

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF102417\  
 Data File : BF099805.D  
 Acq On : 24 Oct 2017 23:14  
 Operator : SJ/JU  
 Sample : I5956-07  
 Misc :  
 ALS Vial : 24 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 EPE-120-1(20-22)

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-BF102317.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         | 5.022       | 496           | 499         | 519          | rBV      | 171697         | 255841        | 14.12%          | 1.789%        |
| 2         | 5.428       | 565           | 568         | 588          | rBV      | 1611727        | 1622314       | 89.55%          | 11.344%       |
| 3         | 6.445       | 737           | 741         | 760          | rBV      | 1607180        | 1604649       | 88.57%          | 11.220%       |
| 4         | 6.575       | 760           | 763         | 773          | rVB      | 1911618        | 1696887       | 93.66%          | 11.865%       |
| 5         | 6.804       | 799           | 802         | 814          | rBB      | 409877         | 354994        | 19.59%          | 2.482%        |
| 6         | 6.957       | 824           | 828         | 844          | rBB      | 1487164        | 1336380       | 73.76%          | 9.344%        |
| 7         | 7.375       | 895           | 899         | 919          | rBV      | 1017386        | 994605        | 54.90%          | 6.954%        |
| 8         | 8.087       | 1017          | 1020        | 1040         | rBV      | 577271         | 465883        | 25.72%          | 3.258%        |
| 9         | 8.645       | 1112          | 1115        | 1127         | rBB      | 82264          | 70900         | 3.91%           | 0.496%        |
| 10        | 9.163       | 1198          | 1203        | 1206         | rBV      | 1985283        | 1811694       | 100.00%         | 12.668%       |
| 11        | 9.557       | 1267          | 1270        | 1277         | rBB      | 30134          | 26724         | 1.48%           | 0.187%        |
| 12        | 9.839       | 1315          | 1318        | 1328         | rBV2     | 548070         | 482927        | 26.66%          | 3.377%        |
| 13        | 10.557      | 1436          | 1440        | 1448         | rBV      | 201057         | 185179        | 10.22%          | 1.295%        |
| 14        | 10.633      | 1450          | 1453        | 1467         | rBV      | 894629         | 815705        | 45.02%          | 5.704%        |
| 15        | 11.333      | 1568          | 1572        | 1580         | rBV      | 457348         | 396840        | 21.90%          | 2.775%        |
| 16        | 12.916      | 1837          | 1841        | 1844         | rBV      | 1342236        | 1108879       | 61.21%          | 7.753%        |
| 17        | 13.798      | 1987          | 1991        | 2003         | rBV2     | 32355          | 43759         | 2.42%           | 0.306%        |
| 18        | 13.980      | 2018          | 2022        | 2030         | rBV      | 347564         | 311041        | 17.17%          | 2.175%        |
| 19        | 15.045      | 2200          | 2203        | 2207         | rBV2     | 17739          | 21094         | 1.16%           | 0.147%        |
| 20        | 15.421      | 2262          | 2267        | 2268         | rBV      | 32971          | 38944         | 2.15%           | 0.272%        |
| 21        | 15.445      | 2268          | 2271        | 2281         | rVB      | 227289         | 302593        | 16.70%          | 2.116%        |
| 22        | 15.839      | 2333          | 2338        | 2346         | rBV      | 43314          | 66542         | 3.67%           | 0.465%        |
| 23        | 16.327      | 2414          | 2421        | 2428         | rBV      | 42335          | 75459         | 4.17%           | 0.528%        |
| 24        | 16.904      | 2511          | 2519        | 2528         | rBV2     | 33529          | 67539         | 3.73%           | 0.472%        |
| 25        | 17.580      | 2627          | 2634        | 2643         | rVB3     | 28761          | 66745         | 3.68%           | 0.467%        |
| 26        | 18.386      | 2765          | 2771        | 2780         | rVB6     | 14771          | 37679         | 2.08%           | 0.263%        |
| 27        | 19.345      | 2924          | 2934        | 2944         | rBV7     | 11097          | 39875         | 2.20%           | 0.279%        |

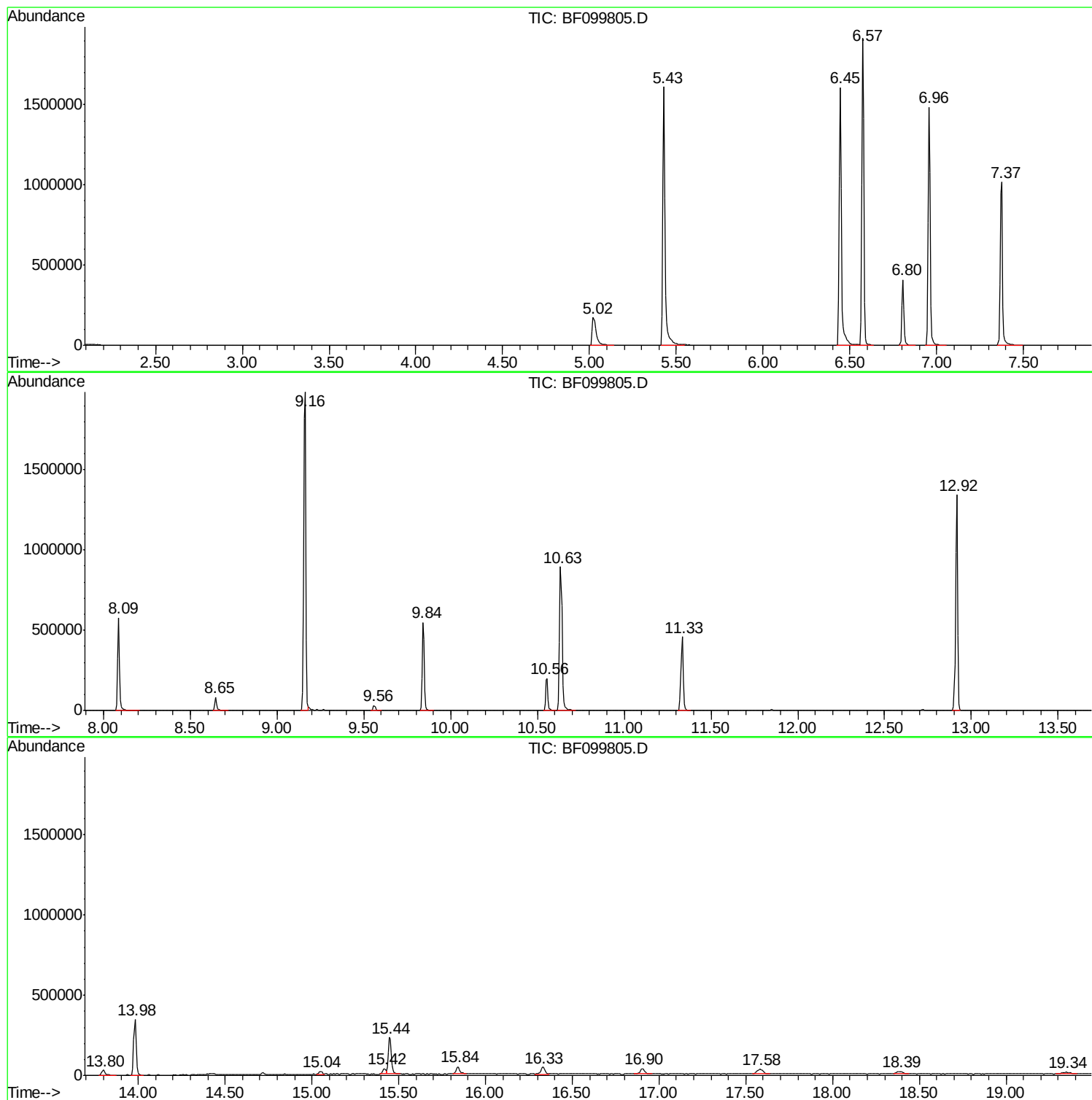
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EPE-120-1(20-22)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF102317.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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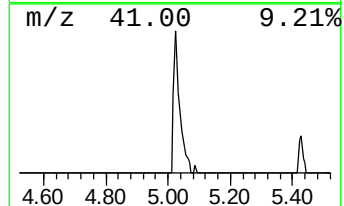
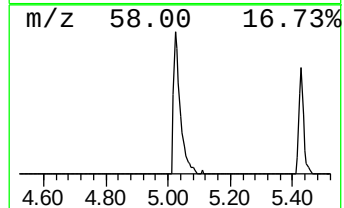
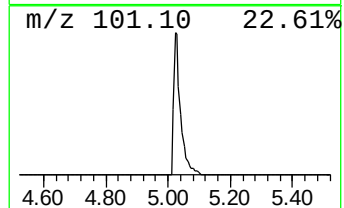
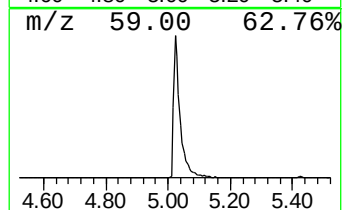
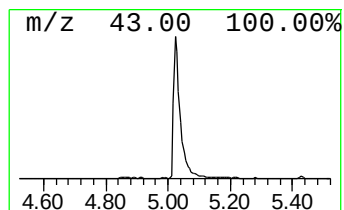
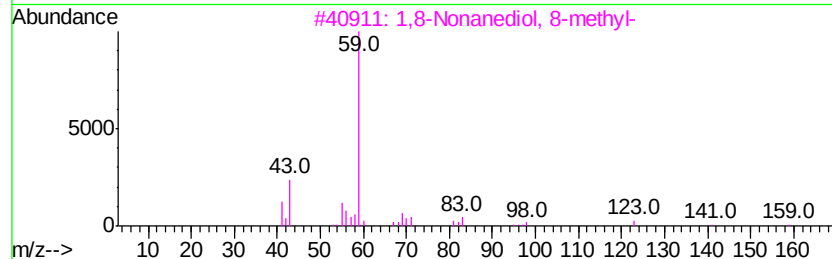
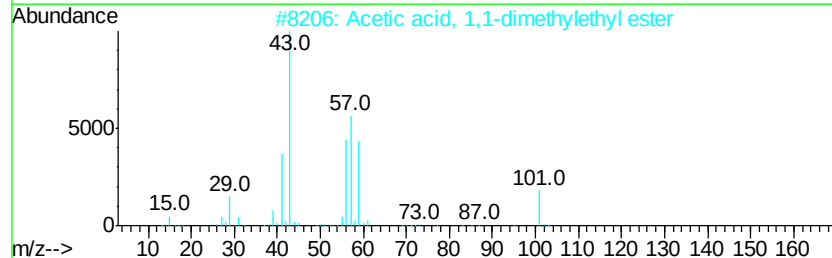
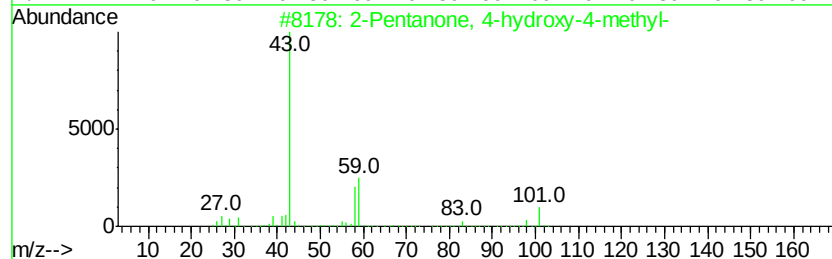
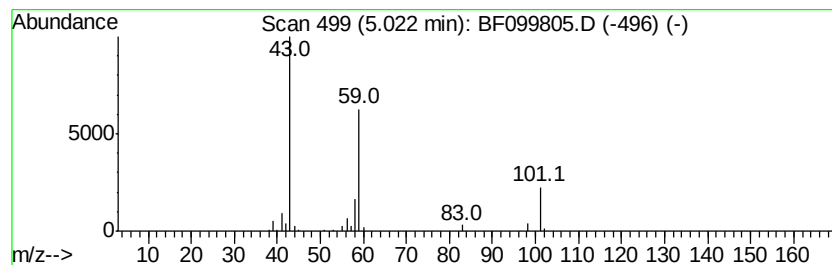
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 5.02 | 14.41 ng | 255841 | 1,4-Dichlorobenzene-d4 | 6.80 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2  | 000123-42-2 | 59   |
| 2    |    |   | Acetic acid, 1,1-dimethylethyl e... | 116 | C6H12O2  | 000540-88-5 | 39   |
| 3    |    |   | 1,8-Nonanediol, 8-methyl-           | 174 | C10H22O2 | 054725-73-4 | 23   |
| 4    |    |   | Acetic acid, cyano-, 1,1-dimethy... | 141 | C7H11NO2 | 001116-98-9 | 17   |
| 5    |    |   | Morpholine, 4-methyl-               | 101 | C5H11NO  | 000109-02-4 | 9    |



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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

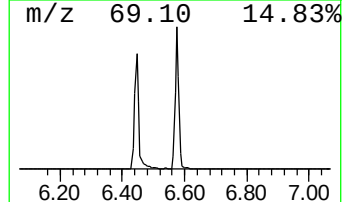
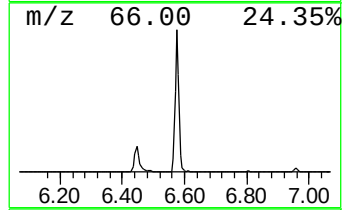
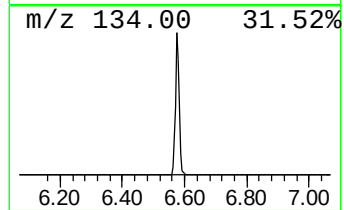
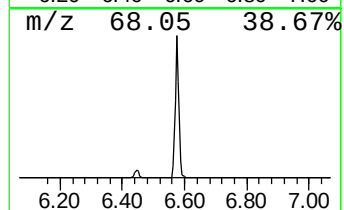
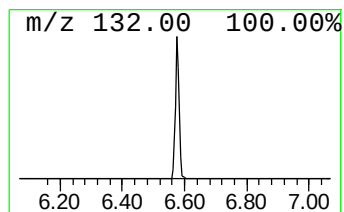
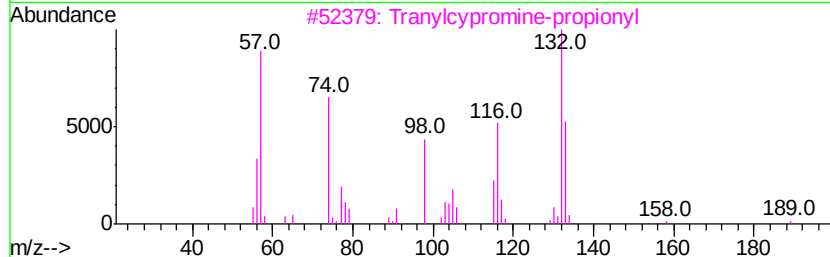
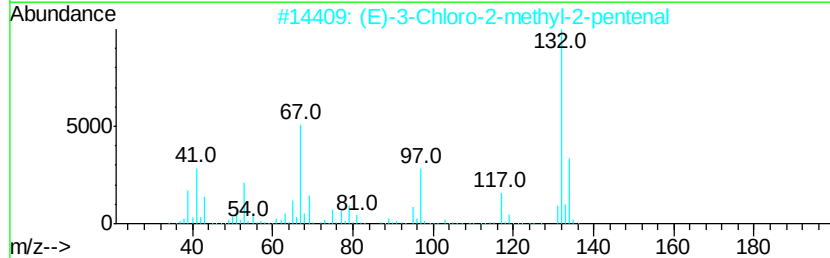
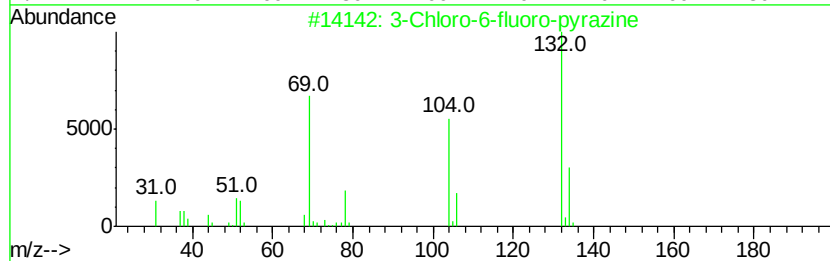
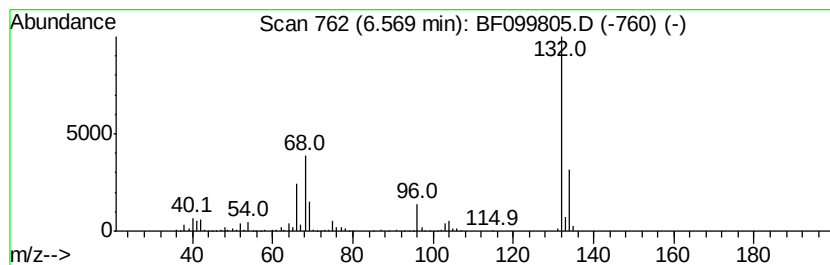
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 unknown6.57 Concentration Rank 1

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.57 | 95.60 ng | 1696890 | 1,4-Dichlorobenzene-d4 | 6.80 |

| Hit# | of | Tentative ID                     | MW  | MolForm   | CAS#         | Qual |
|------|----|----------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine       | 132 | C4H2ClFN2 | 1000146-10-7 | 38   |
| 2    |    | (E)-3-Chloro-2-methyl-2-pentenal | 132 | C6H9ClO   | 031357-76-3  | 35   |
| 3    |    | Tranvlcypropine-propionyl        | 189 | C12H15NO  | 1000123-86-3 | 12   |
| 4    |    | 1H-Indazole, 7-methyl-           | 132 | C8H8N2    | 1000316-00-7 | 10   |
| 5    |    | Xenon                            | 132 | Xe        | 007440-63-3  | 9    |



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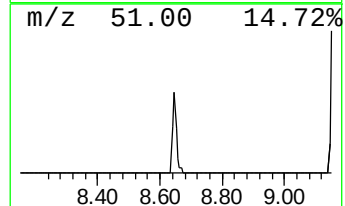
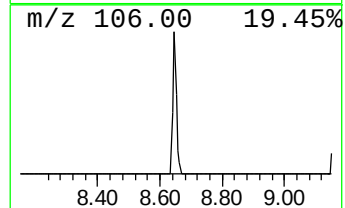
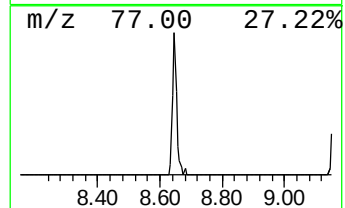
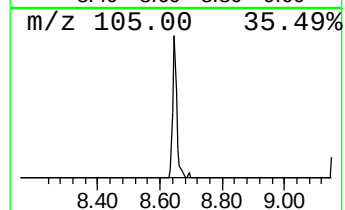
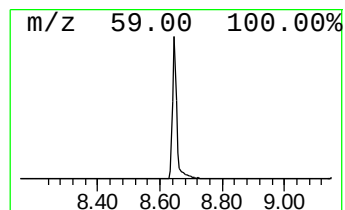
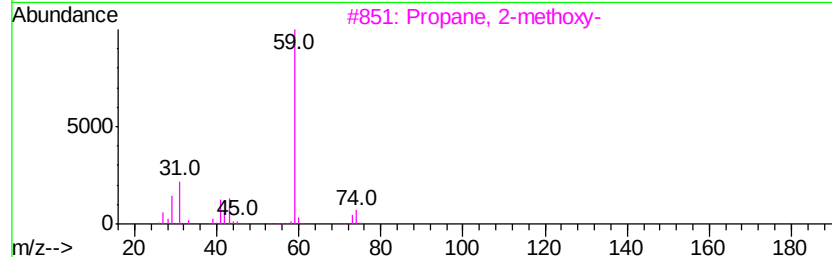
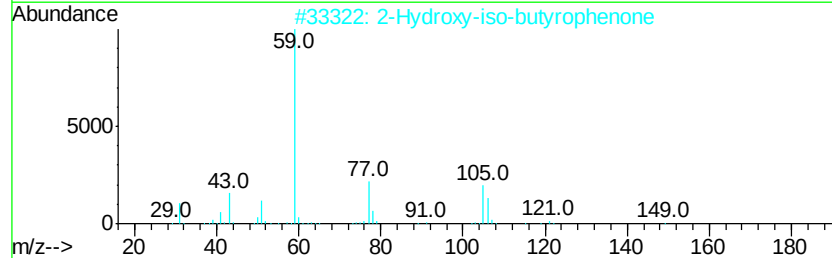
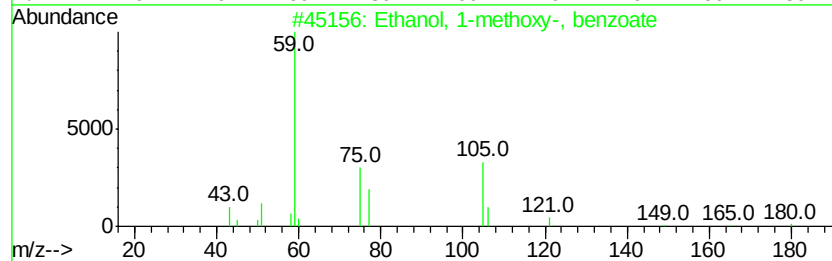
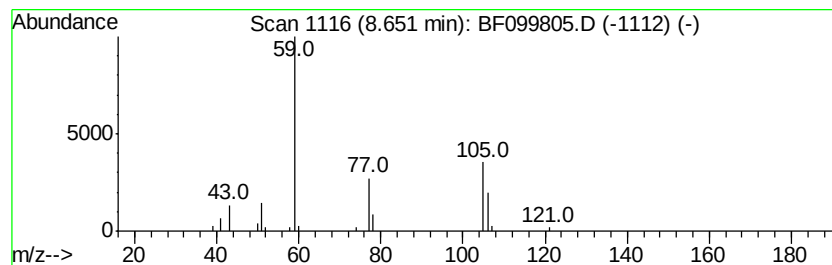
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 Ethanol, 1-methoxy-, benzoate Concentration Rank 9

| R.T. | EstConc | Area  | Relative to ISTD | R.T. |
|------|---------|-------|------------------|------|
| 8.65 | 3.04 ng | 70900 | Napthalene-d8    | 8.09 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Ethanol, 1-methoxy-, benzoate       | 180 | C10H12O3 | 051835-44-0 | 56   |
| 2    |    | 2-Hydroxy-iso-butyrophenone         | 164 | C10H12O2 | 007473-98-5 | 53   |
| 3    |    | Propane, 2-methoxy-                 | 74  | C4H10O   | 000598-53-8 | 9    |
| 4    |    | 1-Propanol, 2-methoxy-              | 90  | C4H10O2  | 001589-47-5 | 9    |
| 5    |    | Propanoic acid, 2-hydroxy-2-meth... | 118 | C5H10O3  | 002110-78-3 | 9    |



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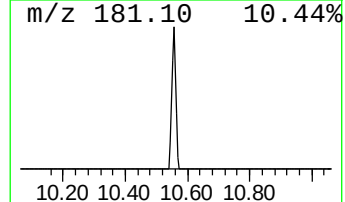
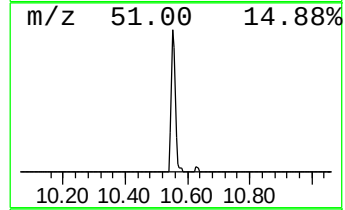
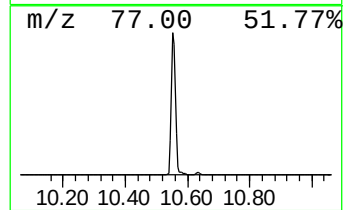
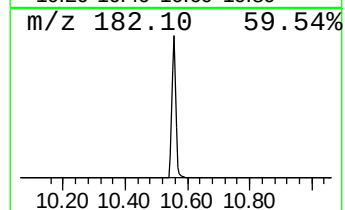
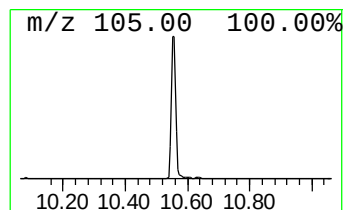
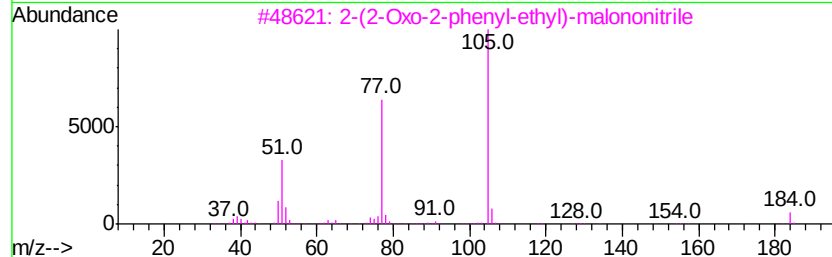
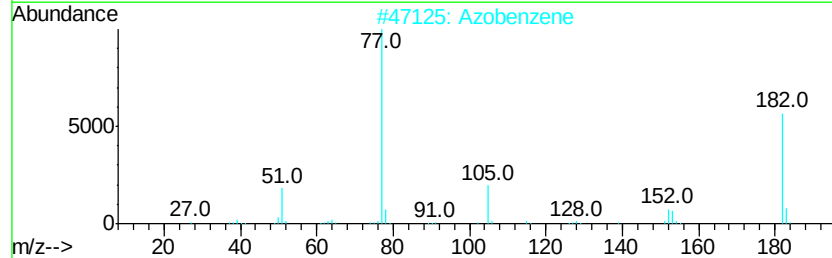
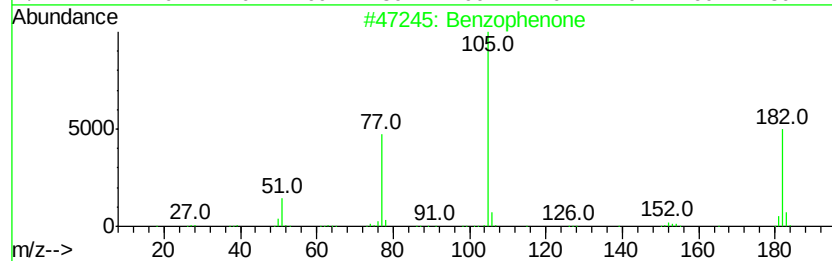
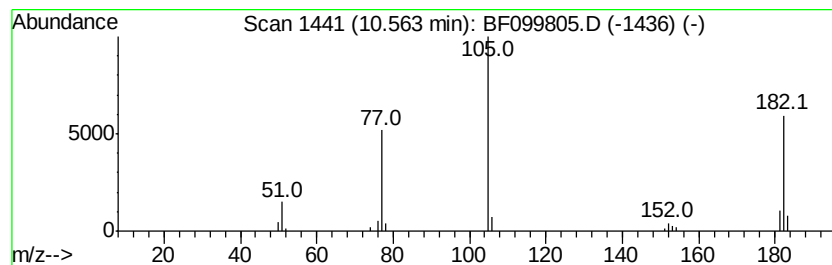
Instrument :  
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ClientSampleId :  
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Peak Number 5 Benzophenone Concentration Rank 4

| R.T.    | EstConc | Area                                   | Relative to ISTD |          | R.T.         |      |
|---------|---------|----------------------------------------|------------------|----------|--------------|------|
| 10.56   | 7.67 ng | 185179                                 | Acenaphthene-d10 |          | 9.84         |      |
| Hit# of | 5       | Tentative ID                           | MW               | MolForm  | CAS#         | Qual |
| 1       |         | Benzophenone                           | 182              | C13H10O  | 000119-61-9  | 97   |
| 2       |         | Azobenzene                             | 182              | C12H10N2 | 000103-33-3  | 47   |
| 3       |         | 2-(2-Oxo-2-phenyl-ethyl)-malononitrile | 184              | C11H8N2O | 1000296-76-9 | 47   |
| 4       |         | Benzeneacetic acid, .alpha.-oxo-       | 164              | C9H8O3   | 015206-55-0  | 43   |
| 5       |         | Acetophenone, 2-chloro-                | 154              | C8H7ClO  | 000532-27-4  | 43   |



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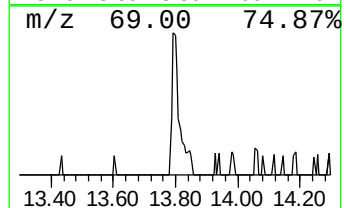
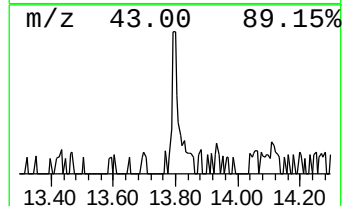
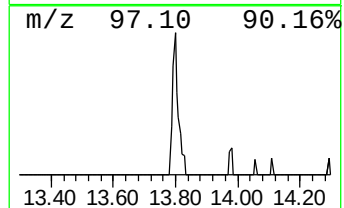
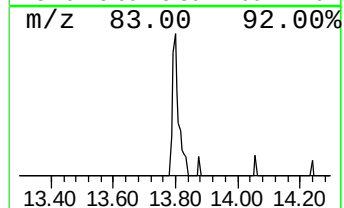
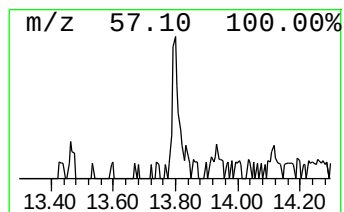
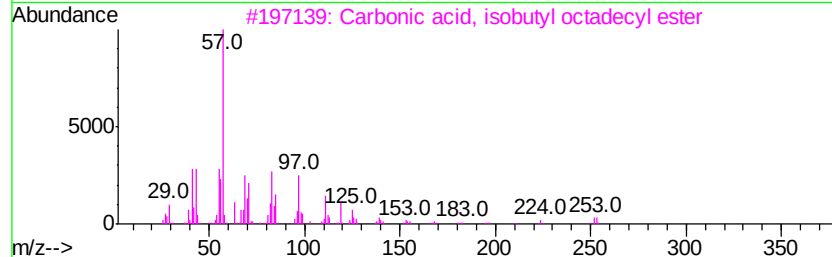
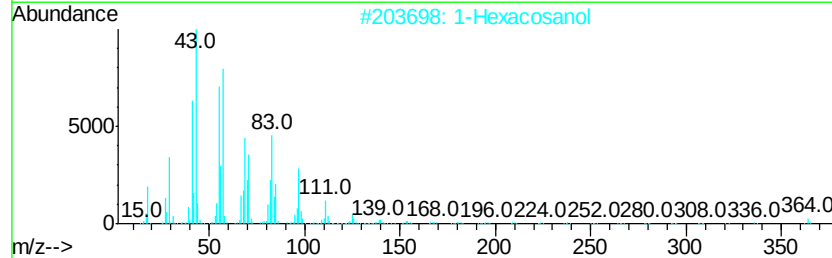
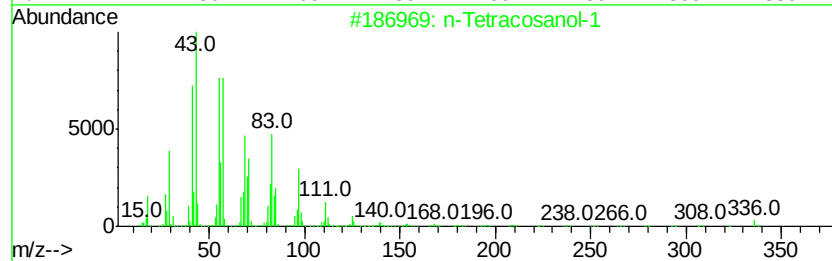
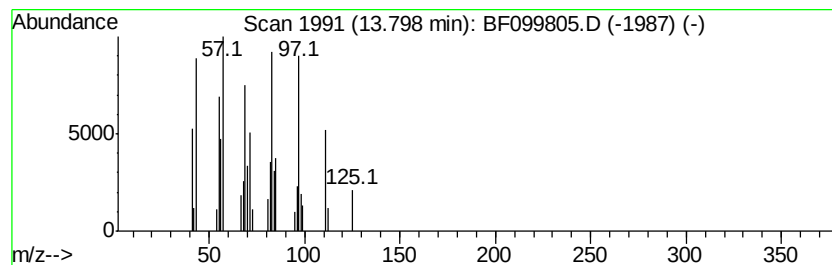
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 n-Tetracosanol-1 Concentration Rank 10

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 13.80 | 2.81 ng | 43759 | Chrysene-d12     | 13.98 |

| Hit# | of | Tentative ID                         | MW  | MolForm   | CAS#         | Qual |
|------|----|--------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | n-Tetracosanol-1                     | 354 | C24H50O   | 000506-51-4  | 68   |
| 2    |    | 1-Hexacosanol                        | 382 | C26H54O   | 000506-52-5  | 68   |
| 3    |    | Carbonic acid, isobutyl octadecyl... | 370 | C23H46O3  | 1000314-61-5 | 52   |
| 4    |    | Cyclopentane, 1,1,3,4-tetramethyl... | 126 | C9H18     | 020309-77-7  | 38   |
| 5    |    | Sulfurous acid, cyclohexylmethyl...  | 402 | C23H46O3S | 1000309-22-4 | 35   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF102417\  
Data File : BF099805.D  
Acq On : 24 Oct 2017 23:14  
Operator : SJ/JU  
Sample : I5956-07  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
EPE-120-1(20-22)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF102317.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

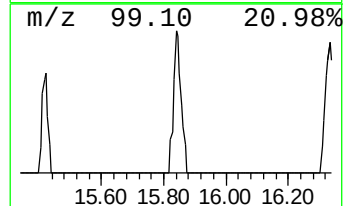
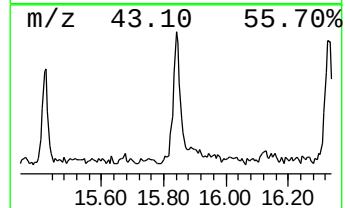
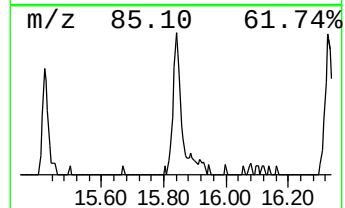
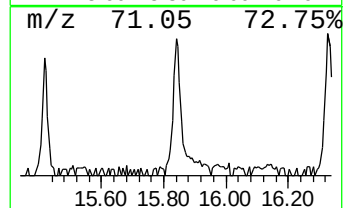
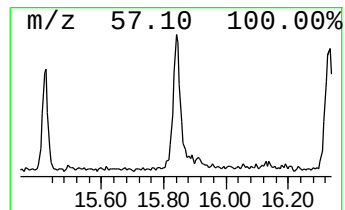
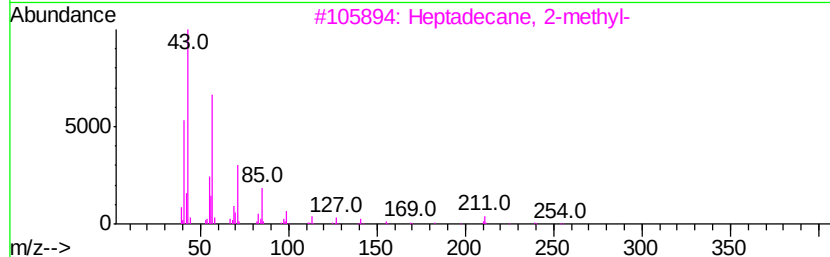
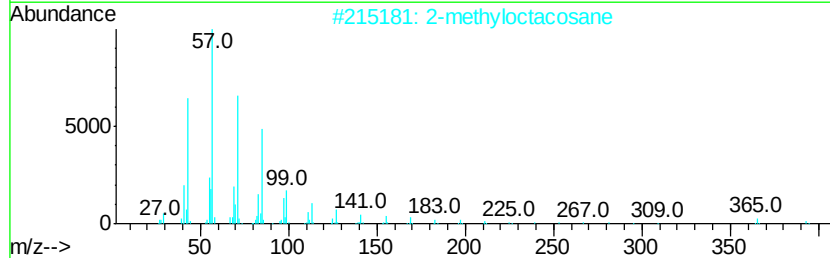
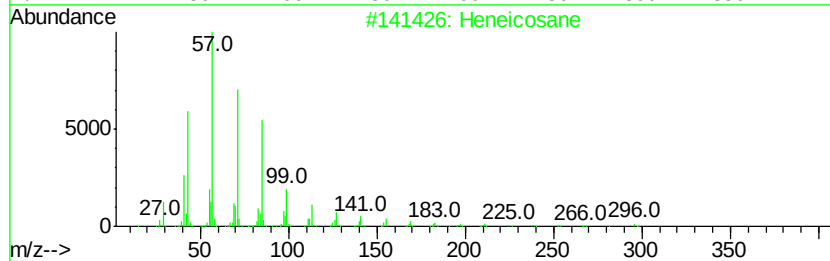
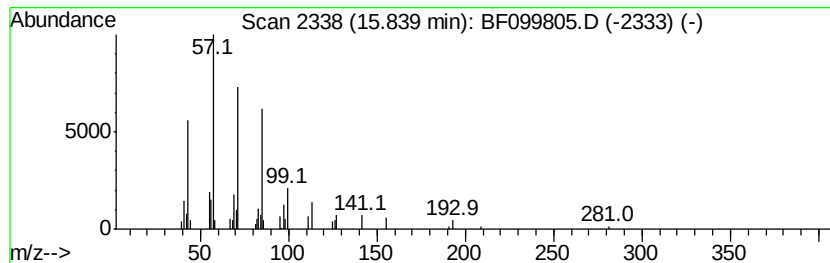
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 Heneicosane Concentration Rank 8

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 15.84 | 4.40 ng | 66542 | Perylene-d12     | 15.44 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#         | Qual |
|------|----|------------------------|-----|---------|--------------|------|
| 1    | 5  | Heneicosane            | 296 | C21H44  | 000629-94-7  | 91   |
| 2    |    | 2-methyloctacosane     | 408 | C29H60  | 1000376-72-8 | 90   |
| 3    |    | Heptadecane, 2-methyl- | 254 | C18H38  | 001560-89-0  | 90   |
| 4    |    | triacontane            | 422 | C30H62  | 000638-68-6  | 90   |
| 5    |    | Eicosane, 10-methyl-   | 296 | C21H44  | 054833-23-7  | 86   |





Data Path : Z:\HPCHEM1\BNA F\DATA\BF102417\  
Data File : BF099805.D  
Acq On : 24 Oct 2017 23:14  
Operator : SJ/JU  
Sample : I5956-07  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

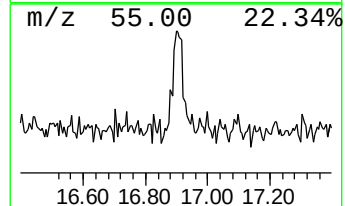
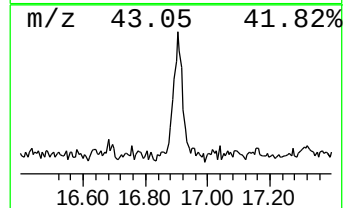
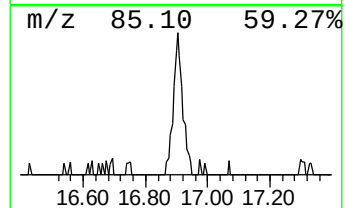
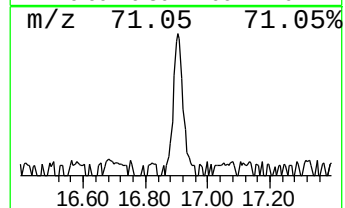
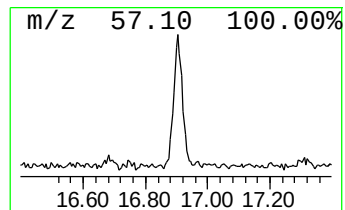
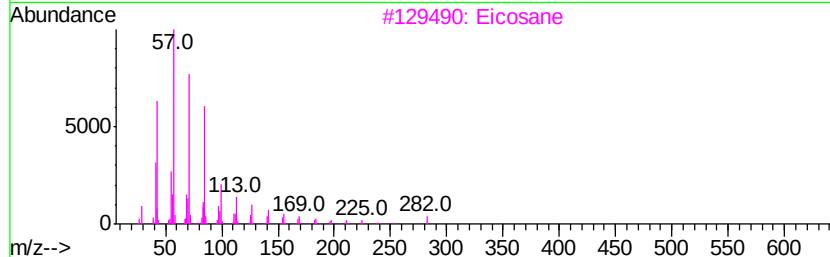
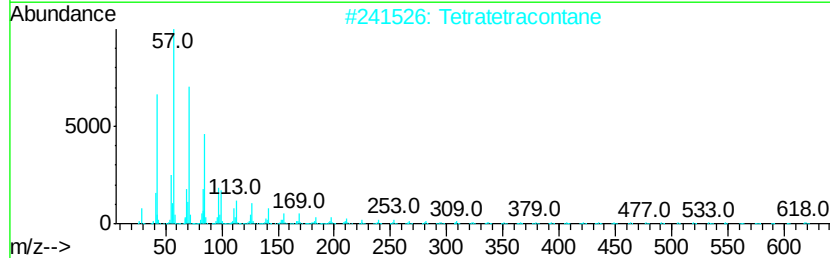
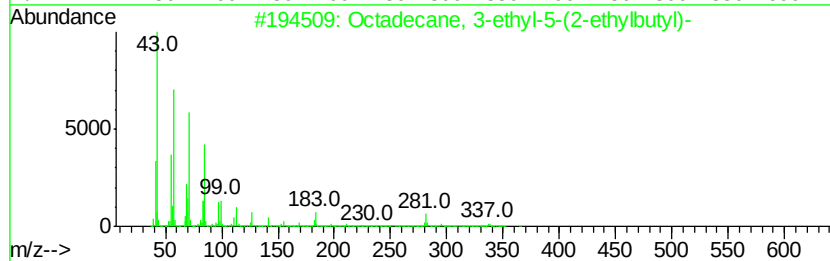
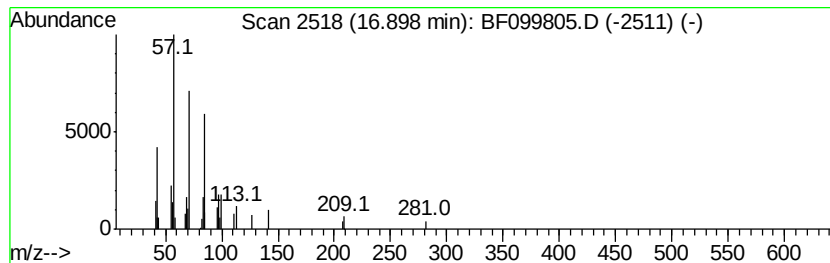
Instrument :  
BNA\_F  
ClientSampleId :  
EPE-120-1(20-22)

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102317.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 9 Octadecane, 3-ethyl-5-(2-ethylbu... Concentration Rank 6

| R.T.      | EstConc                             | Area  | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|-------|------------------|-------------|------|
| 16.90     | 4.46 ng                             | 67539 | Perylene-d12     | 15.44       |      |
| Hit# of 5 | Tentative ID                        | MW    | MolForm          | CAS#        | Qual |
| 1         | Octadecane, 3-ethyl-5-(2-ethylbu... | 366   | C26H54           | 055282-12-7 | 90   |
| 2         | Tetratetracontane                   | 619   | C44H90           | 007098-22-8 | 90   |
| 3         | Eicosane                            | 282   | C20H42           | 000112-95-8 | 87   |
| 4         | Heptacosane                         | 380   | C27H56           | 000593-49-7 | 86   |
| 5         | Hexacosane                          | 366   | C26H54           | 000630-01-3 | 80   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF102417\  
Data File : BF099805.D  
Acq On : 24 Oct 2017 23:14  
Operator : SJ/JU  
Sample : I5956-07  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
EPE-120-1(20-22)

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

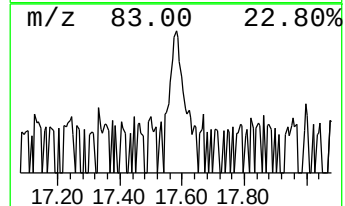
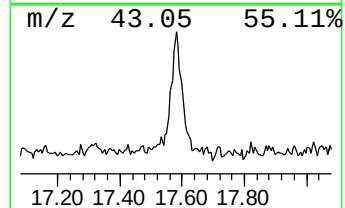
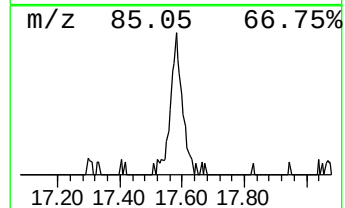
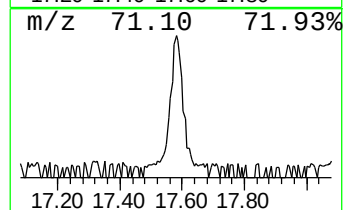
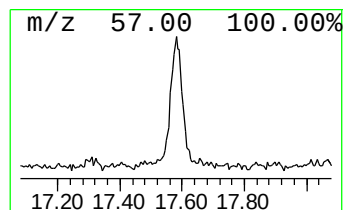
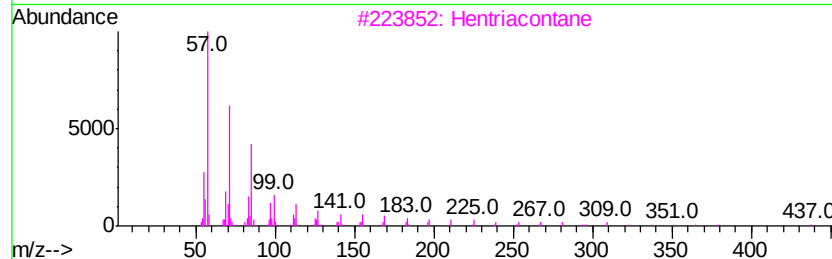
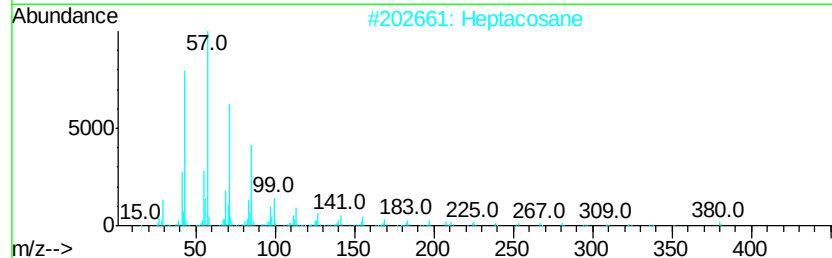
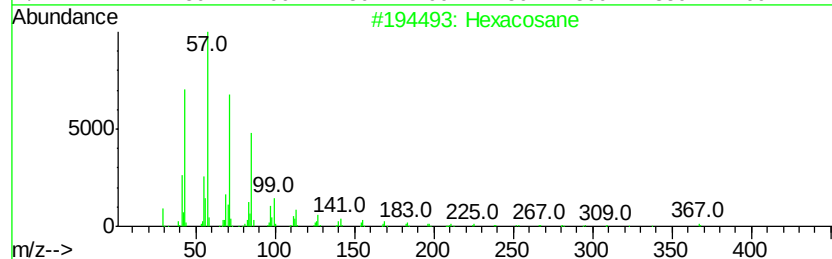
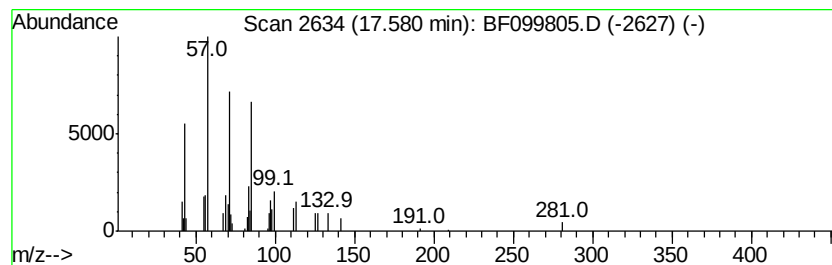
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 Hexacosane Concentration Rank 7

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 17.58 | 4.41 ng | 66745 | Perylene-d12     | 15.44 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm     | CAS#        | Qual |
|------|----|---|-----------------------------|-----|-------------|-------------|------|
| 1    |    |   | Hexacosane                  | 366 | C26H54      | 000630-01-3 | 80   |
| 2    |    |   | Heptacosane                 | 380 | C27H56      | 000593-49-7 | 80   |
| 3    |    |   | Hentriacontane              | 437 | C31H64      | 000630-04-6 | 72   |
| 4    |    |   | Silane, trichlorooctadecyl- | 386 | C18H37Cl3Si | 000112-04-9 | 72   |
| 5    |    |   | Eicosane, 10-methyl-        | 296 | C21H44      | 054833-23-7 | 72   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF102417\  
Data File : BF099805.D  
Acq On : 24 Oct 2017 23:14  
Operator : SJ/JU  
Sample : I5956-07  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
EPE-120-1(20-22)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF102317.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

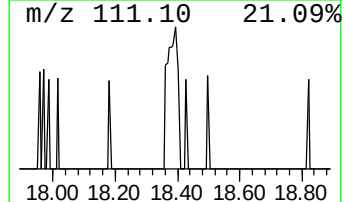
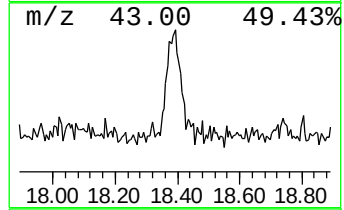
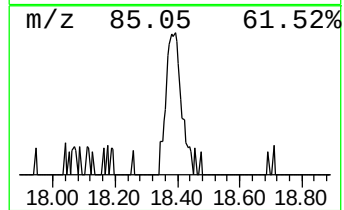
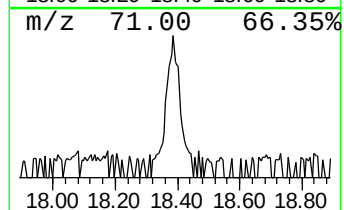
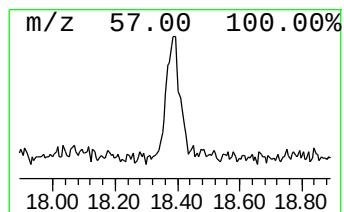
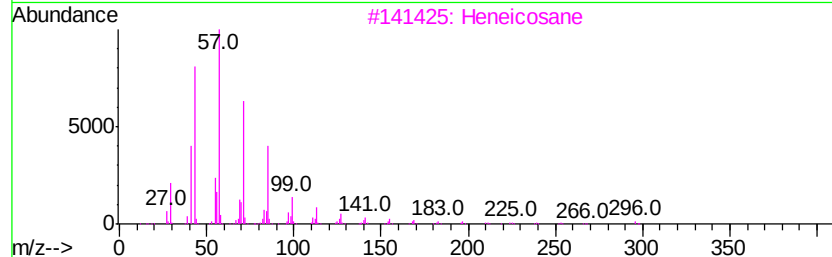
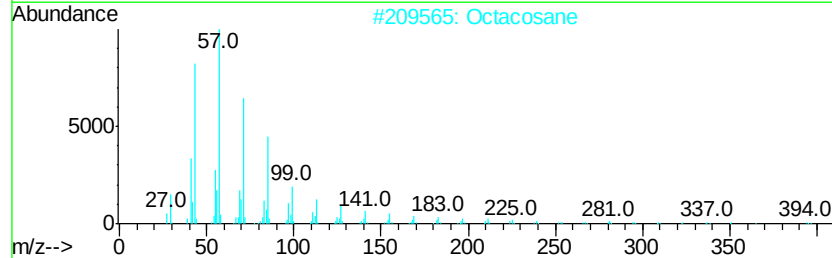
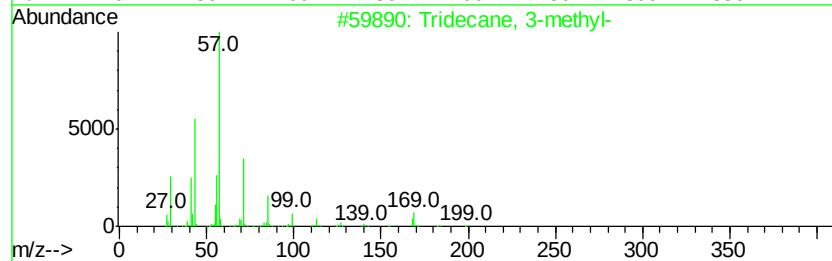
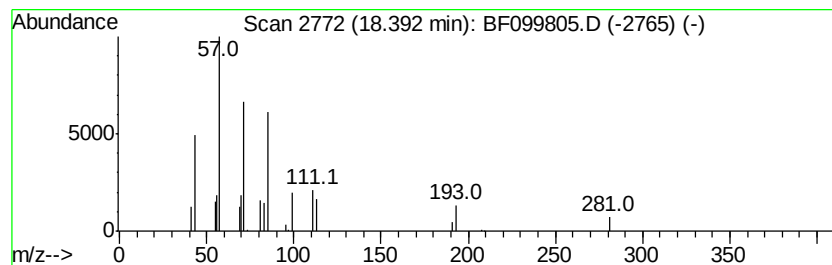
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 Tridecane, 3-methyl- Concentration Rank 12

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 18.39 | 2.49 ng | 37679 | Perylene-d12     | 15.44 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------|-----|---------|-------------|------|
| 1    |    |   | Tridecane, 3-methyl-  | 198 | C14H30  | 006418-41-3 | 80   |
| 2    |    |   | Octacosane            | 394 | C28H58  | 000630-02-4 | 72   |
| 3    |    |   | Heneicosane           | 296 | C21H44  | 000629-94-7 | 72   |
| 4    |    |   | Nonadecane, 2-methyl- | 282 | C20H42  | 001560-86-7 | 72   |
| 5    |    |   | Eicosane, 2-methyl-   | 296 | C21H44  | 001560-84-5 | 72   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF102417\  
Data File : BF099805.D  
Acq On : 24 Oct 2017 23:14  
Operator : SJ/JU  
Sample : I5956-07  
Misc :  
ALS Vial : 24 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
EPE-120-1(20-22)

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

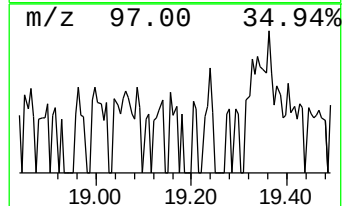
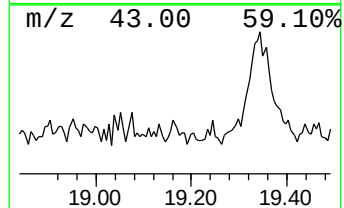
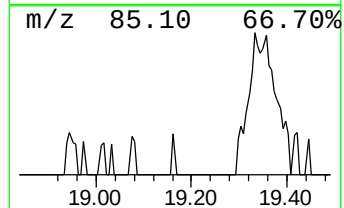
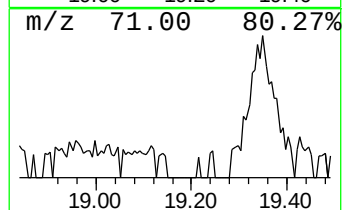
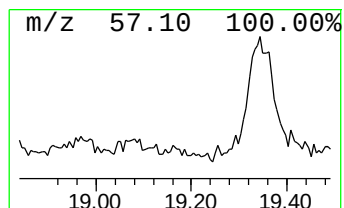
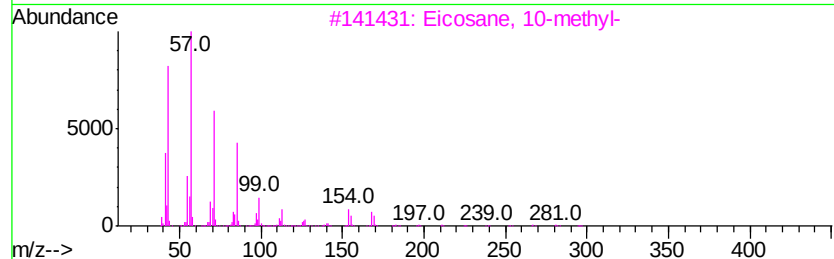
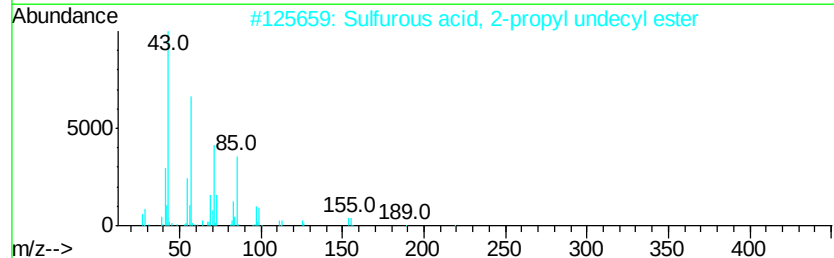
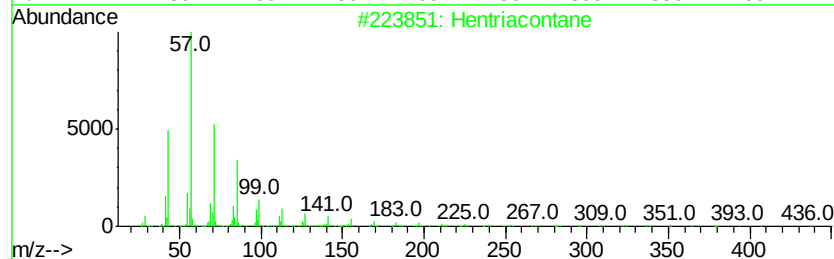
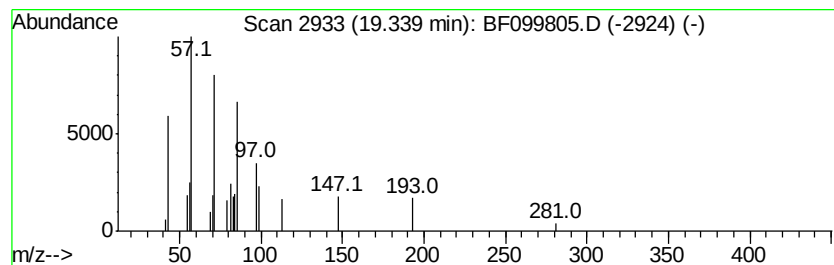
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 12 Hentriacontane Concentration Rank 11

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 19.34 | 2.64 ng | 39875 | Perylene-d12     | 15.44 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Hentriacontane                      | 437 | C31H64    | 000630-04-6  | 59   |
| 2    |    | Sulfurous acid, 2-propyl undecyl... | 278 | C14H30O3S | 1000309-12-2 | 50   |
| 3    |    | Eicosane, 10-methyl-                | 296 | C21H44    | 054833-23-7  | 43   |
| 4    |    | Sulfurous acid, butyl heptadecyl... | 376 | C21H44O3S | 1000309-18-4 | 42   |
| 5    |    | Tetracontane, 3,5,24-trimethyl-     | 605 | C43H88    | 055162-61-3  | 40   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF102417\  
 Data File : BF099805.D  
 Acq On : 24 Oct 2017 23:14  
 Operator : SJ/JU  
 Sample : I5956-07  
 Misc :  
 ALS Vial : 24 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 EPE-120-1(20-22)

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF102317.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 |
| 2-Pentanone, 4-hy... | 5.02  | 14.4    | ng    | 255841   | 1                     | 6.80  | 354994 | 20.0 |
| unknown6.57          | 6.57  | 95.6    | ng    | 1696890  | 1                     | 6.80  | 354994 | 20.0 |
| Ethanol, 1-methox... | 8.65  | 3.0     | ng    | 70900    | 2                     | 8.09  | 465883 | 20.0 |
| Benzophenone         | 10.56 | 7.7     | ng    | 185179   | 3                     | 9.84  | 482927 | 20.0 |
| n-Tetracosanol-1     | 13.80 | 2.8     | ng    | 43759    | 5                     | 13.98 | 311041 | 20.0 |
| Heneicosane          | 15.84 | 4.4     | ng    | 66542    | 6                     | 15.44 | 302593 | 20.0 |
| Octadecane, 3-eth... | 16.90 | 4.5     | ng    | 67539    | 6                     | 15.44 | 302593 | 20.0 |
| Hexacosane           | 17.58 | 4.4     | ng    | 66745    | 6                     | 15.44 | 302593 | 20.0 |
| Tridecane, 3-methyl- | 18.39 | 2.5     | ng    | 37679    | 6                     | 15.44 | 302593 | 20.0 |
| Hentriacontane       | 19.34 | 2.6     | ng    | 39875    | 6                     | 15.44 | 302593 | 20.0 |