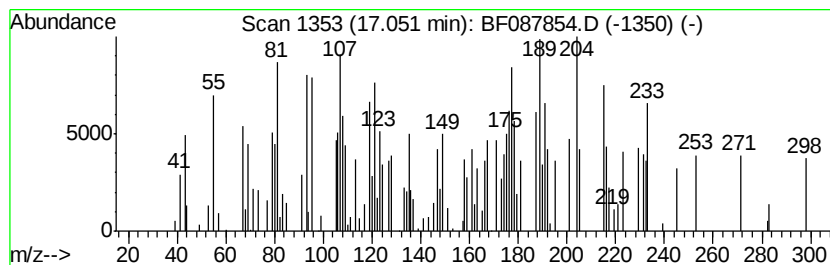
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 unknown17.05 Concentration Rank 19

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 17.05 | 2.63 ng | 88895 | Perylene-d12     | 15.36 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------------------|-----|---------|--------------|------|
| 1    | 5  | 1H-Cycloprop[elazulene, 1a,2,3,5... | 204 | C15H24  | 021747-46-6  | 35   |
| 2    |    | 1,2,4,8-Tetramethylbicyclo[6.3.0... | 204 | C15H24  | 137235-51-9  | 35   |
| 3    |    | N-cyclopropyl-9-phenanthrenamine    | 233 | C17H15N | 348579-19-1  | 30   |
| 4    |    | 1,1,4a-Trimethyl-5,6-dimethylene... | 204 | C15H24  | 1000193-60-8 | 30   |
| 5    |    | Thujopsene-(I2)                     | 204 | C15H24  | 1000152-20-4 | 25   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060916\  
Data File : BF087854.D  
Acq On : 9 Jun 2016 16:02  
Operator : UM/SJ  
Sample : H3399-16 5X  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

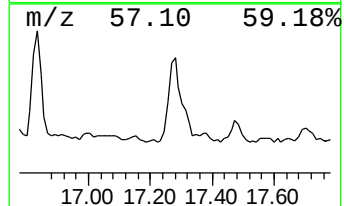
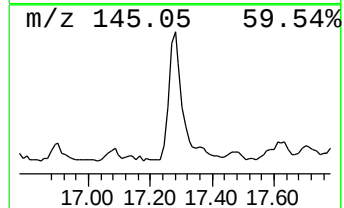
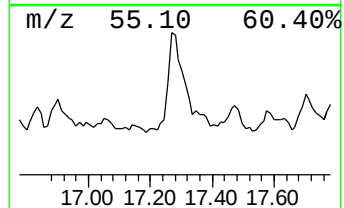
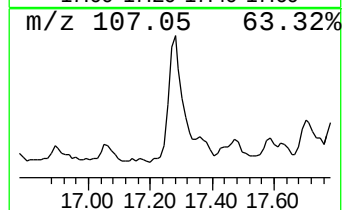
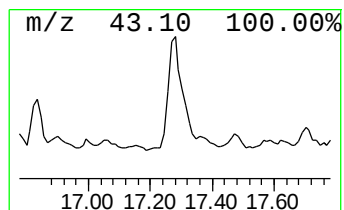
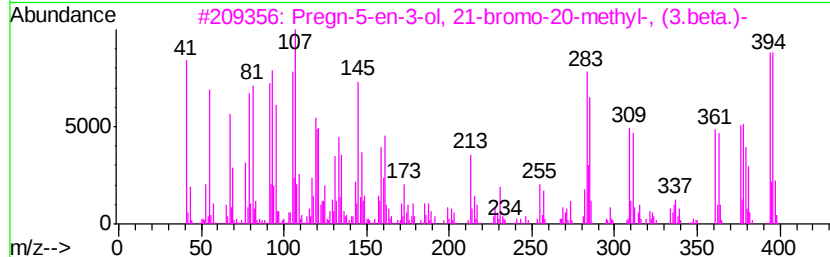
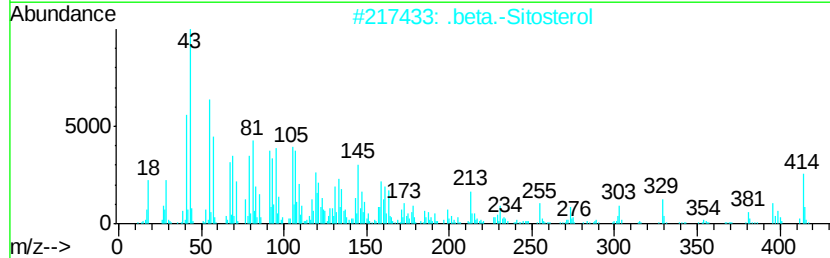
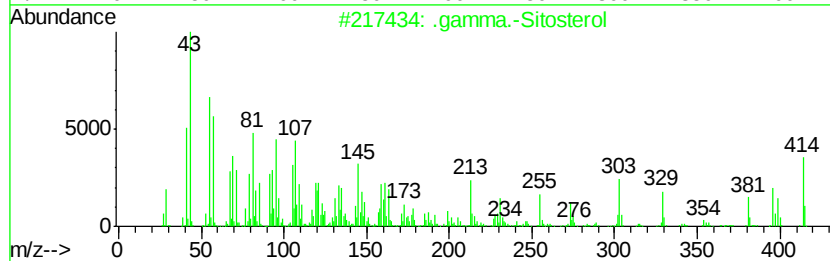
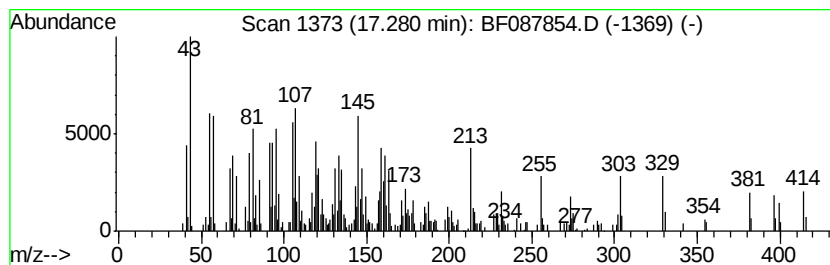
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TIC Integration Parameters: LSCINT.P

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Peak Number 12 .gamma.-Sitosterol Concentration Rank 2

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 17.28 | 28.33 ng | 959231 | Perylene-d12     | 15.36 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | .gamma.-Sitosterol                  | 414 | C29H50O   | 000083-47-6  | 99   |
| 2    |    | .beta.-Sitosterol                   | 414 | C29H50O   | 000083-46-5  | 98   |
| 3    |    | Pregn-5-en-3-ol, 21-bromo-20-met... | 394 | C22H35BrO | 055103-80-5  | 95   |
| 4    |    | 5-Cholestene-3-ol, 24-methyl-       | 400 | C28H48O   | 1000214-17-4 | 53   |
| 5    |    | Cholest-5-ene, 3-ethoxy-, (3.bet... | 414 | C29H50O   | 000986-19-6  | 42   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060916\  
Data File : BF087854.D  
Acq On : 9 Jun 2016 16:02  
Operator : UM/SJ  
Sample : H3399-16 5X  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

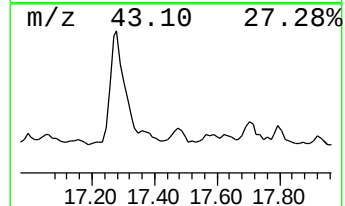
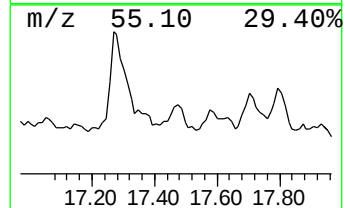
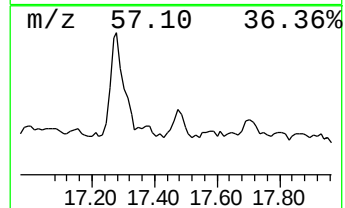
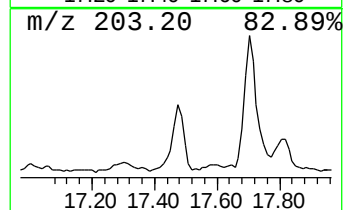
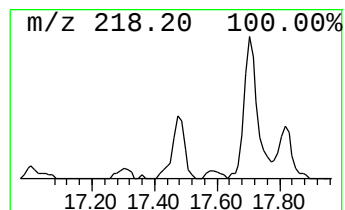
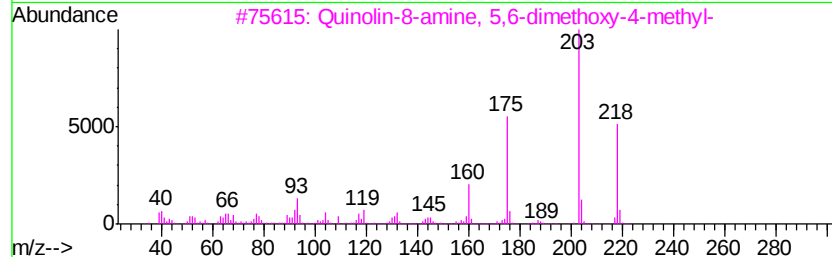
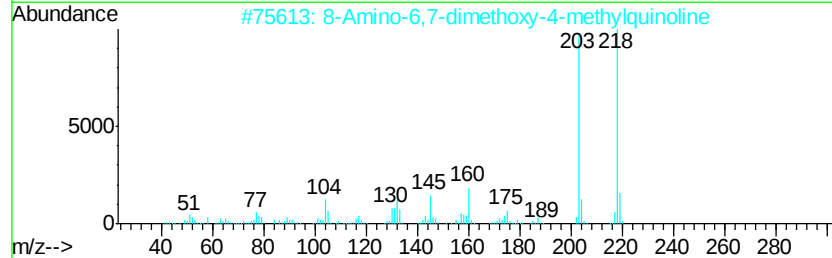
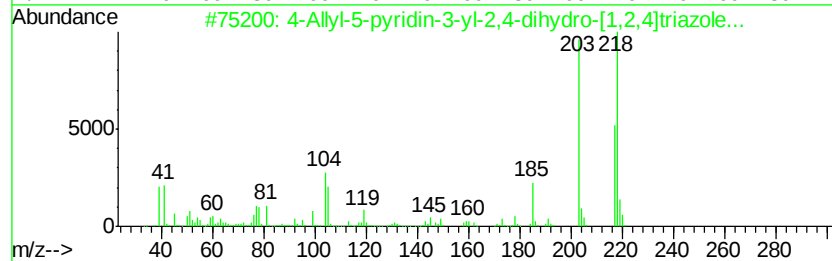
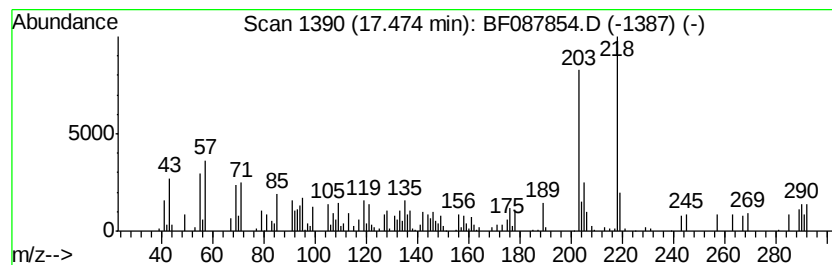
Instrument :  
BNA\_F  
ClientSampleId :  
SS-35

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 13 4-Allyl-5-pyridin-3-yl-2,4-... Concentration Rank 13

| R.T.    | EstConc                             | Area                              | Relative to ISTD | R.T.         |              |      |
|---------|-------------------------------------|-----------------------------------|------------------|--------------|--------------|------|
| 17.47   | 4.55 ng                             | 153979                            | Perylene-d12     | 15.36        |              |      |
| Hit# of | 5                                   | Tentative ID                      | MW               | MolForm      | CAS#         | Qual |
| 1       | 4                                   | Allvl-5-pyridin-3-vl-2,4-dihyd... | 218              | C10H10N4S    | 1000300-40-5 | 52   |
| 2       | 8                                   | Amino-6,7-dimethoxy-4-methylqu... | 218              | C12H14N2O2   | 1000213-07-2 | 52   |
| 3       | Quinolin-8-amine, 5,6-dimethoxy-... | 218                               | C12H14N2O2       | 064992-95-6  | 50           |      |
| 4       | 3-Ethylphenol, O-trifluoroacetyl-   | 218                               | C10H9F3O2        | 1000374-27-3 | 49           |      |
| 5       | l-Valine, n-pentafluoropropionyl... | 291                               | C10H14F5N03      | 1000320-87-8 | 47           |      |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060916\  
Data File : BF087854.D  
Acq On : 9 Jun 2016 16:02  
Operator : UM/SJ  
Sample : H3399-16 5X  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SS-35

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

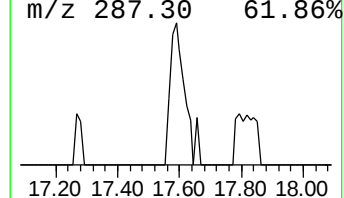
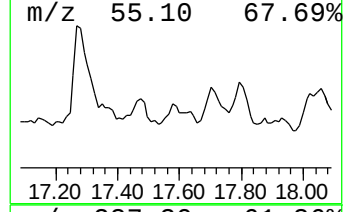
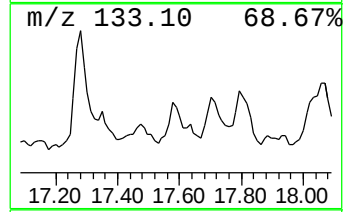
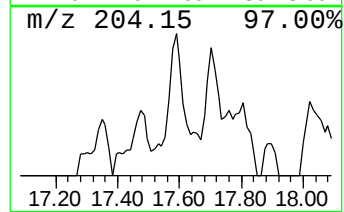
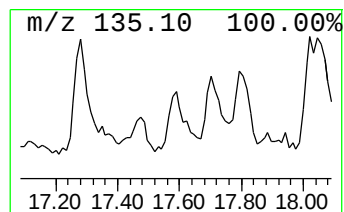
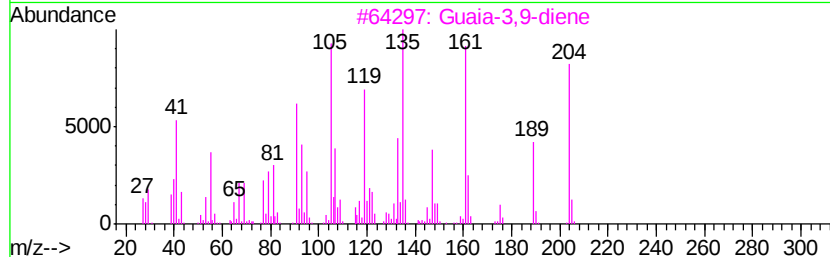
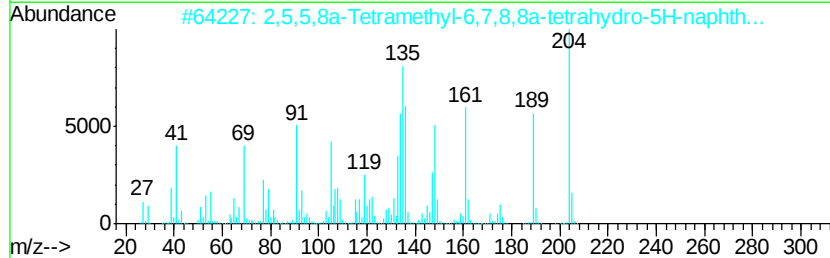
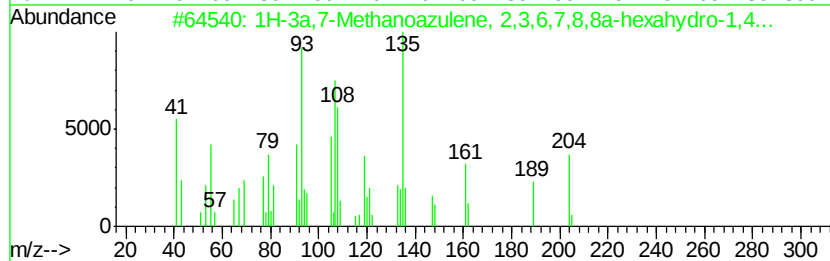
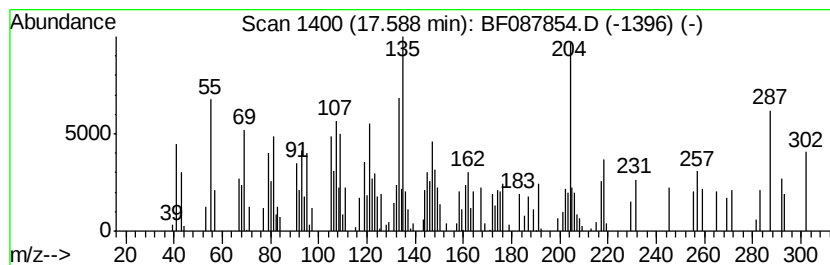
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TIC Integration Parameters: LSCINT.P

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Peak Number 14 unknown17.59 Concentration Rank 16

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 17.59 | 4.00 ng | 135550 | Perylene-d12     | 15.36 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 1H-3a,7-Methanoazulene, 2,3,6,7,... | 204 | C15H24    | 000560-32-7  | 38   |
| 2    |    | 2,5,5,8a-Tetramethyl-6,7,8,8a-te... | 204 | C14H200   | 124957-09-1  | 35   |
| 3    |    | Guaia-3,9-diene                     | 204 | C15H24    | 000489-83-8  | 25   |
| 4    |    | 2H-Cyclopentacyclooctene, 4,5,6,... | 204 | C15H24    | 1000221-85-8 | 25   |
| 5    |    | Quinoline, 8-methoxy-5-nitro-       | 204 | C10H8N2O3 | 1000298-88-7 | 25   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060916\  
Data File : BF087854.D  
Acq On : 9 Jun 2016 16:02  
Operator : UM/SJ  
Sample : H3399-16 5X  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SS-35

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

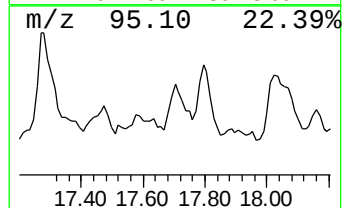
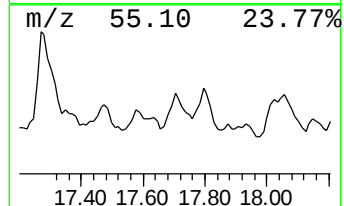
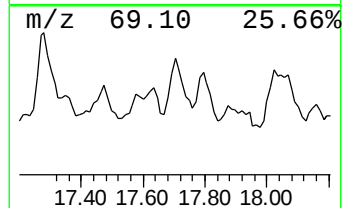
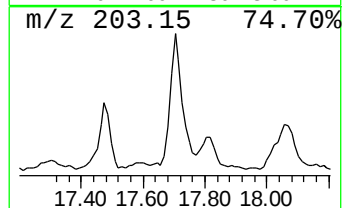
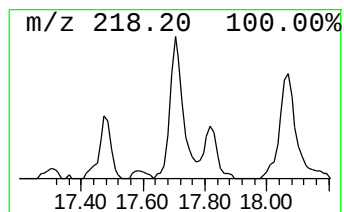
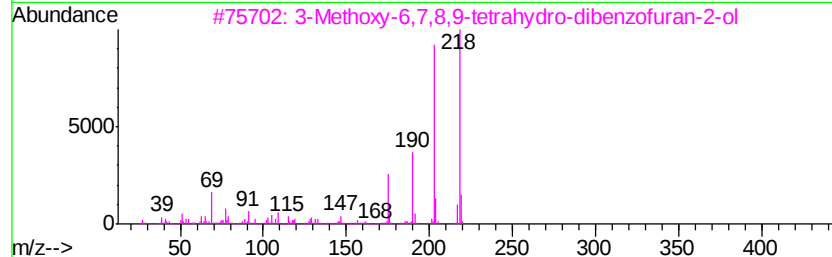
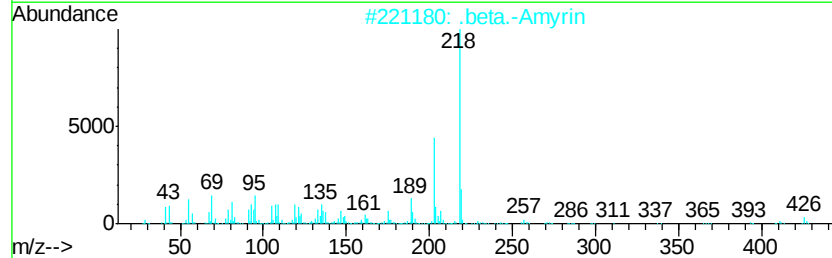
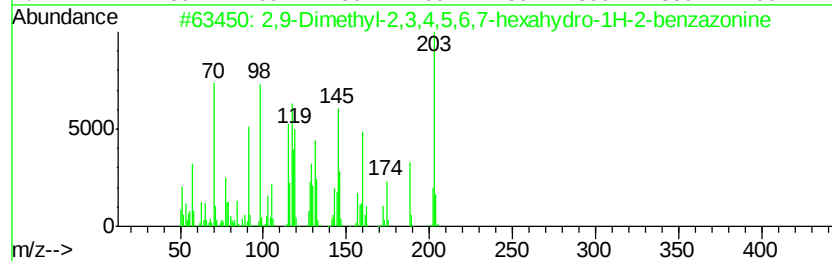
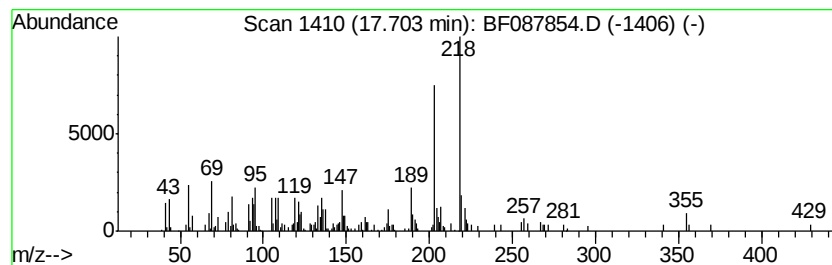
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 15 2,9-Dimethyl-2,3,4,5,6,7-he... Concentration Rank 6

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 17.70 | 11.08 ng | 375299 | Perylene-d12     | 15.36 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | 2,9-Dimethyl-2,3,4,5,6,7-hexahyd... | 203 | C14H21N  | 077581-11-4  | 68   |
| 2    |    | .beta.-Amyrin                       | 426 | C30H50O  | 000559-70-6  | 53   |
| 3    |    | 3-Methoxy-6,7,8,9-tetrahydro-dib... | 218 | C13H14O3 | 1000186-57-9 | 52   |
| 4    |    | 2H-1-Benzopyran-2-one, 6-acetyl-... | 260 | C14H12O5 | 129812-31-3  | 50   |
| 5    |    | 4,4,6a,6b,8a,11,11,14b-Octamethy... | 424 | C30H48O  | 1000194-62-4 | 49   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060916\  
Data File : BF087854.D  
Acq On : 9 Jun 2016 16:02  
Operator : UM/SJ  
Sample : H3399-16 5X  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampled :  
SS-35

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

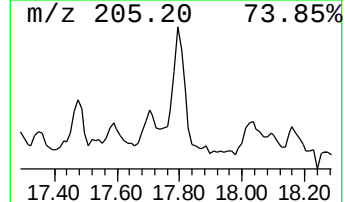
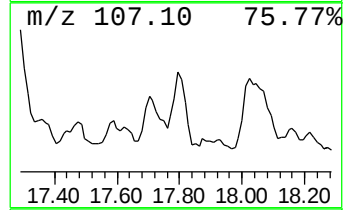
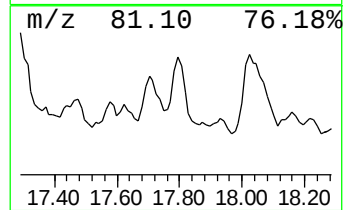
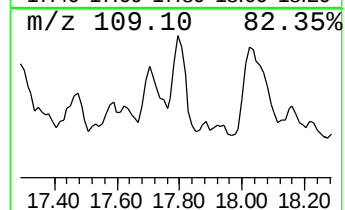
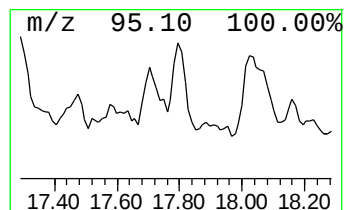
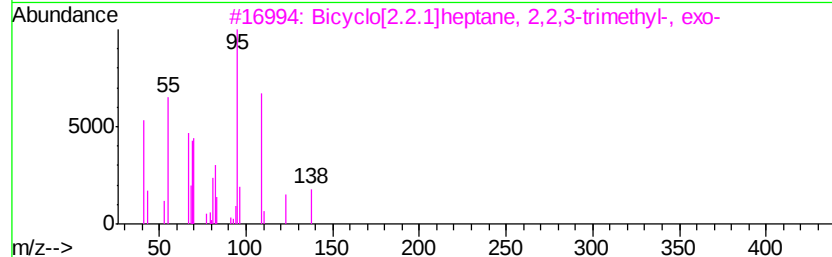
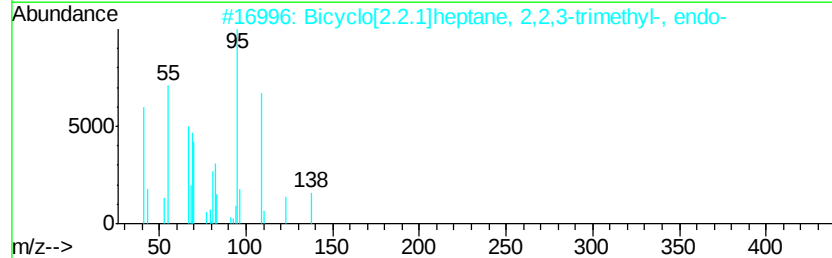
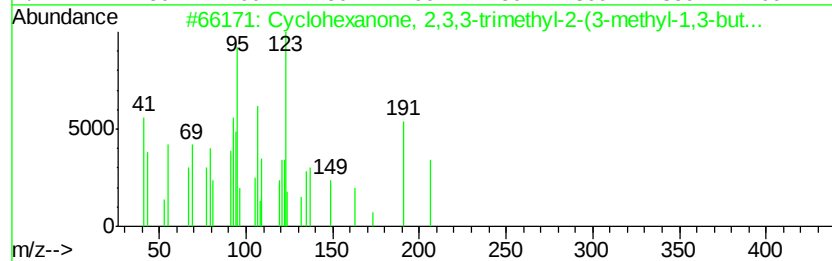
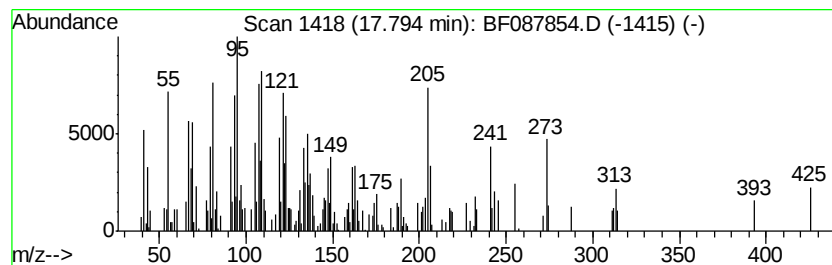
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 unknown17.79 Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 17.79 | 8.92 ng | 301963 | Perylene-d12     | 15.36 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------------------|-----|---------|--------------|------|
| 1    | 5  | Cyclohexanone, 2,3,3-trimethyl-2... | 206 | C14H22O | 069296-90-8  | 35   |
| 2    |    | Bicyclo[2.2.1]heptane, 2,2,3-tri... | 138 | C10H18  | 020536-40-7  | 35   |
| 3    |    | Bicyclo[2.2.1]heptane, 2,2,3-tri... | 138 | C10H18  | 020536-41-8  | 35   |
| 4    |    | But-3-enal, 2-methyl-4-(2,6,6-tr... | 206 | C14H22O | 104900-59-6  | 27   |
| 5    |    | 1-Oxaspiro[2.5]octane, 5,5-dimet... | 206 | C14H22O | 1000195-92-1 | 25   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060916\  
Data File : BF087854.D  
Acq On : 9 Jun 2016 16:02  
Operator : UM/SJ  
Sample : H3399-16 5X  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

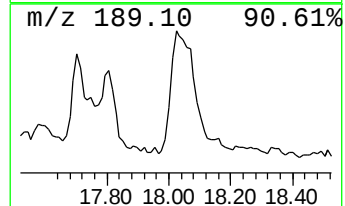
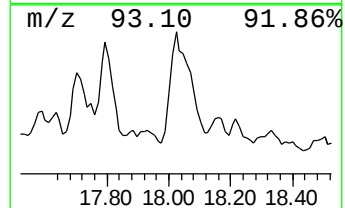
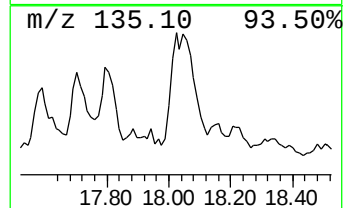
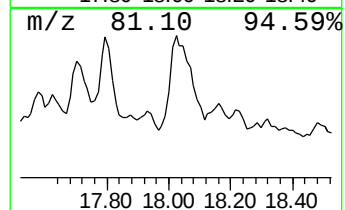
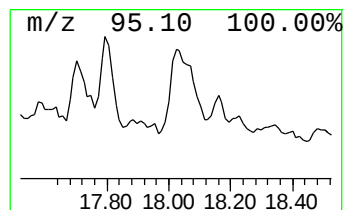
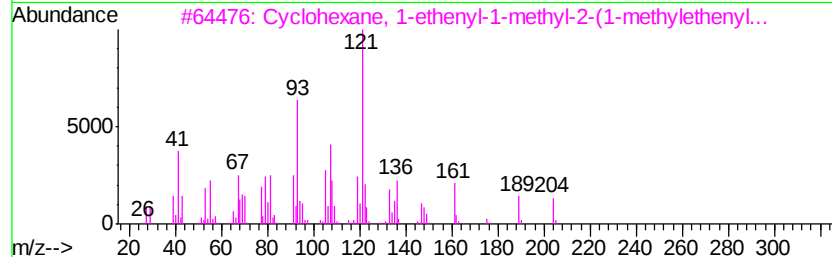
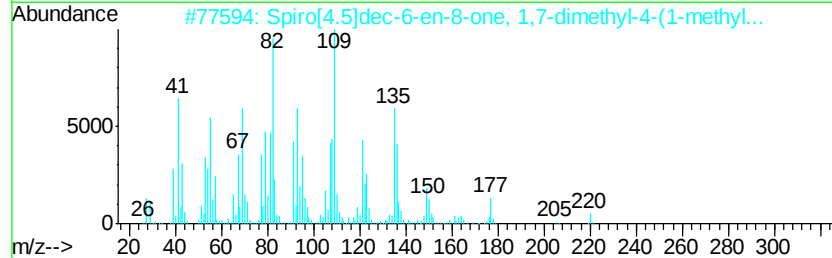
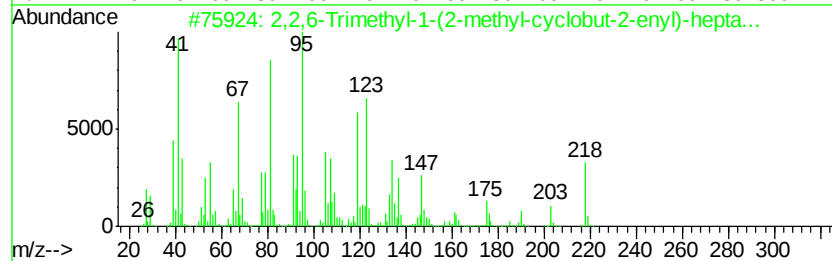
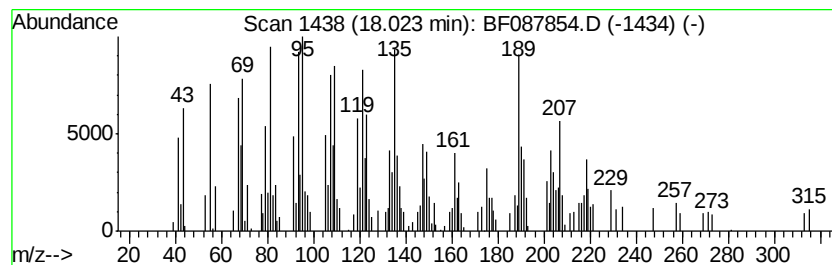
Instrument :  
BNA\_F  
ClientSampled :  
SS-35

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 17 2,2,6-Trimethyl-1-(2-methyl... Concentration Rank 11

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 18.02     | 6.75 ng                             | 228445 | Perylene-d12     | 15.36        |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | 2,2,6-Trimethyl-1-(2-methyl-cycl... | 218    | C15H22O          | 1000188-72-8 | 56   |
| 2         | Spiro[4.5]dec-6-en-8-one, 1,7-di... | 220    | C15H24O          | 039510-36-6  | 53   |
| 3         | Cyclohexane, 1-ethenyl-1-methyl-... | 204    | C15H24           | 003242-08-8  | 46   |
| 4         | Longifolenaldehyde                  | 220    | C15H24O          | 019890-84-7  | 46   |
| 5         | Guaia-9,11-diene                    | 204    | C15H24           | 1000374-19-8 | 42   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060916\  
Data File : BF087854.D  
Acq On : 9 Jun 2016 16:02  
Operator : UM/SJ  
Sample : H3399-16 5X  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SS-35

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

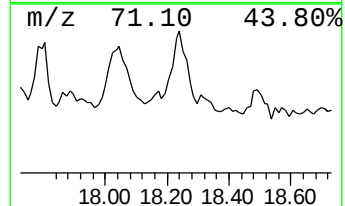
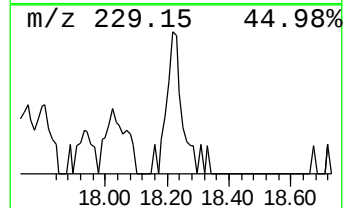
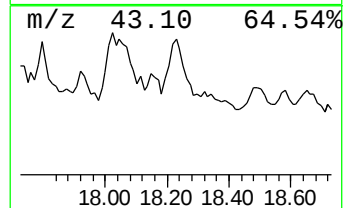
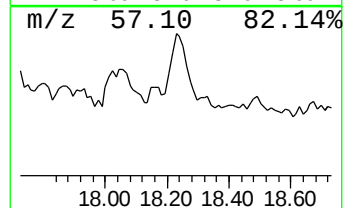
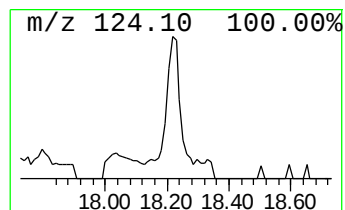
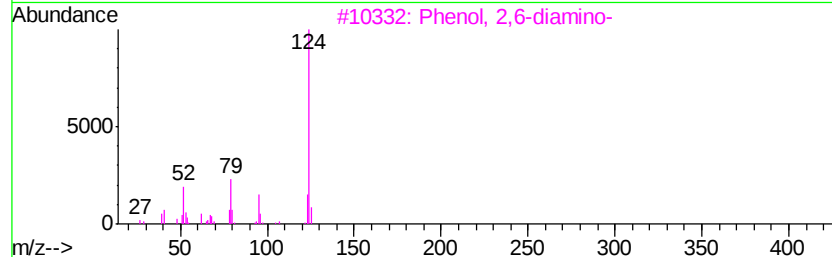
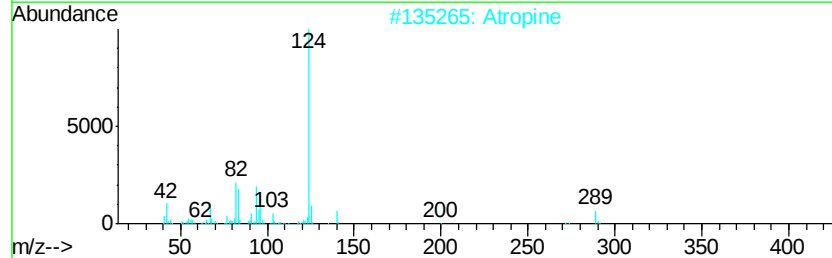
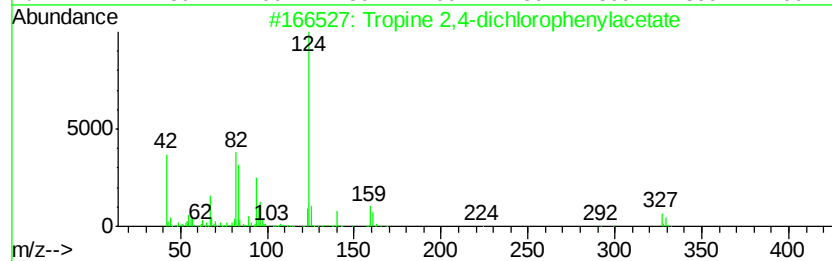
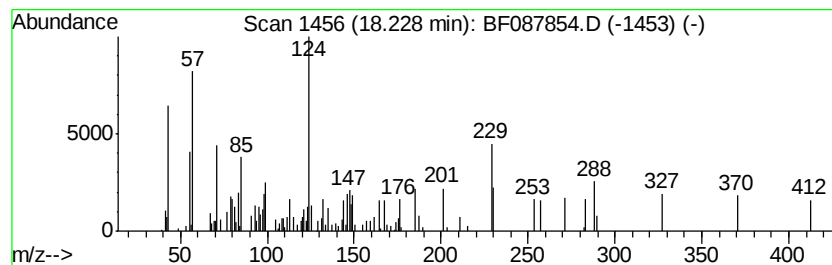
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 unknown18.23 Concentration Rank 18

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 18.23 | 3.09 ng | 104609 | Perylene-d12     | 15.36 |

| Hit# | of | Tentative ID                        | MW  | MolForm      | CAS#         | Qual |
|------|----|-------------------------------------|-----|--------------|--------------|------|
| 1    | 5  | Tropine 2,4-dichlorophenylacetate   | 327 | C16H19Cl2NO2 | 1000212-39-1 | 22   |
| 2    |    | Atropine                            | 289 | C17H23NO3    | 000051-55-8  | 18   |
| 3    |    | Phenol, 2,6-diamino-                | 124 | C6H8N2O      | 022440-82-0  | 18   |
| 4    |    | 1,2,4-Triazin-3-amine, 5,6-dimet... | 124 | C5H8N4       | 017584-12-2  | 18   |
| 5    |    | (2S,4aS,5R,8aR)-2,5-Dimethyldeca... | 167 | C11H21N      | 085956-21-4  | 18   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060916\  
Data File : BF087854.D  
Acq On : 9 Jun 2016 16:02  
Operator : UM/SJ  
Sample : H3399-16 5X  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampled :  
SS-35

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

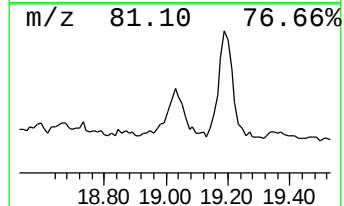
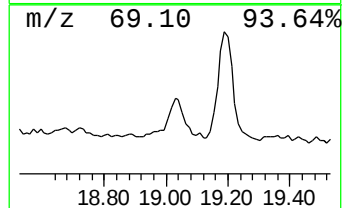
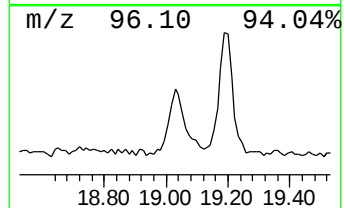
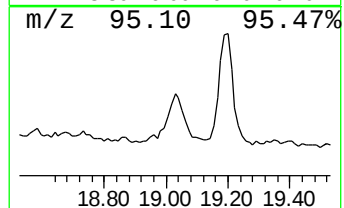
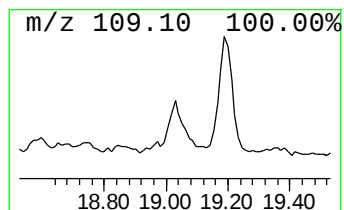
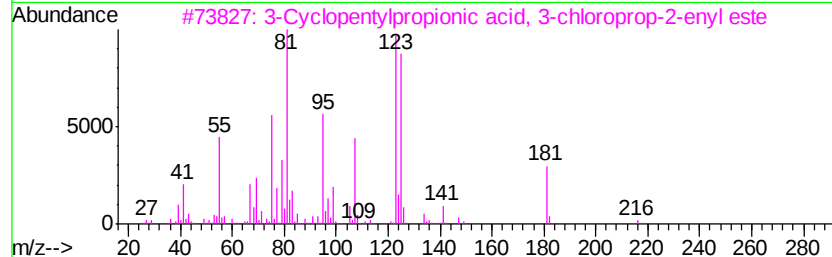
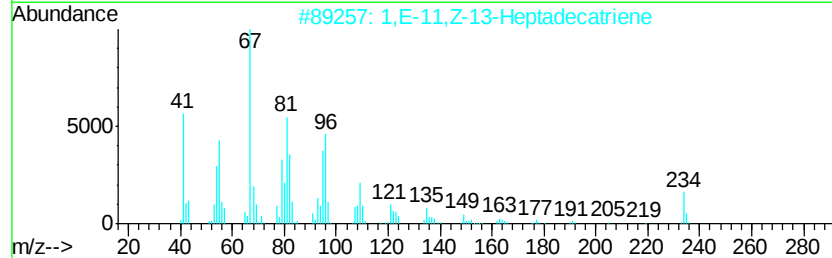
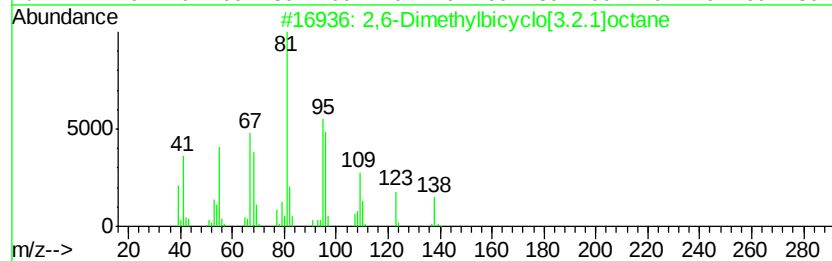
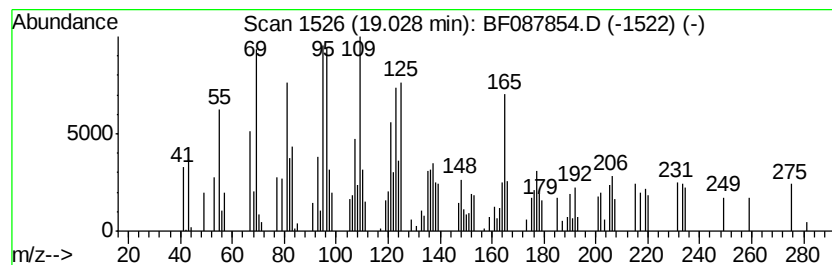
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 19 unknown19.03 Concentration Rank 12

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 19.03 | 4.79 ng | 162357 | Perylene-d12     | 15.36 |

| Hit# | of                                                     | 5   | Tentative ID | MW           | MolForm | CAS# | Qual |
|------|--------------------------------------------------------|-----|--------------|--------------|---------|------|------|
| 1    | 2,6-Dimethylbicyclo[3.2.1]octane                       | 138 | C10H18       | 1000215-28-2 | 30      |      |      |
| 2    | 1,E-11,Z-13-Heptadecatriene                            | 234 | C17H30       | 080625-35-0  | 27      |      |      |
| 3    | 3-Cyclopentylpropionic acid, 3-chloroprop-2-enyl ester | 216 | C11H17ClO2   | 1000292-47-2 | 22      |      |      |
| 4    | 2,5,9-Trimethylcycloundeca-4,8-diene                   | 206 | C14H22O      | 1000195-33-9 | 20      |      |      |
| 5    | 2,3,3-Trimethyl-2-(3-methyl-butyl)-2-butene            | 206 | C14H22O      | 1000193-60-6 | 20      |      |      |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF060916\  
Data File : BF087854.D  
Acq On : 9 Jun 2016 16:02  
Operator : UM/SJ  
Sample : H3399-16 5X  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SS-35

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

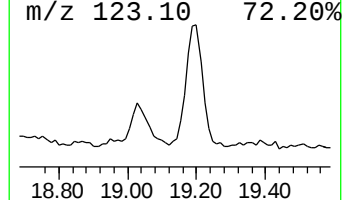
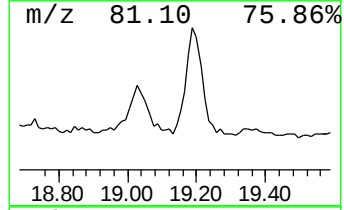
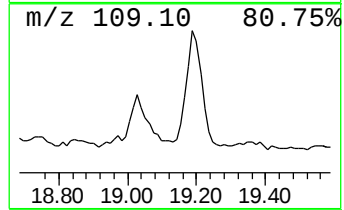
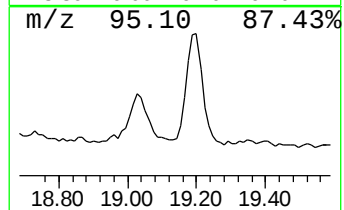
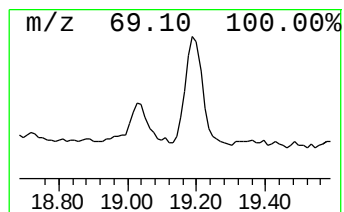
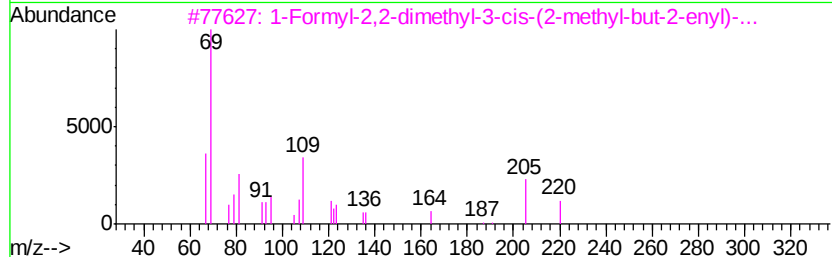
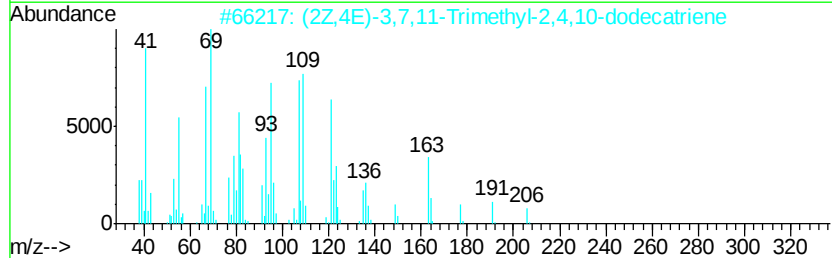
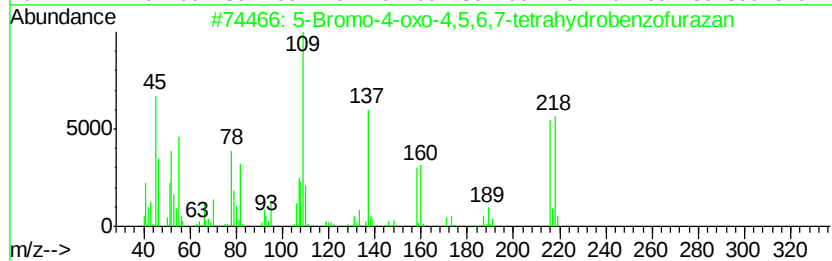
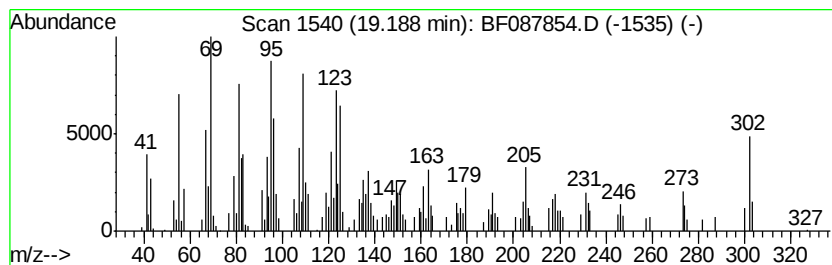
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 5-Bromo-4-oxo-4,5,6,7-tetra... Concentration Rank 5

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 19.19 | 12.24 ng | 414633 | Perylene-d12     | 15.36 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------|-----|------------|--------------|------|
| 1    | 5  | 5-Bromo-4-oxo-4,5,6,7-tetrahydro... | 216 | C6H5BrN2O2 | 300574-36-1  | 59   |
| 2    |    | (2Z,4E)-3,7,11-Trimethyl-2,4,10-... | 206 | C15H26     | 172549-29-0  | 38   |
| 3    |    | 1-Formyl-2,2-dimethyl-3-cis-(2-m... | 220 | C15H24O    | 1000144-10-0 | 38   |
| 4    |    | Bicyclo[2.2.1]heptane, 2,2,3-tri... | 138 | C10H18     | 020536-40-7  | 38   |
| 5    |    | Caparratriene                       | 206 | C15H26     | 1000374-08-7 | 25   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF060916\  
 Data File : BF087854.D  
 Acq On : 9 Jun 2016 16:02  
 Operator : UM/SJ  
 Sample : H3399-16 5X  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SS-35

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060716.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 |
| unknown6.58          | 6.58  | 21.3    | ng    | 759272   | 1                     | 6.82  | 714265 | 20.0 |
| Nonadecane           | 14.43 | 2.5     | ng    | 113190   | 5                     | 13.95 | 889894 | 20.0 |
| Heneicosane          | 15.05 | 11.0    | ng    | 372931   | 6                     | 15.36 | 677238 | 20.0 |
| Benz[e]acephenant... | 15.30 | 4.0     | ng    | 136276   | 6                     | 15.36 | 677238 | 20.0 |
| Heptadecane          | 15.82 | 9.0     | ng    | 304415   | 6                     | 15.36 | 677238 | 20.0 |
| Vitamin E            | 16.06 | 4.5     | ng    | 150866   | 6                     | 15.36 | 677238 | 20.0 |
| unknown16.89         | 16.89 | 3.9     | ng    | 130591   | 6                     | 15.36 | 677238 | 20.0 |
| unknown17.05         | 17.05 | 2.6     | ng    | 88895    | 6                     | 15.36 | 677238 | 20.0 |
| .gamma.-Sitosterol   | 17.28 | 28.3    | ng    | 959231   | 6                     | 15.36 | 677238 | 20.0 |
| 4-Allyl-5-pyridin... | 17.47 | 4.5     | ng    | 153979   | 6                     | 15.36 | 677238 | 20.0 |
| unknown17.59         | 17.59 | 4.0     | ng    | 135550   | 6                     | 15.36 | 677238 | 20.0 |
| 2,9-Dimethyl-2,3,... | 17.70 | 11.1    | ng    | 375299   | 6                     | 15.36 | 677238 | 20.0 |
| unknown17.79         | 17.79 | 8.9     | ng    | 301963   | 6                     | 15.36 | 677238 | 20.0 |
| 2,2,6-Trimethyl-1... | 18.02 | 6.8     | ng    | 228445   | 6                     | 15.36 | 677238 | 20.0 |
| unknown18.23         | 18.23 | 3.1     | ng    | 104609   | 6                     | 15.36 | 677238 | 20.0 |
| unknown19.03         | 19.03 | 4.8     | ng    | 162357   | 6                     | 15.36 | 677238 | 20.0 |
| 5-Bromo-4-oxo-4,5... | 19.19 | 12.2    | ng    | 414633   | 6                     | 15.36 | 677238 | 20.0 |