

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082019\  
 Data File : BM022207.D  
 Acq On : 20 Aug 2019 11:55  
 Operator : HP/JU  
 Sample : K4332-01  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 0617-FIELD-BLANK-38TH-ST

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\_M\METHODS\8270-BM082019.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         | 5.181       | 368           | 373         | 384          | rBV      | 355456         | 506570        | 11.70%          | 2.450%        |
| 2         | 6.758       | 635           | 641         | 657          | rBV      | 207089         | 349380        | 8.07%           | 1.690%        |
| 3         | 7.105       | 693           | 700         | 712          | rBV      | 766685         | 1176990       | 27.18%          | 5.693%        |
| 4         | 7.569       | 773           | 779         | 787          | rBB      | 153174         | 227690        | 5.26%           | 1.101%        |
| 5         | 7.881       | 825           | 832         | 842          | rBB      | 763257         | 1161784       | 26.83%          | 5.619%        |
| 6         | 8.740       | 971           | 978         | 999          | rBV      | 527158         | 908266        | 20.97%          | 4.393%        |
| 7         | 10.346      | 1245          | 1251        | 1268         | rBB      | 158963         | 286037        | 6.60%           | 1.383%        |
| 8         | 12.834      | 1667          | 1674        | 1697         | rBB      | 1662221        | 2414579       | 55.75%          | 11.679%       |
| 9         | 14.228      | 1905          | 1911        | 1922         | rBV      | 271172         | 400508        | 9.25%           | 1.937%        |
| 10        | 15.733      | 2161          | 2167        | 2181         | rBV      | 1638974        | 2161773       | 49.91%          | 10.456%       |
| 11        | 16.992      | 2375          | 2381        | 2392         | rBV      | 321048         | 460936        | 10.64%          | 2.229%        |
| 12        | 17.880      | 2526          | 2532        | 2544         | rBV      | 64506          | 105002        | 2.42%           | 0.508%        |
| 13        | 19.339      | 2775          | 2780        | 2786         | rBV      | 166656         | 192999        | 4.46%           | 0.933%        |
| 14        | 19.633      | 2824          | 2830        | 2836         | rBV      | 3765110        | 4157435       | 95.99%          | 20.108%       |
| 15        | 20.339      | 2940          | 2950        | 2963         | rBV      | 2901720        | 4330962       | 100.00%         | 20.948%       |
| 16        | 20.433      | 2963          | 2966        | 2976         | rVB      | 145394         | 187262        | 4.32%           | 0.906%        |
| 17        | 20.904      | 3042          | 3046        | 3052         | rBV5     | 40620          | 66930         | 1.55%           | 0.324%        |
| 18        | 21.198      | 3091          | 3096        | 3102         | rBV      | 405899         | 507987        | 11.73%          | 2.457%        |
| 19        | 21.298      | 3109          | 3113        | 3116         | rBV      | 38609          | 52766         | 1.22%           | 0.255%        |
| 20        | 22.221      | 3265          | 3270        | 3280         | rVB2     | 264994         | 355883        | 8.22%           | 1.721%        |
| 21        | 22.327      | 3284          | 3288        | 3293         | rVB      | 162416         | 198402        | 4.58%           | 0.960%        |
| 22        | 23.398      | 3464          | 3470        | 3479         | rBV      | 244853         | 465121        | 10.74%          | 2.250%        |

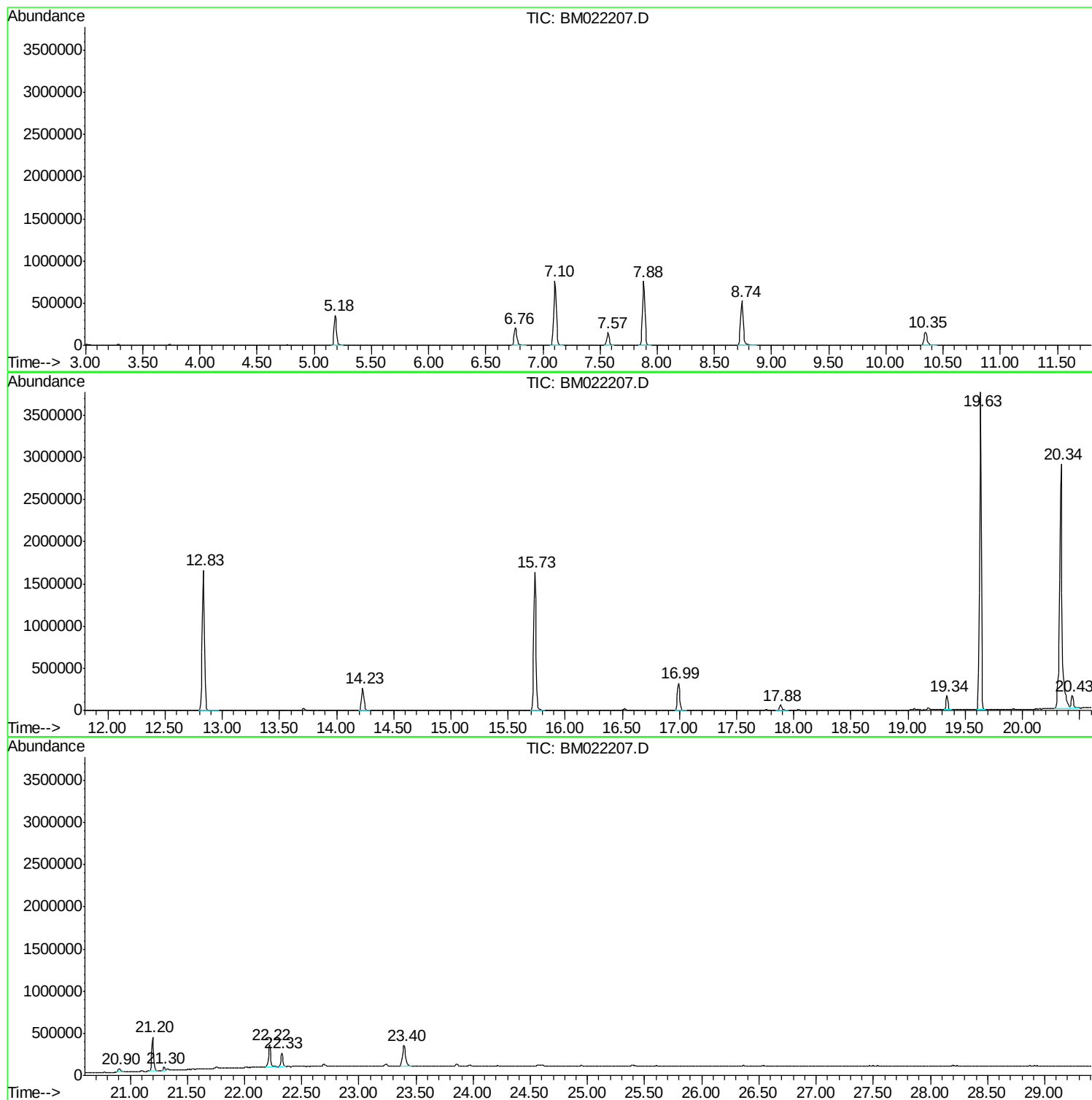
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Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082019\  
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Acq On : 20 Aug 2019 11:55  
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Sample : K4332-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
0617-FIELD-BLANK-38TH-ST

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM082019.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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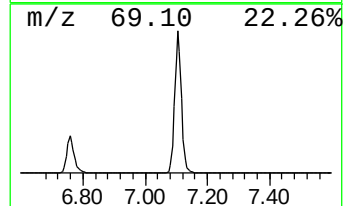
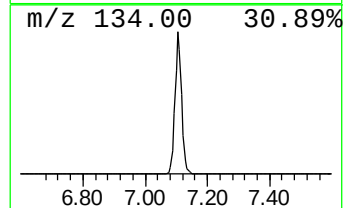
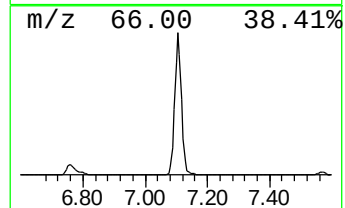
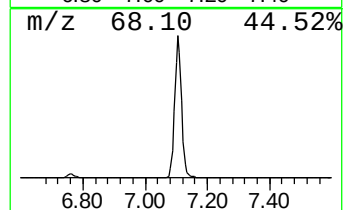
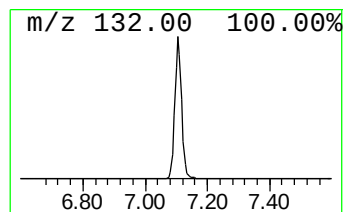
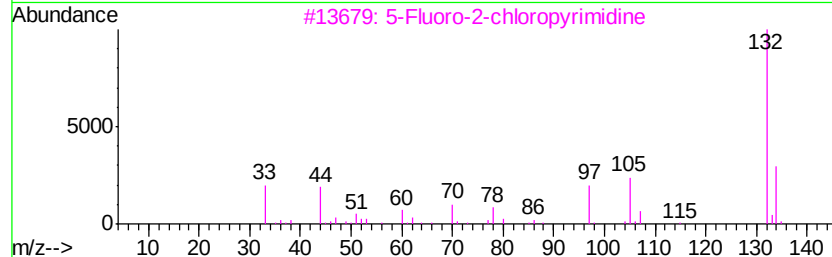
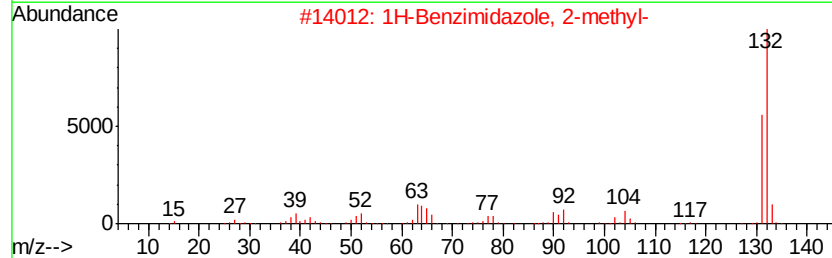
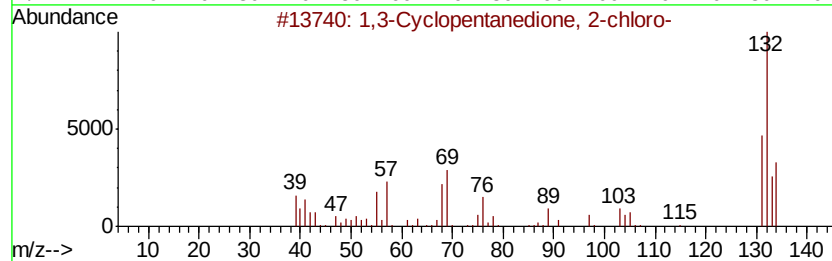
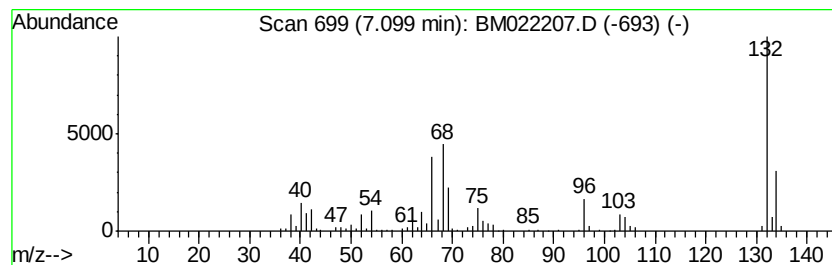
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 unknown7.10 Concentration Rank 2

| R.T. | EstConc   | Area    | Relative to ISTD       | R.T. |
|------|-----------|---------|------------------------|------|
| 7.10 | 103.39 ng | 1176990 | 1,4-Dichlorobenzene-d4 | 7.57 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|-------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | 1,3-Cyclopentanedione, 2-chloro-    | 132 | C5H5ClO2  | 014203-19-1 | 38   |
| 2    |    | 1H-Benzimidazole, 2-methyl-         | 132 | C8H8N2    | 000615-15-6 | 27   |
| 3    |    | 5-Fluoro-2-chloropyrimidine         | 132 | C4H2ClFN2 | 062802-42-0 | 27   |
| 4    |    | 1,3,4-Thiadiazole-2(3H)-thione, ... | 132 | C3H4N2S2  | 029490-19-5 | 27   |
| 5    |    | Benzene, 1,2,4-trifluoro-           | 132 | C6H3F3    | 000367-23-7 | 22   |



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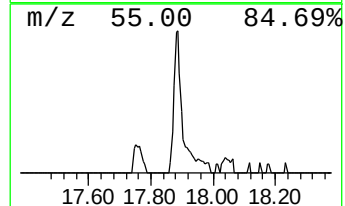
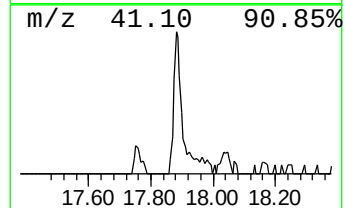
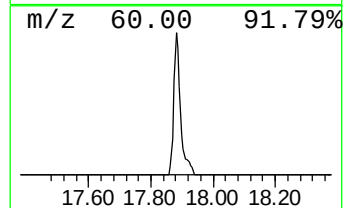
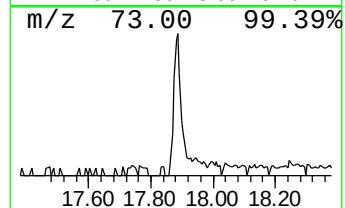
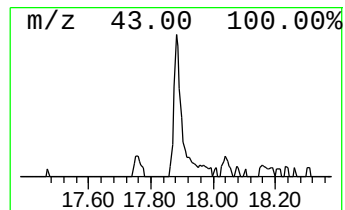
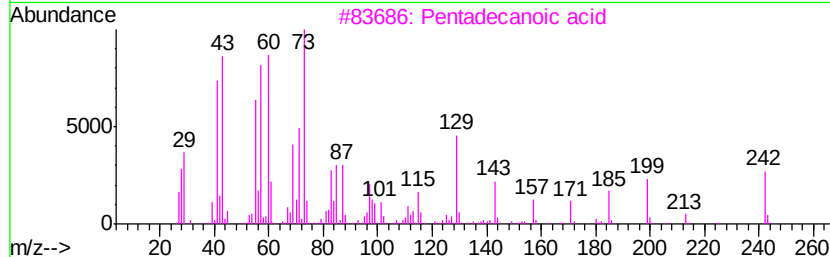
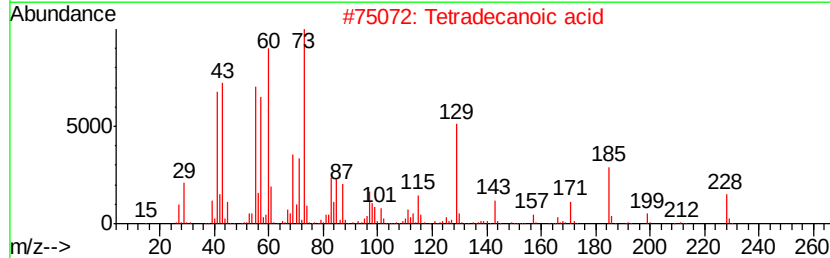
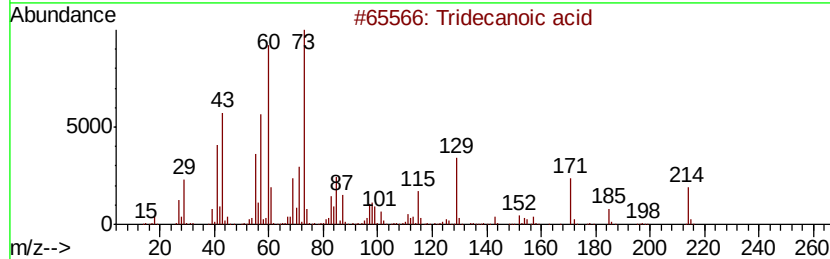
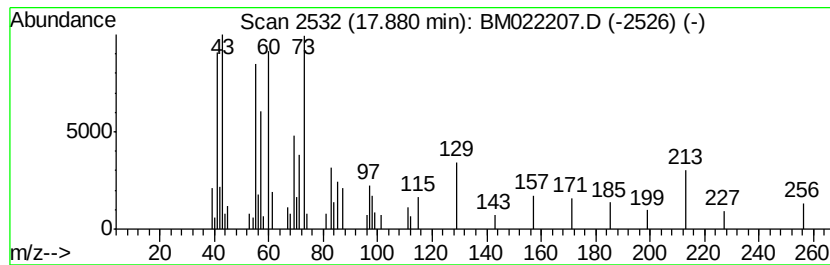
Instrument :  
BNA\_M  
ClientSampleId :  
0617-FIELD-BLANK-38TH-ST

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM082019.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 3 Tridecanoic acid Concentration Rank 8

| R.T.      | EstConc             | Area   | Relative to ISTD | R.T.        |      |
|-----------|---------------------|--------|------------------|-------------|------|
| 17.88     | 4.56 ng             | 105002 | Phenanthrene-d10 | 16.99       |      |
| Hit# of 5 | Tentative ID        | MW     | MolForm          | CAS#        | Qual |
| 1         | Tridecanoic acid    | 214    | C13H26O2         | 000638-53-9 | 72   |
| 2         | Tetradecanoic acid  | 228    | C14H28O2         | 000544-63-8 | 72   |
| 3         | Pentadecanoic acid  | 242    | C15H30O2         | 001002-84-2 | 64   |
| 4         | Undecanoic acid     | 186    | C11H22O2         | 000112-37-8 | 59   |
| 5         | n-Hexadecanoic acid | 256    | C16H32O2         | 000057-10-3 | 58   |



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ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM082019.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

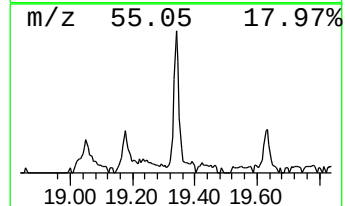
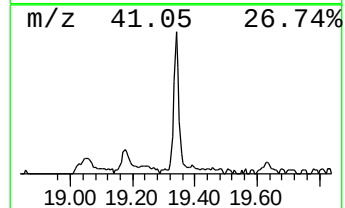
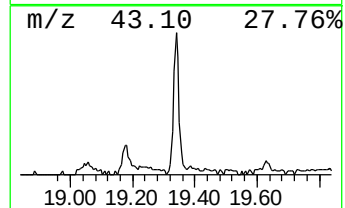
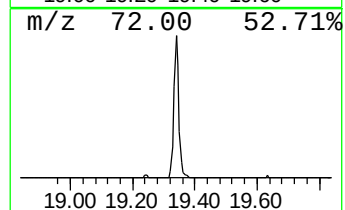
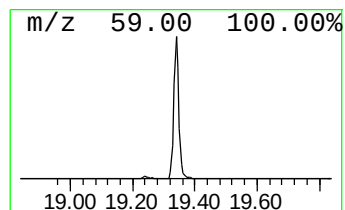
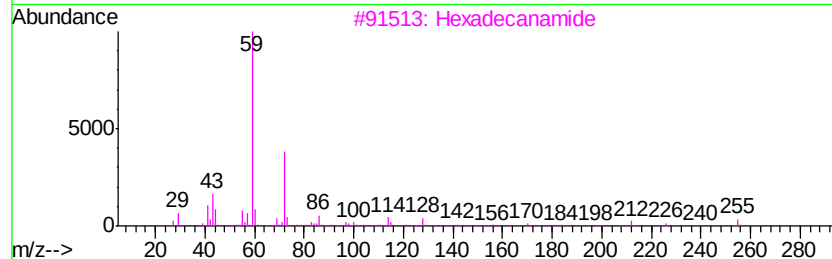
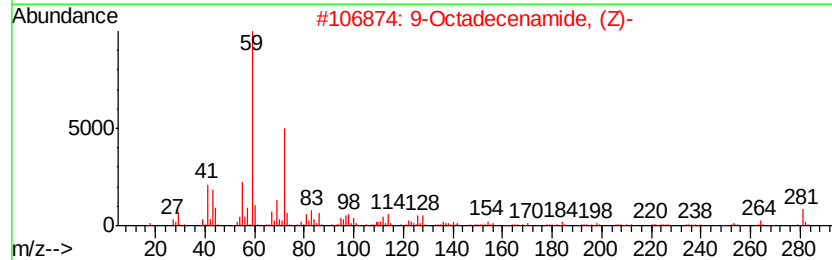
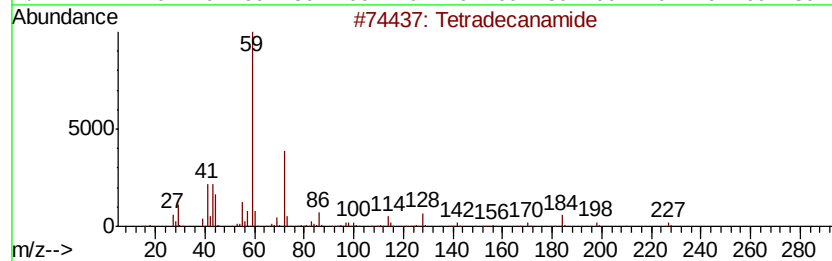
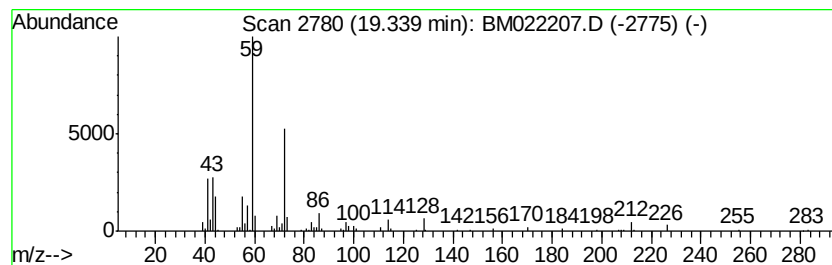
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 Hexadecanamide Concentration Rank 6

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 19.34 | 7.60 ng | 192999 | Chrysene-d12     | 21.20 |

| Hit# | of | Tentative ID               | MW  | MolForm    | CAS#         | Qual |
|------|----|----------------------------|-----|------------|--------------|------|
| 1    | 5  | Tetradecanamide            | 227 | C14H29NO   | 000638-58-4  | 86   |
| 2    |    | 9-Octadecenamide, (Z)-     | 281 | C18H35NO   | 000301-02-0  | 83   |
| 3    |    | Hexadecanamide             | 255 | C16H33NO   | 000629-54-9  | 72   |
| 4    |    | Octadecanamide             | 283 | C18H37NO   | 000124-26-5  | 64   |
| 5    |    | Pentadecanamide, 15-bromo- | 319 | C15H30BrNO | 1000163-86-1 | 64   |



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ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM082019.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

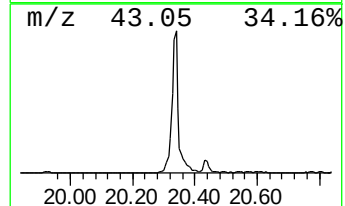
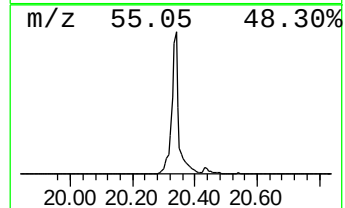
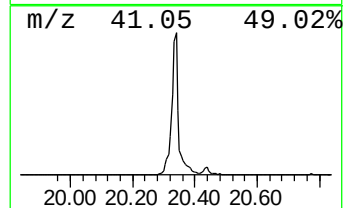
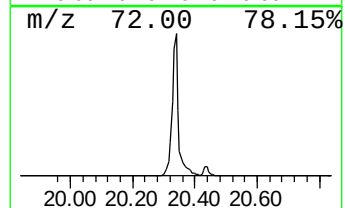
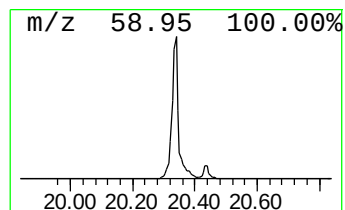
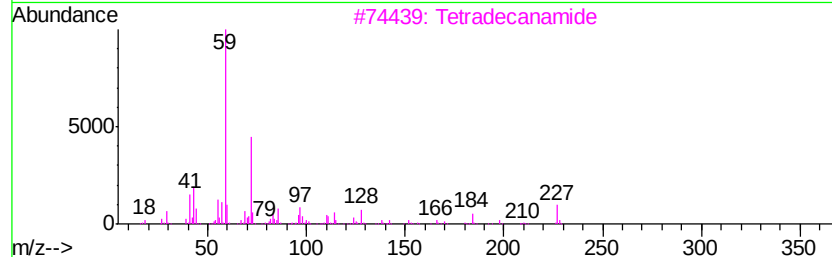
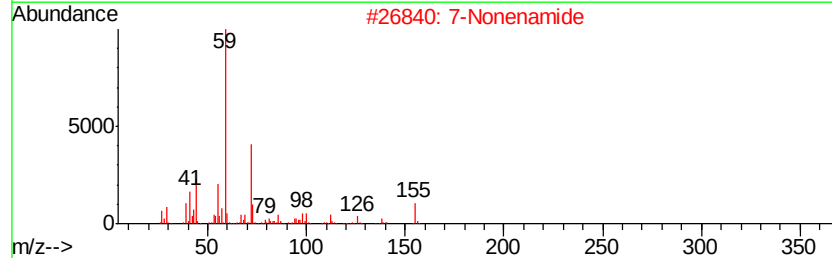
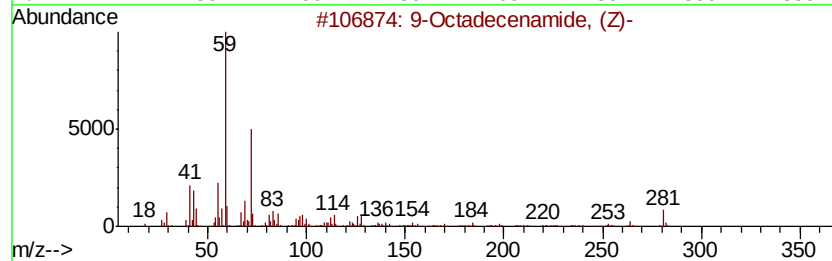
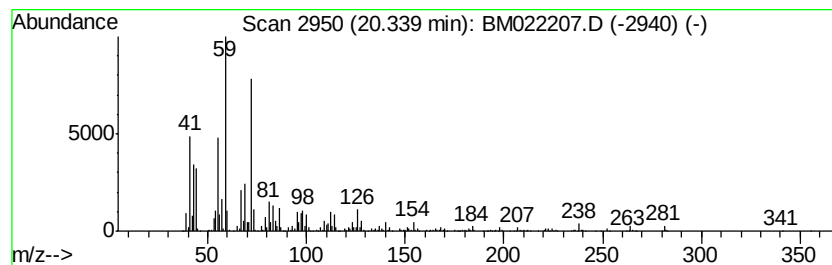
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 9-Octadecenamide, (Z)- Concentration Rank 1

| R.T.  | EstConc   | Area    | Relative to ISTD | R.T.  |
|-------|-----------|---------|------------------|-------|
| 20.34 | 170.52 ng | 4330960 | Chrysene-d12     | 21.20 |

| Hit# of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|---------|---|-------------------------------------|-----|----------|--------------|------|
| 1       |   | 9-Octadecenamide, (Z)-              | 281 | C18H35NO | 000301-02-0  | 94   |
| 2       |   | 7-Nonenamide                        | 155 | C9H17NO  | 090949-53-4  | 47   |
| 3       |   | Tetradecanamide                     | 227 | C14H29NO | 000638-58-4  | 43   |
| 4       |   | Octanamide                          | 143 | C8H17NO  | 000629-01-6  | 38   |
| 5       |   | Methyl 17-methoxy-10-methoxycarb... | 386 | C22H42O5 | 1000110-20-7 | 38   |



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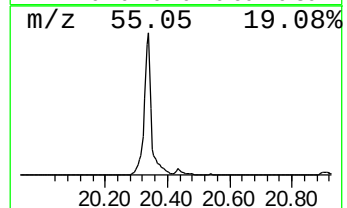
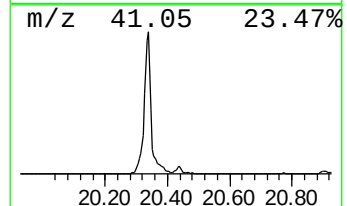
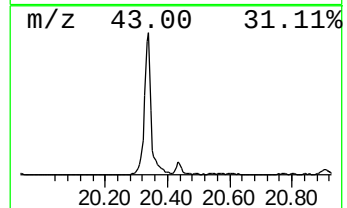
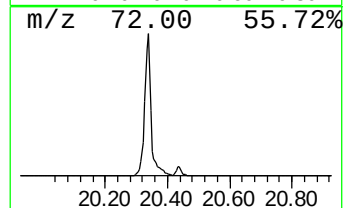
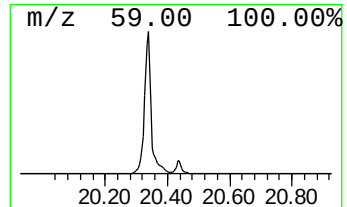
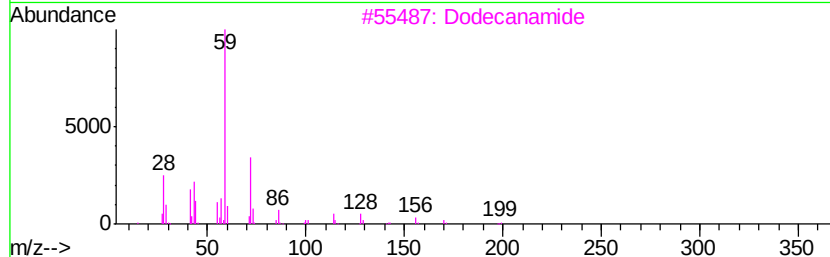
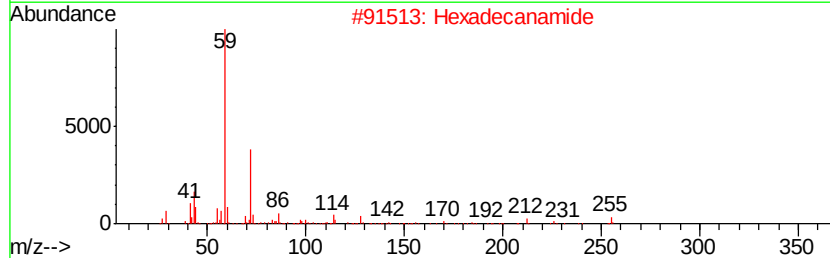
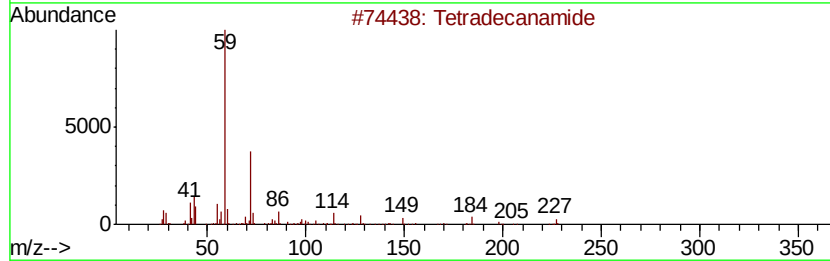
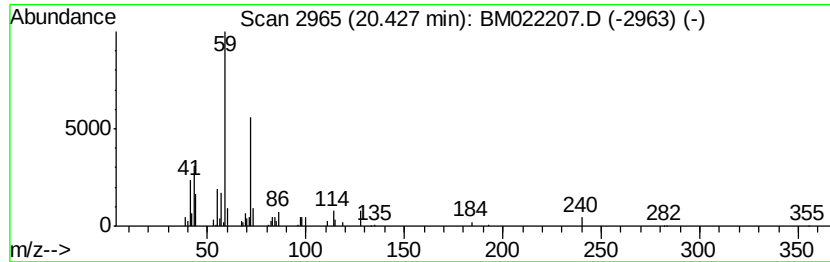
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 Tetradecanamide Concentration Rank 7

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 20.43 | 7.37 ng | 187262 | Chrysene-d12     | 21.20 |

| Hit# | of | Tentative ID           | MW  | MolForm  | CAS#        | Qual |
|------|----|------------------------|-----|----------|-------------|------|
| 1    | 5  | Tetradecanamide        | 227 | C14H29NO | 000638-58-4 | 80   |
| 2    |    | Hexadecanamide         | 255 | C16H33NO | 000629-54-9 | 72   |
| 3    |    | Dodecanamide           | 199 | C12H25NO | 001120-16-7 | 64   |
| 4    |    | Pentanamide, 4-methyl- | 115 | C6H13NO  | 001119-29-5 | 64   |
| 5    |    | 9-Octadecenamide, (Z)- | 281 | C18H35NO | 000301-02-0 | 59   |



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Operator : HP/JU  
Sample : K4332-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

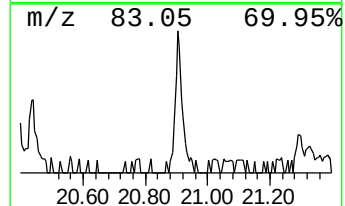
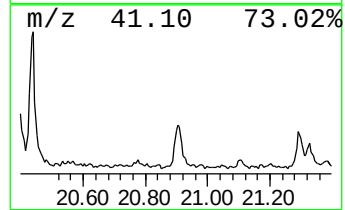
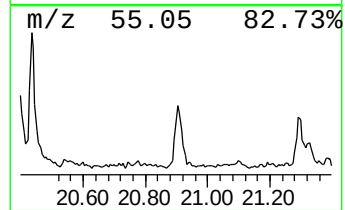
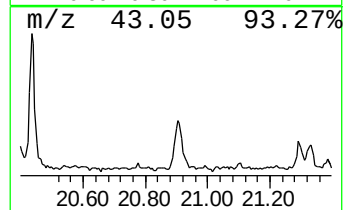
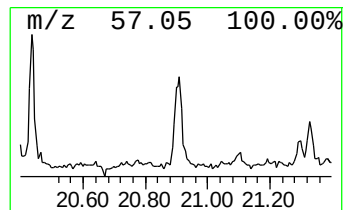
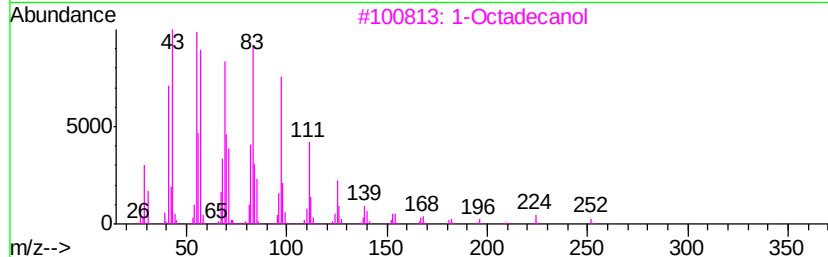
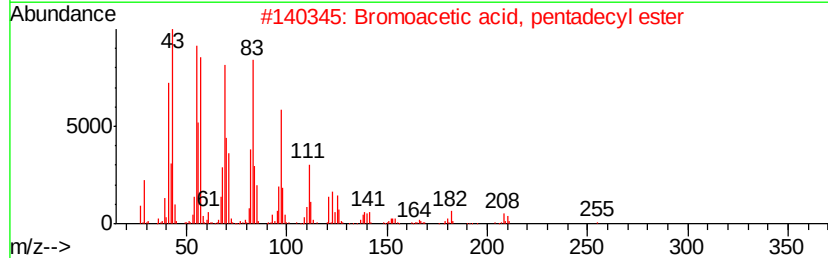
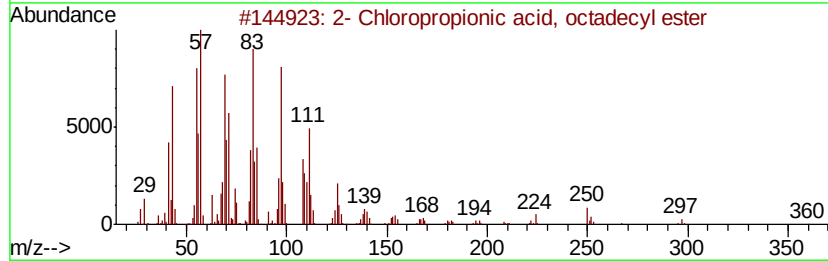
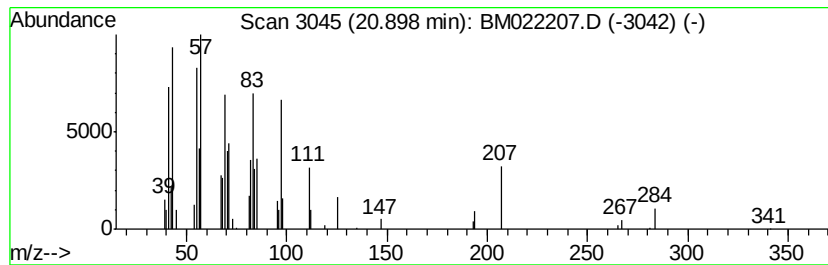
Instrument :  
BNA\_M  
ClientSampleId :  
0617-FIELD-BLANK-38TH-ST

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM082019.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 7 2- Chloropropionic acid, oc... Concentration Rank 9

| R.T.    | EstConc | Area                               | Relative to ISTD | R.T.       |             |      |
|---------|---------|------------------------------------|------------------|------------|-------------|------|
| 20.90   | 2.64 ng | 66930                              | Chrysene-d12     | 21.20      |             |      |
| Hit# of | 5       | Tentative ID                       | MW               | MolForm    | CAS#        | Qual |
| 1       | 2-      | Chloropropionic acid, octadec...   | 360              | C21H41ClO2 | 088104-31-8 | 90   |
| 2       |         | Bromoacetic acid, pentadecyl ester | 348              | C17H33BrO2 | 131143-01-6 | 90   |
| 3       |         | 1-Octadecanol                      | 270              | C18H38O    | 000112-92-5 | 90   |
| 4       | 2-      | Chloropropionic acid, hexadec...   | 332              | C19H37ClO2 | 086711-81-1 | 87   |
| 5       |         | Bromoacetic acid, hexadecyl ester  | 362              | C18H35BrO2 | 005454-48-8 | 87   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082019\  
Data File : BM022207.D  
Acq On : 20 Aug 2019 11:55  
Operator : HP/JU  
Sample : K4332-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

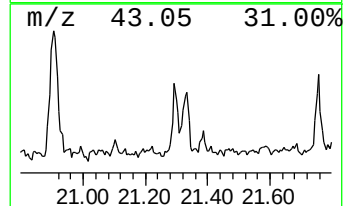
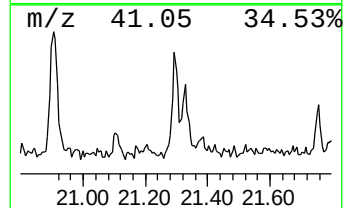
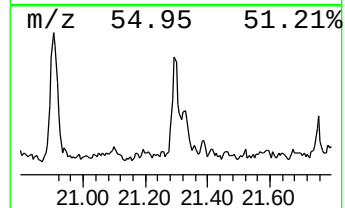
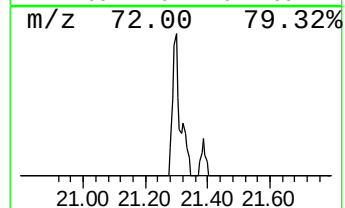
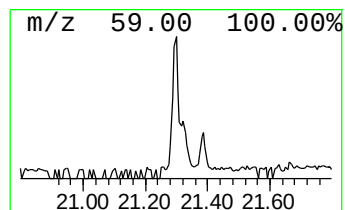
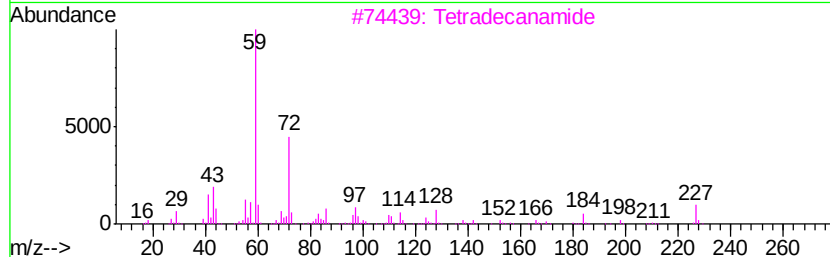
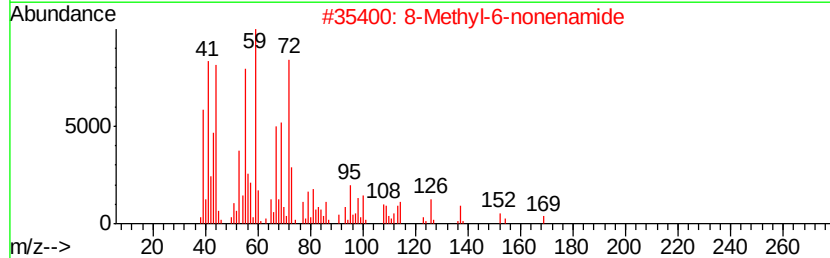
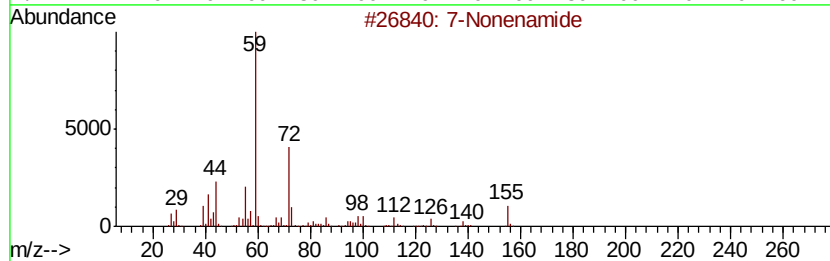
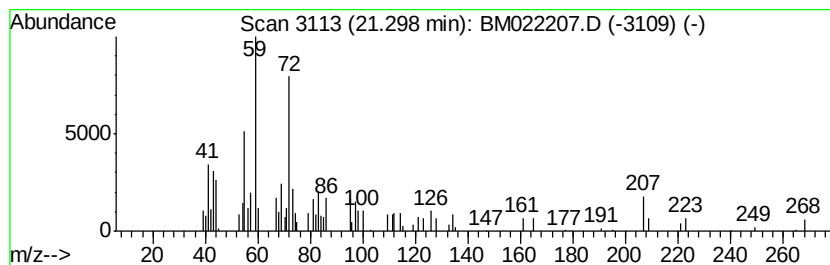
Instrument :  
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ClientSampleId :  
0617-FIELD-BLANK-38TH-ST

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM082019.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 8 7-Nonenamide Concentration Rank 10

| R.T.    | EstConc                             | Area           | Relative to ISTD | R.T.      |
|---------|-------------------------------------|----------------|------------------|-----------|
| 21.30   | 2.08 ng                             | 52766          | Chrysene-d12     | 21.20     |
| Hit# of | 5                                   | Tentative ID   | MW MolForm       | CAS# Qual |
| 1       | 7-Nonenamide                        | 155 C9H17NO    | 090949-53-4      | 78        |
| 2       | 8-Methyl-6-nonenamide               | 169 C10H19NO   | 1000293-20-9     | 53        |
| 3       | Tetradecanamide                     | 227 C14H29NO   | 000638-58-4      | 53        |
| 4       | 1-Propene, 3-[2-(2-methoxyethoxy... | 160 C8H16O3    | 013752-97-1      | 27        |
| 5       | Tridecanoic acid, thiophen-2-ylm... | 322 C18H30N2OS | 1000259-08-5     | 25        |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082019\  
Data File : BM022207.D  
Acq On : 20 Aug 2019 11:55  
Operator : HP/JU  
Sample : K4332-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

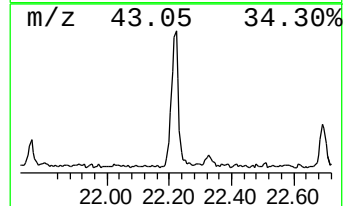
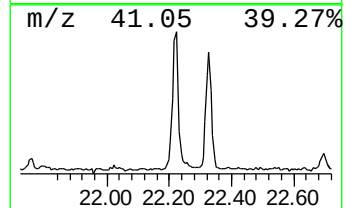
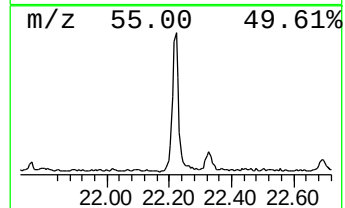
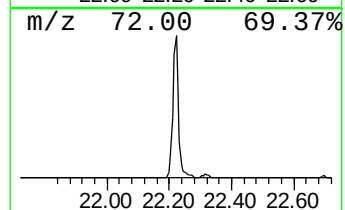
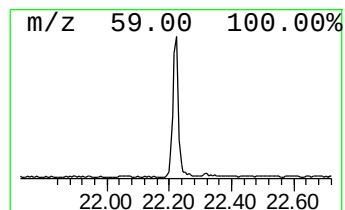
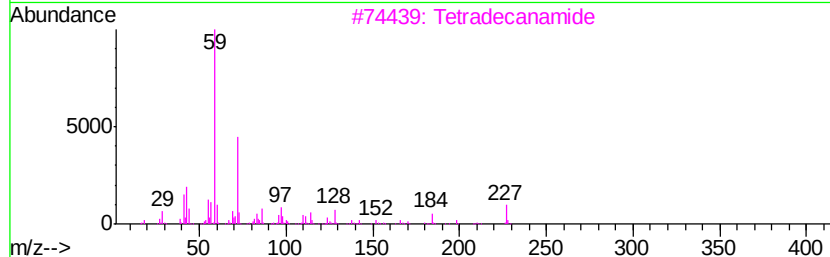
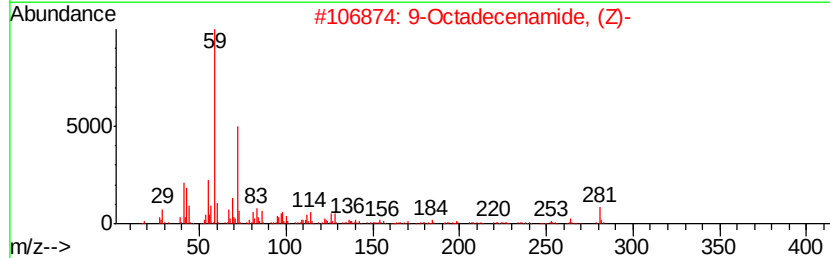
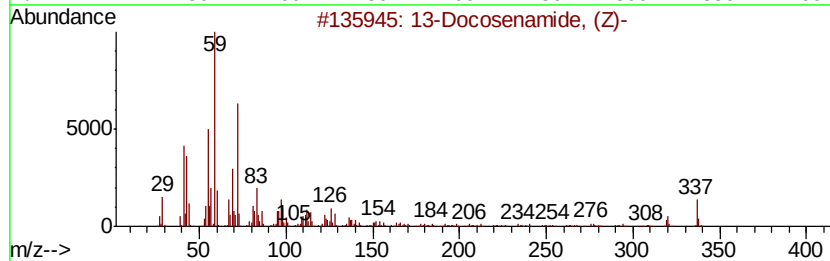
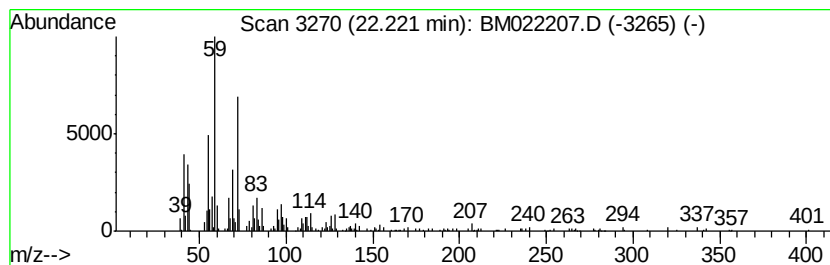
Instrument :  
BNA\_M  
ClientSampleId :  
0617-FIELD-BLANK-38TH-ST

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM082019.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 9 13-Docosenamide, (Z)- Concentration Rank 4

| R.T.    | EstConc  | Area                   | Relative to ISTD | R.T.           |
|---------|----------|------------------------|------------------|----------------|
| 22.22   | 14.01 ng | 355883                 | Chrysene-d12     | 21.20          |
| Hit# of | 5        | Tentative ID           | MW MolForm       | CAS# Qual      |
| 1       |          | 13-Docosenamide, (Z)-  | 337 C22H43NO     | 000112-84-5 76 |
| 2       |          | 9-Octadecenamide, (Z)- | 281 C18H35NO     | 000301-02-0 74 |
| 3       |          | Tetradecanamide        | 227 C14H29NO     | 000638-58-4 64 |
| 4       |          | Octanamide             | 143 C8H17NO      | 000629-01-6 64 |
| 5       |          | Dodecanamide           | 199 C12H25NO     | 001120-16-7 59 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082019\  
Data File : BM022207.D  
Acq On : 20 Aug 2019 11:55  
Operator : HP/JU  
Sample : K4332-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

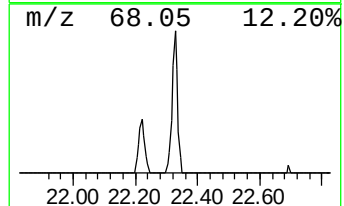
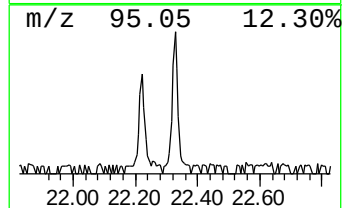
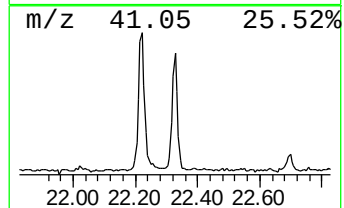
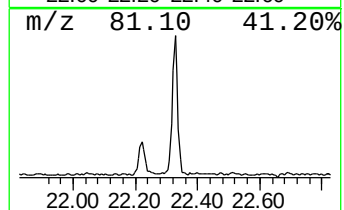
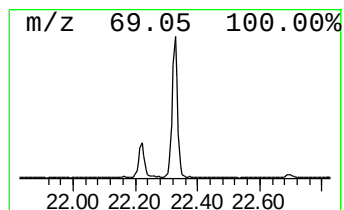
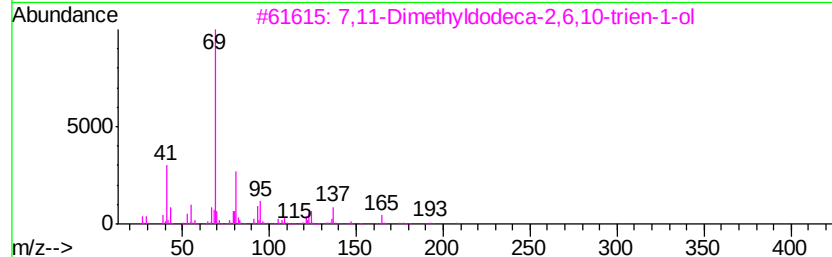
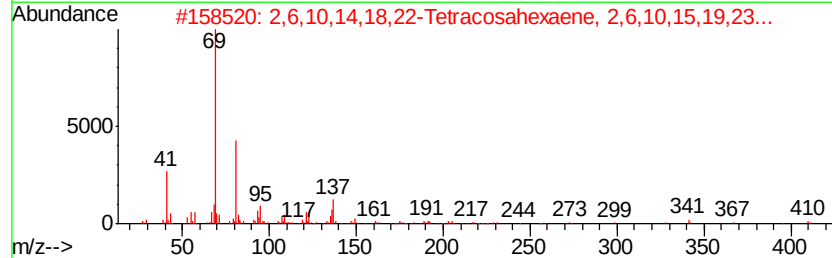
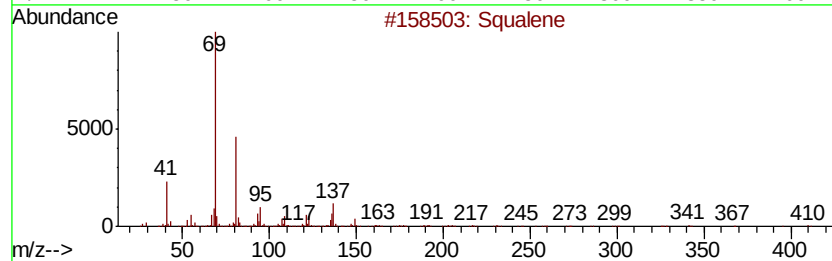
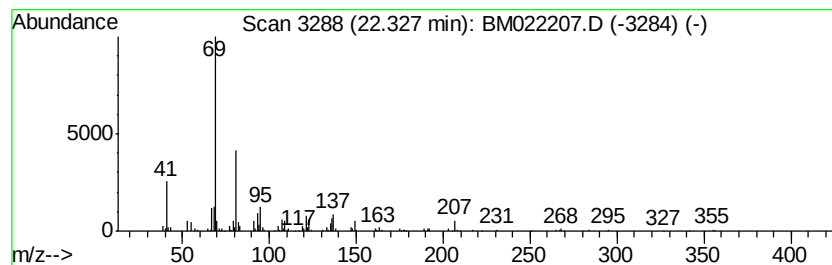
Instrument :  
BNA\_M  
ClientSampleId :  
0617-FIELD-BLANK-38TH-ST

Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM082019.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 10 Squalene Concentration Rank 5

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.           |
|---------|-------------------------------------|--------------|------------------|----------------|
| 22.33   | 8.53 ng                             | 198402       | Perylene-d12     | 23.40          |
| Hit# of | 5                                   | Tentative ID | MW MolForm       | CAS# Qual      |
| 1       | Squalene                            |              | 410 C30H50       | 007683-64-9 90 |
| 2       | 2,6,10,14,18,22-Tetracosahexaene... |              | 410 C30H50       | 000111-02-4 87 |
| 3       | 7,11-Dimethyldodeca-2,6,10-trien... |              | 208 C14H240      | 125529-10-4 81 |
| 4       | 2,6,10,14,18-Pentamethyl-2,6,10,... |              | 342 C25H42       | 075581-03-2 72 |
| 5       | 2,6,10,14-Hexadecatetraenoic aci... |              | 332 C22H36O2     | 024035-35-6 72 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_M\DATA\BM082019\  
Data File : BM022207.D  
Acq On : 20 Aug 2019 11:55  
Operator : HP/JU  
Sample : K4332-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
0617-FIELD-BLANK-38TH-ST

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\8270-BM082019.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

| TIC Top Hit name      | RT    | EstConc | Units | Response | --Internal Standard-- |       |        |      |
|-----------------------|-------|---------|-------|----------|-----------------------|-------|--------|------|
|                       |       |         |       |          | #                     | RT    | Resp   | Conc |
| unknown7.10           | 7.10  | 103.4   | ng    | 1176990  | 1                     | 7.57  | 227690 | 20.0 |
| Tridecanoic acid      | 17.88 | 4.6     | ng    | 105002   | 4                     | 16.99 | 460936 | 20.0 |
| Hexadecanamide        | 19.34 | 7.6     | ng    | 192999   | 5                     | 21.20 | 507987 | 20.0 |
| 9-Octadecenamide, ... | 20.34 | 170.5   | ng    | 4330960  | 5                     | 21.20 | 507987 | 20.0 |
| Tetradecanamide       | 20.43 | 7.4     | ng    | 187262   | 5                     | 21.20 | 507987 | 20.0 |
| 2-Chloropropioni...   | 20.90 | 2.6     | ng    | 66930    | 5                     | 21.20 | 507987 | 20.0 |
| 7-Nonenamide          | 21.30 | 2.1     | ng    | 52766    | 5                     | 21.20 | 507987 | 20.0 |
| 13-Docosenamide, ...  | 22.22 | 14.0    | ng    | 355883   | 5                     | 21.20 | 507987 | 20.0 |
| Squalene              | 22.33 | 8.5     | ng    | 198402   | 6                     | 23.40 | 465121 | 20.0 |