

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP063020\  
 Data File : BP002582.D  
 Acq On : 30 Jun 2020 19:10  
 Operator : CG/JU  
 Sample : L3153-40  
 Misc :  
 ALS Vial : 12 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SP2-K

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\_P\METHODS\8270-BP062920.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         | 5.193       | 385           | 392         | 408          | rBV      | 978600         | 1610610       | 15.37%          | 2.649%        |
| 2         | 6.763       | 653           | 659         | 674          | rVB      | 1183954        | 1988966       | 18.98%          | 3.271%        |
| 3         | 7.569       | 788           | 796         | 806          | rBV      | 268708         | 432540        | 4.13%           | 0.711%        |
| 4         | 7.887       | 843           | 850         | 862          | rBV      | 987127         | 1577594       | 15.06%          | 2.595%        |
| 5         | 8.716       | 983           | 991         | 1008         | rBV      | 811522         | 1380045       | 13.17%          | 2.270%        |
| 6         | 10.346      | 1258          | 1268        | 1285         | rBV      | 392888         | 698444        | 6.67%           | 1.149%        |
| 7         | 11.822      | 1509          | 1519        | 1535         | rBV      | 186679         | 314920        | 3.01%           | 0.518%        |
| 8         | 12.593      | 1639          | 1650        | 1657         | rBV2     | 98433          | 167095        | 1.59%           | 0.275%        |
| 9         | 12.834      | 1684          | 1691        | 1708         | rBV      | 2195424        | 3356546       | 32.03%          | 5.521%        |
| 10        | 13.110      | 1733          | 1738        | 1753         | rVB2     | 160840         | 250024        | 2.39%           | 0.411%        |
| 11        | 13.681      | 1831          | 1835        | 1849         | rVB      | 376157         | 537657        | 5.13%           | 0.884%        |
| 12        | 14.222      | 1914          | 1927        | 1934         | rBV2     | 723827         | 1421488       | 13.57%          | 2.338%        |
| 13        | 14.316      | 1934          | 1943        | 1951         | rVB3     | 75512          | 143807        | 1.37%           | 0.237%        |
| 14        | 14.504      | 1968          | 1975        | 1981         | rBV3     | 106625         | 195541        | 1.87%           | 0.322%        |
| 15        | 15.575      | 2152          | 2157        | 2162         | rBV3     | 123344         | 197728        | 1.89%           | 0.325%        |
| 16        | 15.722      | 2174          | 2182        | 2193         | rBV2     | 2527564        | 3668614       | 35.01%          | 6.034%        |
| 17        | 16.981      | 2383          | 2396        | 2401         | rBV      | 1063519        | 1537142       | 14.67%          | 2.528%        |
| 18        | 17.069      | 2406          | 2411        | 2415         | rBV      | 108211         | 142011        | 1.36%           | 0.234%        |
| 19        | 19.486      | 2817          | 2822        | 2826         | rBV      | 554840         | 677300        | 6.46%           | 1.114%        |
| 20        | 19.569      | 2831          | 2836        | 2845         | rVB      | 5139121        | 6508992       | 62.12%          | 10.706%       |
| 21        | 19.833      | 2876          | 2881        | 2884         | rBV      | 304487         | 356429        | 3.40%           | 0.586%        |
| 22        | 19.975      | 2902          | 2905        | 2910         | rVB      | 185064         | 215985        | 2.06%           | 0.355%        |
| 23        | 20.404      | 2975          | 2978        | 2982         | rVB      | 120262         | 139341        | 1.33%           | 0.229%        |
| 24        | 20.457      | 2983          | 2987        | 2991         | rBV2     | 90942          | 121224        | 1.16%           | 0.199%        |
| 25        | 20.504      | 2991          | 2995        | 3000         | rBV      | 373547         | 440419        | 4.20%           | 0.724%        |
| 26        | 20.663      | 3016          | 3022        | 3027         | rVB      | 7935159        | 10478443      | 100.00%         | 17.235%       |
| 27        | 20.827      | 3046          | 3050        | 3061         | rBV      | 998335         | 1291190       | 12.32%          | 2.124%        |
| 28        | 21.086      | 3089          | 3094        | 3103         | rBV2     | 1732057        | 2160855       | 20.62%          | 3.554%        |
| 29        | 21.263      | 3121          | 3124        | 3132         | rVB      | 114958         | 147676        | 1.41%           | 0.243%        |
| 30        | 21.433      | 3150          | 3153        | 3158         | rVB2     | 144588         | 163133        | 1.56%           | 0.268%        |
| 31        | 21.704      | 3193          | 3199        | 3209         | rBV      | 990392         | 1580567       | 15.08%          | 2.600%        |
| 32        | 22.369      | 3308          | 3312        | 3324         | rVB      | 370368         | 524344        | 5.00%           | 0.862%        |
| 33        | 22.651      | 3355          | 3360        | 3364         | rBV      | 1161732        | 1770525       | 16.90%          | 2.912%        |
| 34        | 22.692      | 3364          | 3367        | 3378         | rVB      | 698144         | 1095931       | 10.46%          | 1.803%        |

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Sample : L3153-40  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SP2-K

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

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

Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\8270-BP062920.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 35 | 23.216 | 3452 | 3456 | 3460 | rVB  | 119923  | 180286  | 1.72%  | 0.297% |
| 36 | 23.280 | 3460 | 3467 | 3476 | rBV  | 1317448 | 2397199 | 22.88% | 3.943% |
| 37 | 23.516 | 3502 | 3507 | 3515 | rVB  | 601602  | 1025551 | 9.79%  | 1.687% |
| 38 | 23.857 | 3559 | 3565 | 3572 | rVB  | 1175573 | 2227717 | 21.26% | 3.664% |
| 39 | 23.933 | 3572 | 3578 | 3591 | rBV  | 447169  | 965745  | 9.22%  | 1.588% |
| 40 | 25.027 | 3758 | 3764 | 3777 | rVB  | 499967  | 1241921 | 11.85% | 2.043% |
| 41 | 25.462 | 3831 | 3838 | 3851 | rVB2 | 369378  | 1041536 | 9.94%  | 1.713% |
| 42 | 25.598 | 3855 | 3861 | 3870 | rVV2 | 191979  | 522538  | 4.99%  | 0.859% |
| 43 | 26.309 | 3973 | 3982 | 3999 | rVB3 | 832930  | 2660512 | 25.39% | 4.376% |
| 44 | 27.045 | 4101 | 4107 | 4112 | rBV4 | 132069  | 347672  | 3.32%  | 0.572% |
| 45 | 27.845 | 4238 | 4243 | 4266 | rVB8 | 219366  | 894160  | 8.53%  | 1.471% |

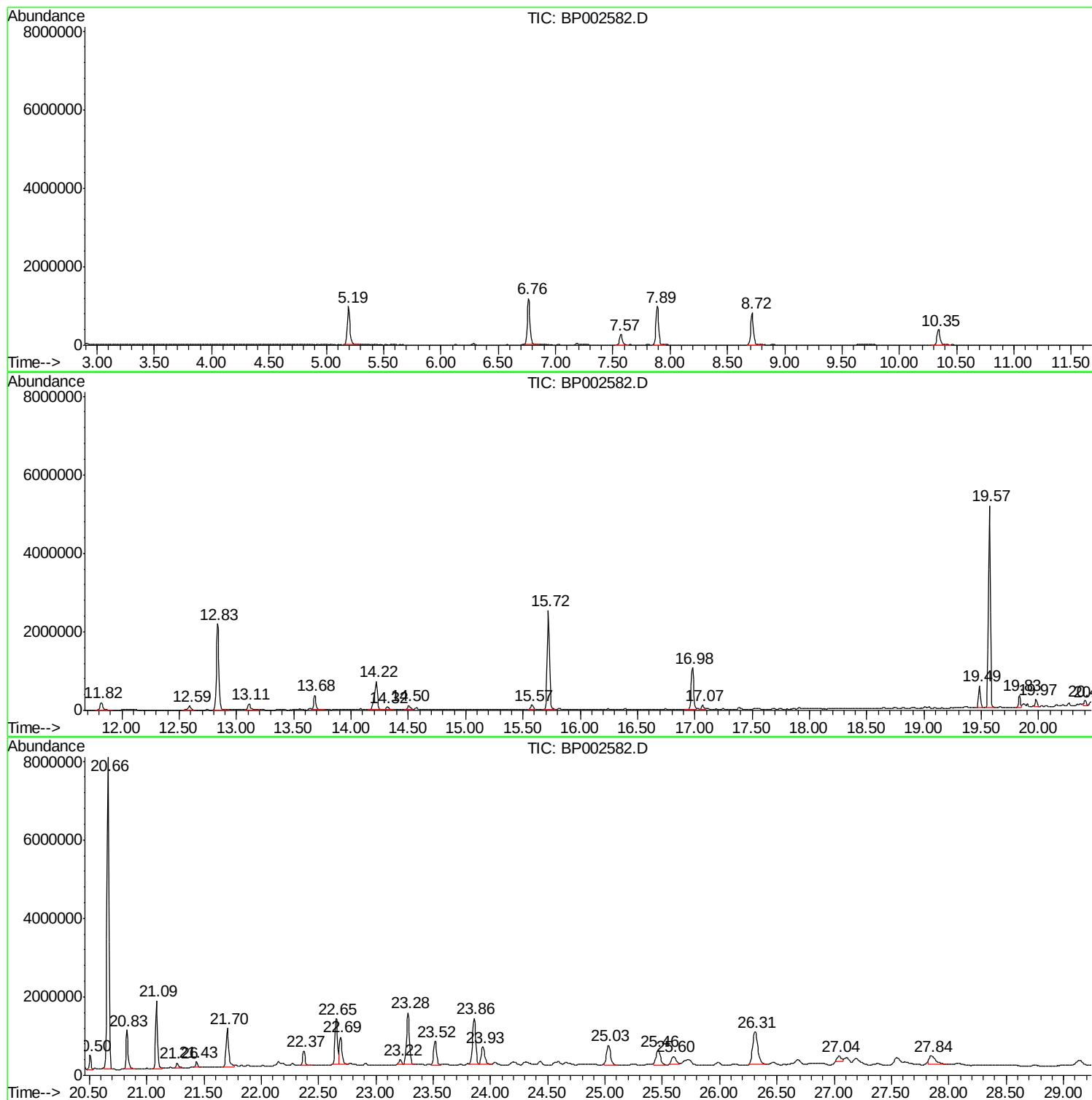
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Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP063020\  
Data File : BP002582.D  
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Sample : L3153-40  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SP2-K

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP063020\  
Data File : BP002582.D  
Acq On : 30 Jun 2020 19:10  
Operator : CG/JU  
Sample : L3153-40  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SP2-K

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

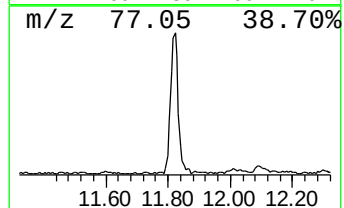
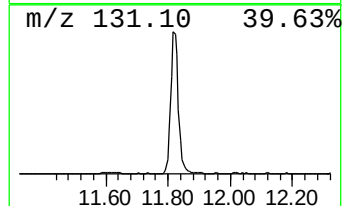
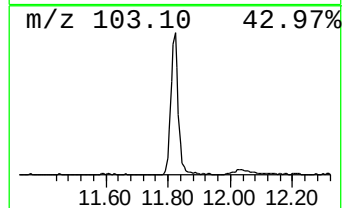
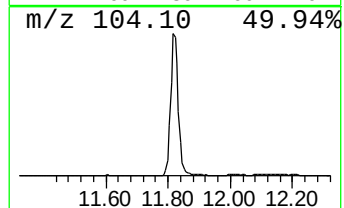
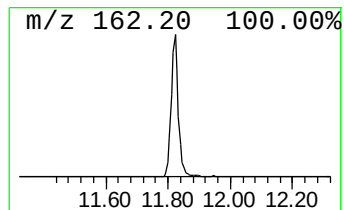
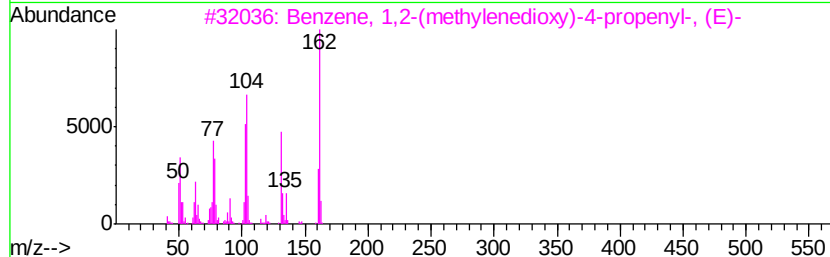
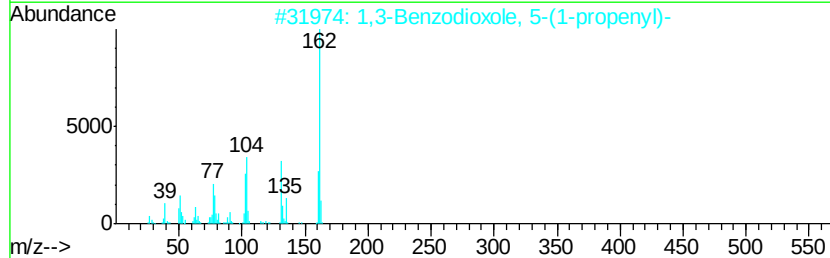
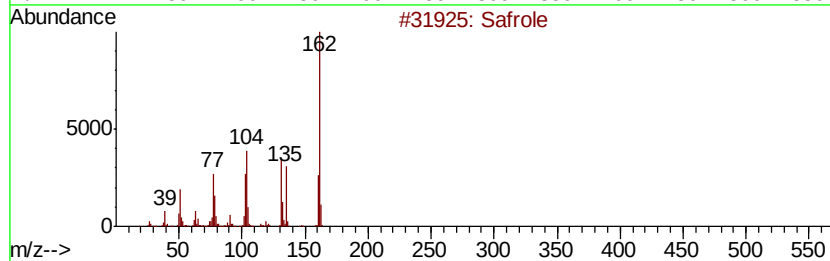
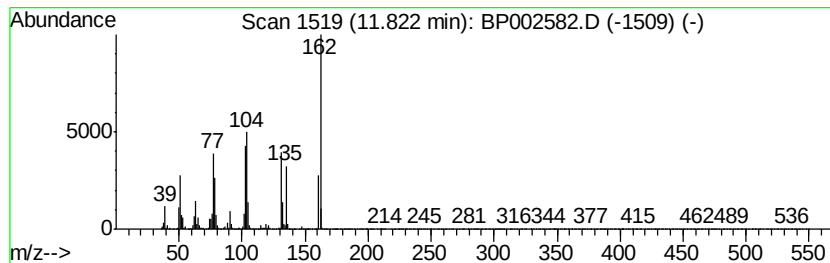
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 Safrole Concentration Rank 10

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.82 | 9.02 ng | 314920 | Naphthalene-d8   | 10.35 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Safrole                             | 162 | C10H10O2 | 000094-59-7 | 98   |
| 2    |    | 1,3-Benzodioxole, 5-(1-propenyl)-   | 162 | C10H10O2 | 000120-58-1 | 97   |
| 3    |    | Benzene, 1,2-(methylenedioxy)-4-... | 162 | C10H10O2 | 004043-71-4 | 96   |
| 4    |    | 1,3-Benzodioxole, 5-(1-propenyl)... | 162 | C10H10O2 | 017627-76-8 | 95   |
| 5    |    | 1H-Isoindole-1,3(2H)-dione, 2-am... | 162 | C8H6N2O2 | 001875-48-5 | 46   |



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BNA\_P  
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SP2-K

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

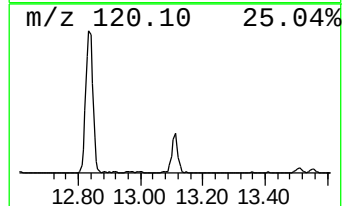
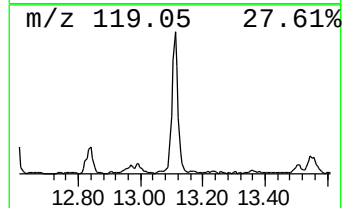
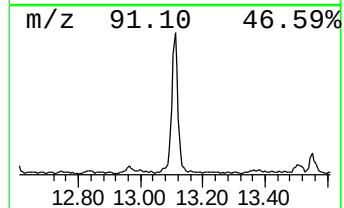
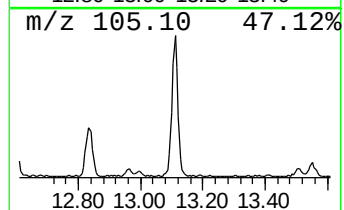
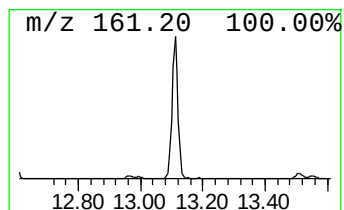
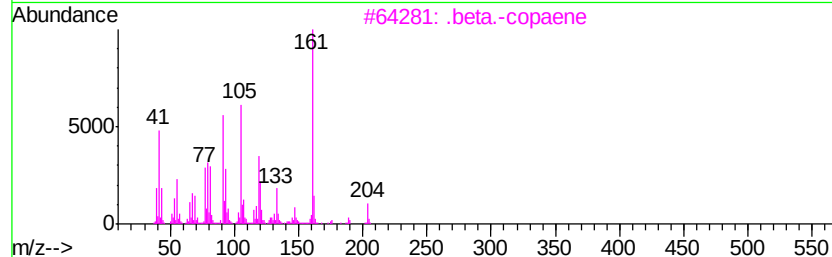
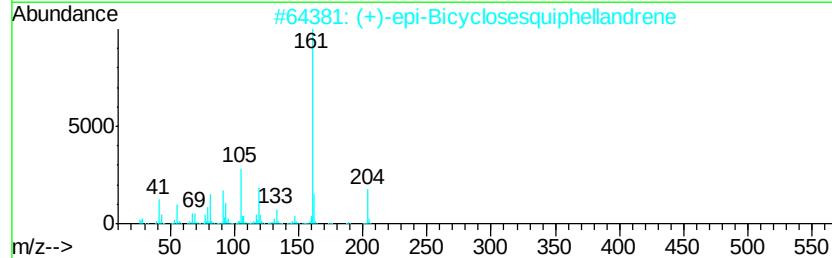
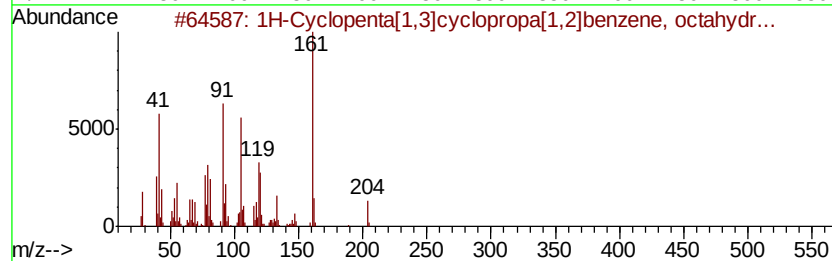
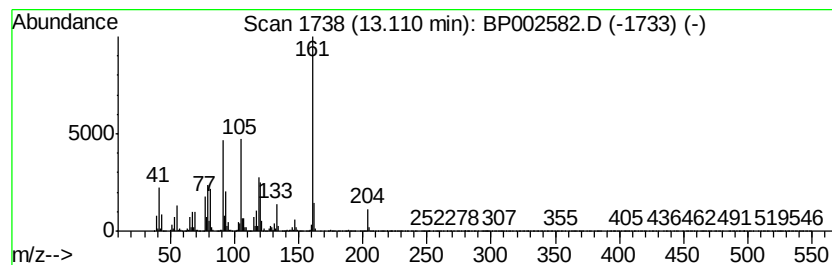
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 1H-Cyclopenta[1,3]cycloprop... Concentration Rank 19

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.11 | 3.52 ng | 250024 | Acenaphthene-d10 | 14.22 |

| Hit# | of | Tentative ID                         | MW  | MolForm | CAS#         | Qual |
|------|----|--------------------------------------|-----|---------|--------------|------|
| 1    | 5  | 1H-Cyclopenta[1,3]cyclopropa[1,2]... | 204 | C15H24  | 013744-15-5  | 98   |
| 2    |    | (+)-epi-Bicyclosquiphellandrene      | 204 | C15H24  | 054274-73-6  | 94   |
| 3    |    | .beta.-copaene                       | 204 | C15H24  | 1000374-18-9 | 94   |
| 4    |    | 1,6-Cyclodecadiene, 1-methyl-5-m...  | 204 | C15H24  | 023986-74-5  | 86   |
| 5    |    | Bicyclo[4.4.0]dec-1-ene, 2-isopr...  | 204 | C15H24  | 150320-52-8  | 81   |



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ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SP2-K

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

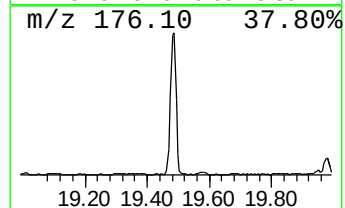
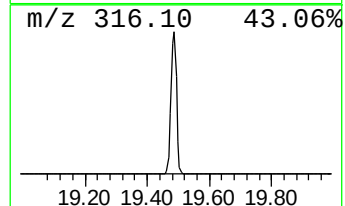
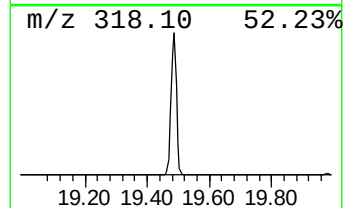
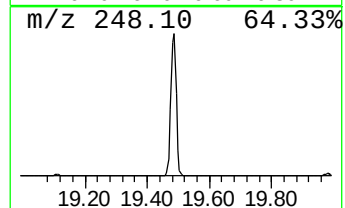
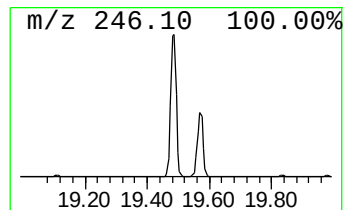
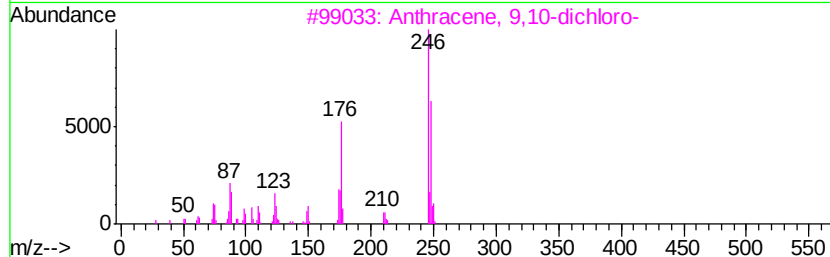
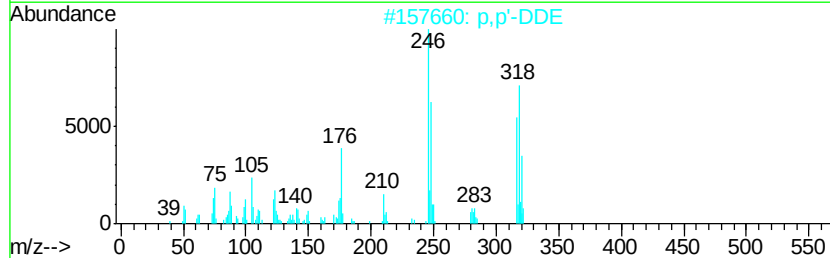
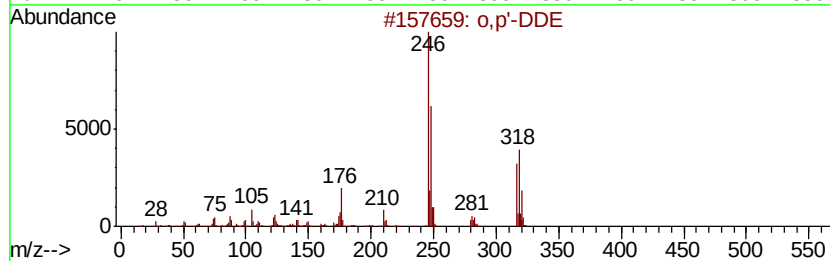
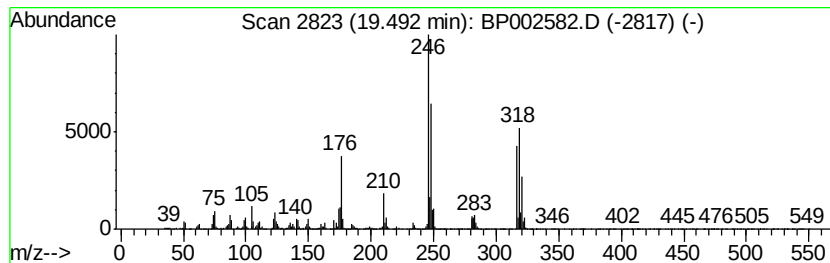
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 o,p'-DDE Concentration Rank 15

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 19.49 | 6.27 ng | 677300 | Chrysene-d12     | 21.09 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | o,p'-DDE                            | 316 | C14H8Cl4 | 003424-82-6 | 99   |
| 2    |    | p,p'-DDE                            | 316 | C14H8Cl4 | 000072-55-9 | 99   |
| 3    |    | Anthracene, 9,10-dichloro-          | 246 | C14H8Cl2 | 000605-48-1 | 97   |
| 4    |    | Benzene, 1,1'-(1,2-dichloro-1,2-... | 316 | C14H8Cl4 | 088970-69-8 | 94   |
| 5    |    | 9H-Fluorene, 9-(dichloromethylene)- | 246 | C14H8Cl2 | 000835-17-6 | 93   |



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SP2-K

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

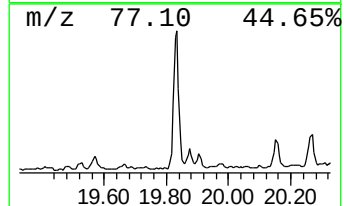
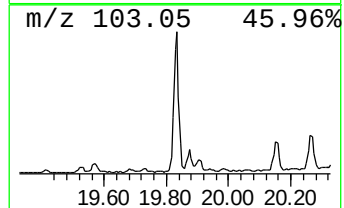
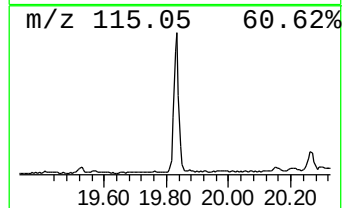
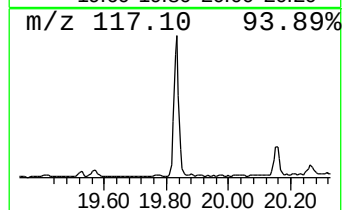
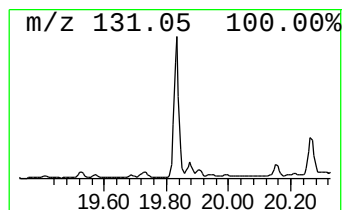
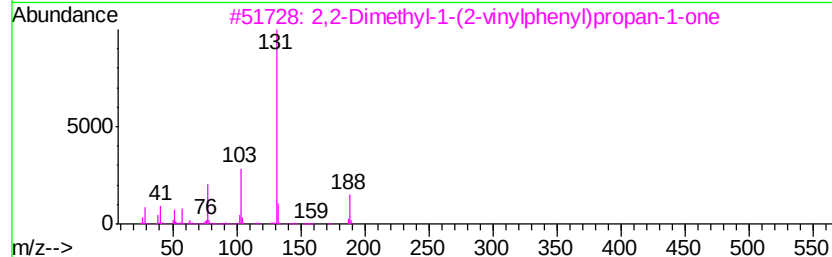
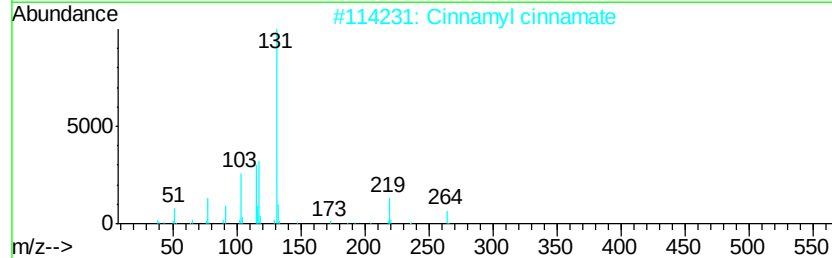
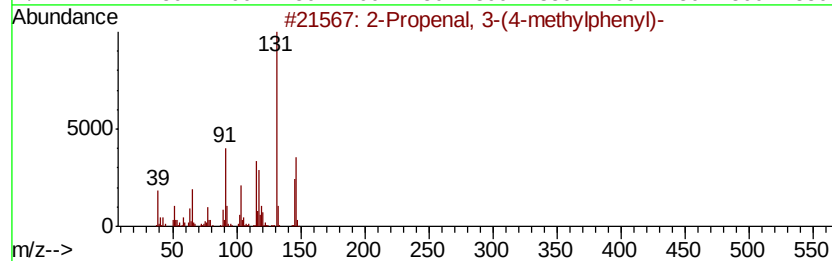
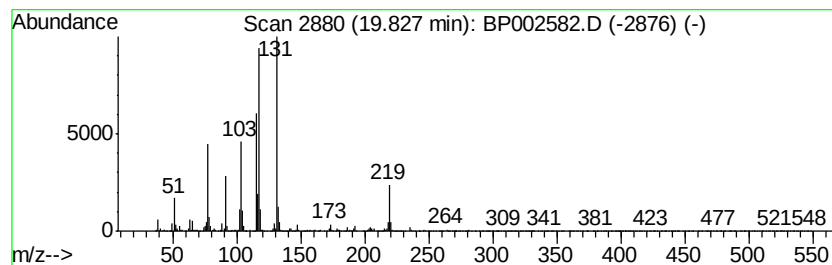
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 2-Propenal, 3-(4-methylphen... Concentration Rank 20

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 19.83 | 3.30 ng | 356429 | Chrysene-d12     | 21.09 |

| Hit# | of | Tentative ID                               | MW  | MolForm   | CAS#         | Qual |
|------|----|--------------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 2-Propenal, 3-(4-methylphenyl)-            | 146 | C10H10O   | 001504-75-2  | 50   |
| 2    |    | Cinnamyl cinnamate                         | 264 | C18H16O2  | 000122-69-0  | 43   |
| 3    |    | 2,2-Dimethyl-1-(2-vinylphenyl)propan-1-one | 188 | C13H16O   | 1000210-99-9 | 43   |
| 4    |    | trans-1-Cinnamoylimidazole                 | 198 | C12H10N2O | 001138-15-4  | 38   |
| 5    |    | 3-Butenoic acid, 2-oxo-4-phenyl-           | 190 | C11H10O3  | 1000142-22-4 | 38   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP063020\  
Data File : BP002582.D  
Acq On : 30 Jun 2020 19:10  
Operator : CG/JU  
Sample : L3153-40  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SP2-K

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

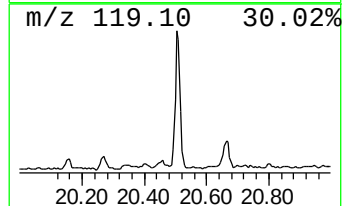
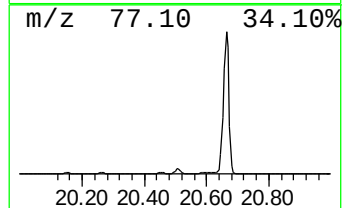
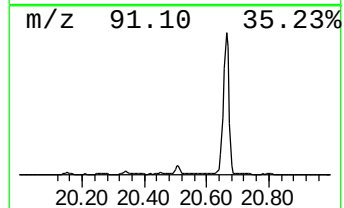
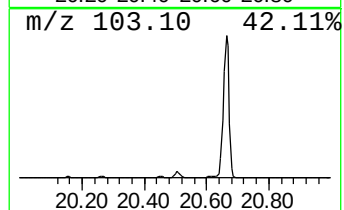
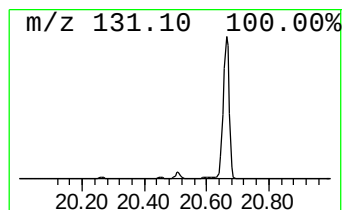
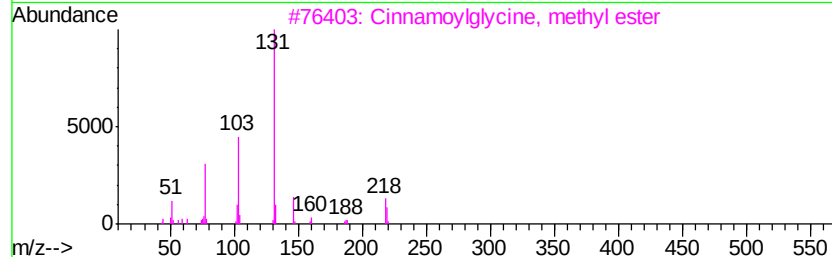
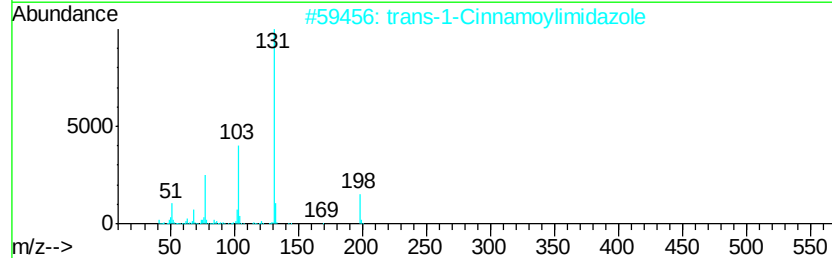
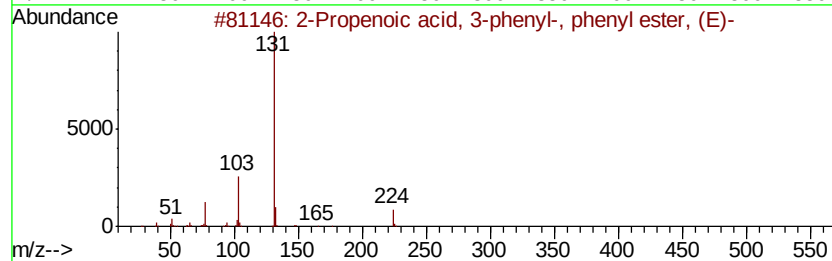
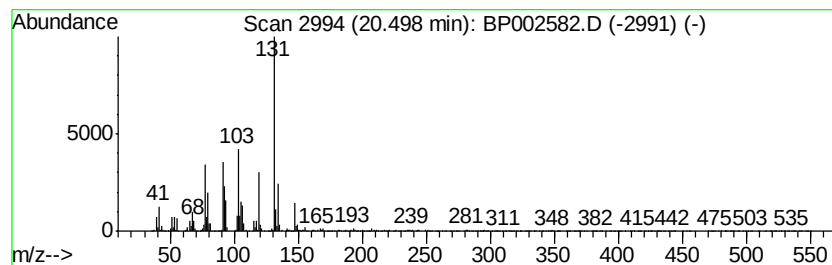
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TIC Integration Parameters: LSCINT.P

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Peak Number 6 unknown20.50 Concentration Rank 18

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 20.50 | 4.08 ng | 440419 | Chrysene-d12     | 21.09 |

| Hit# | of                                  | 5   | Tentative ID | MW           | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|--------------|---------|------|------|
| 1    | 2-Propenoic acid, 3-phenyl-, phe... | 224 | C15H12O2     | 025695-77-6  | 49      |      |      |
| 2    | trans-1-Cinnamoylimidazole          | 198 | C12H10N2O    | 001138-15-4  | 47      |      |      |
| 3    | Cinnamoylglycine, methyl ester      | 219 | C12H13NO3    | 040778-04-9  | 47      |      |      |
| 4    | p-Vinylbenzohydrazide               | 162 | C9H10N2O     | 1000244-74-9 | 47      |      |      |
| 5    | 1-Penten-3-one, 4,4-dimethyl-1-p... | 188 | C13H16O      | 000538-44-3  | 47      |      |      |





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP063020\  
Data File : BP002582.D  
Acq On : 30 Jun 2020 19:10  
Operator : CG/JU  
Sample : L3153-40  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

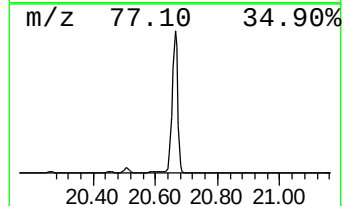
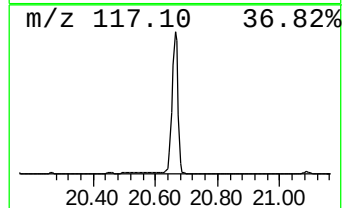
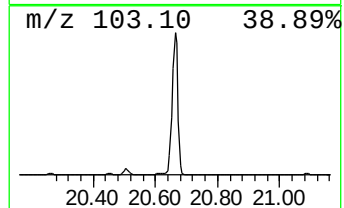
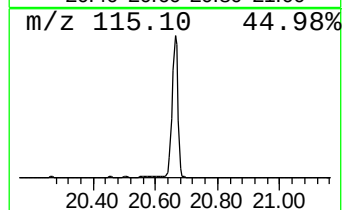
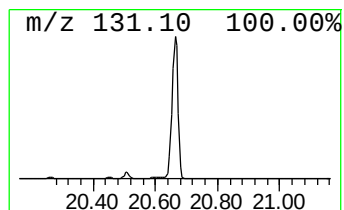
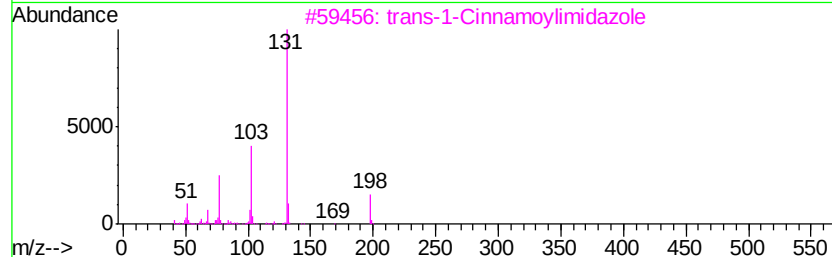
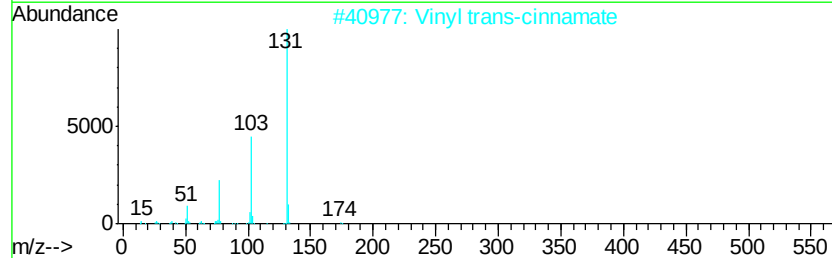
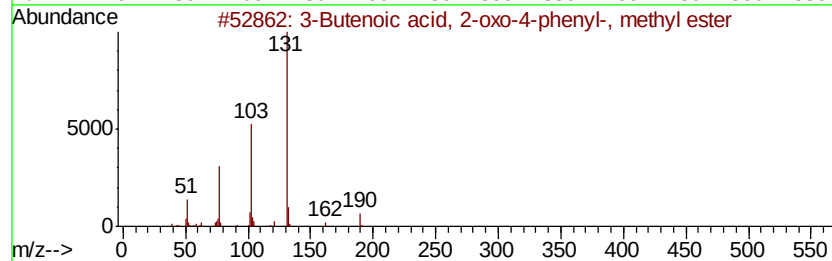
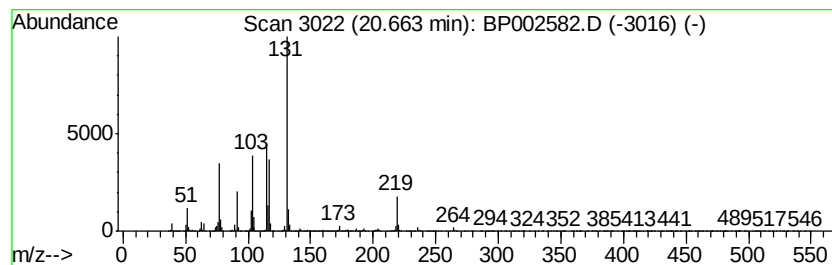
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ClientSampleId :  
SP2-K

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.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 3-Butenoic acid, 2-oxo-4-ph... Concentration Rank 1

| R.T.    | EstConc  | Area                                | Relative to ISTD | R.T.      |              |      |
|---------|----------|-------------------------------------|------------------|-----------|--------------|------|
| 20.66   | 96.98 ng | 10478400                            | Chrysene-d12     | 21.09     |              |      |
| Hit# of | 5        | Tentative ID                        | MW               | MolForm   | CAS#         | Qual |
| 1       |          | 3-Butenoic acid, 2-oxo-4-phenyl-... | 190              | C11H10O3  | 1000142-22-4 | 52   |
| 2       |          | Vinyl trans-cinnamate               | 174              | C11H10O2  | 017719-70-9  | 50   |
| 3       |          | trans-1-Cinnamoylimidazole          | 198              | C12H10N2O | 001138-15-4  | 50   |
| 4       |          | 2-Propenoic acid, 3-phenyl-, phe... | 224              | C15H12O2  | 025695-77-6  | 50   |
| 5       |          | N-(p-Vinylbenzoyl)-l-alanine        | 219              | C12H13NO3 | 139184-60-4  | 50   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP063020\  
Data File : BP002582.D  
Acq On : 30 Jun 2020 19:10  
Operator : CG/JU  
Sample : L3153-40  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

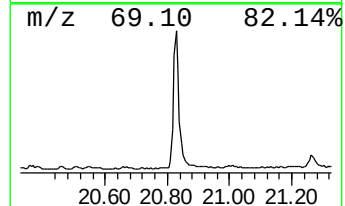
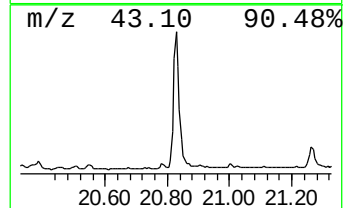
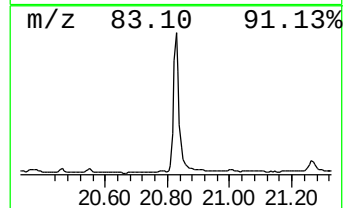
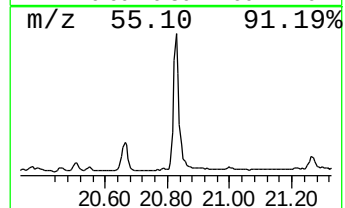
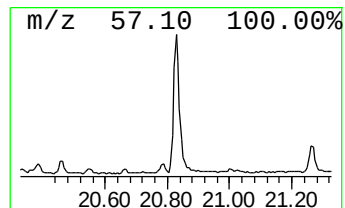
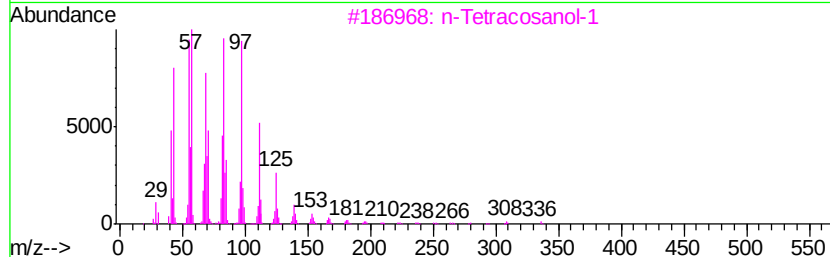
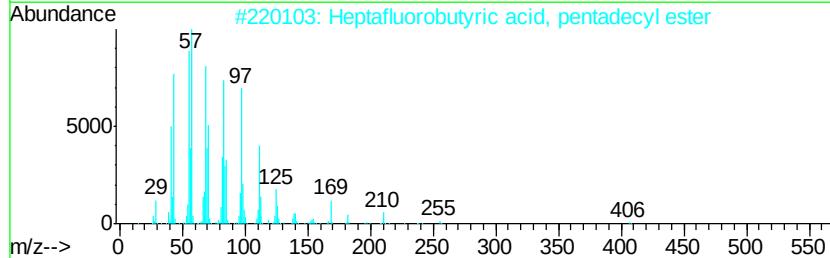
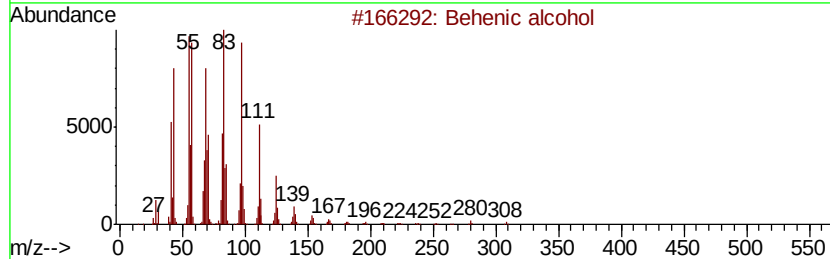
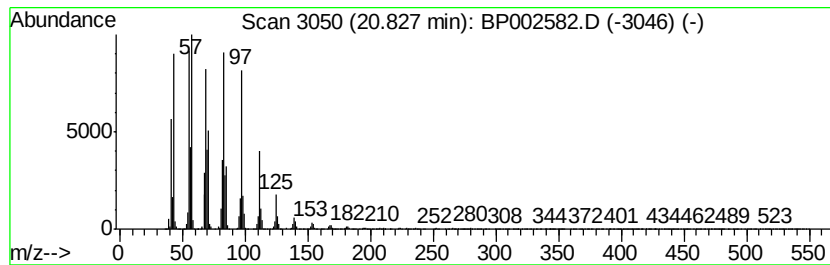
Instrument :  
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ClientSampleId :  
SP2-K

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.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 Behenic alcohol Concentration Rank 7

| R.T.    | EstConc  | Area                                | Relative to ISTD | R.T.        |             |      |
|---------|----------|-------------------------------------|------------------|-------------|-------------|------|
| 20.83   | 11.95 ng | 1291190                             | Chrysene-d12     | 21.09       |             |      |
| Hit# of | 5        | Tentative ID                        | MW               | MolForm     | CAS#        | Qual |
| 1       |          | Behenic alcohol                     | 326              | C22H46O     | 000661-19-8 | 94   |
| 2       |          | Heptafluorobutyric acid, pentade... | 424              | C19H31F7O2  | 959261-23-5 | 94   |
| 3       |          | n-Tetracosanol-1                    | 354              | C24H50O     | 000506-51-4 | 94   |
| 4       |          | 1-Heneicosanol                      | 312              | C21H44O     | 015594-90-8 | 94   |
| 5       |          | Trichloroacetic acid, pentadecyl... | 372              | C17H31Cl3O2 | 074339-53-0 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP063020\  
Data File : BP002582.D  
Acq On : 30 Jun 2020 19:10  
Operator : CG/JU  
Sample : L3153-40  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SP2-K

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

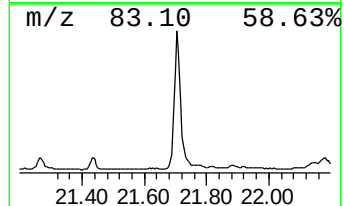
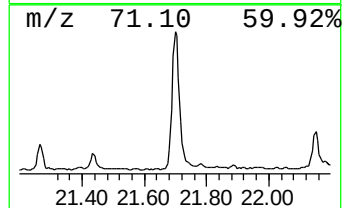
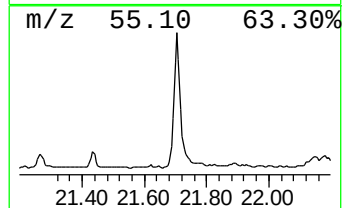
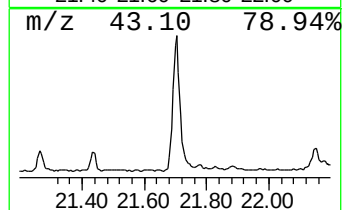
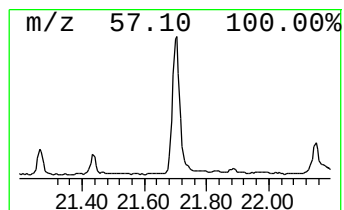
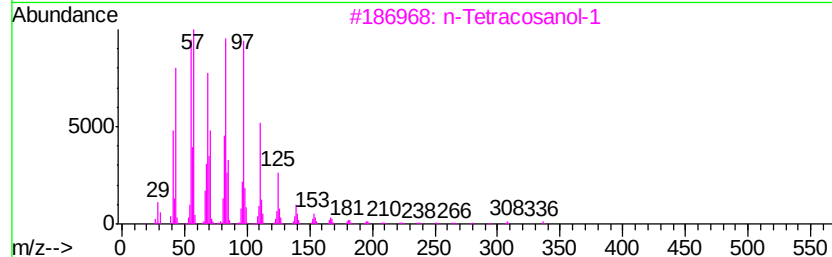
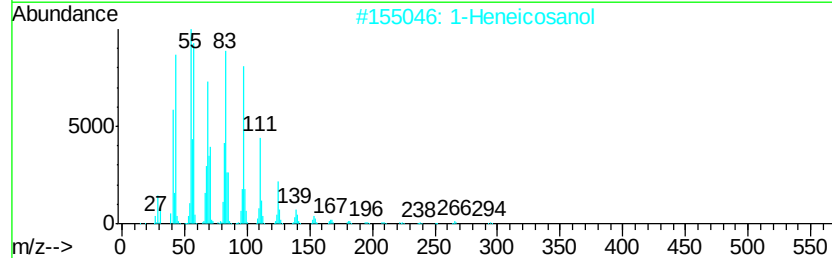
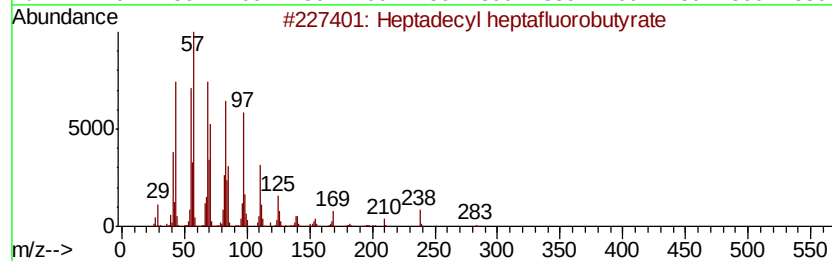
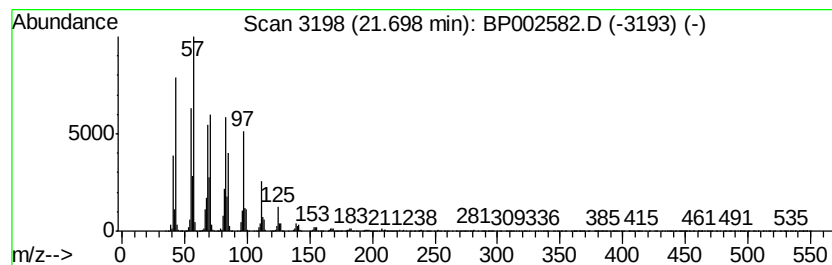
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 Heptadecyl heptafluorobutyrate Concentration Rank 6

| R.T.  | EstConc  | Area    | Relative to ISTD |  | R.T.  |
|-------|----------|---------|------------------|--|-------|
| 21.70 | 14.63 ng | 1580570 | Chrysene-d12     |  | 21.09 |

| Hit# of | 5                           | Tentative ID        | MW  | MolForm    | CAS#        | Qual |
|---------|-----------------------------|---------------------|-----|------------|-------------|------|
| 1       | Heptadecyl                  | heptafluorobutyrate | 452 | C21H35F7O2 | 959085-66-6 | 94   |
| 2       | 1-Heneicosanol              |                     | 312 | C21H44O    | 015594-90-8 | 93   |
| 3       | n-Tetracosanol-1            |                     | 354 | C24H50O    | 000506-51-4 | 93   |
| 4       | Behenic alcohol             |                     | 326 | C22H46O    | 000661-19-8 | 93   |
| 5       | Ethanol, 2-(tetradecyloxy)- |                     | 258 | C16H34O2   | 002136-70-1 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP063020\  
Data File : BP002582.D  
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Operator : CG/JU  
Sample : L3153-40  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

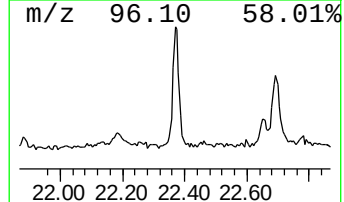
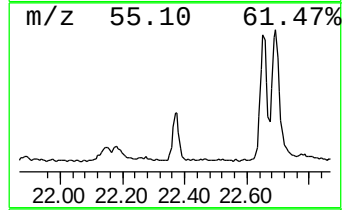
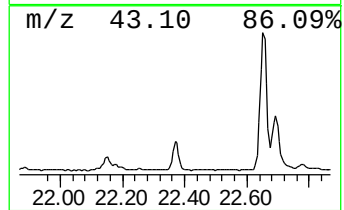
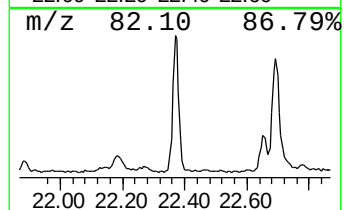
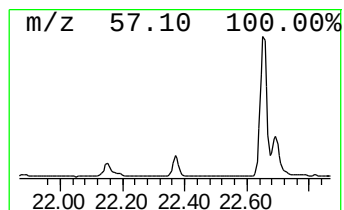
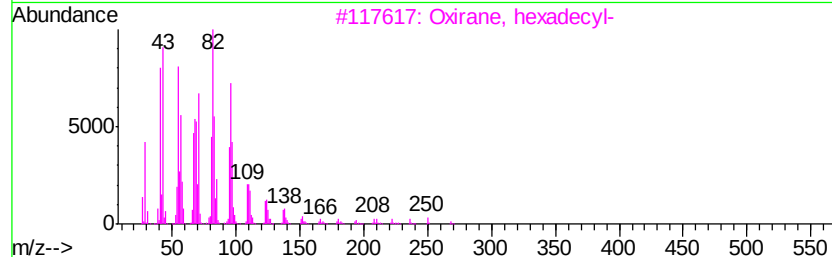
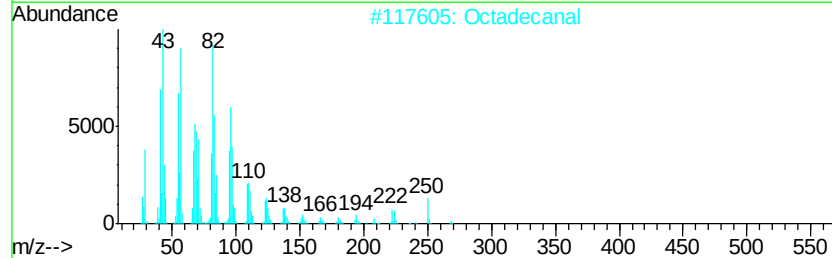
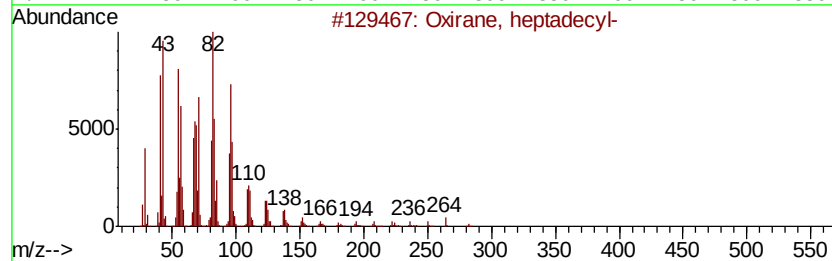
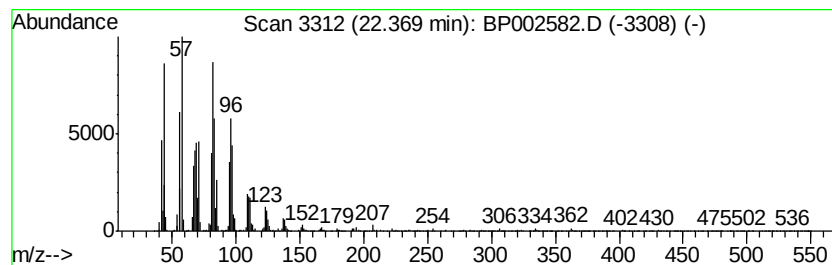
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ClientSampleId :  
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 10 Oxirane, heptadecyl- Concentration Rank 16

| R.T.      | EstConc              | Area   | Relative to ISTD | R.T.        |      |
|-----------|----------------------|--------|------------------|-------------|------|
| 22.37     | 4.37 ng              | 524344 | Perylene-d12     | 23.28       |      |
| Hit# of 5 | Tentative ID         | MW     | MolForm          | CAS#        | Qual |
| 1         | Oxirane, heptadecyl- | 282    | C19H38O          | 067860-04-2 | 91   |
| 2         | Octadecanal          | 268    | C18H36O          | 000638-66-4 | 91   |
| 3         | Oxirane, hexadecyl-  | 268    | C18H36O          | 007390-81-0 | 91   |
| 4         | 1,19-Eicosadiene     | 278    | C20H38           | 014811-95-1 | 90   |
| 5         | 1,13-Tetradecadiene  | 194    | C14H26           | 021964-49-8 | 83   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP063020\  
Data File : BP002582.D  
Acq On : 30 Jun 2020 19:10  
Operator : CG/JU  
Sample : L3153-40  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SP2-K

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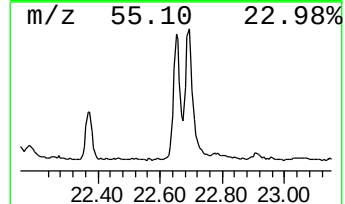
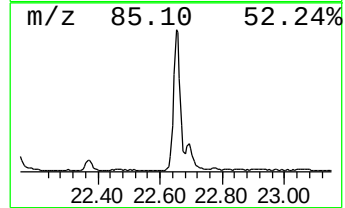
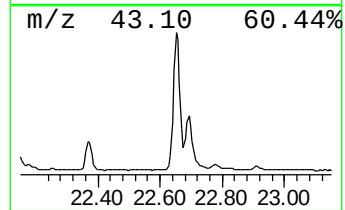
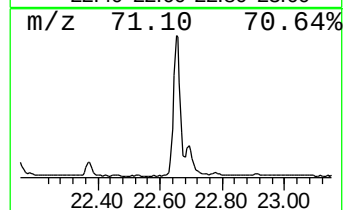
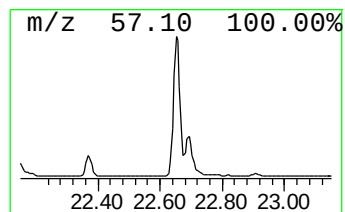
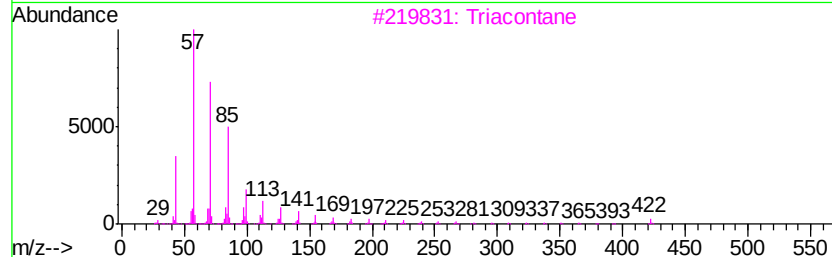
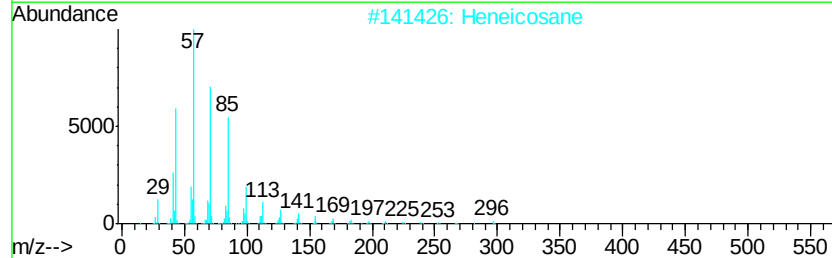
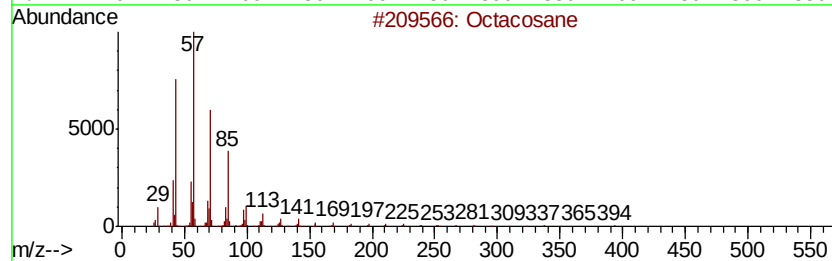
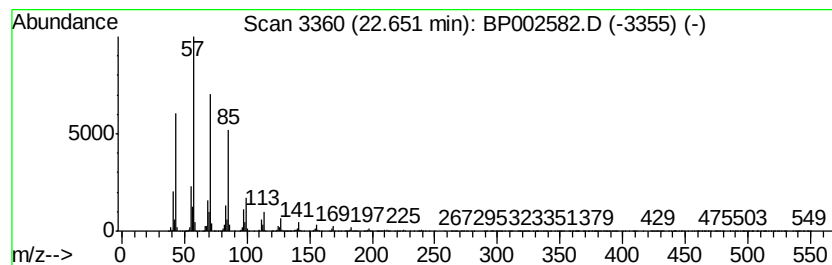
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TIC Integration Parameters: LSCINT.P

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Peak Number 11 Octacosane Concentration Rank 5

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 22.65 | 14.77 ng | 1770530 | Perylene-d12     | 23.28 |

| Hit# | of | Tentative ID       | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------|-----|---------|-------------|------|
| 1    | 5  | Octacosane         | 394 | C28H58  | 000630-02-4 | 98   |
| 2    |    | Heneicosane        | 296 | C21H44  | 000629-94-7 | 97   |
| 3    |    | Triacontane        | 422 | C30H62  | 000638-68-6 | 95   |
| 4    |    | Hentriacontane     | 437 | C31H64  | 000630-04-6 | 95   |
| 5    |    | Eicosane, 7-hexyl- | 366 | C26H54  | 055333-99-8 | 90   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP063020\  
Data File : BP002582.D  
Acq On : 30 Jun 2020 19:10  
Operator : CG/JU  
Sample : L3153-40  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SP2-K

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

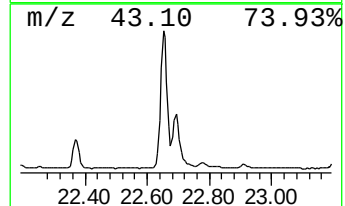
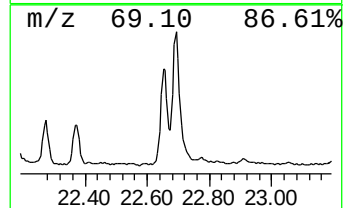
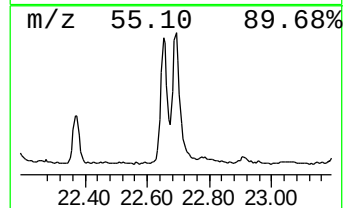
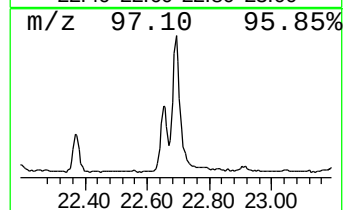
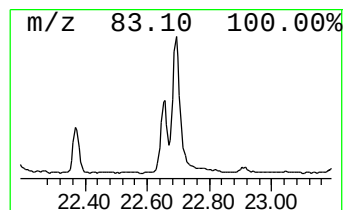
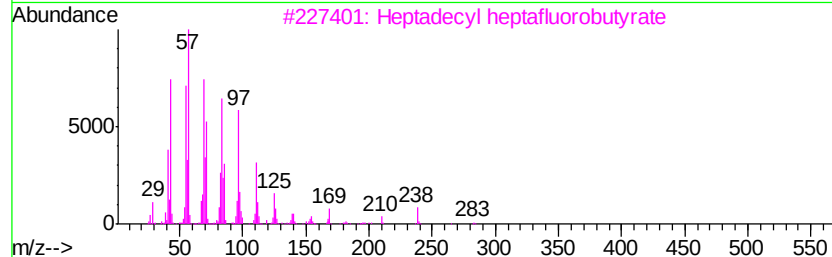
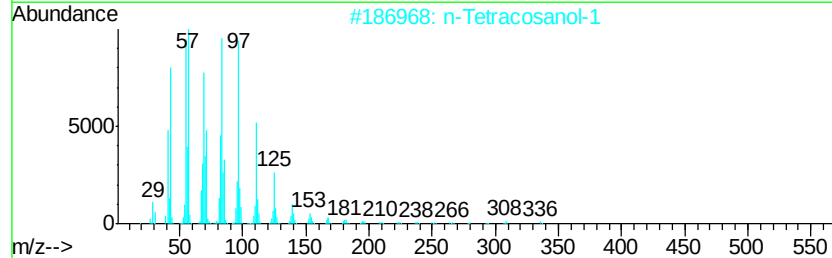
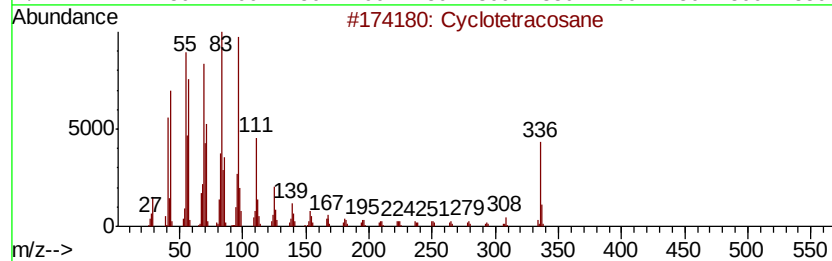
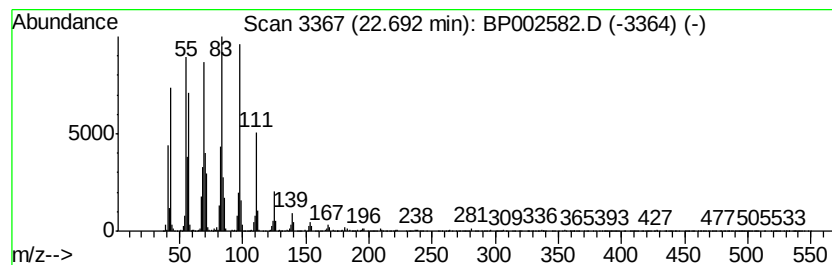
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 12 Cyclotetracosane Concentration Rank 9

| R.T.  | EstConc | Area    | Relative to ISTD | R.T.  |
|-------|---------|---------|------------------|-------|
| 22.69 | 9.14 ng | 1095930 | Perylene-d12     | 23.28 |

| Hit# | of | Tentative ID                   | MW  | MolForm    | CAS#        | Qual |
|------|----|--------------------------------|-----|------------|-------------|------|
| 1    | 5  | Cyclotetracosane               | 336 | C24H48     | 000297-03-0 | 93   |
| 2    |    | n-Tetracosanol-1               | 354 | C24H50O    | 000506-51-4 | 91   |
| 3    |    | Heptadecyl heptafluorobutvrate | 452 | C21H35F7O2 | 959085-66-6 | 90   |
| 4    |    | Octacosanol                    | 410 | C28H58O    | 000557-61-9 | 89   |
| 5    |    | 1-Heneicosyl formate           | 340 | C22H44O2   | 077899-03-7 | 83   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP063020\  
Data File : BP002582.D  
Acq On : 30 Jun 2020 19:10  
Operator : CG/JU  
Sample : L3153-40  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

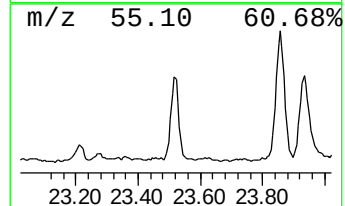
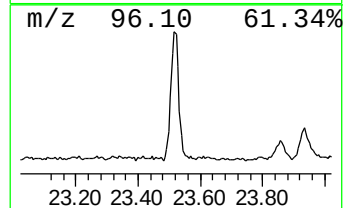
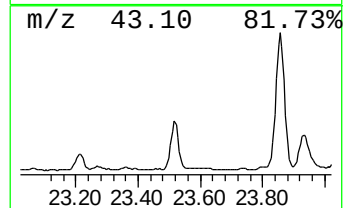
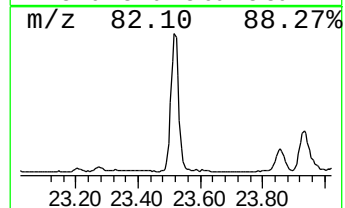
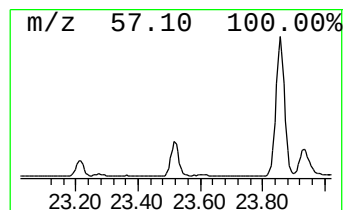
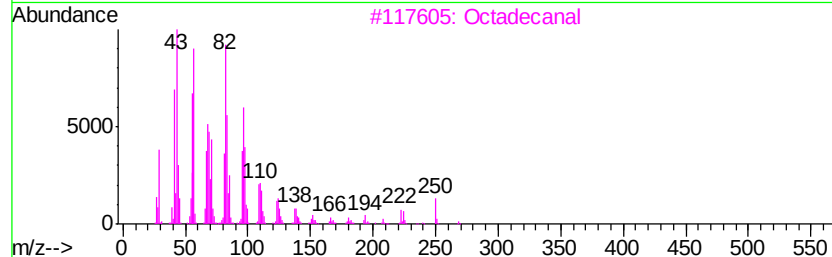
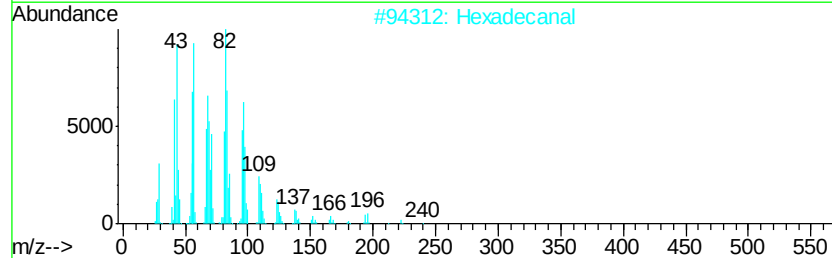
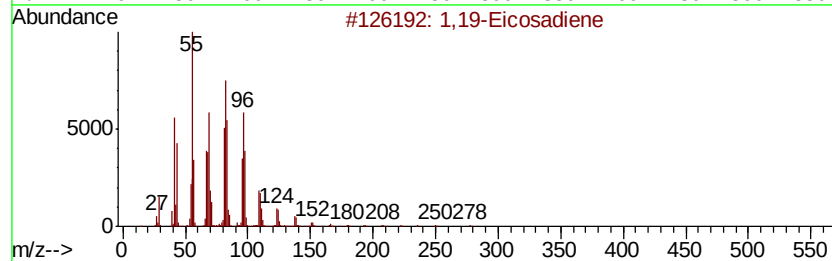
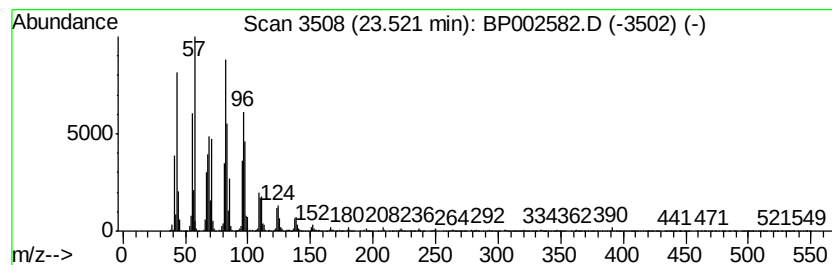
Instrument :  
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ClientSampleId :  
SP2-K

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 13 1,19-Eicosadiene Concentration Rank 12

| R.T.      | EstConc              | Area    | Relative to ISTD | R.T.        |      |
|-----------|----------------------|---------|------------------|-------------|------|
| 23.52     | 8.56 ng              | 1025550 | Perylene-d12     | 23.28       |      |
| Hit# of 5 | Tentative ID         | MW      | MolForm          | CAS#        | Qual |
| 1         | 1,19-Eicosadiene     | 278     | C20H38           | 014811-95-1 | 95   |
| 2         | Hexadecanal          | 240     | C16H32O          | 000629-80-1 | 93   |
| 3         | Octadecanal          | 268     | C18H36O          | 000638-66-4 | 91   |
| 4         | Oxirane, hexadecyl-  | 268     | C18H36O          | 007390-81-0 | 91   |
| 5         | Oxirane, heptadecyl- | 282     | C19H38O          | 067860-04-2 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP063020\  
Data File : BP002582.D  
Acq On : 30 Jun 2020 19:10  
Operator : CG/JU  
Sample : L3153-40  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SP2-K

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

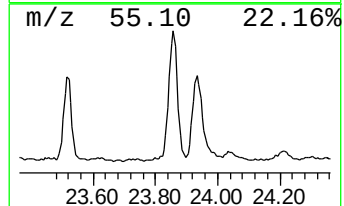
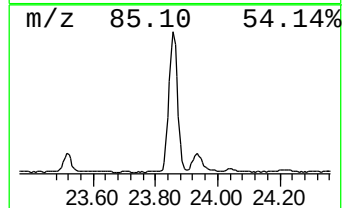
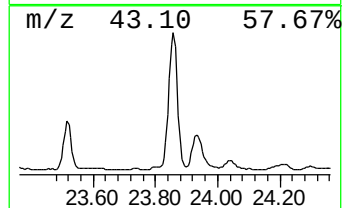
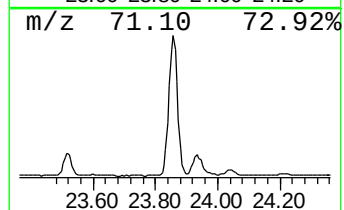
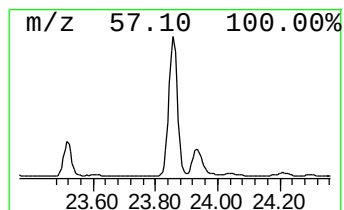
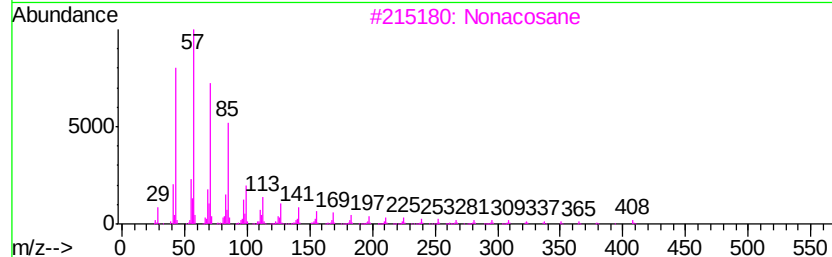
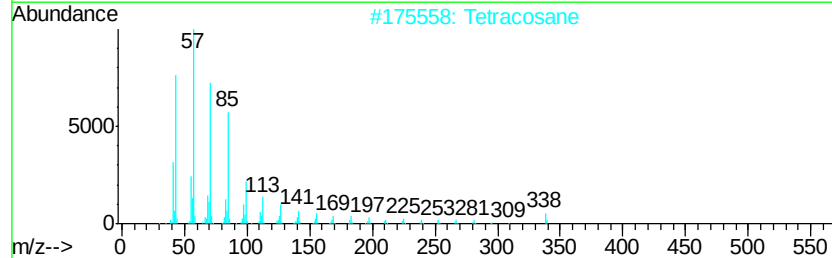
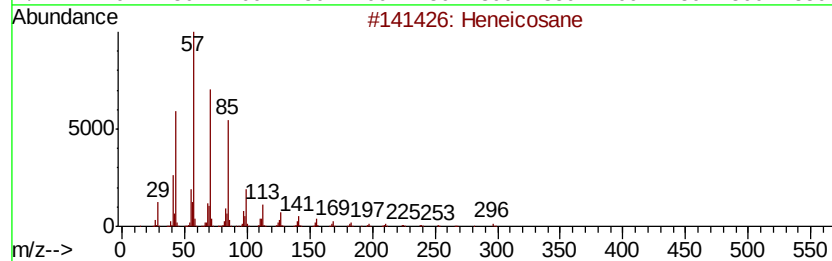
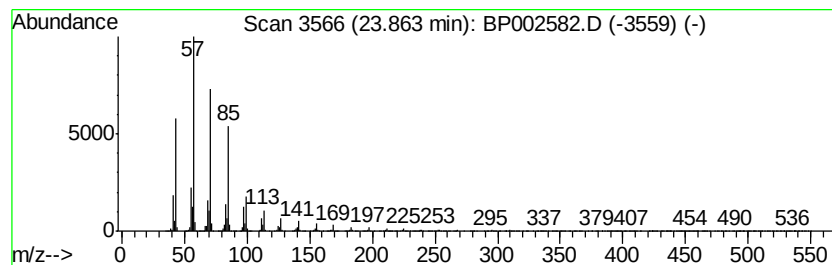
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 14 Heneicosane Concentration Rank 4

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 23.86 | 18.59 ng | 2227720 | Perylene-d12     | 23.28 |

| Hit# of | 5                  | Tentative ID | MW  | MolForm | CAS#         | Qual |
|---------|--------------------|--------------|-----|---------|--------------|------|
| 1       | Heneicosane        |              | 296 | C21H44  | 000629-94-7  | 97   |
| 2       | Tetracosane        |              | 338 | C24H50  | 000646-31-1  | 96   |
| 3       | Nonacosane         |              | 408 | C29H60  | 000630-03-5  | 95   |
| 4       | Triacontane        |              | 422 | C30H62  | 000638-68-6  | 93   |
| 5       | 2-methyloctacosane |              | 408 | C29H60  | 1000376-72-8 | 91   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP063020\  
Data File : BP002582.D  
Acq On : 30 Jun 2020 19:10  
Operator : CG/JU  
Sample : L3153-40  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SP2-K

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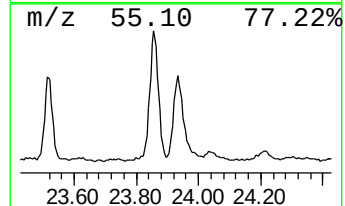
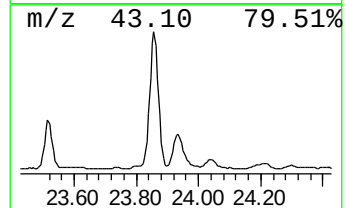
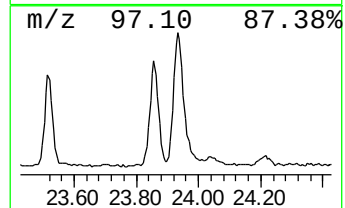
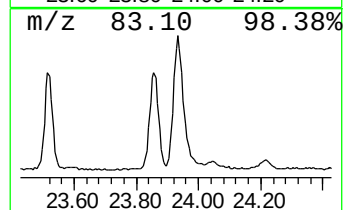
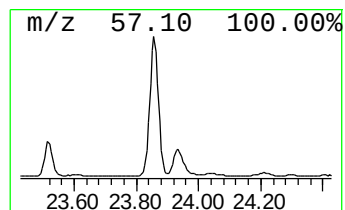
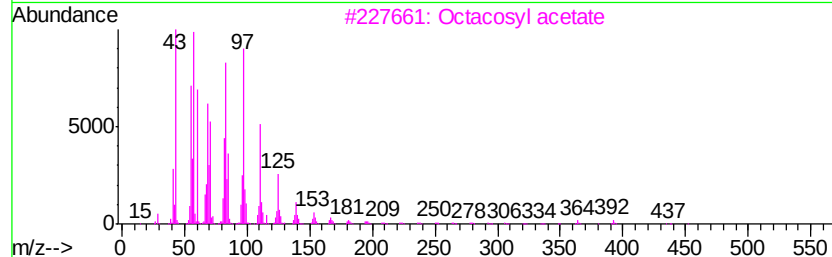
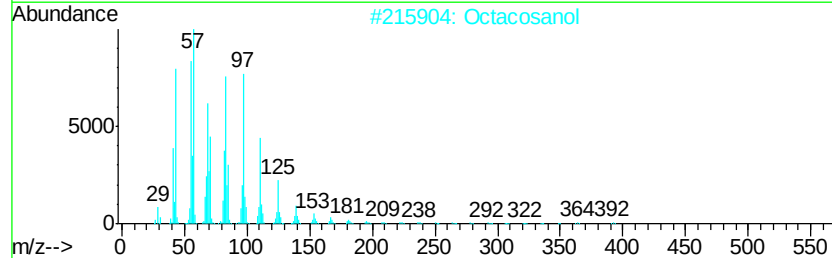
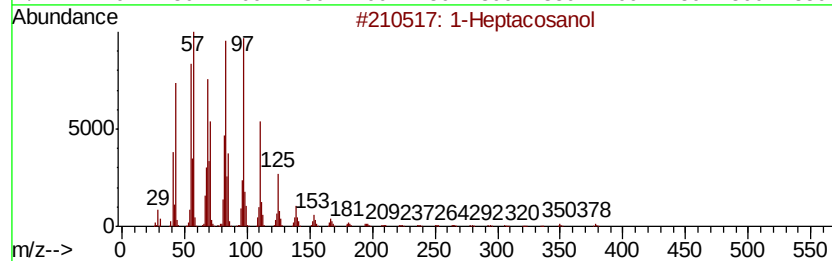
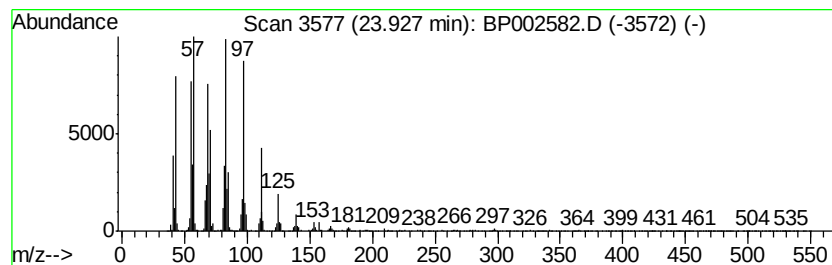
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TIC Integration Parameters: LSCINT.P

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Peak Number 15 1-Heptacosanol Concentration Rank 13

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 23.93 | 8.06 ng | 965745 | Perylene-d12     | 23.28 |

| Hit# | of | Tentative ID                   | MW  | MolForm    | CAS#         | Qual |
|------|----|--------------------------------|-----|------------|--------------|------|
| 1    | 5  | 1-Heptacosanol                 | 396 | C27H56O    | 002004-39-9  | 93   |
| 2    |    | Octacosanol                    | 410 | C28H58O    | 000557-61-9  | 89   |
| 3    |    | Octacosyl acetate              | 452 | C30H60O2   | 018206-97-8  | 86   |
| 4    |    | 1-Heneicosanol                 | 312 | C21H44O    | 015594-90-8  | 81   |
| 5    |    | Tetracosyl heptafluorobutyrate | 550 | C28H49F7O2 | 1000351-83-7 | 78   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP063020\  
Data File : BP002582.D  
Acq On : 30 Jun 2020 19:10  
Operator : CG/JU  
Sample : L3153-40  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SP2-K

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

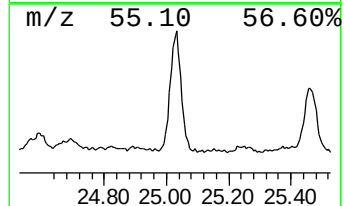
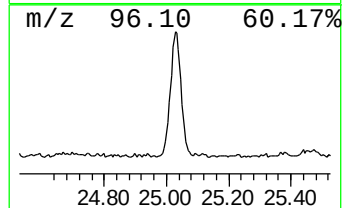
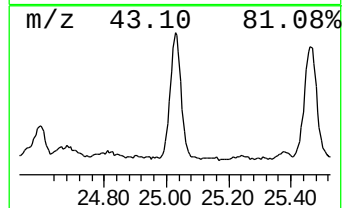
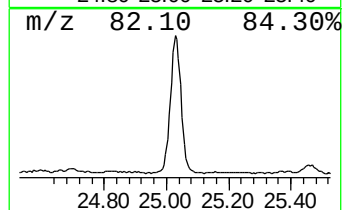
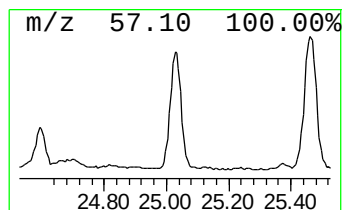
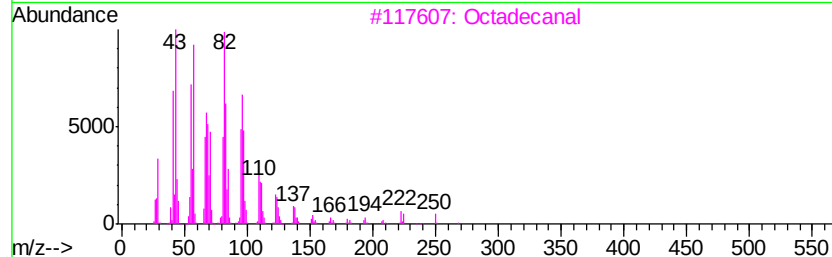
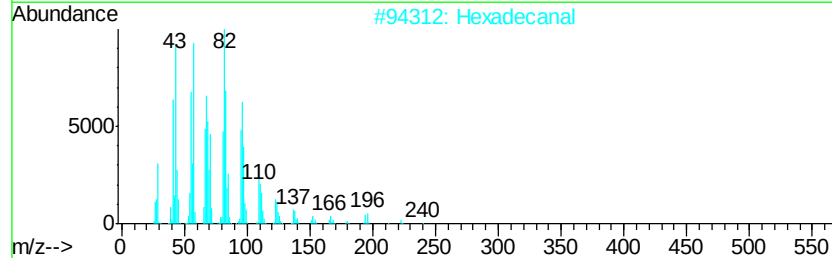
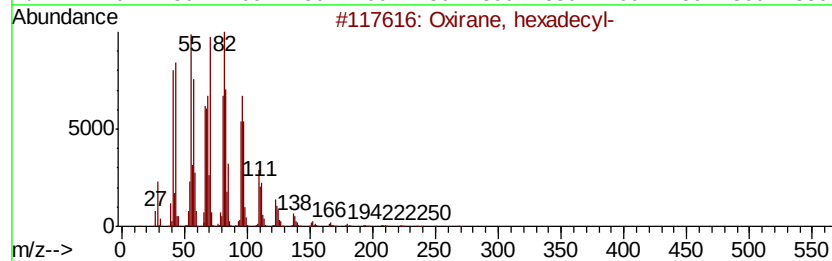
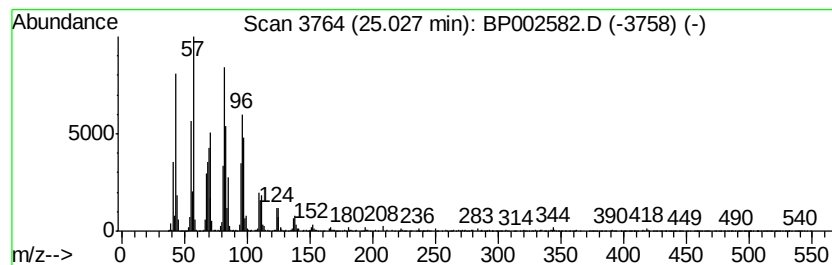
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 16 Oxirane, hexadecyl- Concentration Rank 8

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 25.03 | 10.36 ng | 1241920 | Perylene-d12     | 23.28 |

| Hit# | of | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------|-----|---------|-------------|------|
| 1    | 5  | Oxirane, hexadecyl-  | 268 | C18H36O | 007390-81-0 | 91   |
| 2    |    | Hexadecanal          | 240 | C16H32O | 000629-80-1 | 89   |
| 3    |    | Octadecanal          | 268 | C18H36O | 000638-66-4 | 87   |
| 4    |    | Oxirane, heptadecyl- | 282 | C19H38O | 067860-04-2 | 87   |
| 5    |    | Pentadecanal         | 226 | C15H30O | 002765-11-9 | 87   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP063020\  
Data File : BP002582.D  
Acq On : 30 Jun 2020 19:10  
Operator : CG/JU  
Sample : L3153-40  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SP2-K

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

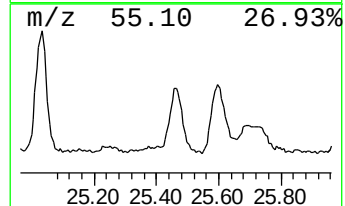
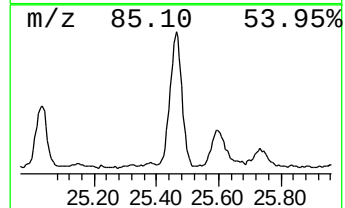
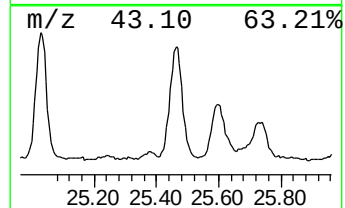
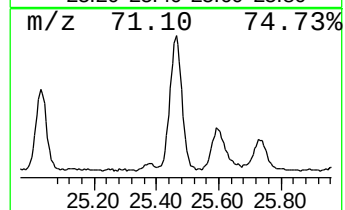
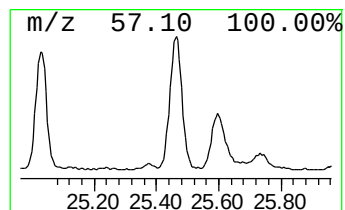
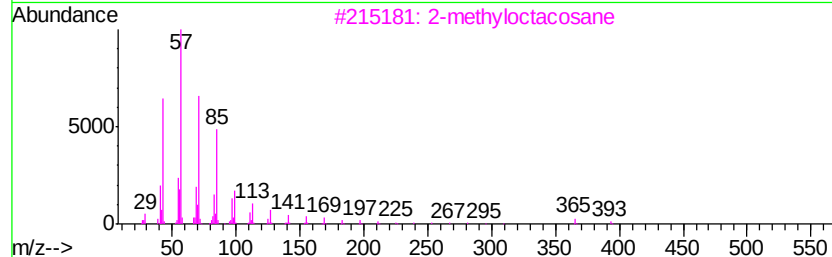
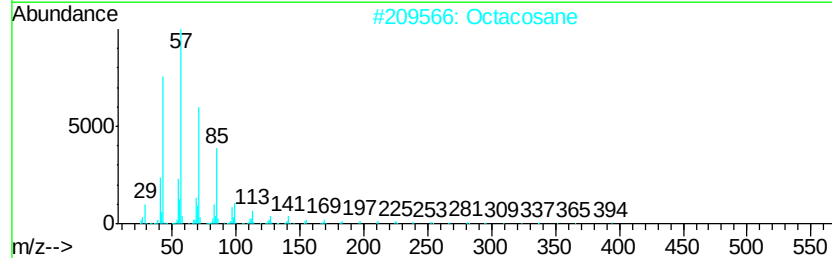
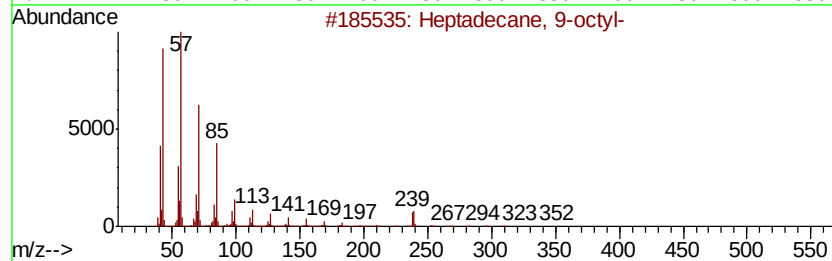
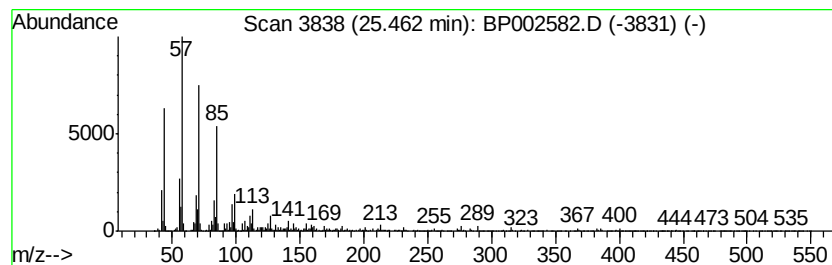
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 17 Heptadecane, 9-octyl- Concentration Rank 11

| R.T.  | EstConc | Area    | Relative to ISTD | R.T.  |
|-------|---------|---------|------------------|-------|
| 25.46 | 8.69 ng | 1041540 | Perylene-d12     | 23.28 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#         | Qual |
|------|----|---|-----------------------|-----|---------|--------------|------|
| 1    |    |   | Heptadecane, 9-octyl- | 352 | C25H52  | 007225-64-1  | 97   |
| 2    |    |   | Octacosane            | 394 | C28H58  | 000630-02-4  | 91   |
| 3    |    |   | 2-methyloctacosane    | 408 | C29H60  | 1000376-72-8 | 91   |
| 4    |    |   | Heneicosane           | 296 | C21H44  | 000629-94-7  | 91   |
| 5    |    |   | Tetracosane           | 338 | C24H50  | 000646-31-1  | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP063020\  
Data File : BP002582.D  
Acq On : 30 Jun 2020 19:10  
Operator : CG/JU  
Sample : L3153-40  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SP2-K

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

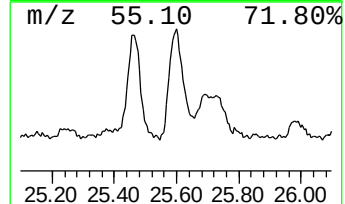
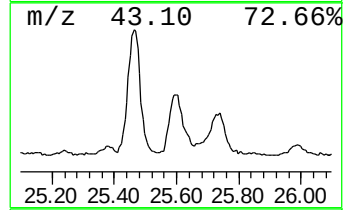
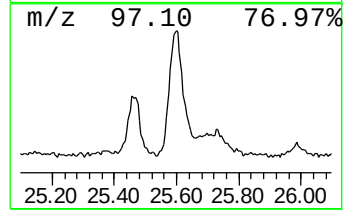
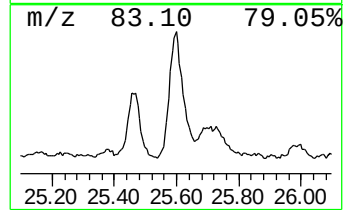
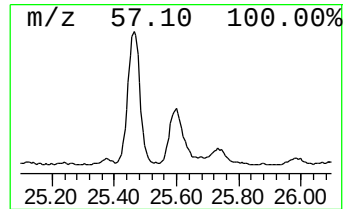
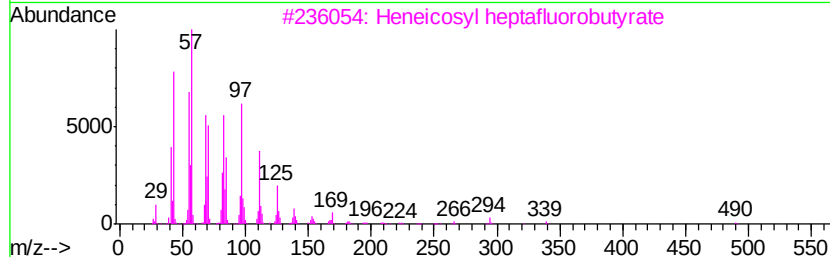
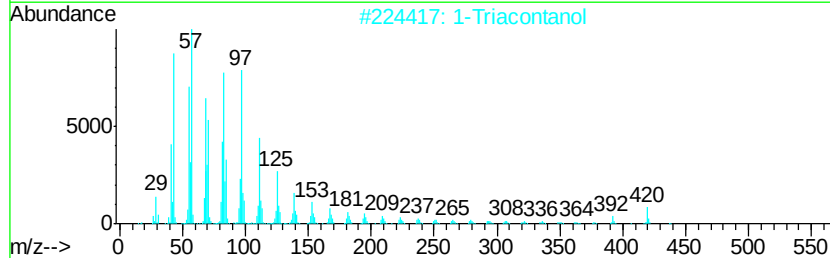
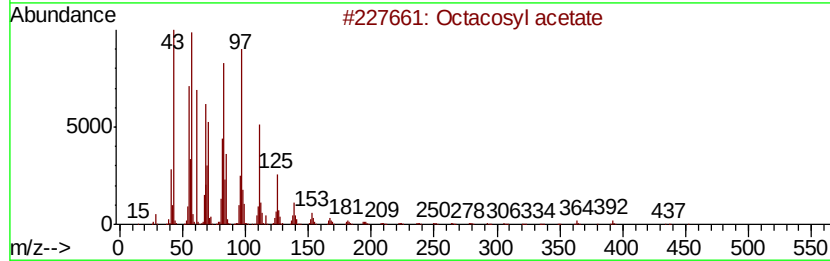
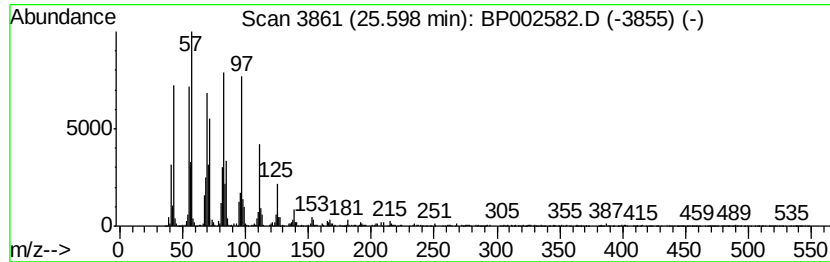
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 Octacosyl acetate Concentration Rank 17

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 25.60 | 4.36 ng | 522538 | Perylene-d12     | 23.28 |

| Hit# | of | Tentative ID                   | MW  | MolForm    | CAS#         | Qual |
|------|----|--------------------------------|-----|------------|--------------|------|
| 1    | 5  | Octacosyl acetate              | 452 | C30H60O2   | 018206-97-8  | 99   |
| 2    |    | 1-Triacontanol                 | 438 | C30H62O    | 000593-50-0  | 96   |
| 3    |    | Heneicosyl heptafluorobutvrate | 508 | C25H43F7O2 | 1000351-83-8 | 94   |
| 4    |    | 1-Heptacosanol                 | 396 | C27H56O    | 002004-39-9  | 93   |
| 5    |    | Heptadecyl heptafluorobutyrate | 452 | C21H35F7O2 | 959085-66-6  | 93   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP063020\  
Data File : BP002582.D  
Acq On : 30 Jun 2020 19:10  
Operator : CG/JU  
Sample : L3153-40  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

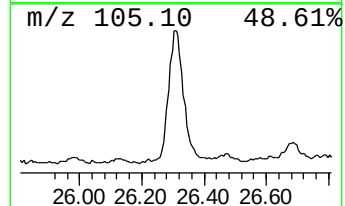
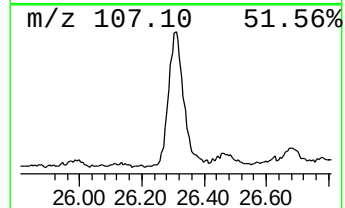
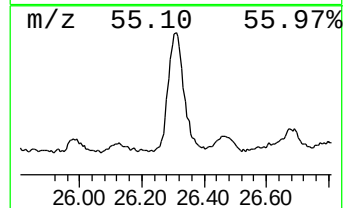
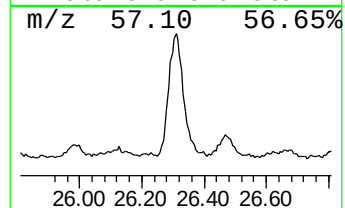
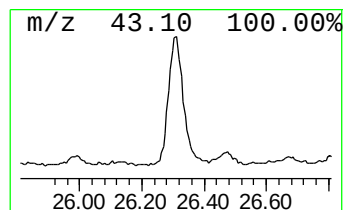
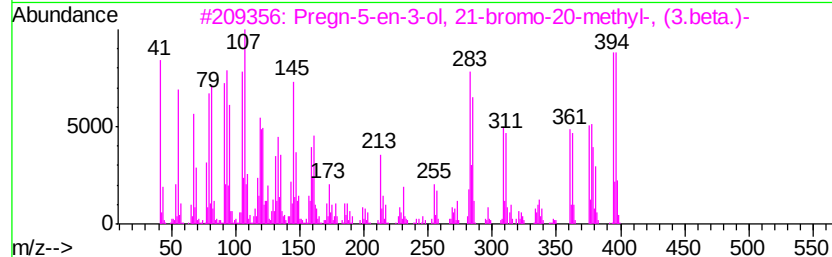
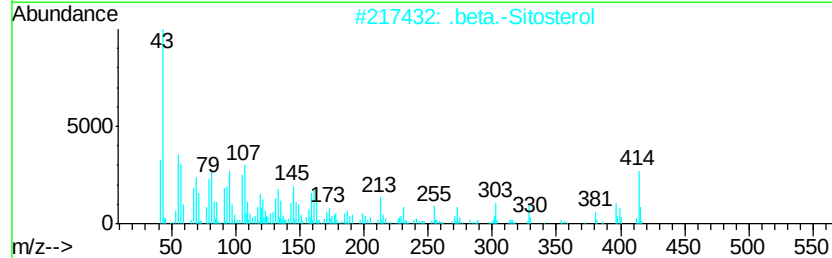
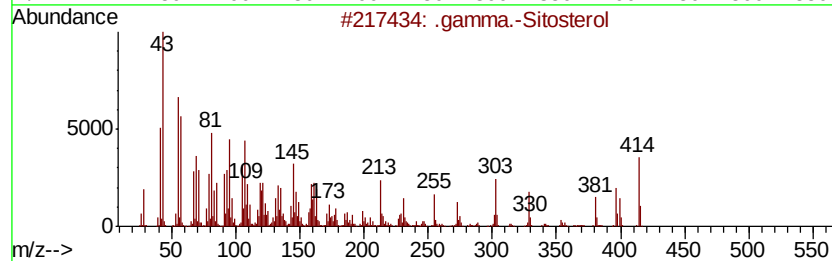
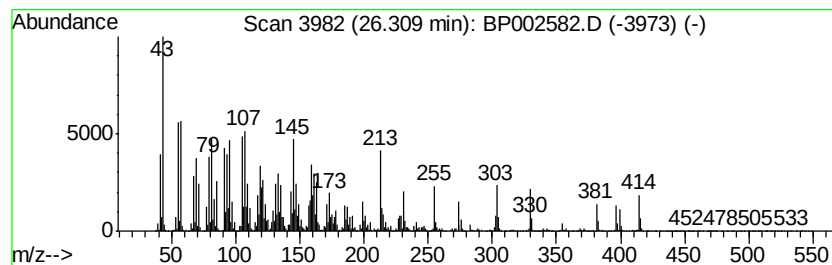
Instrument :  
BNA\_P  
ClientSampleId :  
SP2-K

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.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 19 .gamma.-Sitosterol Concentration Rank 3

| R.T.      | EstConc                             | Area    | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|---------|------------------|-------------|------|
| 26.31     | 22.20 ng                            | 2660510 | Perylene-d12     | 23.28       |      |
| Hit# of 5 | Tentative ID                        | MW      | MolForm          | CAS#        | Qual |
| 1         | .gamma.-Sitosterol                  | 414     | C29H50O          | 000083-47-6 | 99   |
| 2         | .beta.-Sitosterol                   | 414     | C29H50O          | 000083-46-5 | 94   |
| 3         | Pregn-5-en-3-ol, 21-bromo-20-met... | 394     | C22H35BrO        | 055103-80-5 | 93   |
| 4         | Campesterol                         | 400     | C28H48O          | 000474-62-4 | 47   |
| 5         | Ergost-5-en-3-ol, (3.beta.)-        | 400     | C28H48O          | 004651-51-8 | 45   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP063020\  
Data File : BP002582.D  
Acq On : 30 Jun 2020 19:10  
Operator : CG/JU  
Sample : L3153-40  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SP2-K

Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

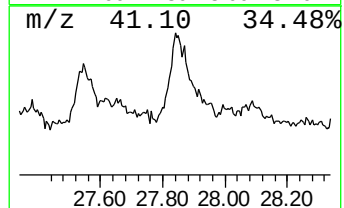
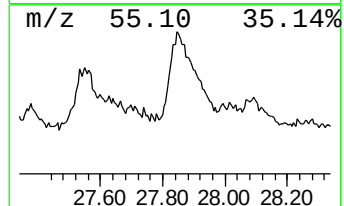
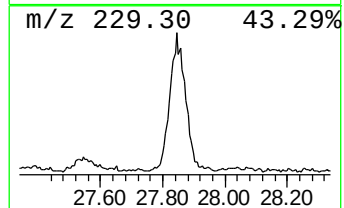
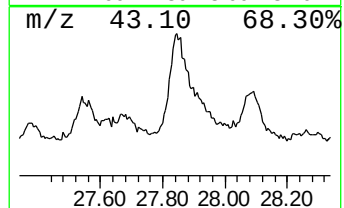
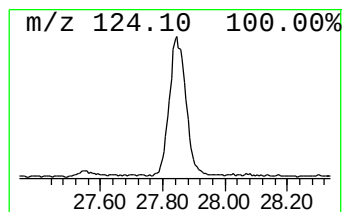
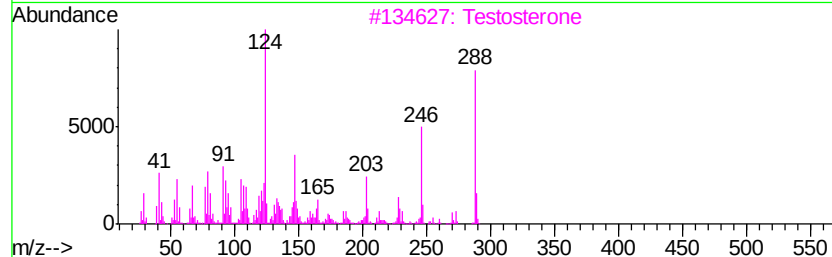
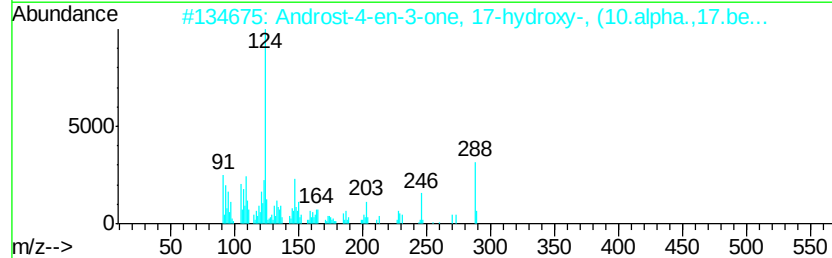
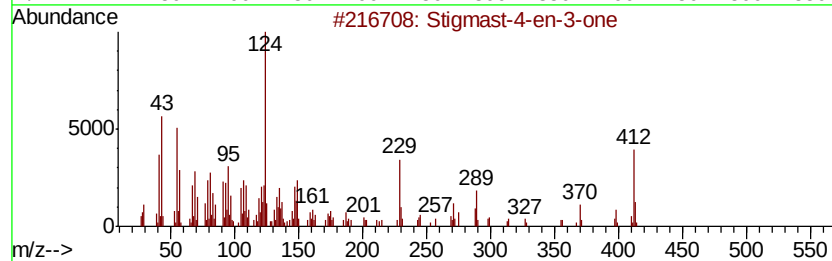
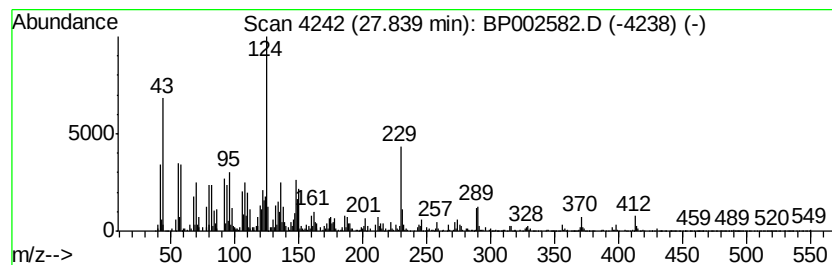
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 Stigmast-4-en-3-one Concentration Rank 14

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 27.84 | 7.46 ng | 894160 | Perylene-d12     | 23.28 |

| Hit# | of | Tentative ID                         | MW  | MolForm    | CAS#         | Qual |
|------|----|--------------------------------------|-----|------------|--------------|------|
| 1    | 5  | Stigmast-4-en-3-one                  | 412 | C29H48O    | 001058-61-3  | 93   |
| 2    |    | Androst-4-en-3-one, 17-hydroxy-, ... | 288 | C19H28O2   | 000604-39-7  | 64   |
| 3    |    | Testosterone                         | 288 | C19H28O2   | 000058-22-0  | 55   |
| 4    |    | Hvoscymamine                         | 289 | C17H23NO3  | 000101-31-5  | 42   |
| 5    |    | Carbonic acid, bis-(2,2,6,6-tetr...  | 370 | C19H34N2O5 | 1000188-97-0 | 41   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_P\DATA\BP063020\  
 Data File : BP002582.D  
 Acq On : 30 Jun 2020 19:10  
 Operator : CG/JU  
 Sample : L3153-40  
 Misc :  
 ALS Vial : 12 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SP2-K

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_P\METHODS\8270-BP062920.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 |
| Safrrole             | 11.82 | 9.0 ng  |       | 314920   | 2                     | 10.35 | 698444  | 20.0 |
| 1H-Cyclopenta[1,3... | 13.11 | 3.5 ng  |       | 250024   | 3                     | 14.22 | 1421490 | 20.0 |
| o,p'-DDE             | 19.49 | 6.3 ng  |       | 677300   | 5                     | 21.09 | 2160860 | 20.0 |
| 2-Propenal, 3-(4-... | 19.83 | 3.3 ng  |       | 356429   | 5                     | 21.09 | 2160860 | 20.0 |
| unknown20.50         | 20.50 | 4.1 ng  |       | 440419   | 5                     | 21.09 | 2160860 | 20.0 |
| 3-Butenoic acid, ... | 20.66 | 97.0 ng |       | 10478400 | 5                     | 21.09 | 2160860 | 20.0 |
| Behenic alcohol      | 20.83 | 11.9 ng |       | 1291190  | 5                     | 21.09 | 2160860 | 20.0 |
| Heptadecyl hepta...  | 21.70 | 14.6 ng |       | 1580570  | 5                     | 21.09 | 2160860 | 20.0 |
| Oxirane, heptadecyl- | 22.37 | 4.4 ng  |       | 524344   | 6                     | 23.28 | 2397200 | 20.0 |
| Octacosane           | 22.65 | 14.8 ng |       | 1770530  | 6                     | 23.28 | 2397200 | 20.0 |
| Cyclotetracosane     | 22.69 | 9.1 ng  |       | 1095930  | 6                     | 23.28 | 2397200 | 20.0 |
| 1,19-Eicosadiene     | 23.52 | 8.6 ng  |       | 1025550  | 6                     | 23.28 | 2397200 | 20.0 |
| Heneicosane          | 23.86 | 18.6 ng |       | 2227720  | 6                     | 23.28 | 2397200 | 20.0 |
| 1-Heptacosanol       | 23.93 | 8.1 ng  |       | 965745   | 6                     | 23.28 | 2397200 | 20.0 |
| Oxirane, hexadecyl-  | 25.03 | 10.4 ng |       | 1241920  | 6                     | 23.28 | 2397200 | 20.0 |
| Heptadecane, 9-oc... | 25.46 | 8.7 ng  |       | 1041540  | 6                     | 23.28 | 2397200 | 20.0 |
| Octacosyl acetate    | 25.60 | 4.4 ng  |       | 522538   | 6                     | 23.28 | 2397200 | 20.0 |
| .gamma.-Sitosterol   | 26.31 | 22.2 ng |       | 2660510  | 6                     | 23.28 | 2397200 | 20.0 |
| Stigmast-4-en-3-one  | 27.84 | 7.5 ng  |       | 894160   | 6                     | 23.28 | 2397200 | 20.0 |