

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP070220\  
 Data File : BP002638.D  
 Acq On : 02 Jul 2020 19:21  
 Operator : CG/JU  
 Sample : L3188-01 5X  
 Misc :  
 ALS Vial : 20 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 TR-04-070120

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       | 386           | 392         | 406          | rBV      | 236216         | 406373        | 1.13%           | 0.322%        |
| 2         | 6.764       | 650           | 659         | 671          | rBV      | 258293         | 485966        | 1.35%           | 0.386%        |
| 3         | 7.569       | 788           | 796         | 807          | rBV      | 607131         | 997946        | 2.78%           | 0.792%        |
| 4         | 7.887       | 840           | 850         | 862          | rBV      | 289712         | 485969        | 1.35%           | 0.386%        |
| 5         | 8.716       | 976           | 991         | 1003         | rBV      | 185024         | 390227        | 1.09%           | 0.310%        |
| 6         | 10.340      | 1259          | 1267        | 1281         | rBV      | 858824         | 1538107       | 4.28%           | 1.221%        |
| 7         | 12.834      | 1685          | 1691        | 1701         | rBV      | 571678         | 871094        | 2.43%           | 0.691%        |
| 8         | 14.216      | 1920          | 1926        | 1933         | rBV2     | 1360217        | 2023748       | 5.63%           | 1.606%        |
| 9         | 15.722      | 2176          | 2182        | 2192         | rVB      | 368179         | 589051        | 1.64%           | 0.467%        |
| 10        | 16.975      | 2388          | 2395        | 2400         | rBV      | 1760868        | 2614556       | 7.28%           | 2.075%        |
| 11        | 17.016      | 2400          | 2402        | 2409         | rVB      | 317395         | 390565        | 1.09%           | 0.310%        |
| 12        | 17.557      | 2491          | 2494        | 2502         | rVB2     | 397276         | 492813        | 1.37%           | 0.391%        |
| 13        | 17.692      | 2511          | 2517        | 2522         | rBV      | 10486317       | 16657246      | 46.38%          | 13.219%       |
| 14        | 18.216      | 2597          | 2606        | 2614         | rVB5     | 254431         | 666463        | 1.86%           | 0.529%        |
| 15        | 18.357      | 2626          | 2630        | 2634         | rVB2     | 367528         | 417407        | 1.16%           | 0.331%        |
| 16        | 18.822      | 2700          | 2709        | 2712         | rVV3     | 12993732       | 35917778      | 100.00%         | 28.504%       |
| 17        | 18.857      | 2712          | 2715        | 2722         | rVV3     | 12311136       | 22302017      | 62.09%          | 17.699%       |
| 18        | 18.916      | 2722          | 2725        | 2729         | rVV      | 836846         | 1070256       | 2.98%           | 0.849%        |
| 19        | 18.969      | 2729          | 2734        | 2738         | rVV      | 7234769        | 9879945       | 27.51%          | 7.841%        |
| 20        | 19.010      | 2738          | 2741        | 2745         | rVB      | 518006         | 617615        | 1.72%           | 0.490%        |
| 21        | 19.316      | 2786          | 2793        | 2797         | rBV3     | 382973         | 816610        | 2.27%           | 0.648%        |
| 22        | 19.363      | 2797          | 2801        | 2805         | rVV      | 465246         | 651437        | 1.81%           | 0.517%        |
| 23        | 19.569      | 2830          | 2836        | 2840         | rBV      | 1422306        | 2068735       | 5.76%           | 1.642%        |
| 24        | 19.828      | 2876          | 2880        | 2882         | rBV2     | 417576         | 575104        | 1.60%           | 0.456%        |
| 25        | 19.851      | 2882          | 2884        | 2886         | rVV2     | 474735         | 524188        | 1.46%           | 0.416%        |
| 26        | 19.875      | 2886          | 2888        | 2890         | rVV      | 434220         | 530820        | 1.48%           | 0.421%        |
| 27        | 19.922      | 2892          | 2896        | 2905         | rVB2     | 1266087        | 2017418       | 5.62%           | 1.601%        |
| 28        | 20.033      | 2911          | 2915        | 2917         | rBV      | 1327004        | 1446197       | 4.03%           | 1.148%        |
| 29        | 20.063      | 2917          | 2920        | 2926         | rVV3     | 991699         | 1514994       | 4.22%           | 1.202%        |
| 30        | 20.122      | 2926          | 2930        | 2937         | rVV2     | 1011123        | 1453743       | 4.05%           | 1.154%        |
| 31        | 20.322      | 2960          | 2964        | 2971         | rVB2     | 1796286        | 2390966       | 6.66%           | 1.897%        |
| 32        | 20.633      | 3010          | 3017        | 3021         | rVB2     | 1051213        | 1999548       | 5.57%           | 1.587%        |
| 33        | 20.957      | 3069          | 3072        | 3083         | rVB      | 1541602        | 1700805       | 4.74%           | 1.350%        |
| 34        | 21.086      | 3088          | 3094        | 3098         | rVV2     | 2623111        | 3805854       | 10.60%          | 3.020%        |

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Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
TR-04-070120

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 | 21.122 | 3098 | 3100 | 3108 | rVB  | 375062  | 449909  | 1.25% | 0.357% |
| 36 | 21.827 | 3216 | 3220 | 3228 | rVB  | 549405  | 744641  | 2.07% | 0.591% |
| 37 | 22.645 | 3352 | 3359 | 3369 | rVB3 | 318297  | 987289  | 2.75% | 0.783% |
| 38 | 23.186 | 3447 | 3451 | 3460 | rVB2 | 273450  | 715198  | 1.99% | 0.568% |
| 39 | 23.280 | 3461 | 3467 | 3473 | rBV2 | 1530434 | 2801962 | 7.80% | 2.224% |

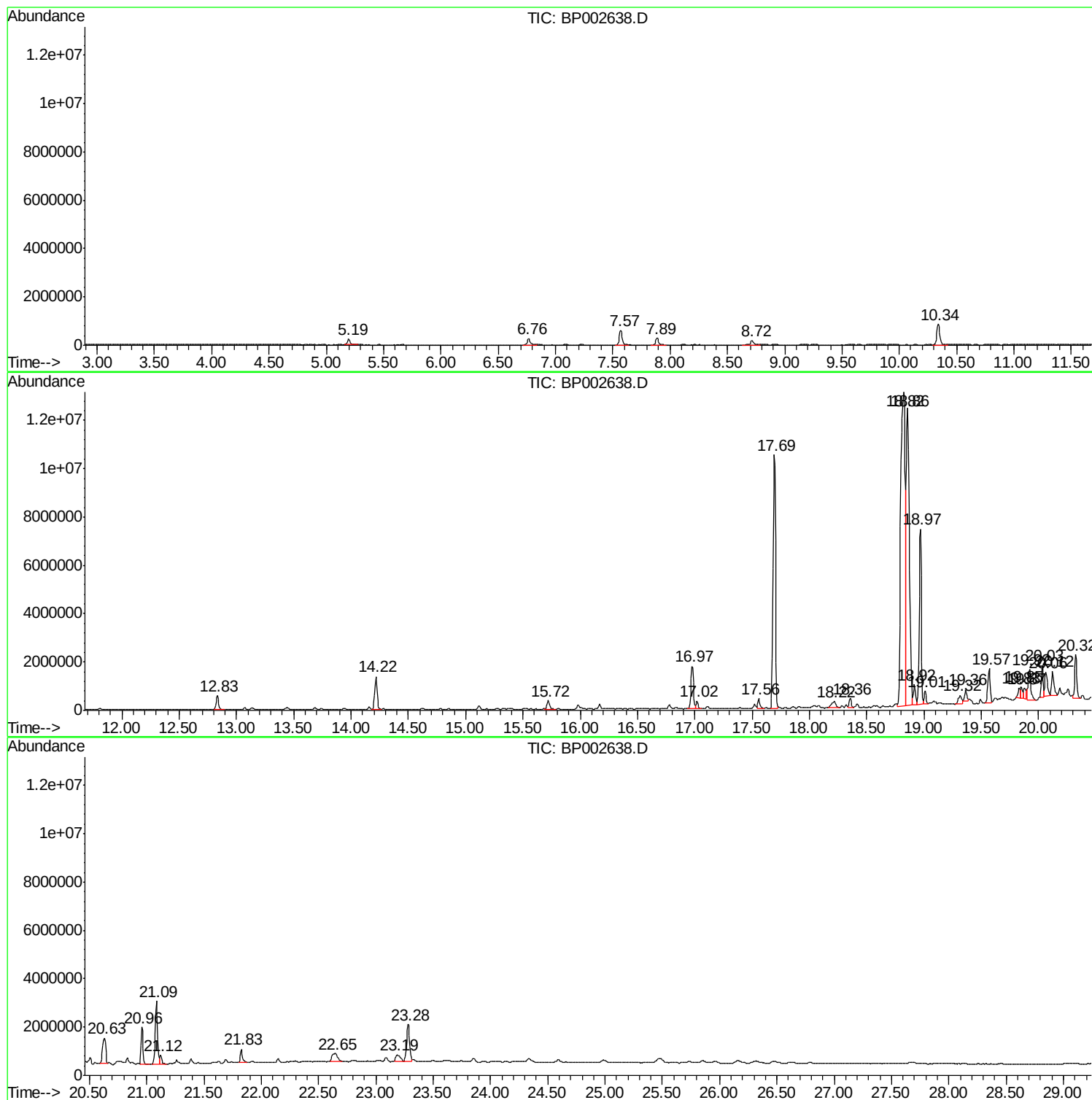
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP062920.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
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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

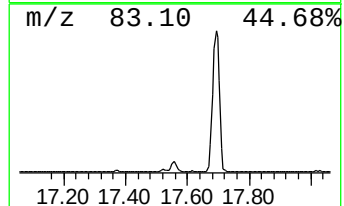
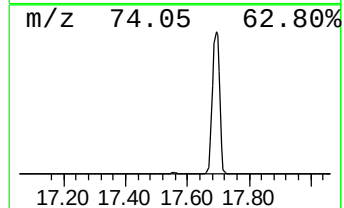
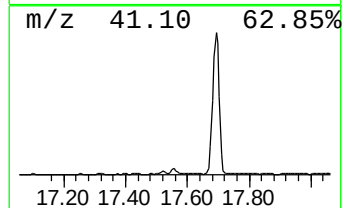
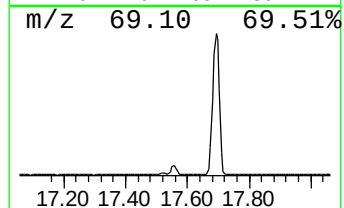
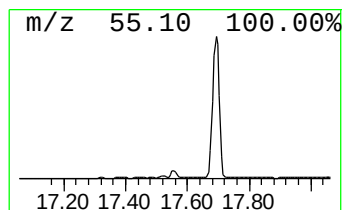
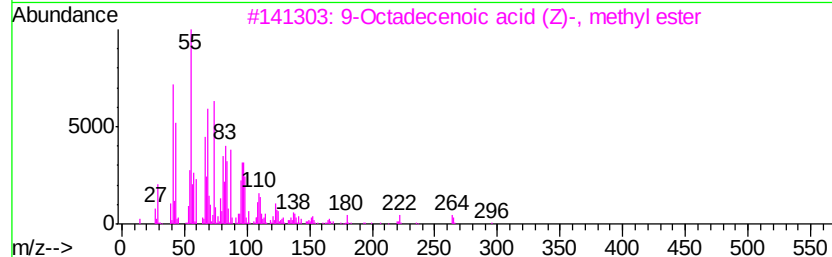
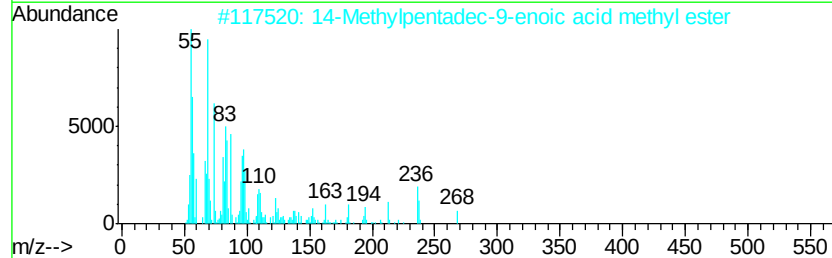
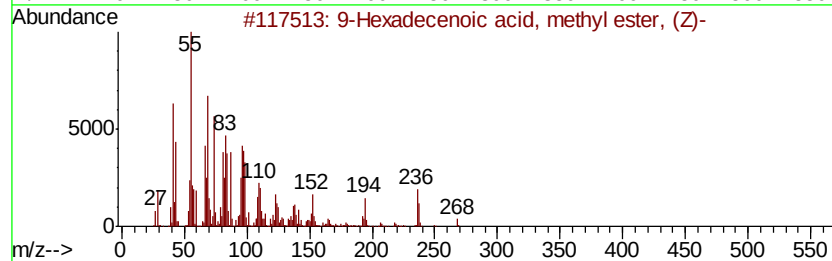
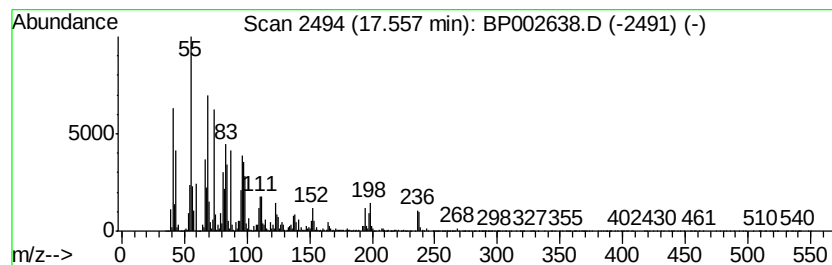
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 9-Hexadecenoic acid, methyl... Concentration Rank 15

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 17.56 | 3.77 ng | 492813 | Phenanthrene-d10 | 16.97 |

| Hit# of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|---------|---|-------------------------------------|-----|----------|--------------|------|
| 1       |   | 9-Hexadecenoic acid, methyl este... | 268 | C17H32O2 | 001120-25-8  | 99   |
| 2       |   | 14-Methylpentadec-9-enoic acid m... | 268 | C17H32O2 | 1000365-89-7 | 95   |
| 3       |   | 9-Octadecenoic acid (Z)-, methyl... | 296 | C19H36O2 | 000112-62-9  | 93   |
| 4       |   | Methyl hexadec-9-enoate             | 268 | C17H32O2 | 010030-74-7  | 91   |
| 5       |   | 13-Borabicyclo[7.3.0]tridecane, ... | 236 | C15H29BO | 1000156-41-7 | 83   |



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Misc :  
ALS Vial : 20 Sample Multiplier: 1

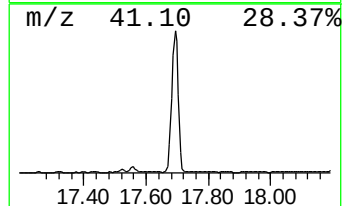
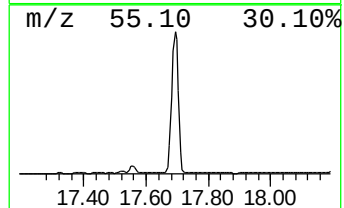
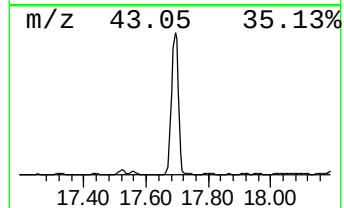
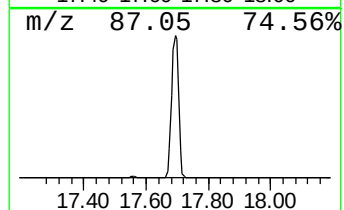
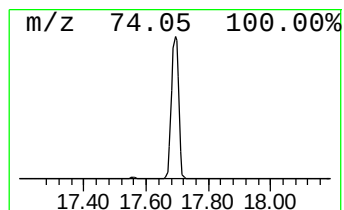
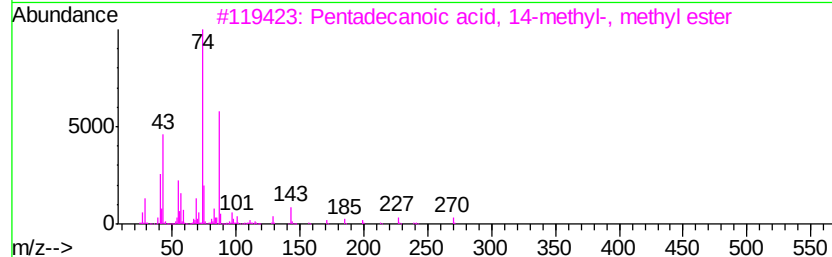
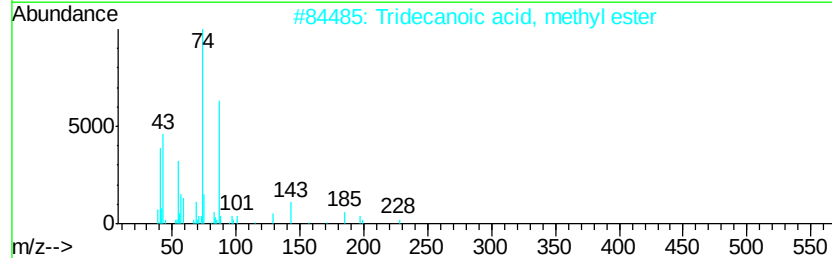
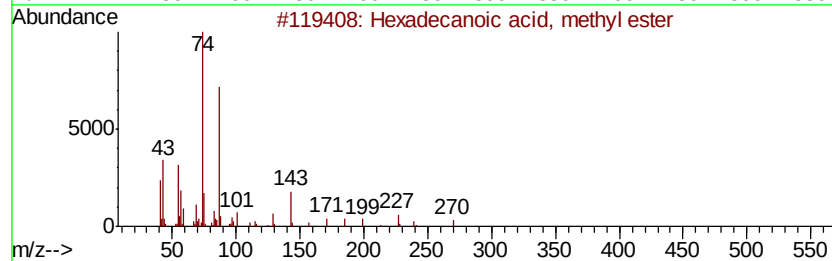
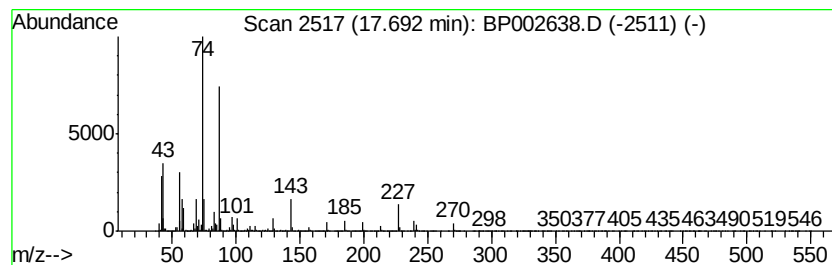
Instrument :  
BNA\_P  
ClientSampleId :  
TR-04-070120

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 Hexadecanoic acid, methyl e... Concentration Rank 3

| R.T.    | EstConc   | Area                                | Relative to ISTD | R.T.     |             |      |
|---------|-----------|-------------------------------------|------------------|----------|-------------|------|
| 17.69   | 127.42 ng | 16657200                            | Phenanthrene-d10 | 16.97    |             |      |
| Hit# of | 5         | Tentative ID                        | MW               | MolForm  | CAS#        | Qual |
| 1       |           | Hexadecanoic acid, methyl ester     | 270              | C17H34O2 | 000112-39-0 | 97   |
| 2       |           | Tridecanoic acid, methyl ester      | 228              | C14H28O2 | 001731-88-0 | 92   |
| 3       |           | Pentadecanoic acid, 14-methyl-, ... | 270              | C17H34O2 | 005129-60-2 | 91   |
| 4       |           | Methyl tetradecanoate               | 242              | C15H30O2 | 000124-10-7 | 87   |
| 5       |           | Hexadecanoic acid, 15-methyl-, m... | 284              | C18H36O2 | 006929-04-0 | 87   |



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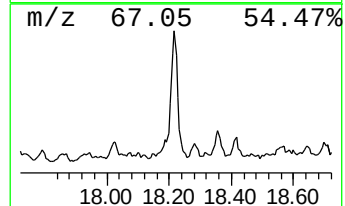
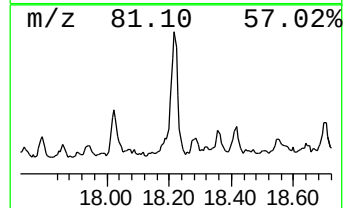
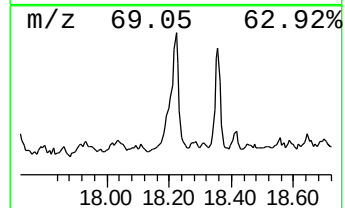
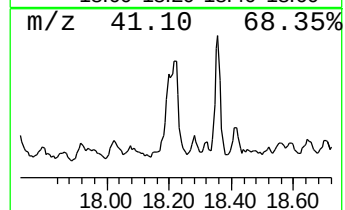
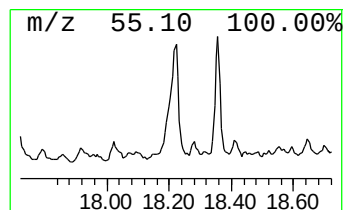
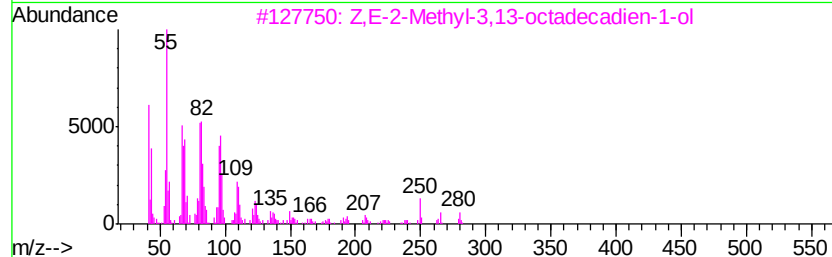
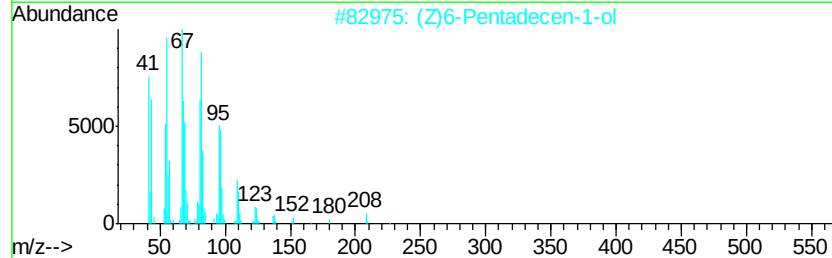
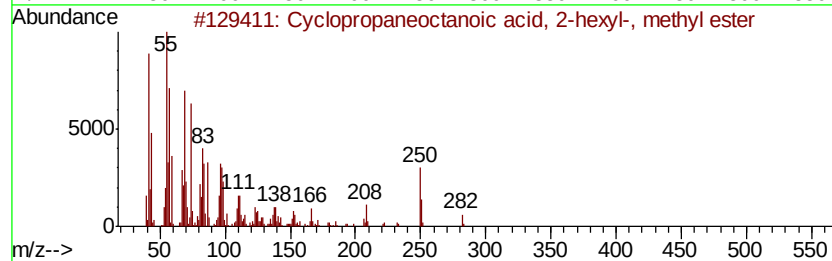
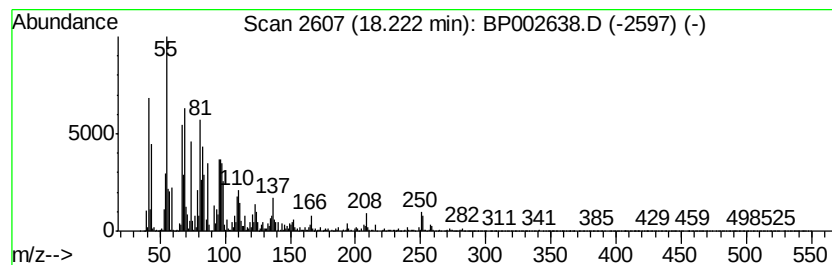
Instrument :  
BNA\_P  
ClientSampleId :  
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 4 Cyclopropaneoctanoic acid, ... Concentration Rank 13

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 18.22     | 5.10 ng                             | 666463 | Phenanthrene-d10 | 16.97        |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | Cyclopropaneoctanoic acid, 2-hex... | 282    | C18H34O2         | 010152-61-1  | 89   |
| 2         | (Z)6-Pentadecen-1-ol                | 226    | C15H30O          | 068797-95-5  | 64   |
| 3         | Z,E-2-Methyl-3,13-octadecadien-1-ol | 280    | C19H36O          | 1000131-10-5 | 64   |
| 4         | 5-Octadecyne                        | 250    | C18H34           | 071899-42-8  | 55   |
| 5         | Z,Z-2,15-Octadecadien-1-ol acetate  | 308    | C20H36O2         | 1000130-95-1 | 48   |



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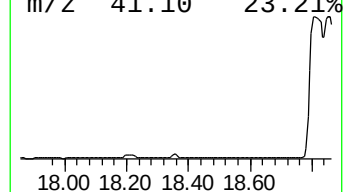
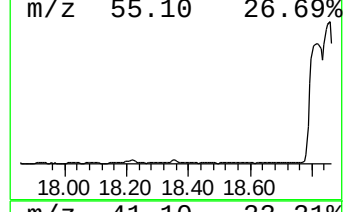
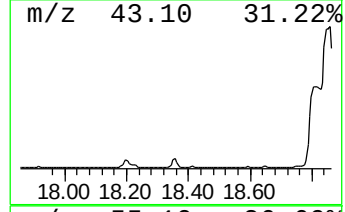
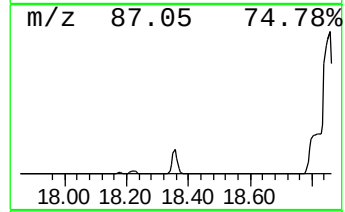
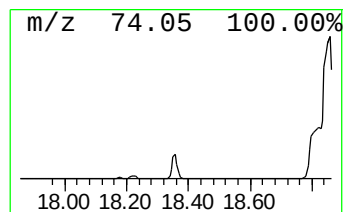
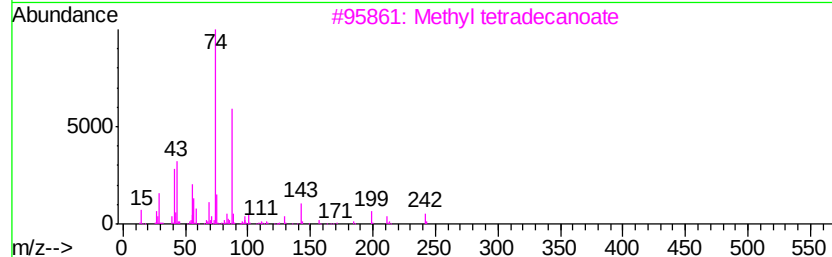
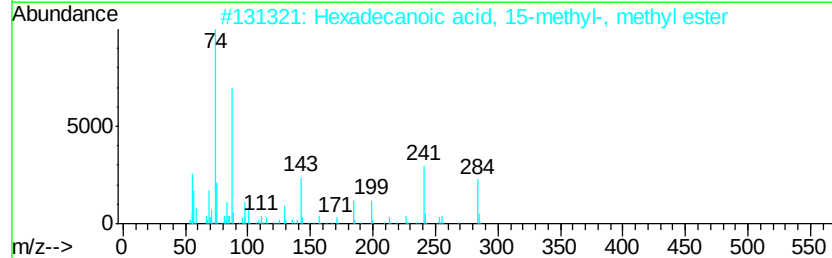
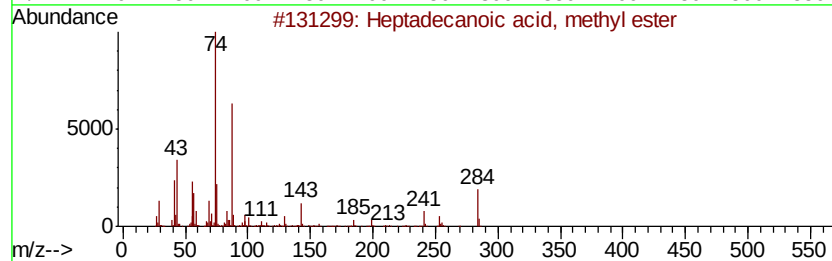
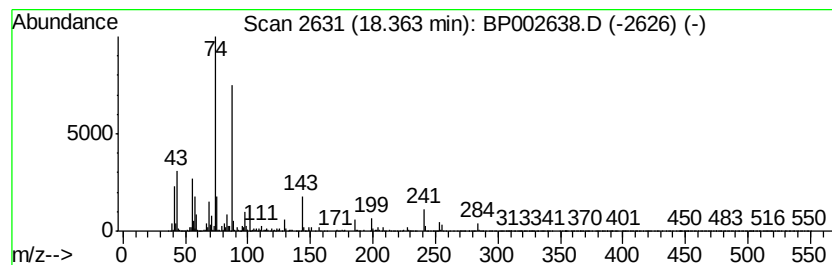
TIC Library : C:\DATABASE\NIST11.L  
 TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
 Peak Number 5 Heptadecanoic acid, methyl ... Concentration Rank 16

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 18.36 | 3.19 ng | 417407 | Phenanthrene-d10 | 16.97 |

| Hit# | of | Tentative ID                                | MW  | MolForm  | CAS#        | Qual |
|------|----|---------------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Heptadecanoic acid, methyl ester            | 284 | C18H36O2 | 001731-92-6 | 97   |
| 2    |    | Hexadecanoic acid, 15-methyl-, methyl ester | 284 | C18H36O2 | 006929-04-0 | 95   |
| 3    |    | Methyl tetradecanoate                       | 242 | C15H30O2 | 000124-10-7 | 91   |
| 4    |    | Tridecanoic acid, 12-methyl-, methyl ester  | 242 | C15H30O2 | 005129-58-8 | 90   |
| 5    |    | Hexadecanoic acid, methyl ester             | 270 | C17H34O2 | 000112-39-0 | 87   |



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 ClientSampleId :  
 TR-04-070120

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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

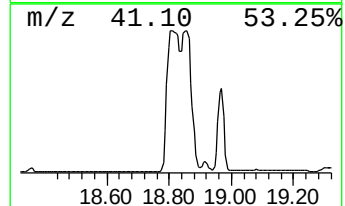
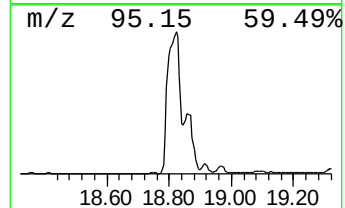
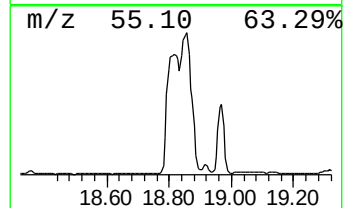
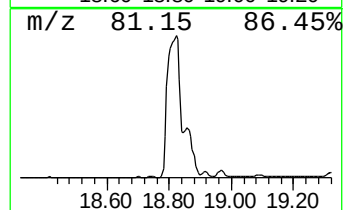
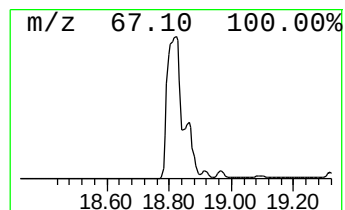
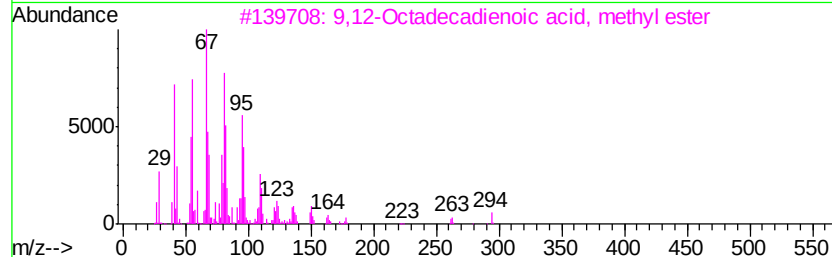
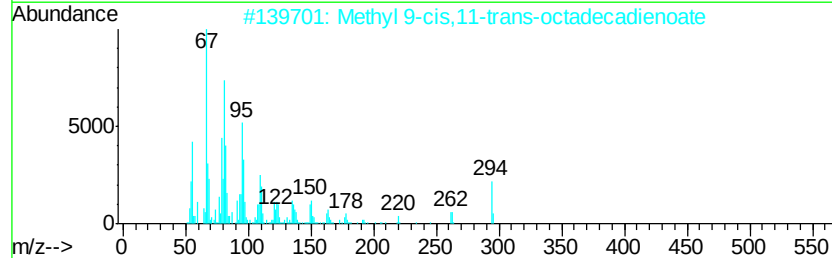
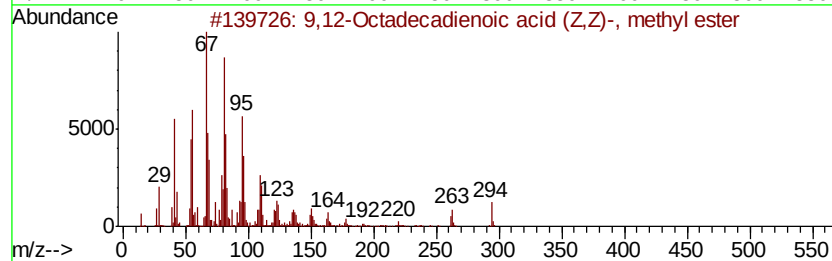
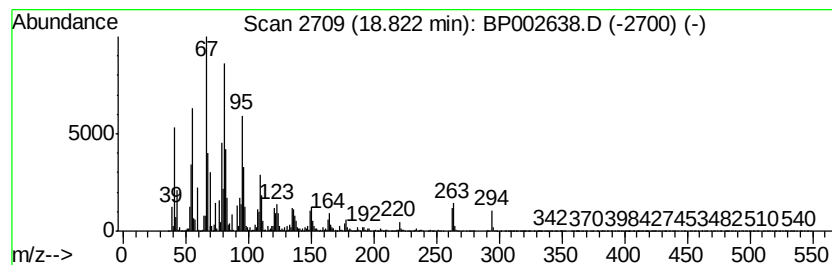
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 TIC Integration Parameters: LSCINT.P

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 Peak Number 6 9,12-Octadecadienoic acid (... Concentration Rank 1

| R.T.  | EstConc   | Area     | Relative to ISTD | R.T.  |
|-------|-----------|----------|------------------|-------|
| 18.82 | 274.75 ng | 35917800 | Phenanthrene-d10 | 16.97 |

| Hit# of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|---------|---|-------------------------------------|-----|----------|--------------|------|
| 1       |   | 9,12-Octadecadienoic acid (Z,Z)-... | 294 | C19H34O2 | 000112-63-0  | 99   |
| 2       |   | Methyl 9-cis,11-trans-octadecadi... | 294 | C19H34O2 | 1000336-44-0 | 99   |
| 3       |   | 9,12-Octadecadienoic acid, methy... | 294 | C19H34O2 | 002462-85-3  | 99   |
| 4       |   | 9,12-Octadecadienoic acid, methy... | 294 | C19H34O2 | 002566-97-4  | 99   |
| 5       |   | Methyl 10-trans,12-cis-octadecad... | 294 | C19H34O2 | 1000336-44-2 | 99   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP070220\  
 Data File : BP002638.D  
 Acq On : 02 Jul 2020 19:21  
 Operator : CG/JU  
 Sample : L3188-01 5X  
 Misc :  
 ALS Vial : 20 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 TR-04-070120

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

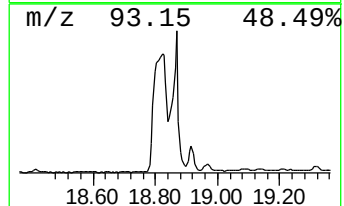
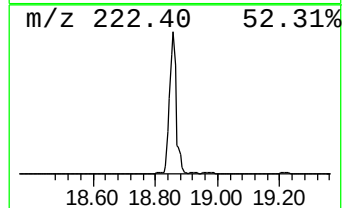
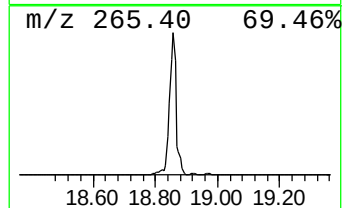
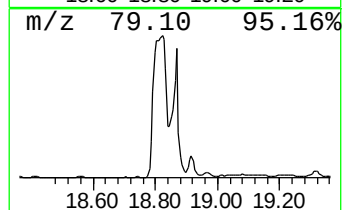
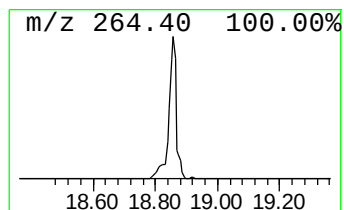
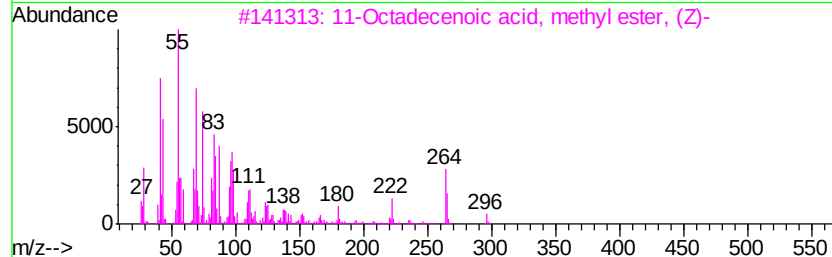
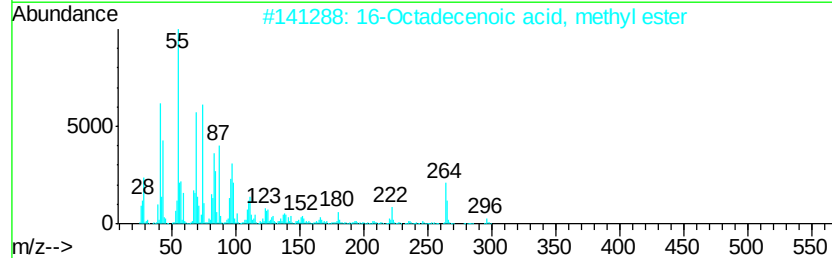
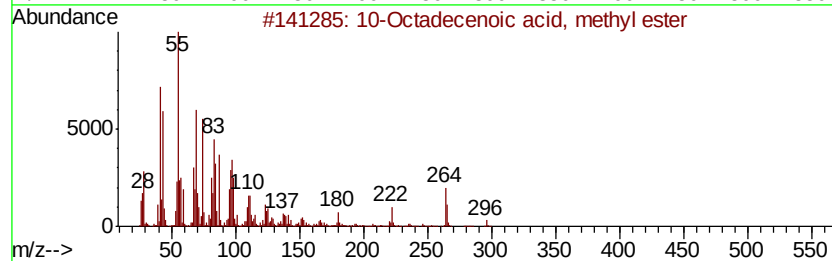
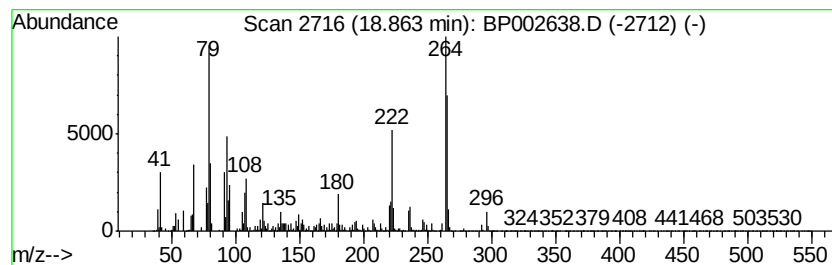
TIC Library : C:\DATABASE\NIST11.L  
 TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
 Peak Number 7 10-Octadecenoic acid, methyl ester Concentration Rank 2

| R.T.  | EstConc   | Area     | Relative to ISTD | R.T.  |
|-------|-----------|----------|------------------|-------|
| 18.86 | 170.60 ng | 22302000 | Phenanthrene-d10 | 16.97 |

| Hit# | of | Tentative ID                           | MW  | MolForm  | CAS#         | Qual |
|------|----|----------------------------------------|-----|----------|--------------|------|
| 1    | 5  | 10-Octadecenoic acid, methyl ester     | 296 | C19H36O2 | 013481-95-3  | 91   |
| 2    |    | 16-Octadecenoic acid, methyl ester     | 296 | C19H36O2 | 056554-49-5  | 91   |
| 3    |    | 11-Octadecenoic acid, methyl ester     | 296 | C19H36O2 | 001937-63-9  | 90   |
| 4    |    | cis-13-Octadecenoic acid, methyl ester | 296 | C19H36O2 | 1000333-58-3 | 86   |
| 5    |    | 14-Octadecenoic acid, methyl ester     | 296 | C19H36O2 | 056554-48-4  | 81   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP070220\  
Data File : BP002638.D  
Acq On : 02 Jul 2020 19:21  
Operator : CG/JU  
Sample : L3188-01 5X  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
TR-04-070120

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

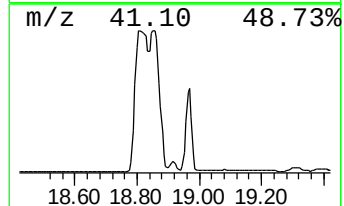
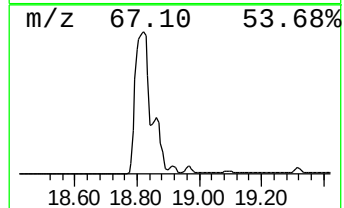
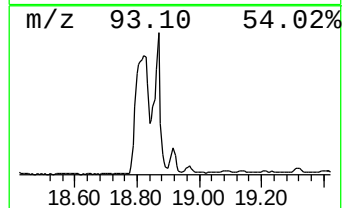
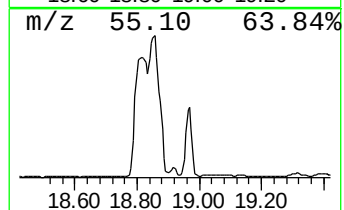
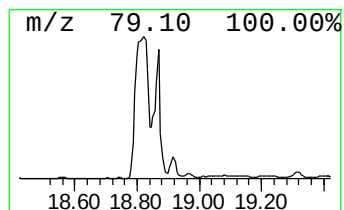
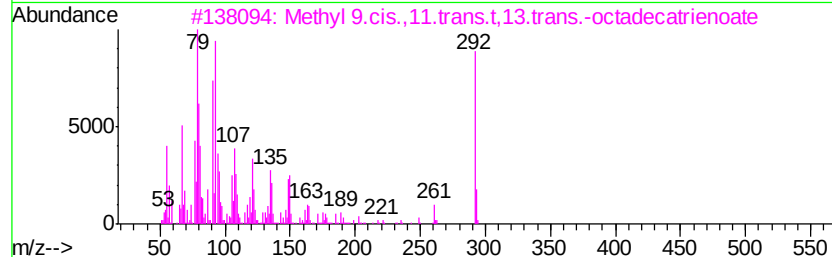
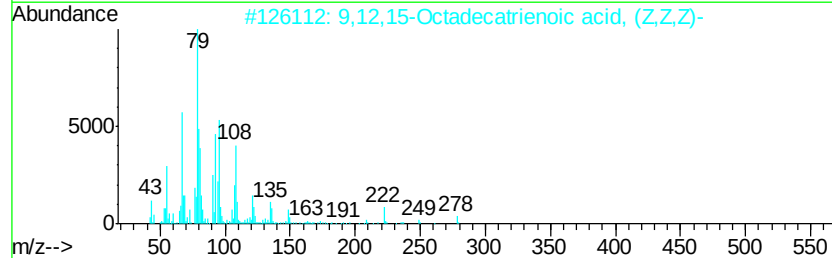
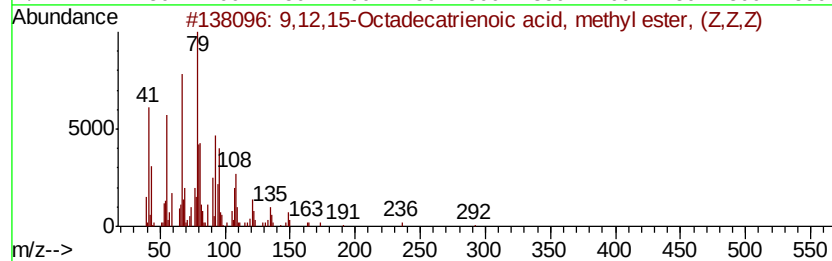
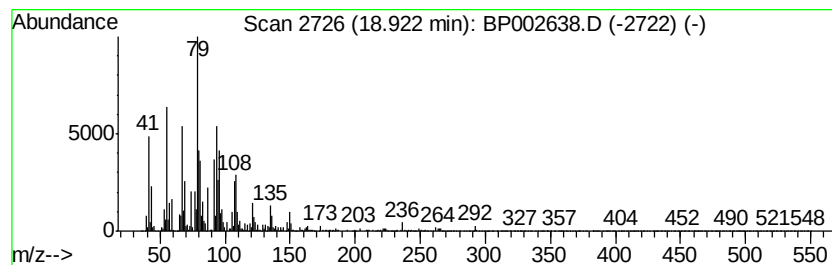
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 9,12,15-Octadecatrienoic ac... Concentration Rank 9

| R.T.  | EstConc | Area    | Relative to ISTD | R.T.  |
|-------|---------|---------|------------------|-------|
| 18.92 | 8.19 ng | 1070260 | Phenanthrene-d10 | 16.97 |

| Hit# | of | 5 | Tentative ID                         | MW  | MolForm  | CAS#         | Qual |
|------|----|---|--------------------------------------|-----|----------|--------------|------|
| 1    |    |   | 9,12,15-Octadecatrienoic acid, m...  | 292 | C19H32O2 | 000301-00-8  | 99   |
| 2    |    |   | 9,12,15-Octadecatrienoic acid, (...) | 278 | C18H30O2 | 000463-40-1  | 90   |
| 3    |    |   | Methyl 9.cis.,11.trans.t,13.trans... | 292 | C19H32O2 | 1000336-42-6 | 87   |
| 4    |    |   | 3-Dodecen-1-yne, (Z)-                | 164 | C12H20   | 025091-24-1  | 76   |
| 5    |    |   | cis,cis,cis-7,10,13-Hexadecatrienal  | 234 | C16H26O  | 056797-43-4  | 64   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP070220\  
Data File : BP002638.D  
Acq On : 02 Jul 2020 19:21  
Operator : CG/JU  
Sample : L3188-01 5X  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
TR-04-070120

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

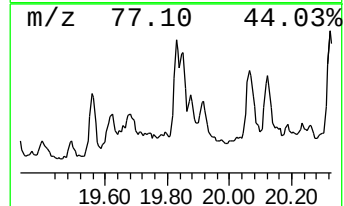
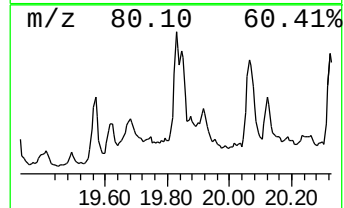
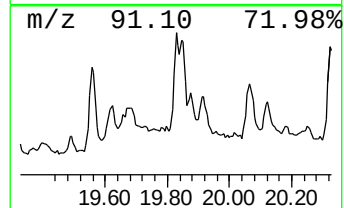
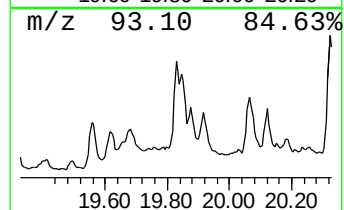
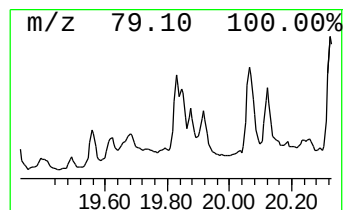
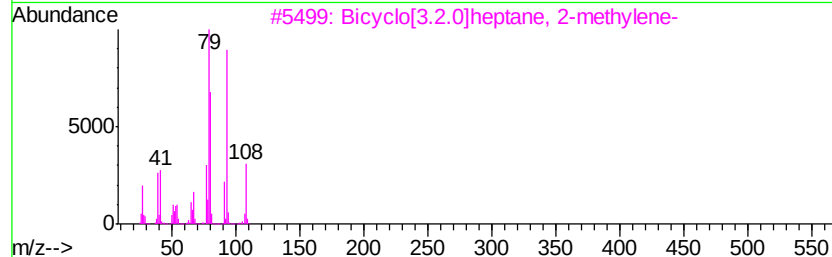
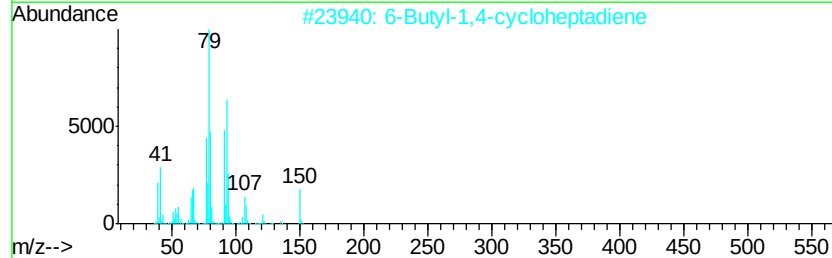
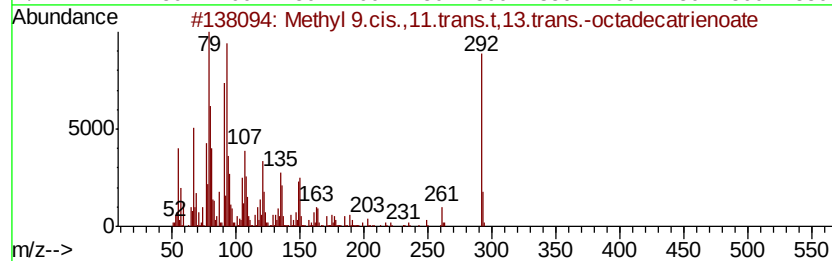
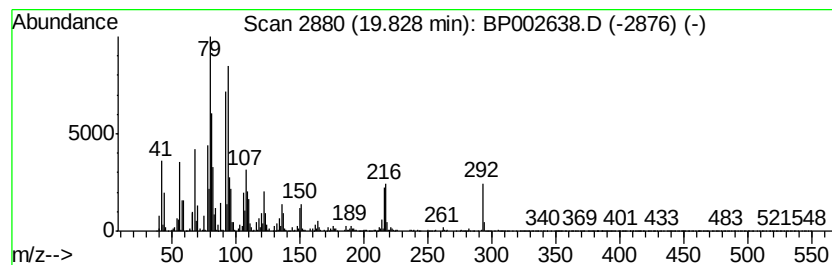
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 Methyl 9.cis.,11.trans.t,13... Concentration Rank 17

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 19.83 | 3.02 ng | 575104 | Chrysene-d12     | 21.09 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | Methyl 9.cis.,11.trans.t,13.tran... | 292 | C19H32O2 | 1000336-42-6 | 99   |
| 2    |    | 6-Butyl-1,4-cycloheptadiene         | 150 | C11H18   | 022735-58-6  | 70   |
| 3    |    | Bicyclo[3.2.0]heptane, 2-methylene- | 108 | C8H12    | 089243-94-7  | 50   |
| 4    |    | Methyl 6-cis,9-cis,11-trans-octa... | 292 | C19H32O2 | 1000336-37-7 | 50   |
| 5    |    | Bicyclo[3.1.1]heptane, 6,6-dimet... | 136 | C10H16   | 016022-04-1  | 49   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP070220\  
Data File : BP002638.D  
Acq On : 02 Jul 2020 19:21  
Operator : CG/JU  
Sample : L3188-01 5X  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

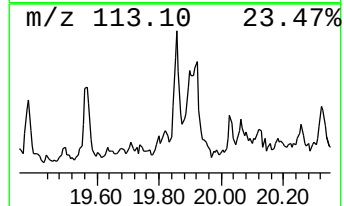
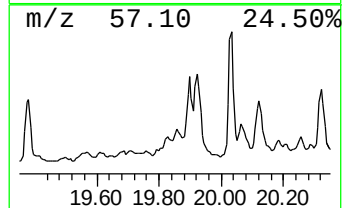
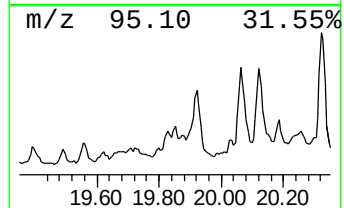
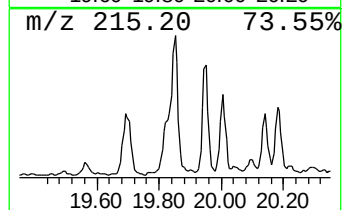
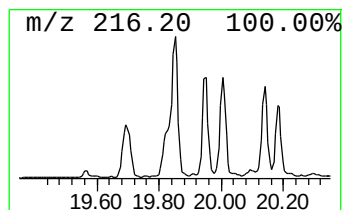
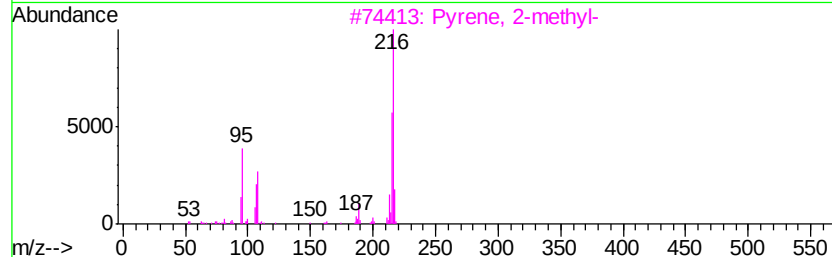
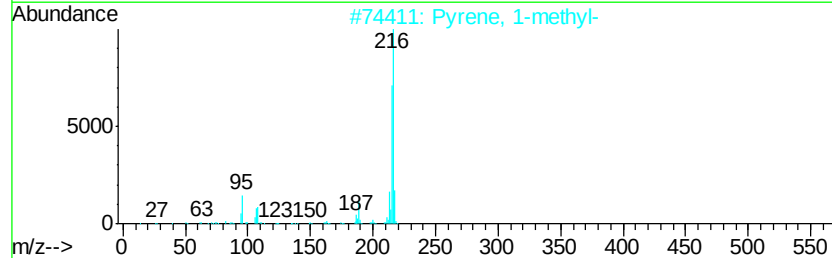
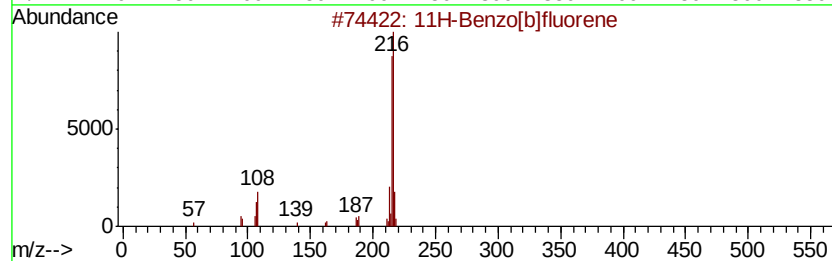
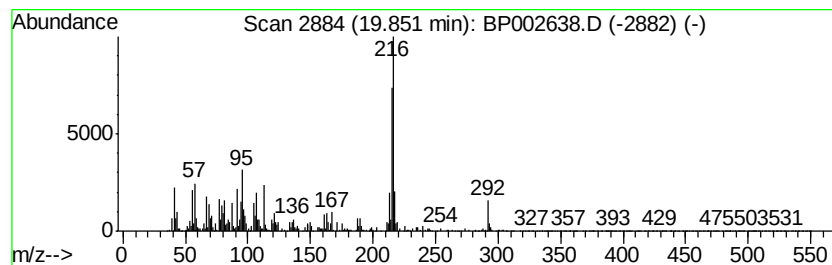
Instrument :  
BNA\_P  
ClientSampleId :  
TR-04-070120

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

\*\*\*\*\*  
Peak Number 10 11H-Benzo[b]fluorene Concentration Rank 19

| R.T.    | EstConc | Area                 | Relative to ISTD |         | R.T.        |      |
|---------|---------|----------------------|------------------|---------|-------------|------|
| 19.85   | 2.75 ng | 524188               | Chrysene-d12     |         | 21.09       |      |
| Hit# of | 5       | Tentative ID         | MW               | MolForm | CAS#        | Qual |
| 1       |         | 11H-Benzo[b]fluorene | 216              | C17H12  | 000243-17-4 | 92   |
| 2       |         | Pyrene, 1-methyl-    | 216              | C17H12  | 002381-21-7 | 91   |
| 3       |         | Pyrene, 2-methyl-    | 216              | C17H12  | 003442-78-2 | 90   |
| 4       |         | 7H-Benzo[c]fluorene  | 216              | C17H12  | 000205-12-9 | 76   |
| 5       |         | 11H-Benzo[a]fluorene | 216              | C17H12  | 000238-84-6 | 68   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP070220\  
Data File : BP002638.D  
Acq On : 02 Jul 2020 19:21  
Operator : CG/JU  
Sample : L3188-01 5X  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
TR-04-070120

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

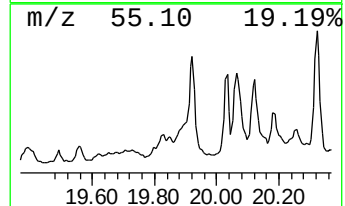
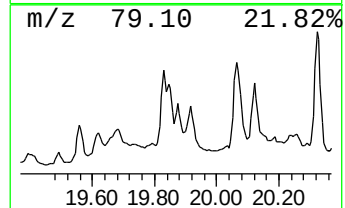
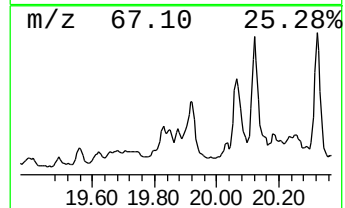
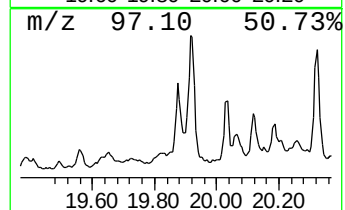
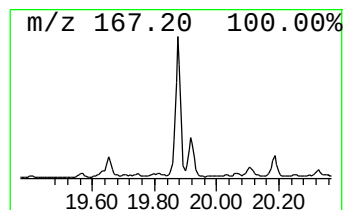
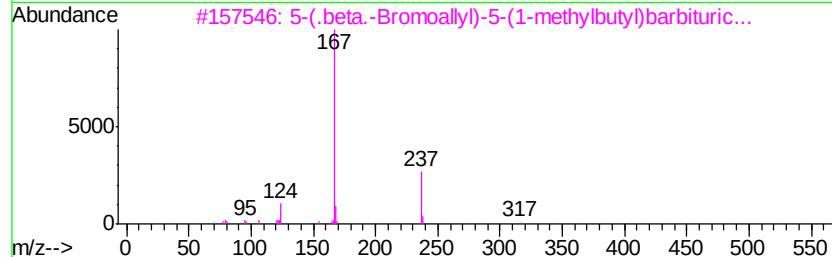
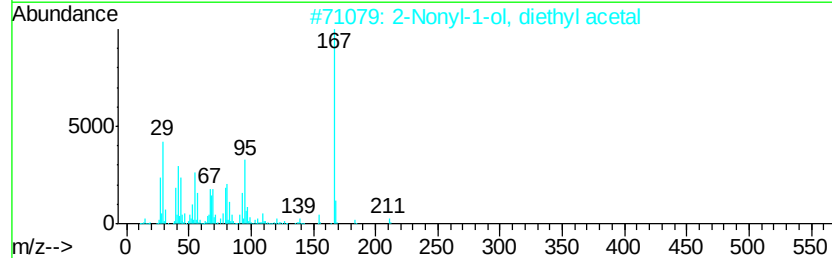
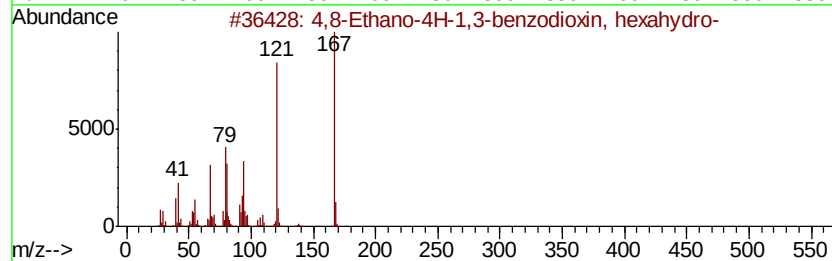
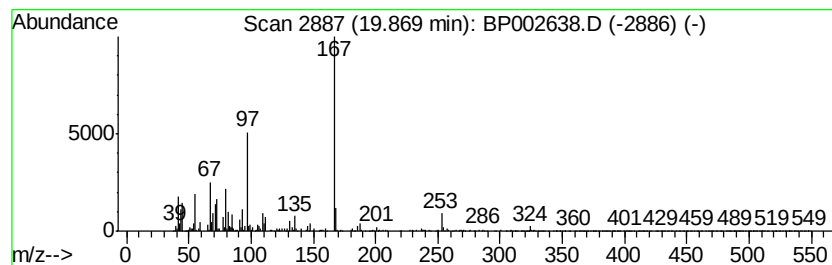
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 unknown19.87 Concentration Rank 18

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 19.87 | 2.79 ng | 530820 | Chrysene-d12     | 21.09 |

| Hit# | of | Tentative ID                        | MW  | MolForm      | CAS#         | Qual |
|------|----|-------------------------------------|-----|--------------|--------------|------|
| 1    | 5  | 4,8-Ethano-4H-1,3-benzodioxin, h... | 168 | C10H16O2     | 057790-44-0  | 47   |
| 2    |    | 2-Nonyl-1-ol, diethyl acetal        | 212 | C13H24O2     | 1000343-44-6 | 42   |
| 3    |    | 5-(.beta.-Bromoallyl)-5-(1-methv... | 316 | C12H17BrN2O3 | 001216-40-6  | 35   |
| 4    |    | 2-Methyl-7H-1,3,4-thiadiazolo[3,... | 167 | C6H5N3OS     | 056347-07-0  | 35   |
| 5    |    | 6H-Purine-6-thione, 9-amino-1,9-... | 167 | C5H5N5S      | 020914-56-1  | 35   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP070220\  
Data File : BP002638.D  
Acq On : 02 Jul 2020 19:21  
Operator : CG/JU  
Sample : L3188-01 5X  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
TR-04-070120

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

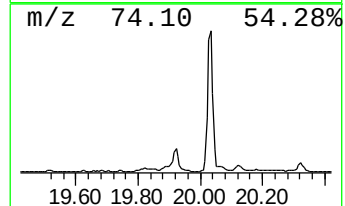
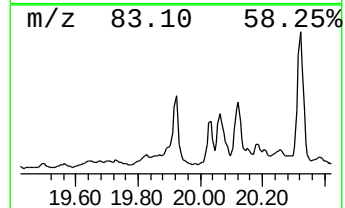
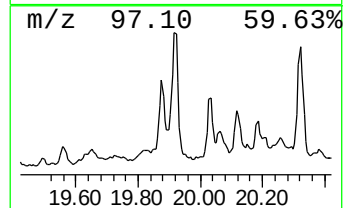
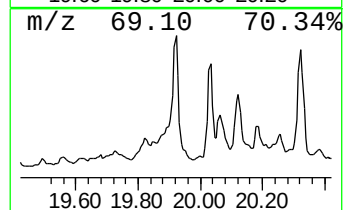
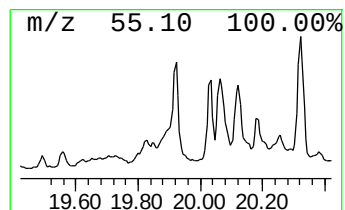
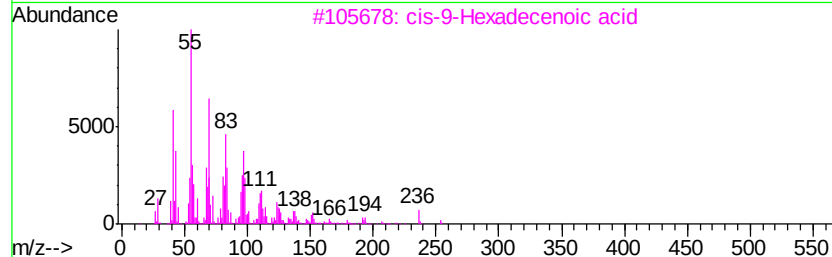
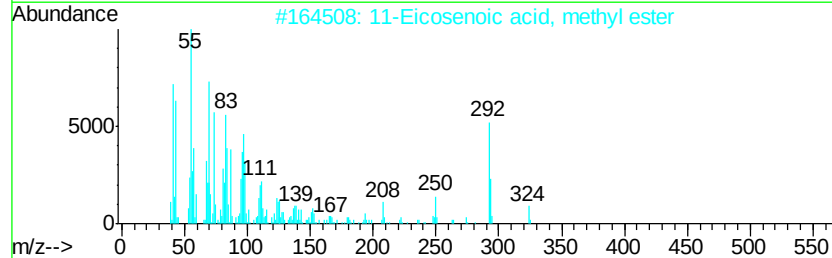
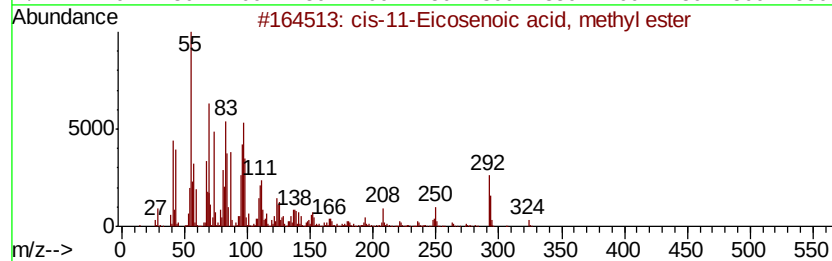
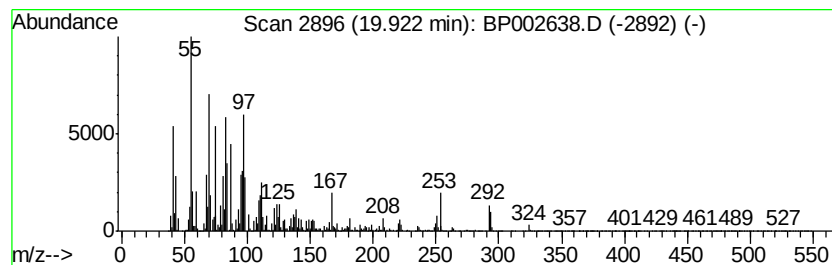
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TIC Integration Parameters: LSCINT.P

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Peak Number 12 cis-11-Eicosenoic acid, met... Concentration Rank 5

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 19.92 | 10.60 ng | 2017420 | Chrysene-d12     | 21.09 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | cis-11-Eicosenoic acid, methyl e... | 324 | C21H40O2 | 1000333-63-8 | 90   |
| 2    |    | 11-Eicosenoic acid, methyl ester    | 324 | C21H40O2 | 003946-08-5  | 62   |
| 3    |    | cis-9-Hexadecenoic acid             | 254 | C16H30O2 | 1000333-19-5 | 55   |
| 4    |    | 14-Octadecenoic acid, methyl ester  | 296 | C19H36O2 | 056554-48-4  | 55   |
| 5    |    | trans-13-Octadecenoic acid, meth... | 296 | C19H36O2 | 1000333-61-3 | 53   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP070220\  
Data File : BP002638.D  
Acq On : 02 Jul 2020 19:21  
Operator : CG/JU  
Sample : L3188-01 5X  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

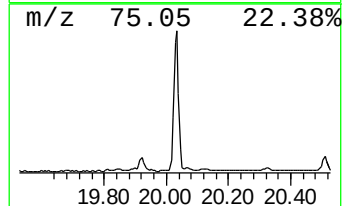
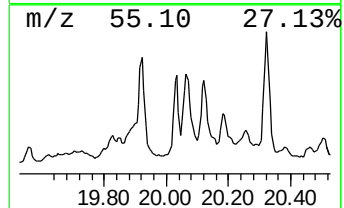
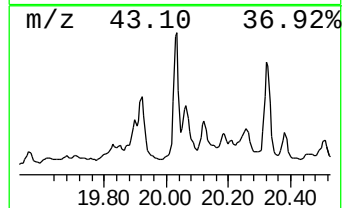
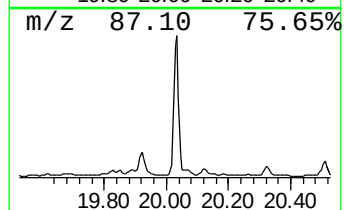
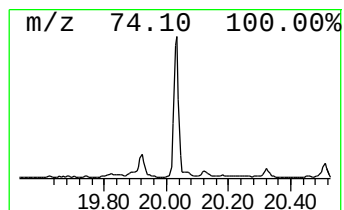
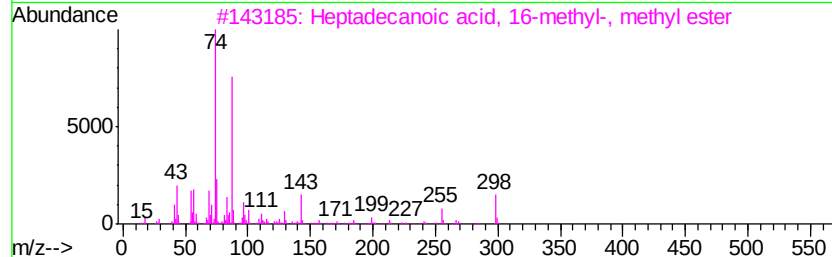
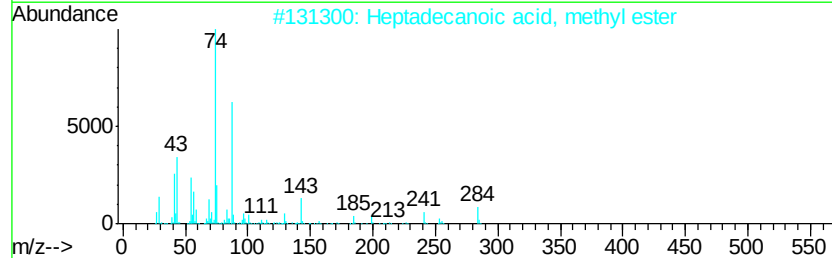
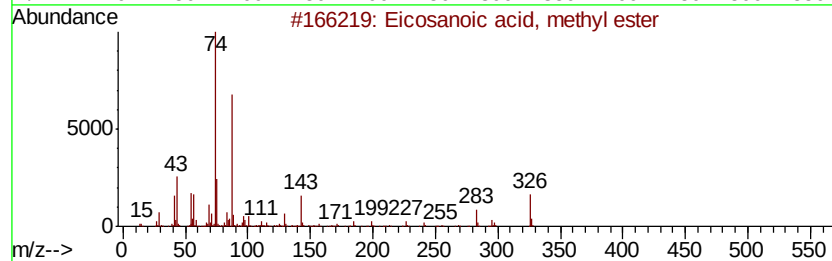
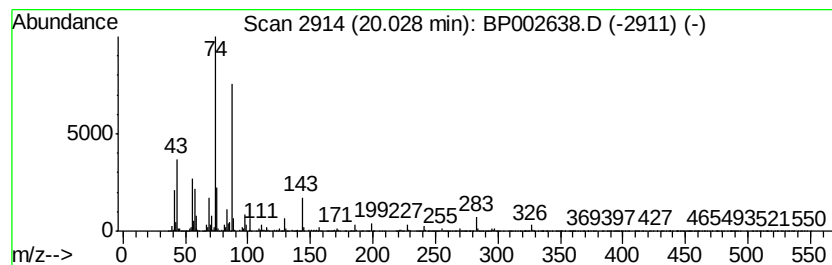
Instrument :  
BNA\_P  
ClientSampleId :  
TR-04-070120

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

\*\*\*\*\*  
Peak Number 13 Eicosanoic acid, methyl ester Concentration Rank 12

| R.T.    | EstConc                             | Area         | Relative to ISTD | R.T.         |      |      |
|---------|-------------------------------------|--------------|------------------|--------------|------|------|
| 20.03   | 7.60 ng                             | 1446200      | Chrysene-d12     | 21.09        |      |      |
| Hit# of | 5                                   | Tentative ID | MW               | MolForm      | CAS# | Qual |
| 1       | Eicosanoic acid, methyl ester       | 326          | C21H42O2         | 001120-28-1  | 98   |      |
| 2       | Heptadecanoic acid, methyl ester    | 284          | C18H36O2         | 001731-92-6  | 91   |      |
| 3       | Heptadecanoic acid, 16-methyl-, ... | 298          | C19H38O2         | 005129-61-3  | 91   |      |
| 4       | Methyl 18-methylnonadecanoate       | 326          | C21H42O2         | 1000352-20-6 | 91   |      |
| 5       | Hexadecanoic acid, 15-methyl-, m... | 284          | C18H36O2         | 006929-04-0  | 90   |      |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP070220\  
Data File : BP002638.D  
Acq On : 02 Jul 2020 19:21  
Operator : CG/JU  
Sample : L3188-01 5X  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
TR-04-070120

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

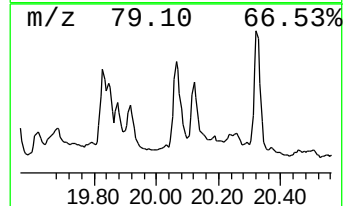
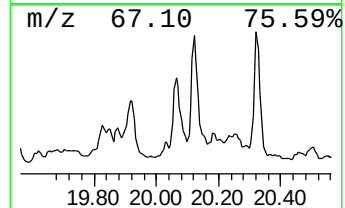
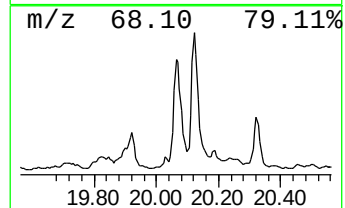
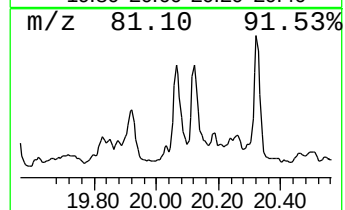
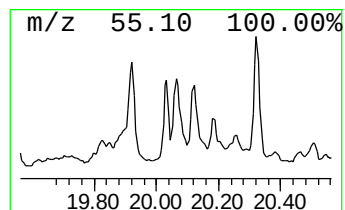
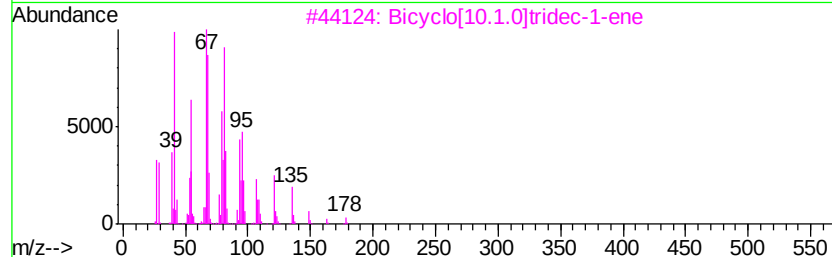
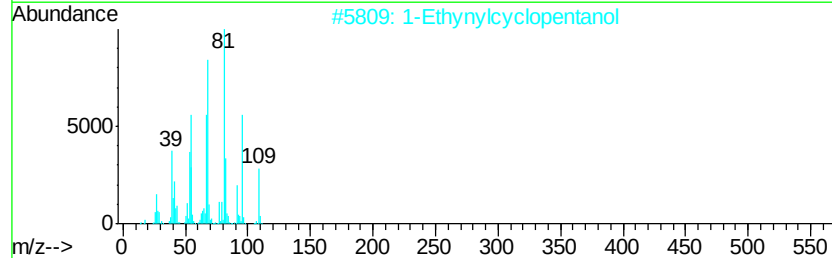
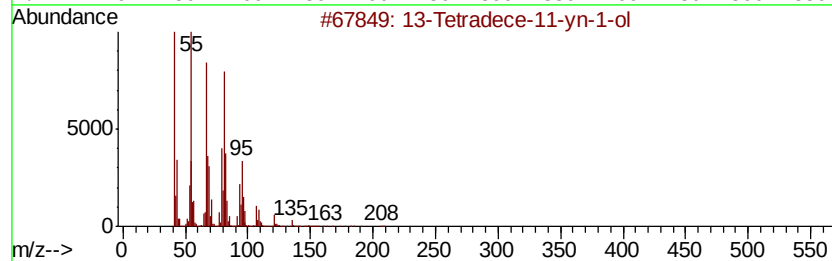
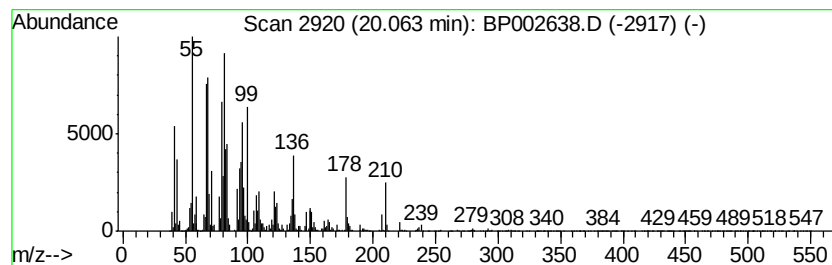
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 14 13-Tetradec-11-yn-1-ol Concentration Rank 10

| R.T.  | EstConc | Area    | Relative to ISTD | R.T.  |
|-------|---------|---------|------------------|-------|
| 20.06 | 7.96 ng | 1514990 | Chrysene-d12     | 21.09 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------------------|-----|---------|--------------|------|
| 1    | 5  | 13-Tetradec-11-vn-1-ol              | 208 | C14H24O | 1000131-00-4 | 78   |
| 2    |    | 1-Ethynylcyclopentanol              | 110 | C7H10O  | 017356-19-3  | 58   |
| 3    |    | Bicyclo[10.1.0]tridec-1-ene         | 178 | C13H22  | 054766-91-5  | 52   |
| 4    |    | Bicyclo[4.1.0]heptane, 3-methyl-    | 110 | C8H14   | 041977-47-3  | 52   |
| 5    |    | 3-Methyl-cis-3a,4,7,7a-tetrahydr... | 136 | C10H16  | 1000144-04-4 | 50   |





Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP070220\  
Data File : BP002638.D  
Acq On : 02 Jul 2020 19:21  
Operator : CG/JU  
Sample : L3188-01 5X  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
TR-04-070120

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

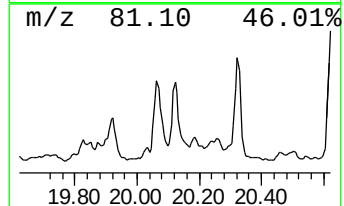
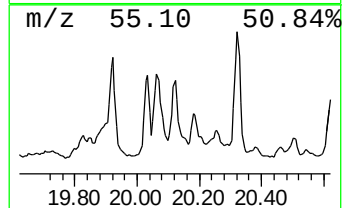
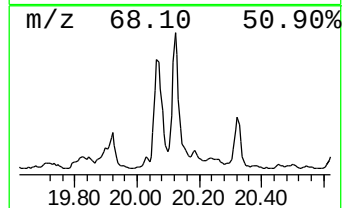
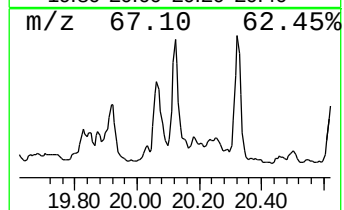
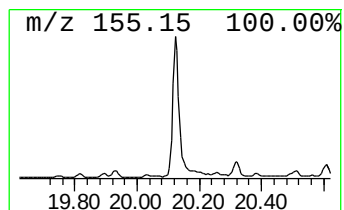
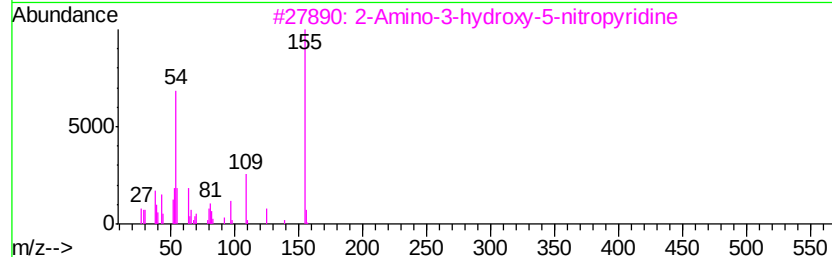
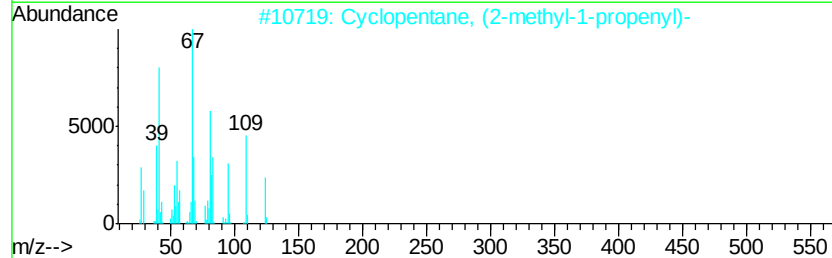
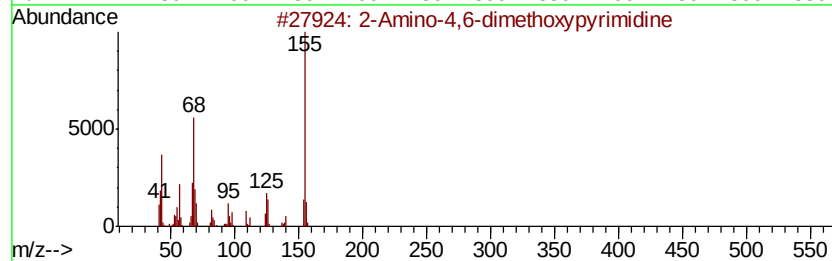
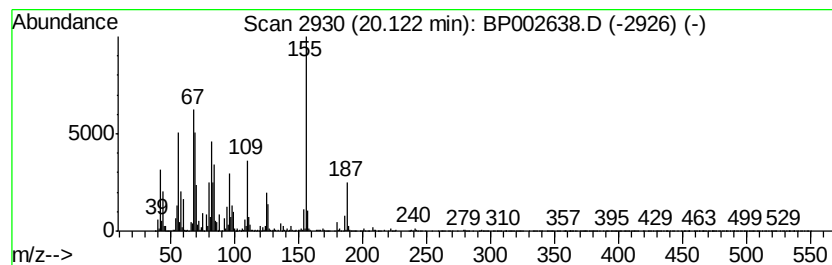
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 15 unknown20.12 Concentration Rank 11

| R.T.  | EstConc | Area    | Relative to ISTD | R.T.  |
|-------|---------|---------|------------------|-------|
| 20.12 | 7.64 ng | 1453740 | Chrysene-d12     | 21.09 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | 2-Amino-4,6-dimethoxypyrimidine     | 155 | C6H9N3O2 | 036315-01-2  | 47   |
| 2    |    |   | Cyclopentane, (2-methyl-1-propen... | 124 | C9H16    | 053366-57-7  | 46   |
| 3    |    |   | 2-Amino-3-hydroxy-5-nitropyridine   | 155 | C5H5N3O3 | 1000289-29-0 | 35   |
| 4    |    |   | 9-Hexadecyn-1-ol                    | 238 | C16H30O  | 1000342-40-3 | 30   |
| 5    |    |   | 1,2-Cyclohexanedicarboxylic acid... | 348 | C20H28O5 | 1000339-66-9 | 27   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP070220\  
Data File : BP002638.D  
Acq On : 02 Jul 2020 19:21  
Operator : CG/JU  
Sample : L3188-01 5X  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
TR-04-070120

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

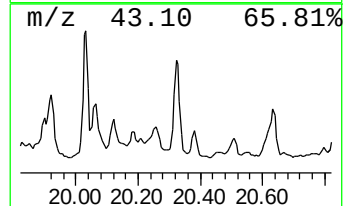
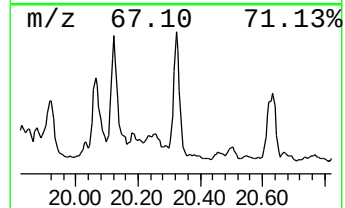
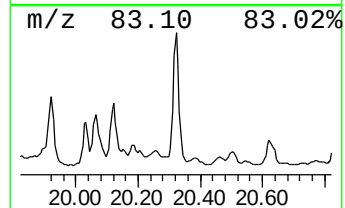
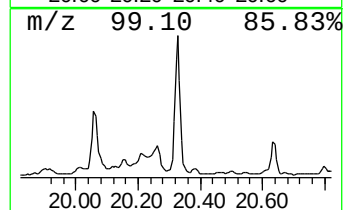
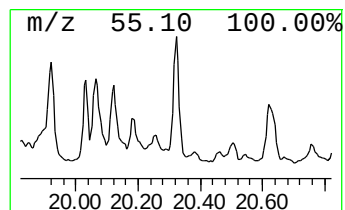
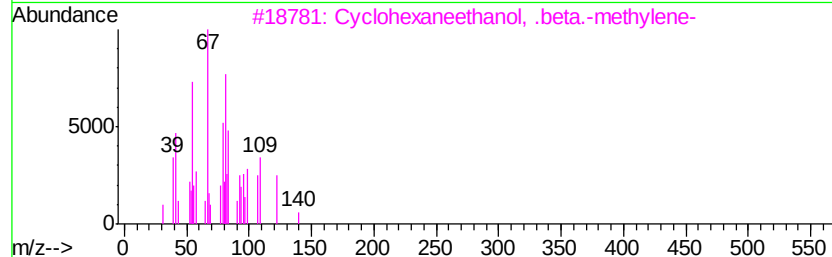
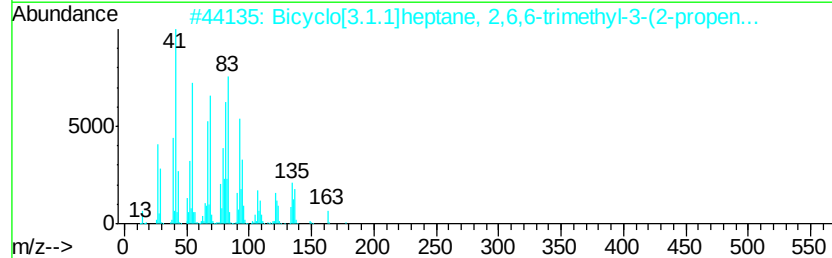
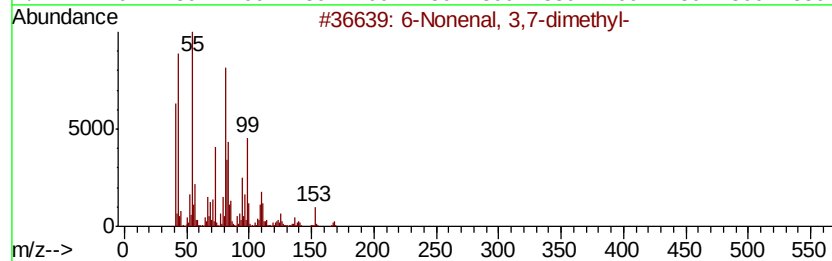
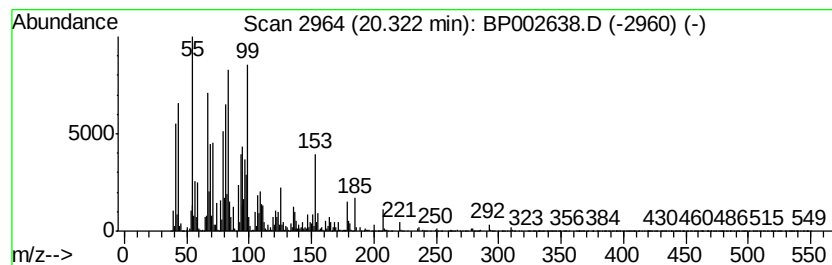
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 16 unknown20.32 Concentration Rank 4

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 20.32 | 12.56 ng | 2390970 | Chrysene-d12     | 21.09 |

| Hit# | of                                  | 5 | Tentative ID | MW  | MolForm | CAS#         | Qual |
|------|-------------------------------------|---|--------------|-----|---------|--------------|------|
| 1    | 6-Nonenal, 3,7-dimethyl-            |   |              | 168 | C11H20O | 1000132-43-9 | 47   |
| 2    | Bicyclo[3.1.1]heptane, 2,6,6-tri... |   |              | 178 | C13H22  | 050746-55-9  | 27   |
| 3    | Cyclohexaneethanol, .beta.-methv... |   |              | 140 | C9H16O  | 007697-55-4  | 25   |
| 4    | 1,3-dicyclohexylpropene             |   |              | 206 | C15H26  | 1000214-96-8 | 15   |
| 5    | Silane, dimethyldi-2-propenyl-      |   |              | 140 | C8H16Si | 001113-12-8  | 14   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP070220\  
Data File : BP002638.D  
Acq On : 02 Jul 2020 19:21  
Operator : CG/JU  
Sample : L3188-01 5X  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
TR-04-070120

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

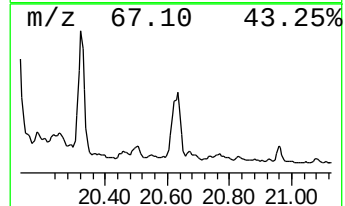
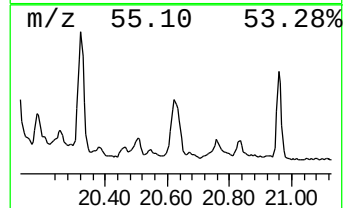
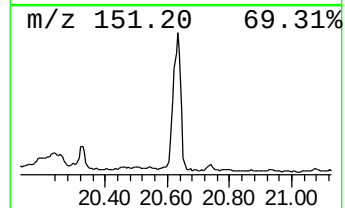
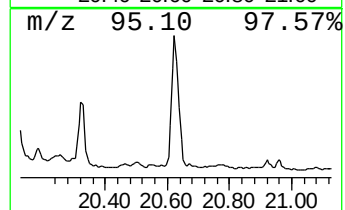
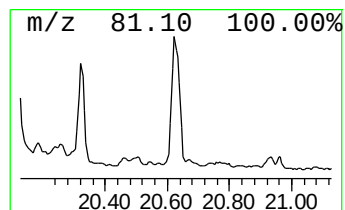
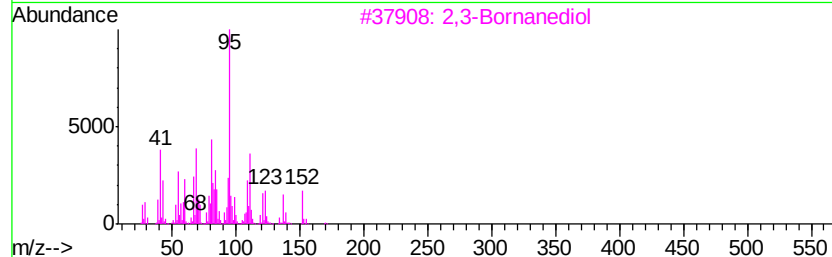
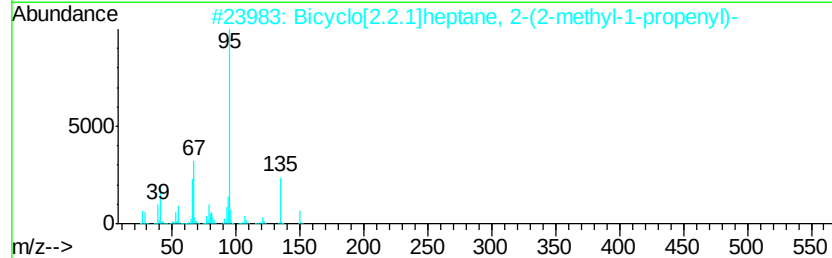
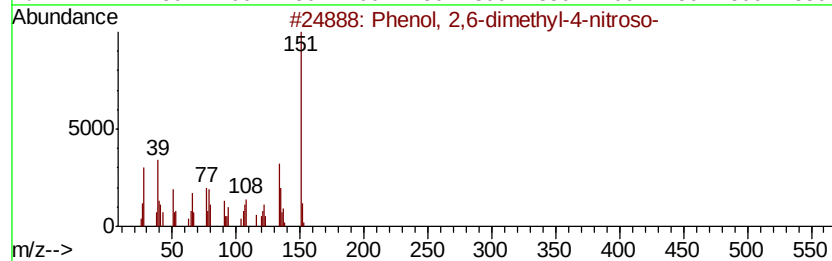
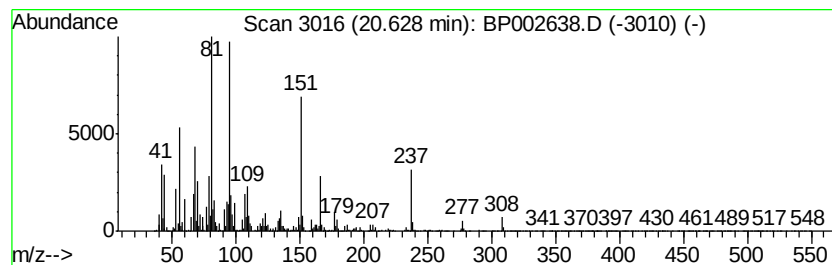
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 17 unknown20.63 Concentration Rank 6

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 20.63 | 10.51 ng | 1999550 | Chrysene-d12     | 21.09 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | Phenol, 2,6-dimethyl-4-nitroso-     | 151 | C8H9NO2  | 013331-93-6  | 35   |
| 2    |    | Bicyclo[2.2.1]heptane, 2-(2-meth... | 150 | C11H18   | 061142-27-6  | 30   |
| 3    |    | 2,3-Bornanediol                     | 170 | C10H18O2 | 025696-56-4  | 25   |
| 4    |    | 2-Cyanomethyl-hexahydro-pyrimidi... | 175 | C8H9N5   | 1000185-88-3 | 15   |
| 5    |    | Daucol                              | 238 | C15H26O2 | 000887-08-1  | 15   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP070220\  
 Data File : BP002638.D  
 Acq On : 02 Jul 2020 19:21  
 Operator : CG/JU  
 Sample : L3188-01 5X  
 Misc :  
 ALS Vial : 20 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 TR-04-070120

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

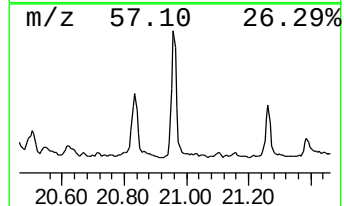
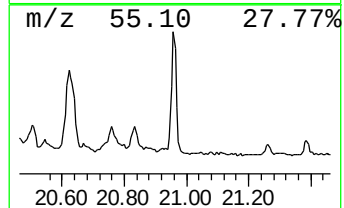
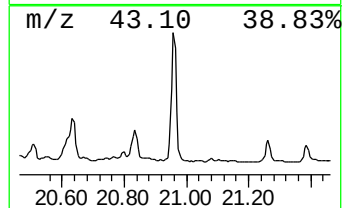
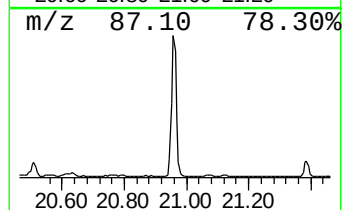
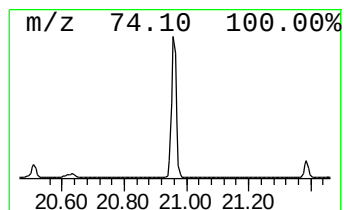
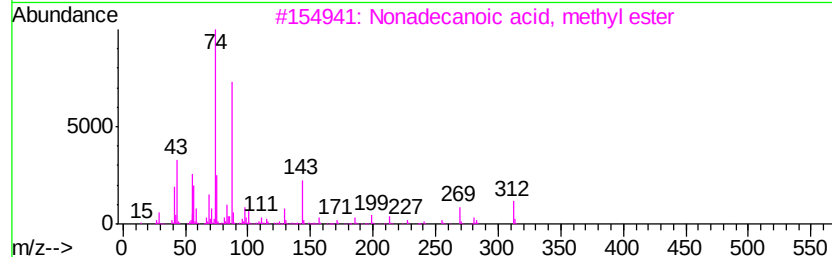
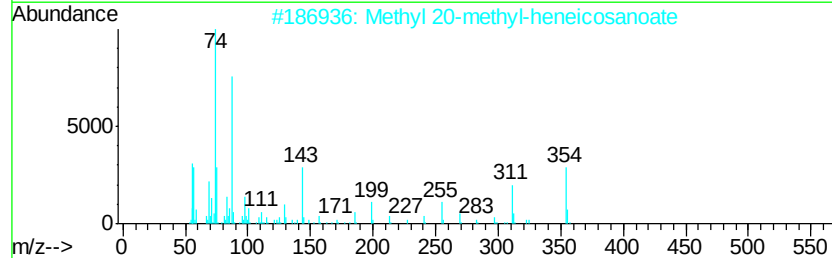
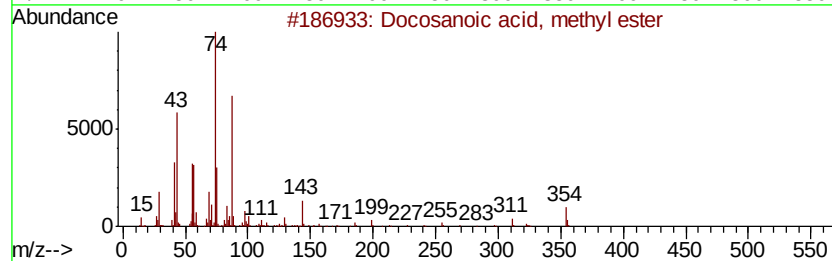
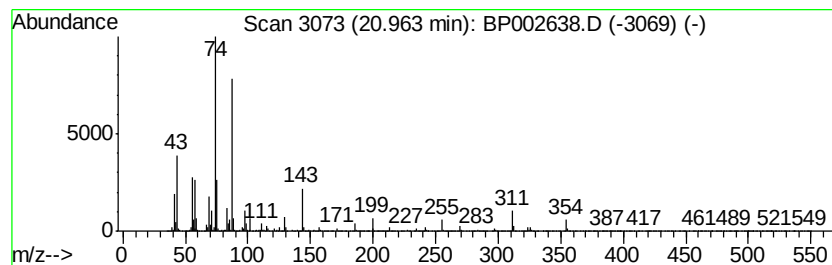
TIC Library : C:\DATABASE\NIST11.L  
 TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
 Peak Number 18 Docosanoic acid, methyl ester Concentration Rank 8

| R.T.  | EstConc | Area    | Relative to ISTD | R.T.  |
|-------|---------|---------|------------------|-------|
| 20.96 | 8.94 ng | 1700810 | Chrysene-d12     | 21.09 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | Docosanoic acid, methyl ester       | 354 | C23H46O2 | 000929-77-1  | 99   |
| 2    |    | Methyl 20-methyl-heneicosanoate     | 354 | C23H46O2 | 1000336-47-4 | 98   |
| 3    |    | Nonadecanoic acid, methyl ester     | 312 | C20H40O2 | 001731-94-8  | 90   |
| 4    |    | Methyl stearate                     | 298 | C19H38O2 | 000112-61-8  | 87   |
| 5    |    | Hexadecanoic acid, 15-methyl-, m... | 284 | C18H36O2 | 006929-04-0  | 83   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP070220\  
Data File : BP002638.D  
Acq On : 02 Jul 2020 19:21  
Operator : CG/JU  
Sample : L3188-01 5X  
Misc :  
ALS Vial : 20 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
TR-04-070120

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

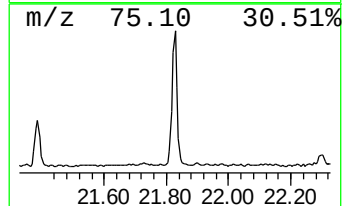
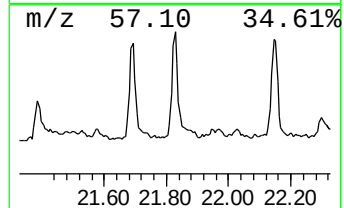
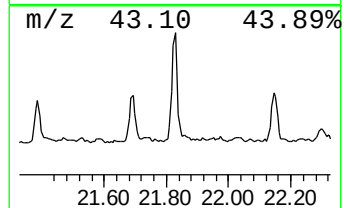
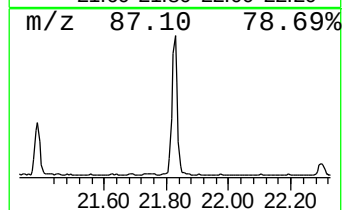
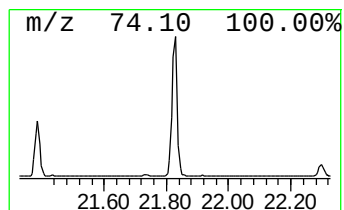
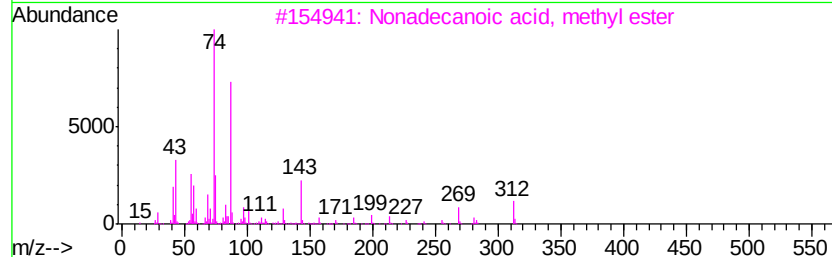
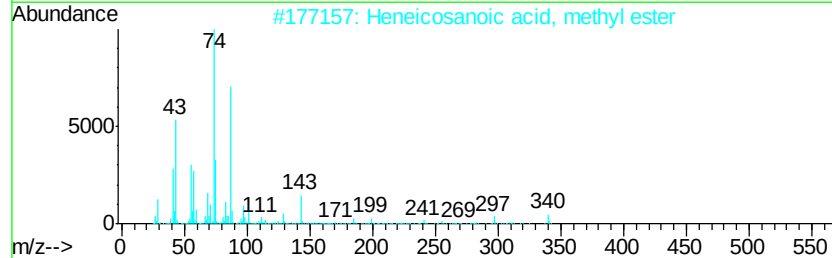
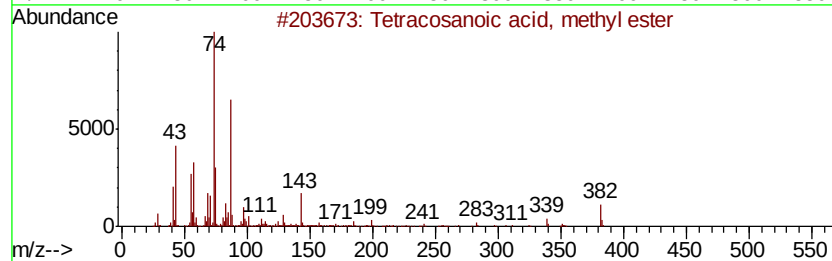
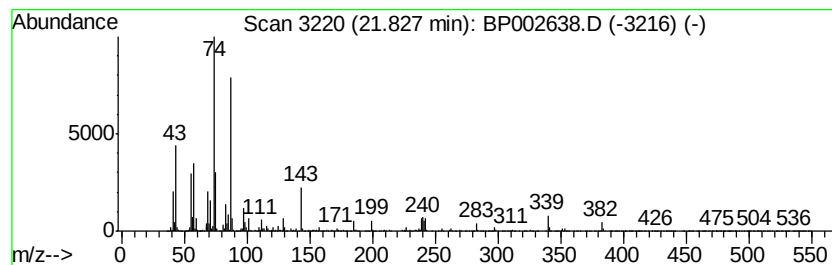
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 Tetracosanoic acid, methyl ... Concentration Rank 14

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 21.83 | 3.91 ng | 744641 | Chrysene-d12     | 21.09 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Tetracosanoic acid, methyl ester    | 382 | C25H50O2 | 002442-49-1 | 98   |
| 2    |    |   | Heneicosanoic acid, methyl ester    | 340 | C22H44O2 | 006064-90-0 | 93   |
| 3    |    |   | Nonadecanoic acid, methyl ester     | 312 | C20H40O2 | 001731-94-8 | 87   |
| 4    |    |   | Methyl stearate                     | 298 | C19H38O2 | 000112-61-8 | 86   |
| 5    |    |   | Heptadecanoic acid, 16-methyl-, ... | 298 | C19H38O2 | 005129-61-3 | 83   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_P\DATA\BP070220\  
 Data File : BP002638.D  
 Acq On : 02 Jul 2020 19:21  
 Operator : CG/JU  
 Sample : L3188-01 5X  
 Misc :  
 ALS Vial : 20 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 TR-04-070120

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 |
| 9-Hexadecenoic ac...   | 17.56 | 3.8     | ng    | 492813   | 4                     | 16.97 | 2614560 | 20.0 |
| Hexadecanoic acid...   | 17.69 | 127.4   | ng    | 16657200 | 4                     | 16.97 | 2614560 | 20.0 |
| Cyclopropaneoctan...   | 18.22 | 5.1     | ng    | 666463   | 4                     | 16.97 | 2614560 | 20.0 |
| Heptadecanoic aci...   | 18.36 | 3.2     | ng    | 417407   | 4                     | 16.97 | 2614560 | 20.0 |
| 9,12-Octadecadien...   | 18.82 | 274.8   | ng    | 35917800 | 4                     | 16.97 | 2614560 | 20.0 |
| 10-Octadecenoic a...   | 18.86 | 170.6   | ng    | 22302000 | 4                     | 16.97 | 2614560 | 20.0 |
| 9,12,15-Octadecat...   | 18.92 | 8.2     | ng    | 1070260  | 4                     | 16.97 | 2614560 | 20.0 |
| Methyl 9.cis.,11...    | 19.83 | 3.0     | ng    | 575104   | 5                     | 21.09 | 3805850 | 20.0 |
| 11H-Benzo[b]fluorene   | 19.85 | 2.8     | ng    | 524188   | 5                     | 21.09 | 3805850 | 20.0 |
| unknown19.87           | 19.87 | 2.8     | ng    | 530820   | 5                     | 21.09 | 3805850 | 20.0 |
| cis-11-Eicosenoic...   | 19.92 | 10.6    | ng    | 2017420  | 5                     | 21.09 | 3805850 | 20.0 |
| Eicosanoic acid, ...   | 20.03 | 7.6     | ng    | 1446200  | 5                     | 21.09 | 3805850 | 20.0 |
| 13-Tetradecene-11-y... | 20.06 | 8.0     | ng    | 1514990  | 5                     | 21.09 | 3805850 | 20.0 |
| unknown20.12           | 20.12 | 7.6     | ng    | 1453740  | 5                     | 21.09 | 3805850 | 20.0 |
| unknown20.32           | 20.32 | 12.6    | ng    | 2390970  | 5                     | 21.09 | 3805850 | 20.0 |
| unknown20.63           | 20.63 | 10.5    | ng    | 1999550  | 5                     | 21.09 | 3805850 | 20.0 |
| Docosanoic acid, ...   | 20.96 | 8.9     | ng    | 1700810  | 5                     | 21.09 | 3805850 | 20.0 |
| Tetracosanoic aci...   | 21.83 | 3.9     | ng    | 744641   | 5                     | 21.09 | 3805850 | 20.0 |