

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF051519\  
 Data File : BF114284.D  
 Acq On : 15 May 2019 14:48  
 Operator : JU/SJ  
 Sample : K2837-01  
 Misc :  
 ALS Vial : 8 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SOIL-COMP

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

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

Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF051019.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 2.110       | 3             | 4           | 23           | rVB      | 1072741        | 1223102       | 24.45%          | 2.497%        |
| 2         | 5.481       | 573           | 577         | 594          | rBV      | 4394401        | 4706192       | 94.07%          | 9.609%        |
| 3         | 6.492       | 745           | 749         | 767          | rBV      | 4521507        | 5003021       | 100.00%         | 10.216%       |
| 4         | 6.622       | 767           | 771         | 775          | rBV      | 4783211        | 4753358       | 95.01%          | 9.706%        |
| 5         | 6.845       | 806           | 809         | 813          | rBV      | 1507147        | 1284238       | 25.67%          | 2.622%        |
| 6         | 7.004       | 832           | 836         | 840          | rBV      | 4094554        | 3510599       | 70.17%          | 7.168%        |
| 7         | 7.410       | 900           | 905         | 917          | rBV      | 3192146        | 3098807       | 61.94%          | 6.327%        |
| 8         | 8.128       | 1023          | 1027        | 1044         | rBV      | 1758677        | 1643800       | 32.86%          | 3.356%        |
| 9         | 9.204       | 1206          | 1210        | 1213         | rBV      | 5440187        | 4727322       | 94.49%          | 9.653%        |
| 10        | 9.598       | 1273          | 1277        | 1289         | rBV      | 568295         | 578435        | 11.56%          | 1.181%        |
| 11        | 9.886       | 1322          | 1326        | 1336         | rVB      | 1884599        | 1821842       | 36.41%          | 3.720%        |
| 12        | 10.674      | 1456          | 1460        | 1476         | rBV      | 3521320        | 3339598       | 66.75%          | 6.819%        |
| 13        | 11.368      | 1574          | 1578        | 1589         | rBV      | 2068270        | 2011912       | 40.21%          | 4.108%        |
| 14        | 11.892      | 1664          | 1667        | 1680         | rBV2     | 229324         | 274991        | 5.50%           | 0.561%        |
| 15        | 12.680      | 1797          | 1801        | 1808         | rBV2     | 39445          | 60998         | 1.22%           | 0.125%        |
| 16        | 12.951      | 1843          | 1847        | 1851         | rBV      | 4947281        | 4756608       | 95.07%          | 9.712%        |
| 17        | 13.521      | 1938          | 1944        | 1946         | rBV2     | 41751          | 54296         | 1.09%           | 0.111%        |
| 18        | 13.851      | 1996          | 2000        | 2015         | rBV      | 400355         | 659810        | 13.19%          | 1.347%        |
| 19        | 14.015      | 2024          | 2028        | 2042         | rVB      | 1844940        | 1884057       | 37.66%          | 3.847%        |
| 20        | 14.174      | 2051          | 2055        | 2060         | rBV      | 113547         | 109209        | 2.18%           | 0.223%        |
| 21        | 14.492      | 2104          | 2109        | 2112         | rBV      | 183405         | 202357        | 4.04%           | 0.413%        |
| 22        | 14.798      | 2157          | 2161        | 2166         | rBV      | 235141         | 262521        | 5.25%           | 0.536%        |
| 23        | 15.139      | 2215          | 2219        | 2225         | rBV      | 284814         | 304309        | 6.08%           | 0.621%        |
| 24        | 15.509      | 2277          | 2282        | 2291         | rBV      | 1490987        | 2086083       | 41.70%          | 4.260%        |
| 25        | 15.951      | 2352          | 2357        | 2364         | rVB      | 215031         | 324415        | 6.48%           | 0.662%        |
| 26        | 16.456      | 2438          | 2443        | 2454         | rVB2     | 121563         | 221802        | 4.43%           | 0.453%        |
| 27        | 17.045      | 2540          | 2543        | 2549         | rVB      | 41179          | 70728         | 1.41%           | 0.144%        |

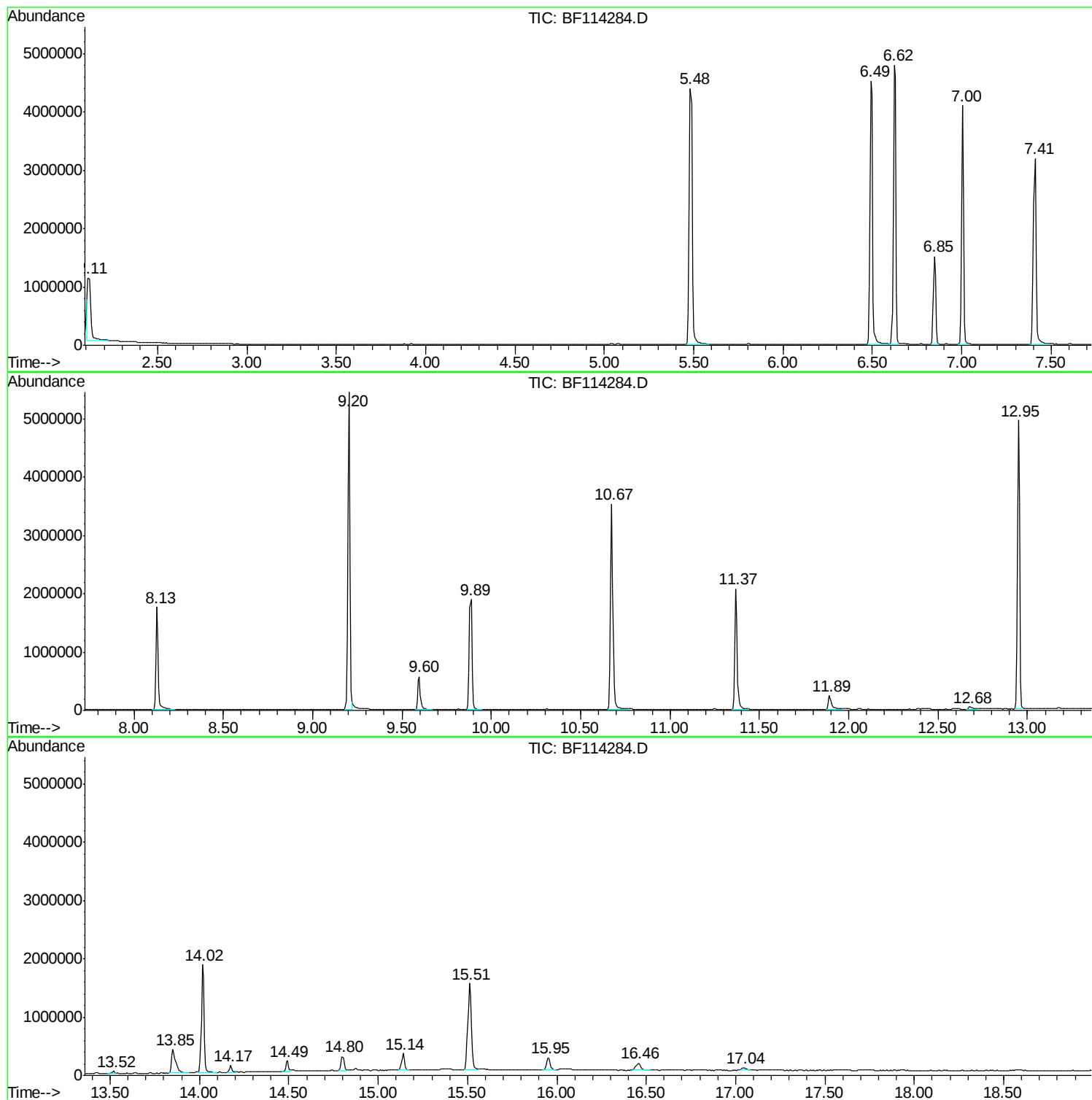
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF051019.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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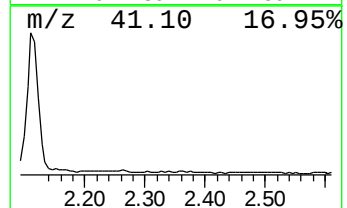
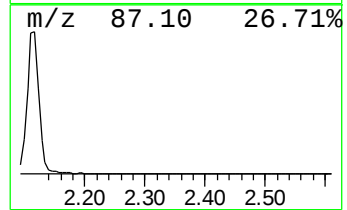
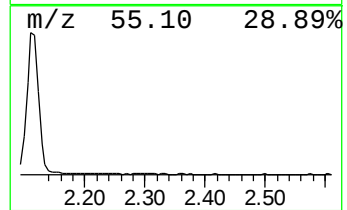
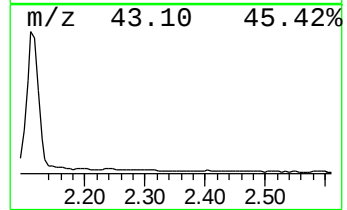
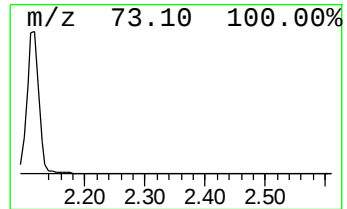
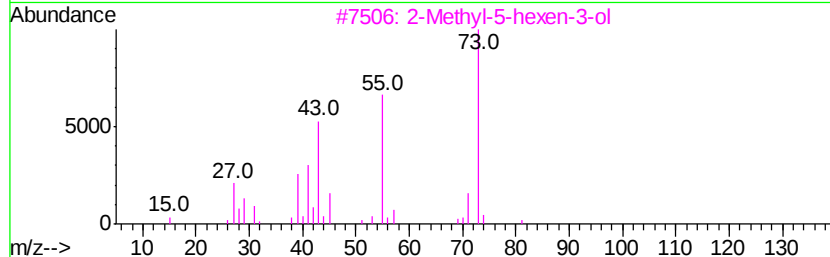
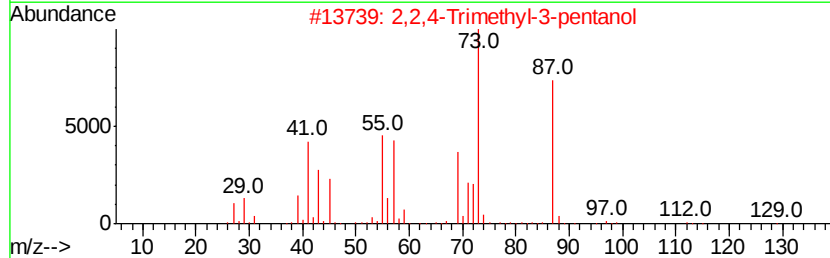
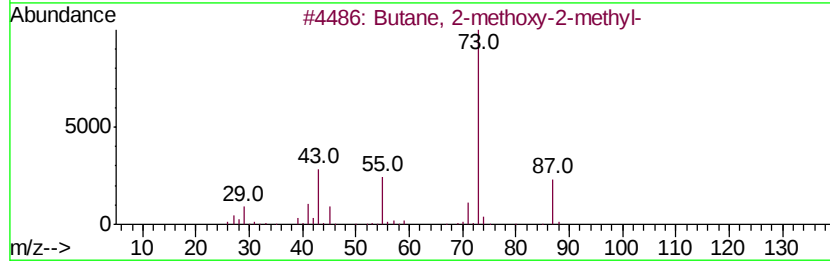
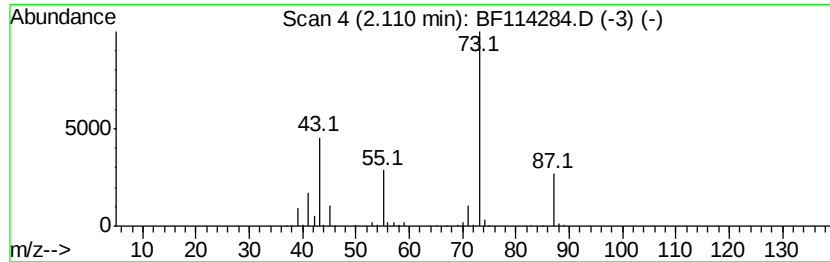
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF051019.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 2.11 | 19.05 ng | 1223100 | 1,4-Dichlorobenzene-d4 | 6.85 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | Butane, 2-methoxy-2-methyl- | 102 | C6H14O  | 000994-05-8 | 83   |
| 2    |    | 2,2,4-Trimethyl-3-pentanol  | 130 | C8H18O  | 005162-48-1 | 39   |
| 3    |    | 2-Methyl-5-hexen-3-ol       | 114 | C7H14O  | 032815-70-6 | 38   |
| 4    |    | 3-Pentanol, 2,4-dimethyl-   | 116 | C7H16O  | 000600-36-2 | 37   |
| 5    |    | 1,3-Dioxolane, 2-methyl-    | 88  | C4H8O2  | 000497-26-7 | 25   |



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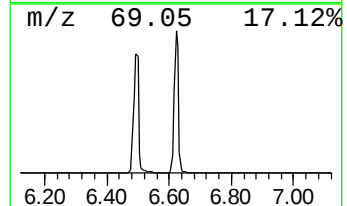
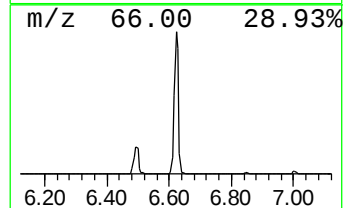
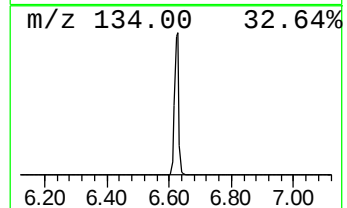
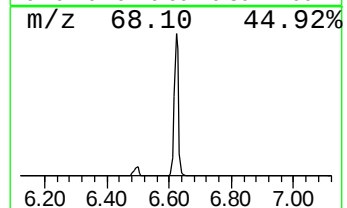
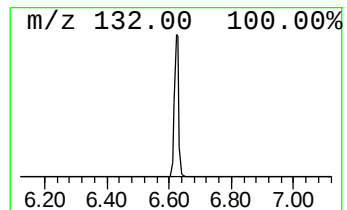
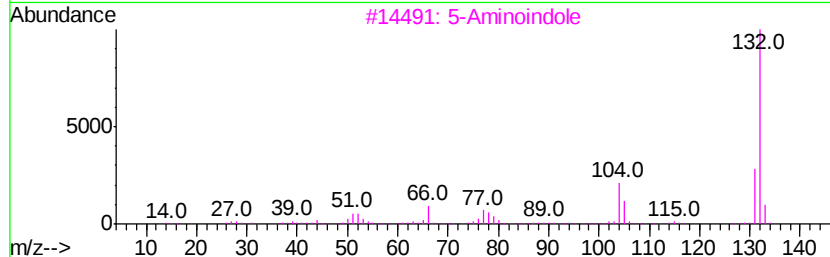
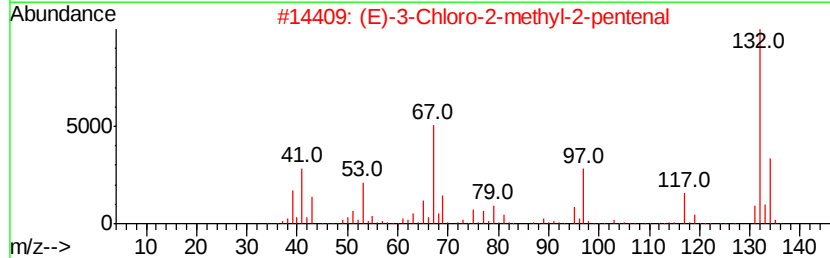
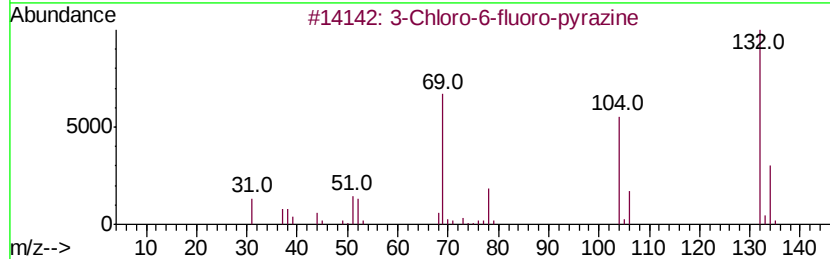
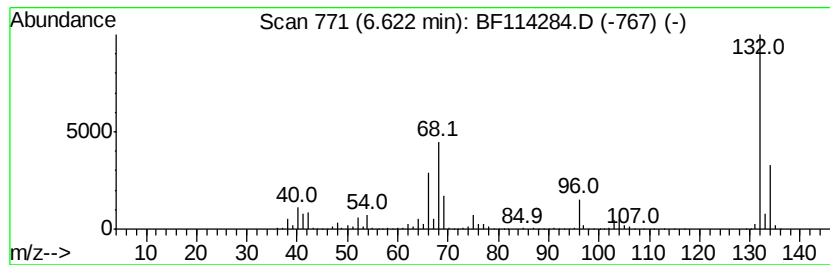
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF051019.M  
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TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 unknown6.62 Concentration Rank 1

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.62 | 74.03 ng | 4753360 | 1,4-Dichlorobenzene-d4 | 6.85 |

| Hit# | of | Tentative ID                     | MW  | MolForm   | CAS#         | Qual |
|------|----|----------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine       | 132 | C4H2ClFN2 | 1000146-10-7 | 27   |
| 2    |    | (E)-3-Chloro-2-methyl-2-pentenal | 132 | C6H9ClO   | 031357-76-3  | 16   |
| 3    |    | 5-Aminoindole                    | 132 | C8H8N2    | 005192-03-0  | 12   |
| 4    |    | Tranvlcvpromine-propionyl        | 189 | C12H15NO  | 1000123-86-3 | 10   |
| 5    |    | 1,2,4-Trifluorobenzene           | 132 | C6H3F3    | 000367-23-7  | 9    |



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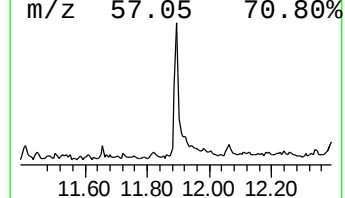
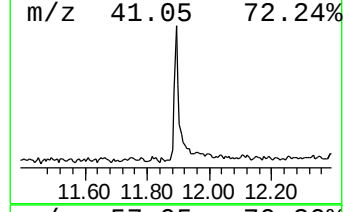
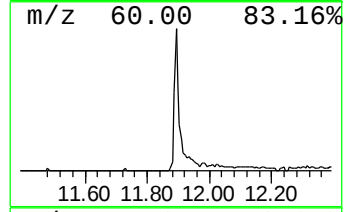
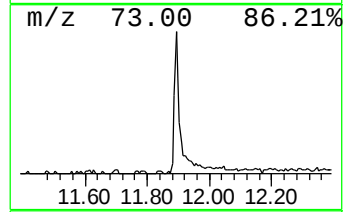
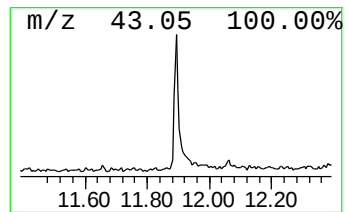
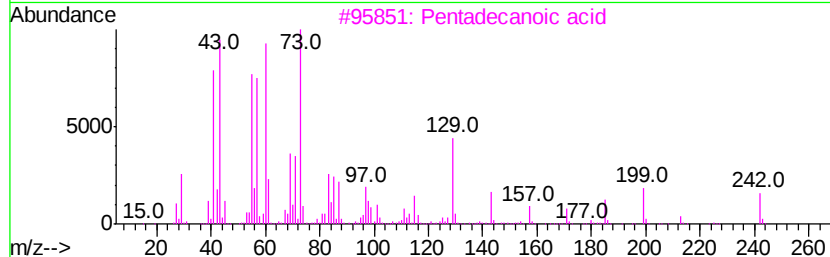
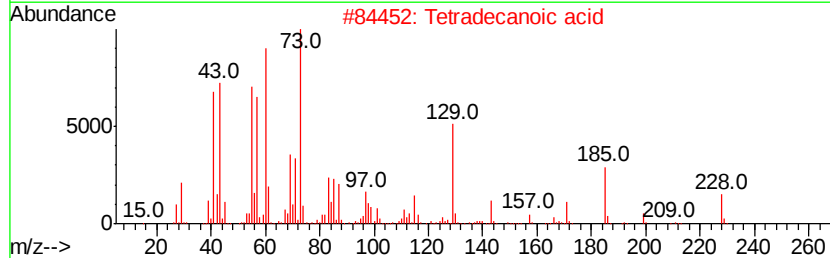
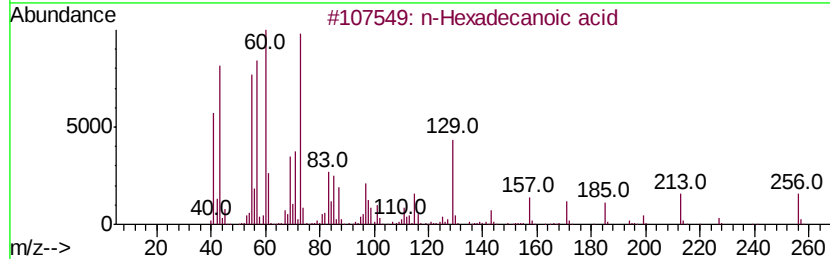
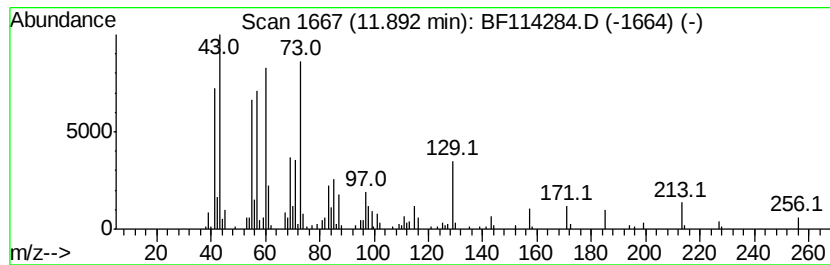
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 n-Hexadecanoic acid Concentration Rank 7

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.89 | 2.73 ng | 274991 | Phenanthrene-d10 | 11.37 |

| Hit# of | 5 | Tentative ID        | MW  | MolForm  | CAS#        | Qual |
|---------|---|---------------------|-----|----------|-------------|------|
| 1       |   | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 99   |
| 2       |   | Tetradecanoic acid  | 228 | C14H28O2 | 000544-63-8 | 95   |
| 3       |   | Pentadecanoic acid  | 242 | C15H30O2 | 001002-84-2 | 90   |
| 4       |   | Tridecanoic acid    | 214 | C13H26O2 | 000638-53-9 | 72   |
| 5       |   | n-Decanoic acid     | 172 | C10H20O2 | 000334-48-5 | 53   |



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SOIL-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF051019.M  
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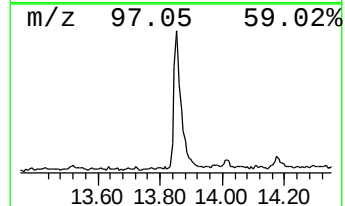
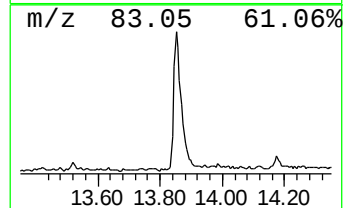
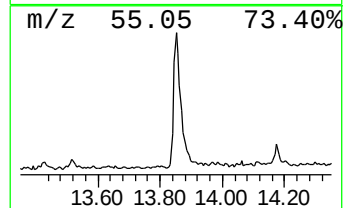
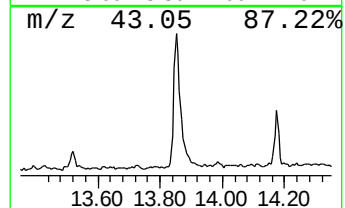
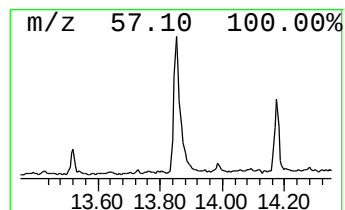
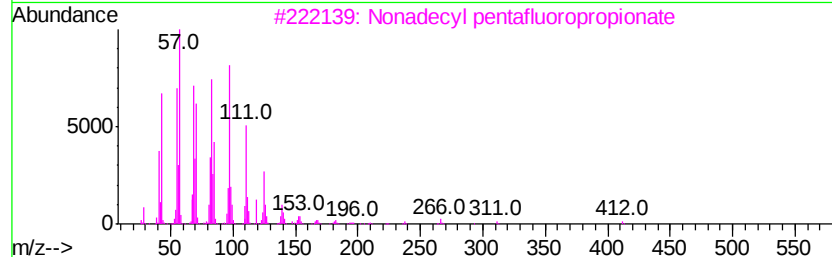
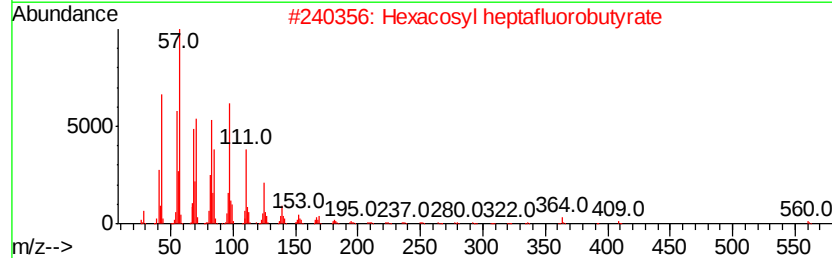
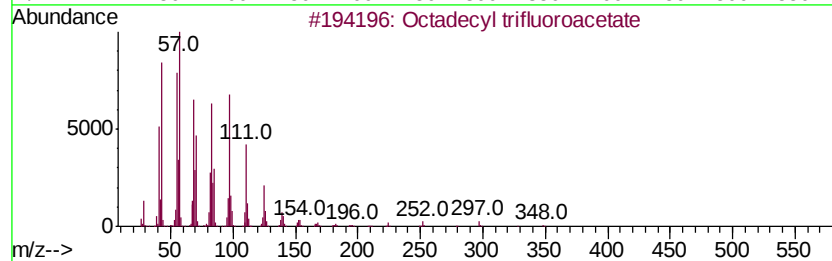
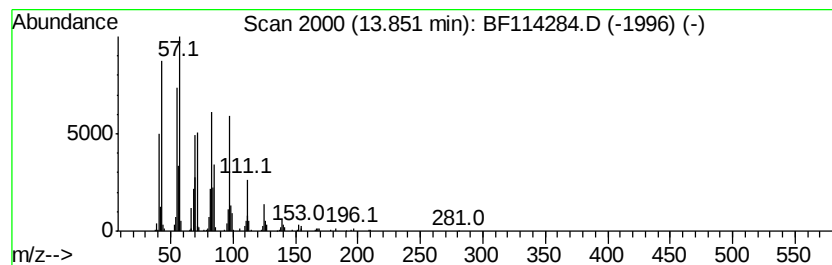
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Peak Number 5 Octadecyl trifluoroacetate Concentration Rank 4

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.85 | 7.00 ng | 659810 | Chrysene-d12     | 14.02 |

| Hit# of | 5 | Tentative ID                    | MW  | MolForm    | CAS#         | Qual |
|---------|---|---------------------------------|-----|------------|--------------|------|
| 1       |   | Octadecyl trifluoroacetate      | 366 | C20H37F3O2 | 079392-43-1  | 91   |
| 2       |   | Hexacosyl heptafluorobutyrate   | 578 | C30H53F7O2 | 1000351-83-3 | 91   |
| 3       |   | Nonadecyl pentafluoropropionate | 430 | C22H39F5O2 | 1000351-88-8 | 91   |
| 4       |   | Heneicosyl heptafluorobutyrate  | 508 | C25H43F7O2 | 1000351-83-8 | 90   |
| 5       |   | Octacosyl trifluoroacetate      | 506 | C30H57F3O2 | 1000351-74-9 | 87   |



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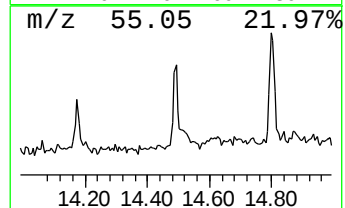
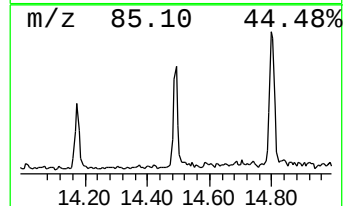
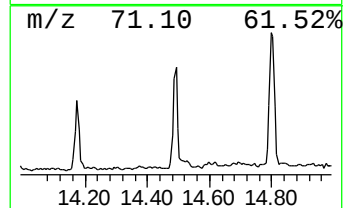
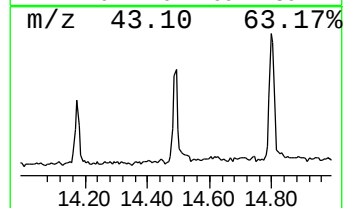
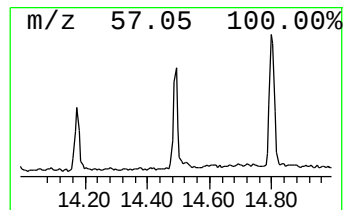
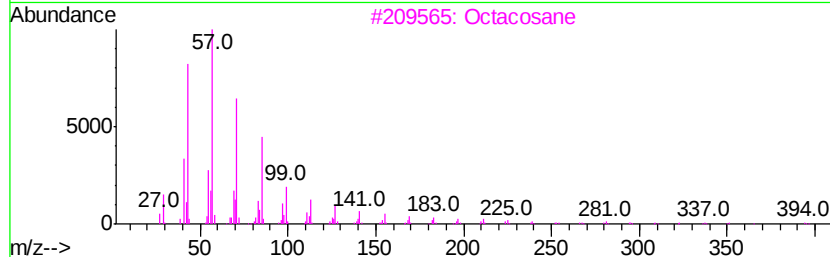
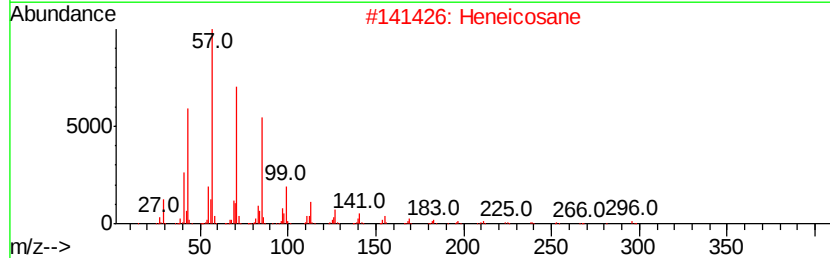
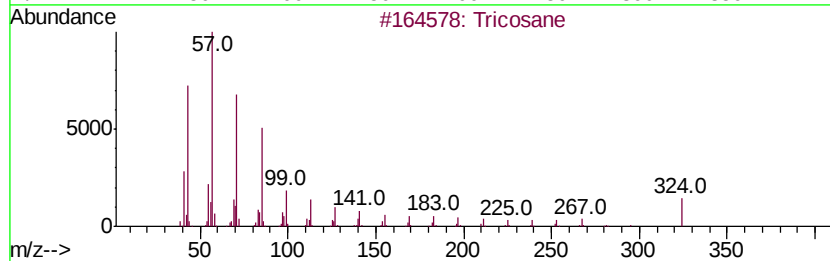
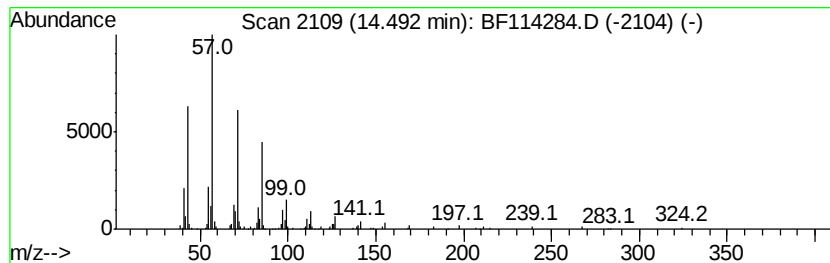
Instrument :  
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ClientSampleId :  
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 6 Tricosane Concentration Rank 9

| R.T.    | EstConc     | Area         | Relative to ISTD | R.T.      |
|---------|-------------|--------------|------------------|-----------|
| 14.49   | 2.15 ng     | 202357       | Chrysene-d12     | 14.02     |
| Hit# of | 5           | Tentative ID | MW MolForm       | CAS# Qual |
| 1       | Tricosane   | 324 C23H48   | 000638-67-5      | 93        |
| 2       | Heneicosane | 296 C21H44   | 000629-94-7      | 91        |
| 3       | Octacosane  | 394 C28H58   | 000630-02-4      | 91        |
| 4       | Heptacosane | 380 C27H56   | 000593-49-7      | 91        |
| 5       | Tetracosane | 338 C24H50   | 000646-31-1      | 91        |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF051519\  
Data File : BF114284.D  
Acq On : 15 May 2019 14:48  
Operator : JU/SJ  
Sample : K2837-01  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SOIL-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF051019.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

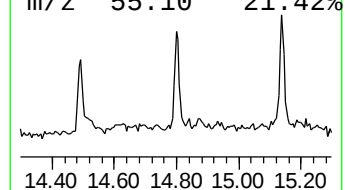
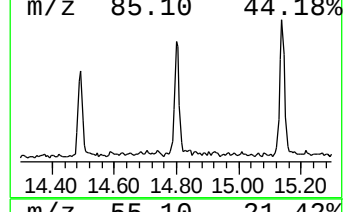
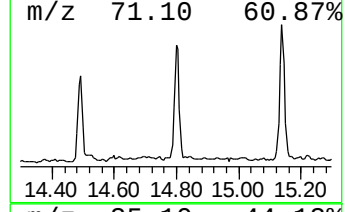
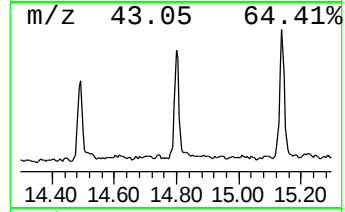
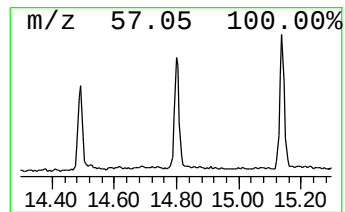
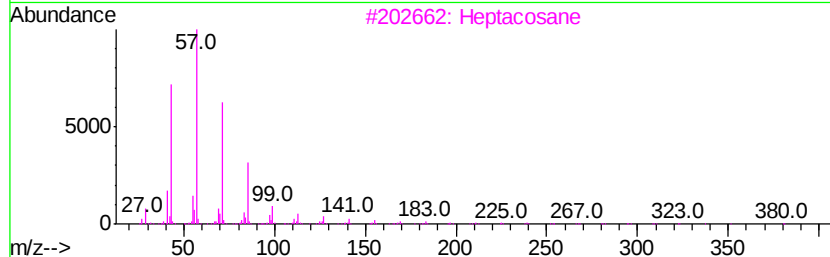
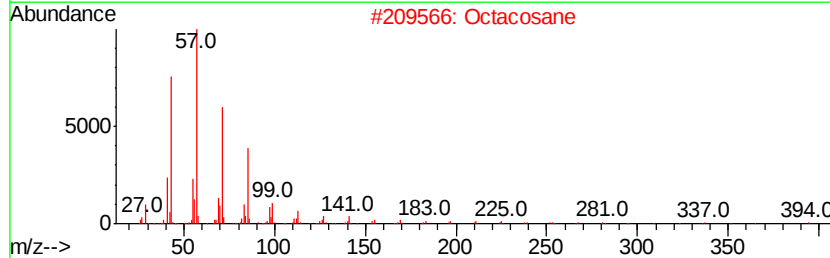
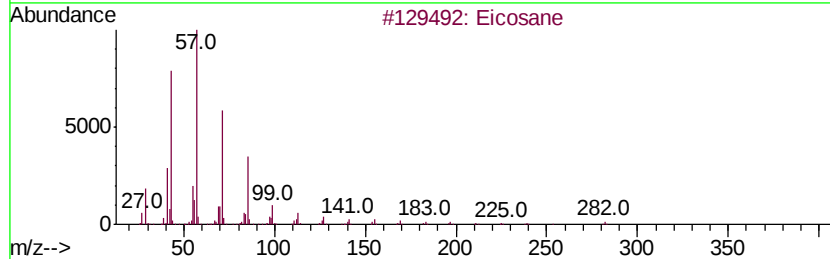
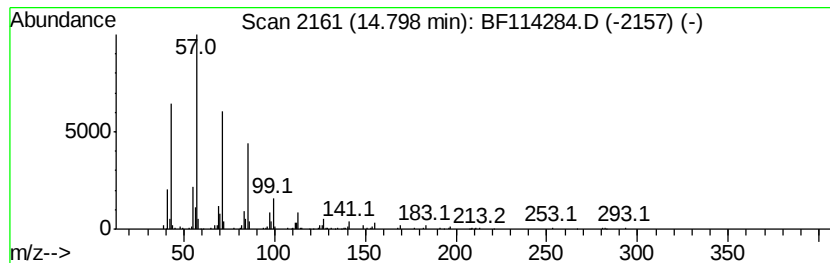
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 Eicosane Concentration Rank 8

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 14.80 | 2.52 ng | 262521 | Perylene-d12     | 15.51 |

| Hit# | of | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|--------------|-----|---------|-------------|------|
| 1    | 5  | Eicosane     | 282 | C20H42  | 000112-95-8 | 95   |
| 2    |    | Octacosane   | 394 | C28H58  | 000630-02-4 | 91   |
| 3    |    | Heptacosane  | 380 | C27H56  | 000593-49-7 | 91   |
| 4    |    | Octadecane   | 254 | C18H38  | 000593-45-3 | 91   |
| 5    |    | Pentadecane  | 212 | C15H32  | 000629-62-9 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF051519\  
Data File : BF114284.D  
Acq On : 15 May 2019 14:48  
Operator : JU/SJ  
Sample : K2837-01  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SOIL-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF051019.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

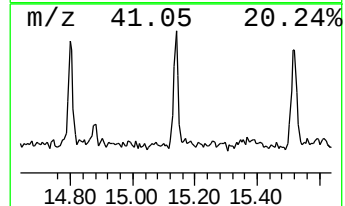
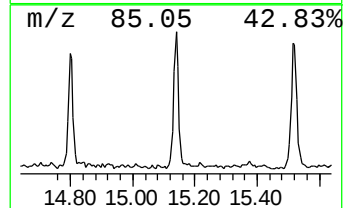
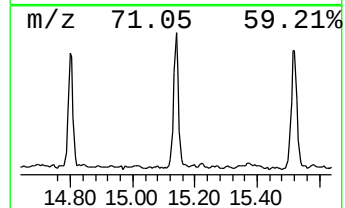
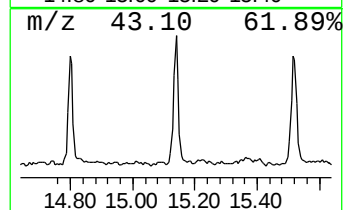
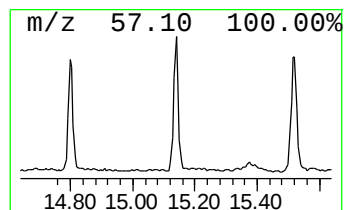
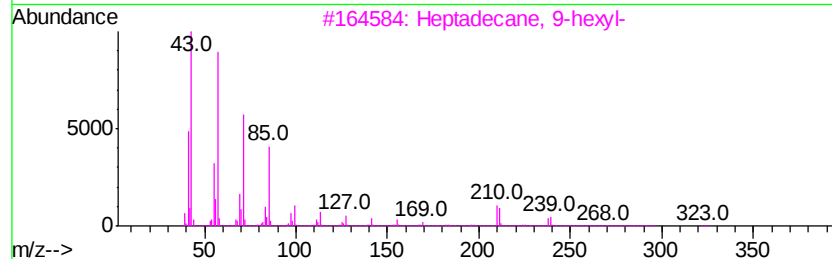
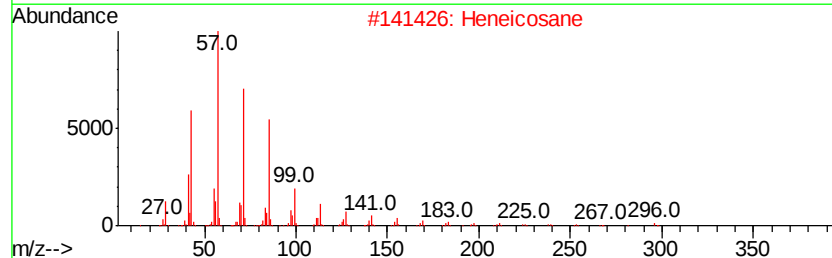
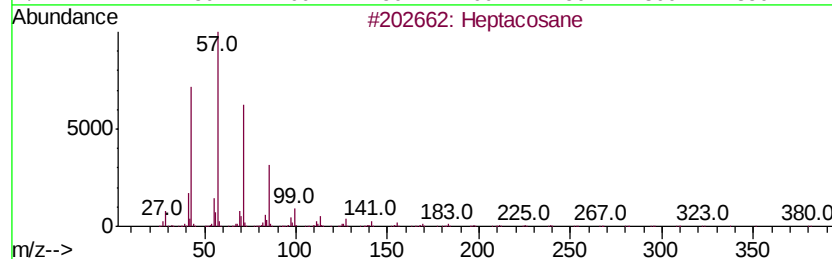
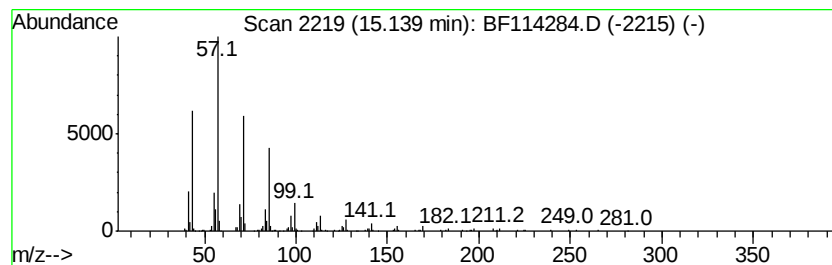
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 Heptacosane Concentration Rank 6

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 15.14 | 2.92 ng | 304309 | Perylene-d12     | 15.51 |

| Hit# | of | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------------|-----|---------|-------------|------|
| 1    | 5  | Heptacosane                      | 380 | C27H56  | 000593-49-7 | 91   |
| 2    |    | Heneicosane                      | 296 | C21H44  | 000629-94-7 | 91   |
| 3    |    | Heptadecane, 9-hexyl-            | 324 | C23H48  | 055124-79-3 | 90   |
| 4    |    | Heneicosane, 11-(1-ethylpropyl)- | 366 | C26H54  | 055282-11-6 | 90   |
| 5    |    | Octadecane, 2-methyl-            | 268 | C19H40  | 001560-88-9 | 90   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF051519\  
Data File : BF114284.D  
Acq On : 15 May 2019 14:48  
Operator : JU/SJ  
Sample : K2837-01  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SOIL-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF051019.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

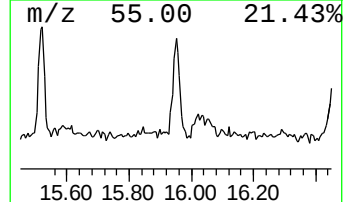
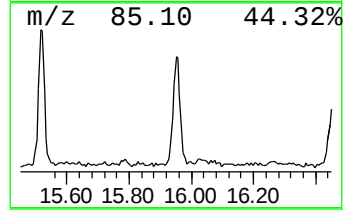
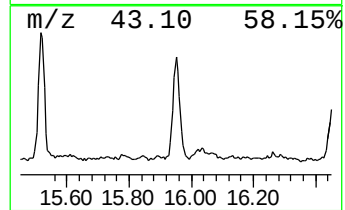
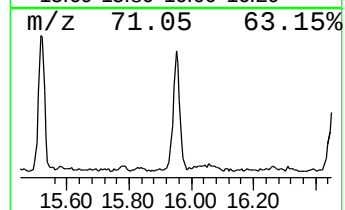
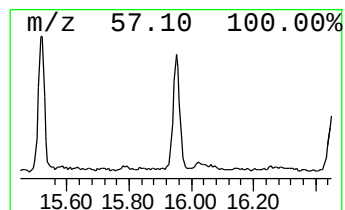
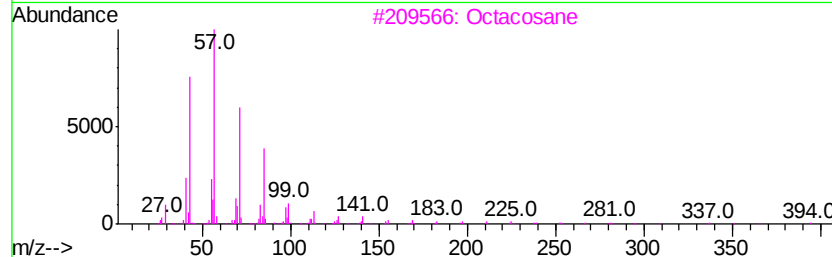
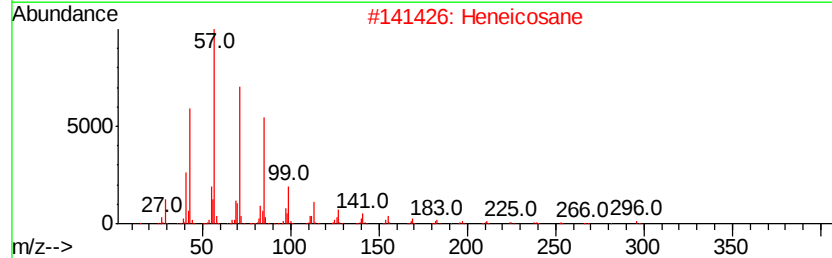
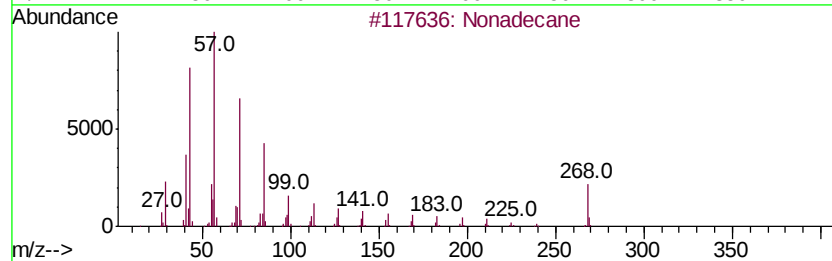
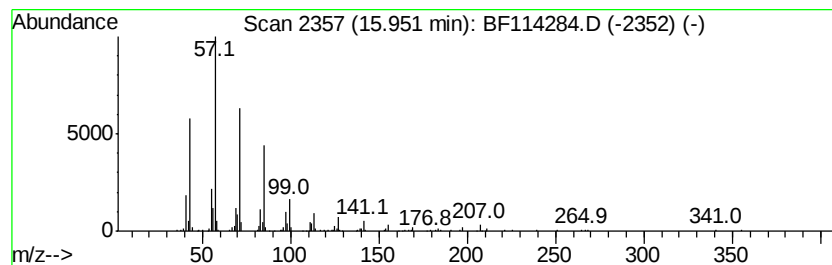
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 Nonadecane Concentration Rank 5

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 15.95 | 3.11 ng | 324415 | Perylene-d12     | 15.51 |

| Hit# | of | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|--------------|-----|---------|-------------|------|
| 1    | 5  | Nonadecane   | 268 | C19H40  | 000629-92-5 | 93   |
| 2    |    | Heneicosane  | 296 | C21H44  | 000629-94-7 | 91   |
| 3    |    | Octacosane   | 394 | C28H58  | 000630-02-4 | 91   |
| 4    |    | Docosane     | 310 | C22H46  | 000629-97-0 | 91   |
| 5    |    | Hexacosane   | 366 | C26H54  | 000630-01-3 | 90   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF051519\  
Data File : BF114284.D  
Acq On : 15 May 2019 14:48  
Operator : JU/SJ  
Sample : K2837-01  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SOIL-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF051019.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

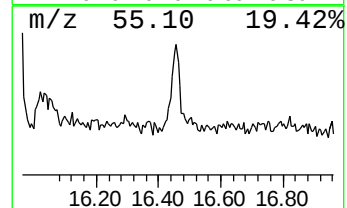
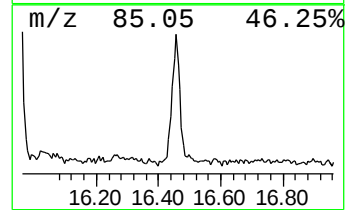
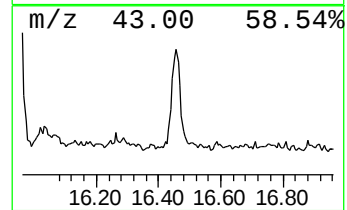
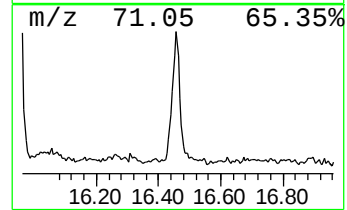
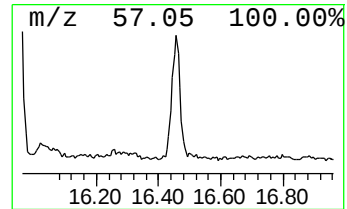
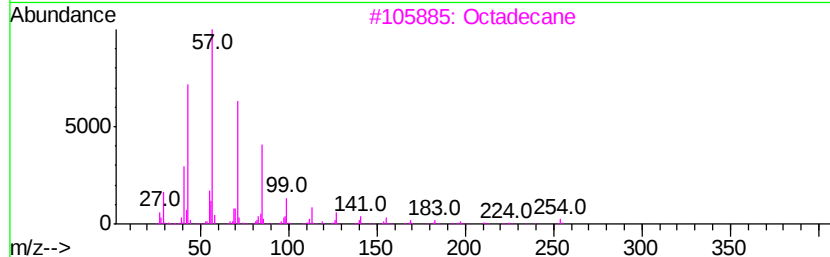
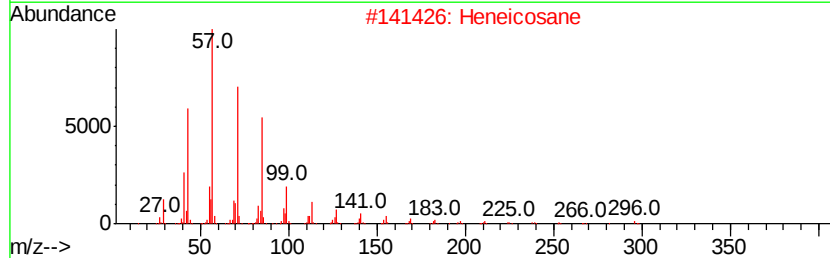
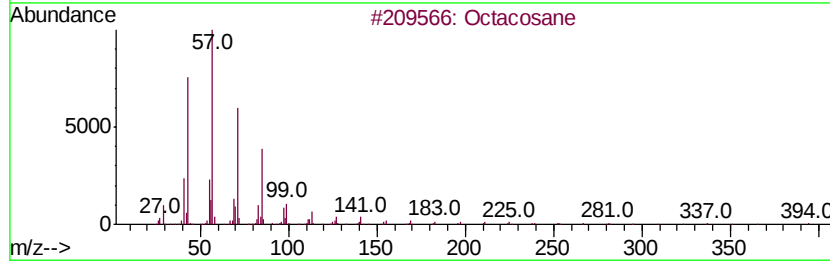
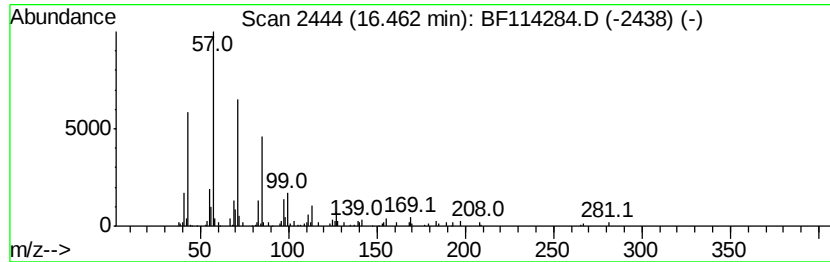
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 Octacosane Concentration Rank 10

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 16.46 | 2.13 ng | 221802 | Perylene-d12     | 15.51 |

| Hit# | of | 5 | Tentative ID        | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------|-----|---------|-------------|------|
| 1    |    |   | Octacosane          | 394 | C28H58  | 000630-02-4 | 90   |
| 2    |    |   | Heneicosane         | 296 | C21H44  | 000629-94-7 | 87   |
| 3    |    |   | Octadecane          | 254 | C18H38  | 000593-45-3 | 87   |
| 4    |    |   | Octadecane, 1-iodo- | 380 | C18H37I | 000629-93-6 | 87   |
| 5    |    |   | Heptadecane         | 240 | C17H36  | 000629-78-7 | 86   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF051519\  
Data File : BF114284.D  
Acq On : 15 May 2019 14:48  
Operator : JU/SJ  
Sample : K2837-01  
Misc :  
ALS Vial : 8 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SOIL-COMP

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF051019.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 |
| Butane, 2-methoxy... | 2.11  | 19.1    | ng    | 1223100  | 1                     | 6.85  | 1284240 | 20.0 |
| unknown6.62          | 6.62  | 74.0    | ng    | 4753360  | 1                     | 6.85  | 1284240 | 20.0 |
| n-Hexadecanoic acid  | 11.89 | 2.7     | ng    | 274991   | 4                     | 11.37 | 2011910 | 20.0 |
| Octadecyl trifluo... | 13.85 | 7.0     | ng    | 659810   | 5                     | 14.02 | 1884060 | 20.0 |
| Tricosane            | 14.49 | 2.1     | ng    | 202357   | 5                     | 14.02 | 1884060 | 20.0 |
| Eicosane             | 14.80 | 2.5     | ng    | 262521   | 6                     | 15.51 | 2086080 | 20.0 |
| Heptacosane          | 15.14 | 2.9     | ng    | 304309   | 6                     | 15.51 | 2086080 | 20.0 |
| Nonadecane           | 15.95 | 3.1     | ng    | 324415   | 6                     | 15.51 | 2086080 | 20.0 |
| Octacosane           | 16.46 | 2.1     | ng    | 221802   | 6                     | 15.51 | 2086080 | 20.0 |