

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP012522\  
 Data File : BP008913.D  
 Acq On : 25 Jan 2022 18:16  
 Operator : CG/JU  
 Sample : N1233-01  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 RT3876

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\8270E-BP011422.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP008913.D\data.ms

| 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         | 3.040       | 5             | 9           | 25           | rVB      | 436756         | 607811        | 4.82%           | 0.786%        |
| 2         | 5.422       | 407           | 414         | 432          | rBV      | 3261086        | 5179192       | 41.06%          | 6.698%        |
| 3         | 7.016       | 677           | 685         | 709          | rBV      | 2982679        | 5245220       | 41.58%          | 6.783%        |
| 4         | 7.834       | 817           | 824         | 834          | rBV      | 819238         | 1345402       | 10.67%          | 1.740%        |
| 5         | 8.993       | 1013          | 1021        | 1040         | rBV      | 1907534        | 3429480       | 27.19%          | 4.435%        |
| 6         | 10.634      | 1291          | 1300        | 1315         | rBV      | 961081         | 1783964       | 14.14%          | 2.307%        |
| 7         | 13.098      | 1712          | 1719        | 1734         | rBV      | 5444360        | 8337186       | 66.09%          | 10.782%       |
| 8         | 14.475      | 1945          | 1953        | 1960         | rBV      | 1524671        | 2317645       | 18.37%          | 2.997%        |
| 9         | 15.969      | 2200          | 2207        | 2227         | rBV      | 4023473        | 6245048       | 49.51%          | 8.076%        |
| 10        | 17.234      | 2415          | 2422        | 2427         | rBV      | 1849395        | 2760183       | 21.88%          | 3.570%        |
| 11        | 17.275      | 2427          | 2429        | 2435         | rVB      | 139054         | 165881        | 1.32%           | 0.215%        |
| 12        | 17.910      | 2532          | 2537        | 2543         | rBV      | 1531831        | 1770873       | 14.04%          | 2.290%        |
| 13        | 18.128      | 2570          | 2574        | 2582         | rBV2     | 333832         | 539381        | 4.28%           | 0.698%        |
| 14        | 19.010      | 2719          | 2724        | 2726         | rBV      | 4006020        | 4493324       | 35.62%          | 5.811%        |
| 15        | 19.045      | 2726          | 2730        | 2734         | rVB2     | 6625650        | 8084050       | 64.09%          | 10.455%       |
| 16        | 19.169      | 2747          | 2751        | 2757         | rBV      | 720340         | 818466        | 6.49%           | 1.058%        |
| 17        | 19.245      | 2759          | 2764        | 2772         | rVV      | 356774         | 547029        | 4.34%           | 0.707%        |
| 18        | 19.363      | 2781          | 2784        | 2793         | rVB2     | 92614          | 148423        | 1.18%           | 0.192%        |
| 19        | 19.598      | 2820          | 2824        | 2832         | rVB      | 231599         | 391297        | 3.10%           | 0.506%        |
| 20        | 19.792      | 2852          | 2857        | 2863         | rBV      | 9996883        | 12614193      | 100.00%         | 16.313%       |
| 21        | 20.128      | 2911          | 2914        | 2921         | rVB3     | 166200         | 207268        | 1.64%           | 0.268%        |
| 22        | 21.039      | 3065          | 3069        | 3076         | rBV      | 941453         | 1294266       | 10.26%          | 1.674%        |
| 23        | 21.322      | 3112          | 3117        | 3122         | rBV      | 2619190        | 3595450       | 28.50%          | 4.650%        |
| 24        | 21.439      | 3133          | 3137        | 3142         | rBV      | 1172102        | 1657636       | 13.14%          | 2.144%        |
| 25        | 23.739      | 3521          | 3528        | 3536         | rVB2     | 1818340        | 3747086       | 29.71%          | 4.846%        |

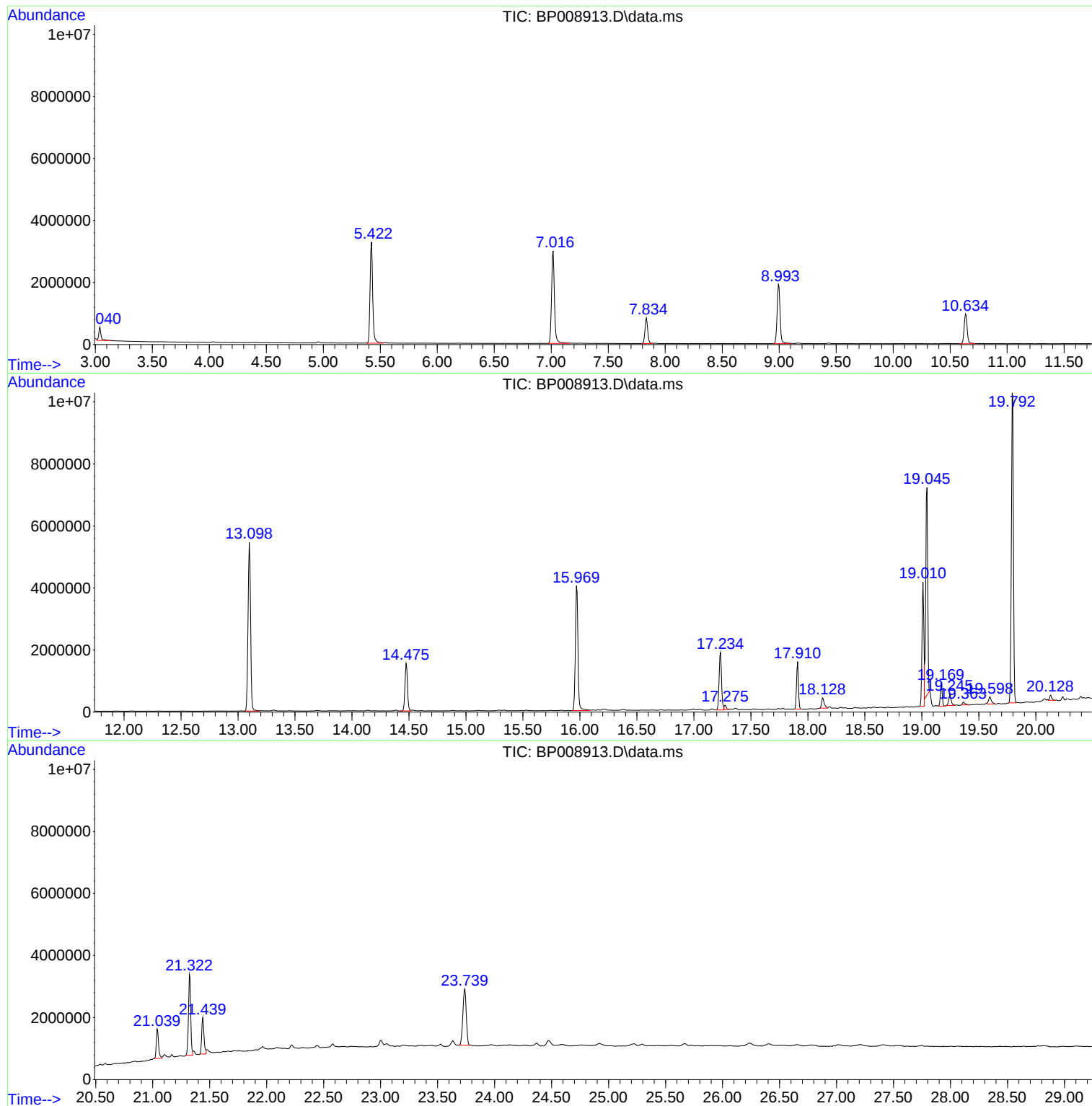
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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP011422.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P



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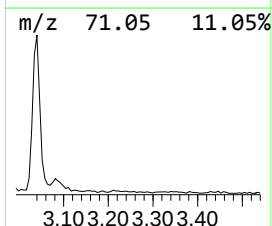
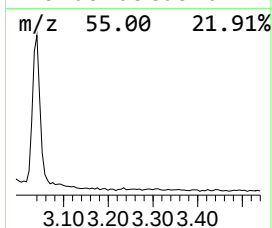
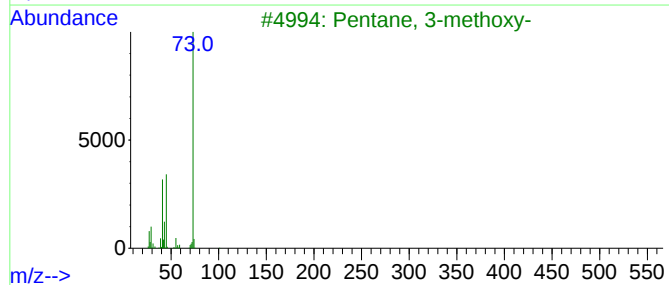
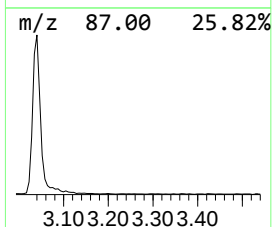
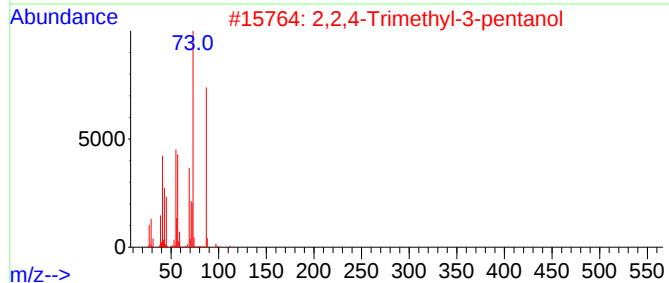
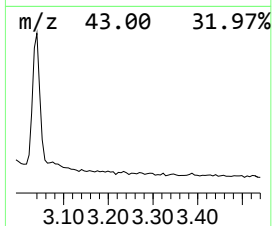
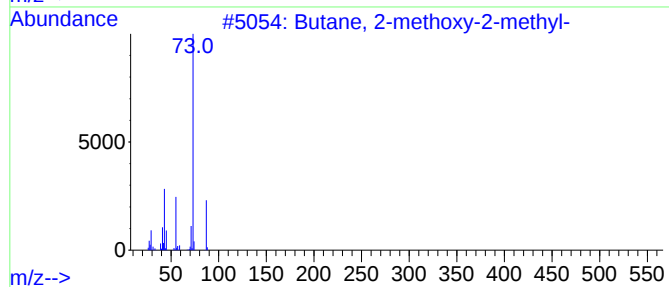
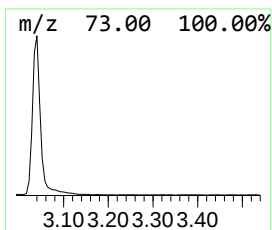
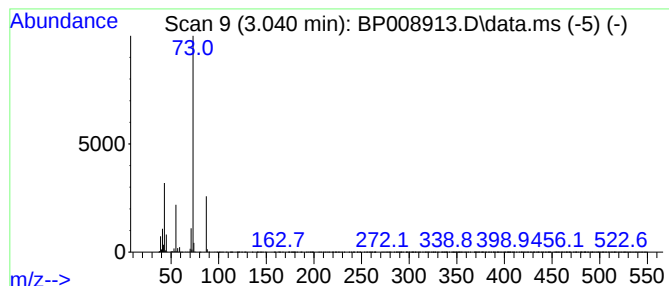
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 5

| R.T.  | EstConc | Area   | Relative to ISTD       | R.T.  |
|-------|---------|--------|------------------------|-------|
| 3.040 | 9.04 ng | 607811 | 1,4-Dichlorobenzene-d4 | 7.834 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Butane, 2-methoxy-2-methyl- | 102 | C6H14O  | 000994-05-8 | 90   |
| 2    |    |   | 2,2,4-Trimethyl-3-pentanol  | 130 | C8H18O  | 005162-48-1 | 43   |
| 3    |    |   | Pentane, 3-methoxy-         | 102 | C6H14O  | 036839-67-5 | 23   |
| 4    |    |   | Silane, tetramethyl-        | 88  | C4H12Si | 000075-76-3 | 9    |
| 5    |    |   | N-Ethylformamide            | 73  | C3H7NO  | 000627-45-2 | 9    |



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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP011422.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

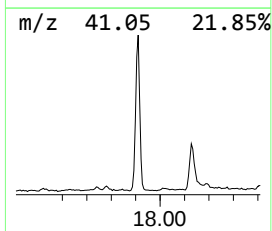
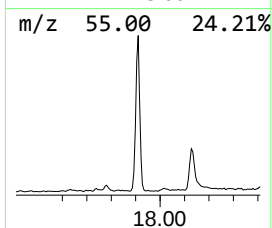
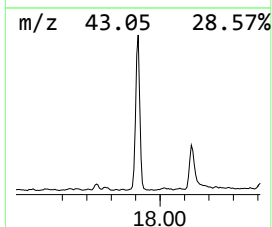
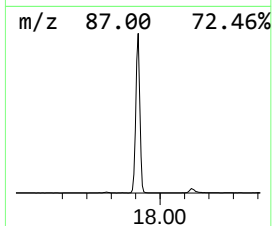
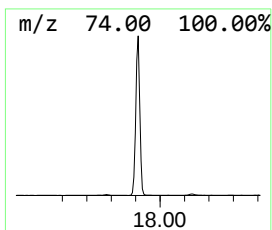
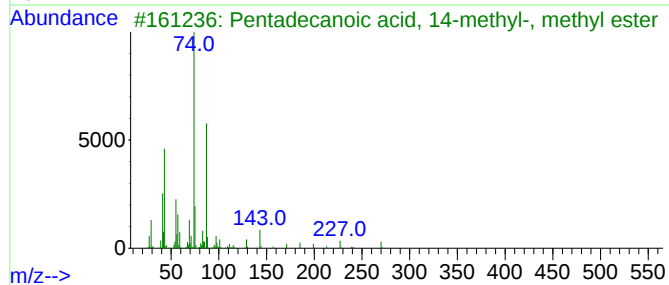
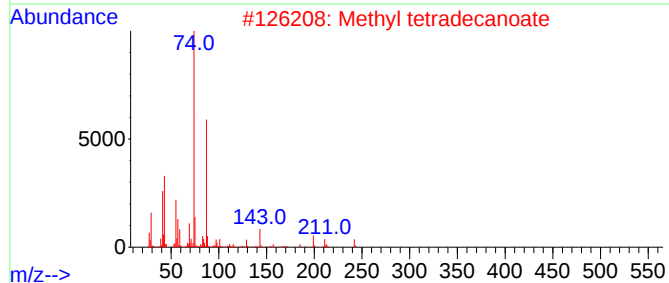
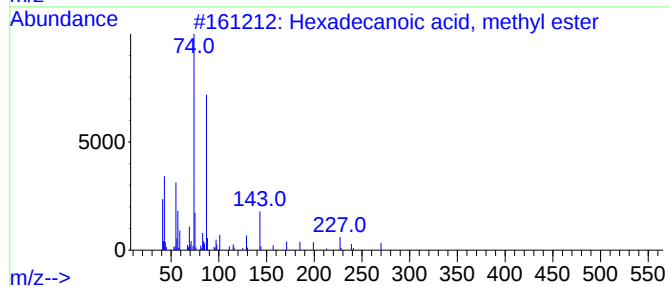
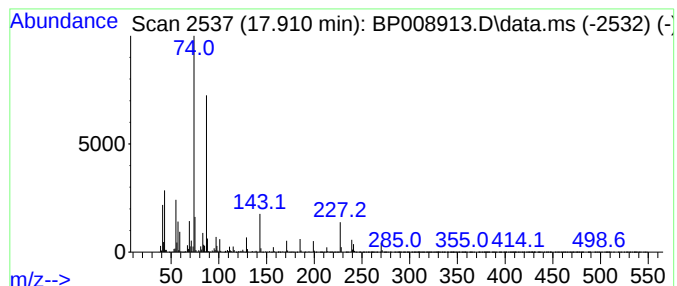
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 Hexadecanoic acid, methyl e... Concentration Rank 3

| R.T.   | EstConc  | Area    | Relative to ISTD | R.T.   |
|--------|----------|---------|------------------|--------|
| 17.910 | 12.83 ng | 1770870 | Phenanthrene-d10 | 17.234 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Hexadecanoic acid, methyl ester     | 270 | C17H34O2 | 000112-39-0 | 98   |
| 2    |    |   | Methyl tetradecanoate               | 242 | C15H30O2 | 000124-10-7 | 93   |
| 3    |    |   | Pentadecanoic acid, 14-methyl-, ... | 270 | C17H34O2 | 005129-60-2 | 93   |
| 4    |    |   | Tridecanoic acid, methyl ester      | 228 | C14H28O2 | 001731-88-0 | 89   |
| 5    |    |   | Decanoic acid, methyl ester         | 186 | C11H22O2 | 000110-42-9 | 81   |



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ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
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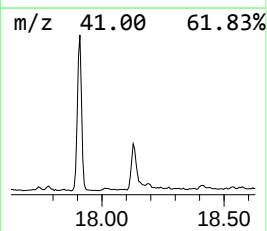
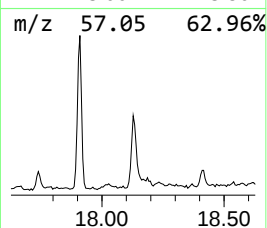
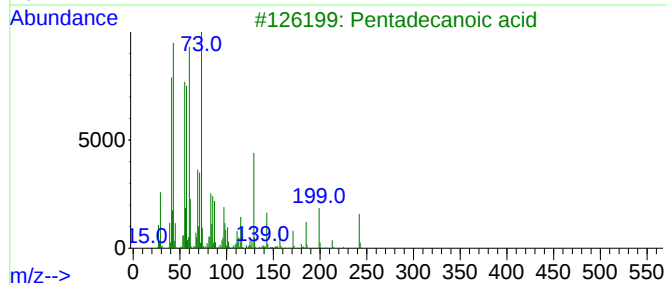
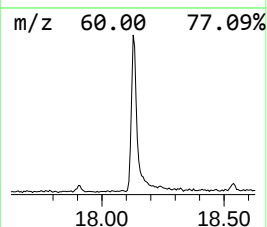
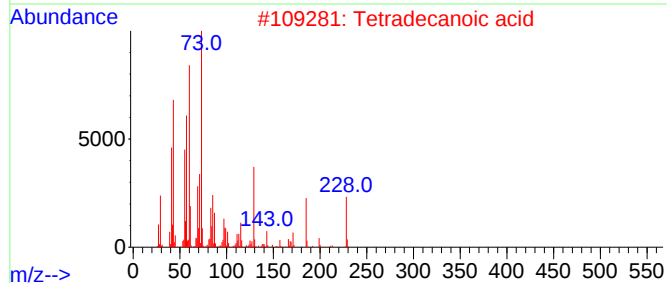
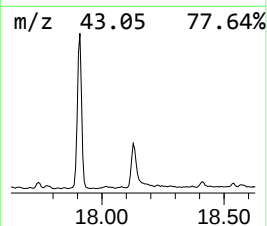
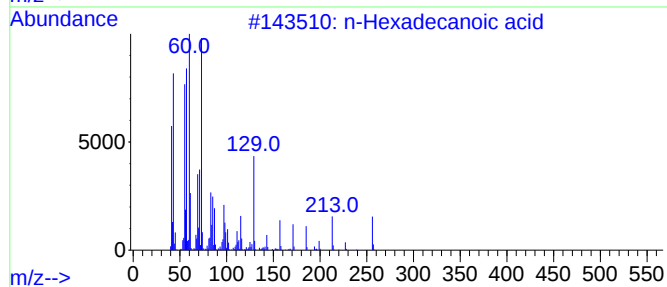
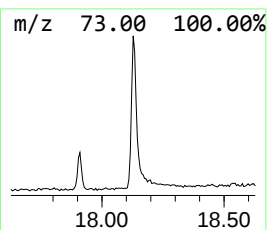
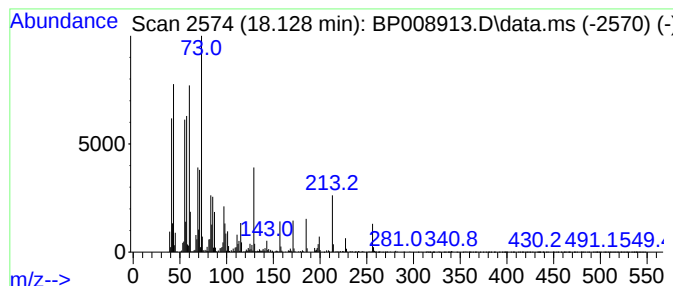
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP011422.M  
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TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 n-Hexadecanoic acid Concentration Rank 8

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 18.128 | 3.91 ng | 539381 | Phenanthrene-d10 | 17.234 |

| Hit# | of | 5 | Tentative ID        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|---------------------|-----|----------|-------------|------|
| 1    |    |   | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 99   |
| 2    |    |   | Tetradecanoic acid  | 228 | C14H28O2 | 000544-63-8 | 89   |
| 3    |    |   | Pentadecanoic acid  | 242 | C15H30O2 | 001002-84-2 | 81   |
| 4    |    |   | Octadecanoic acid   | 284 | C18H36O2 | 000057-11-4 | 76   |
| 5    |    |   | Tridecanoic acid    | 214 | C13H26O2 | 000638-53-9 | 76   |



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Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
RT3876

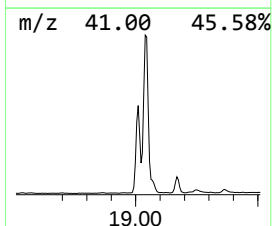
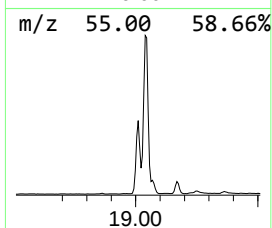
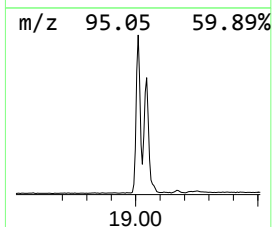
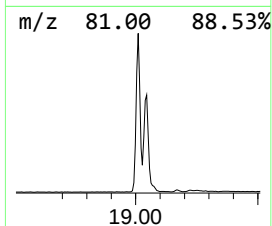
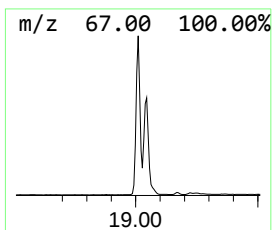
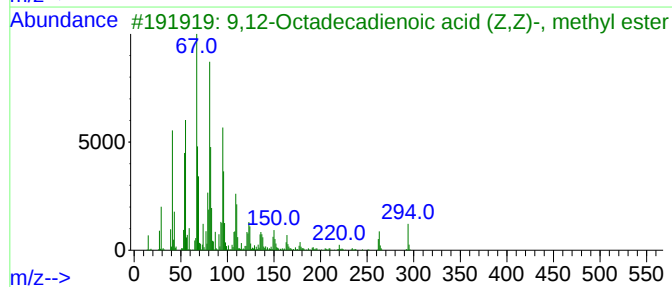
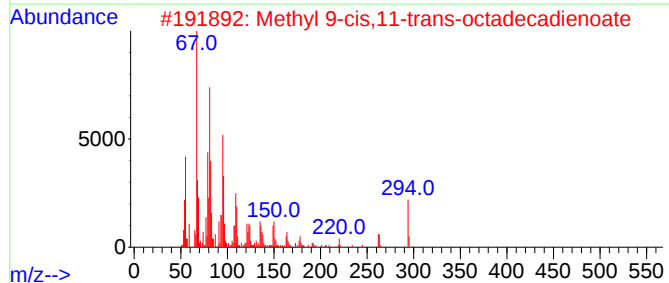
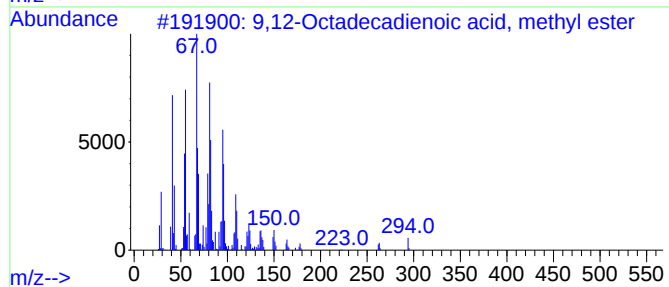
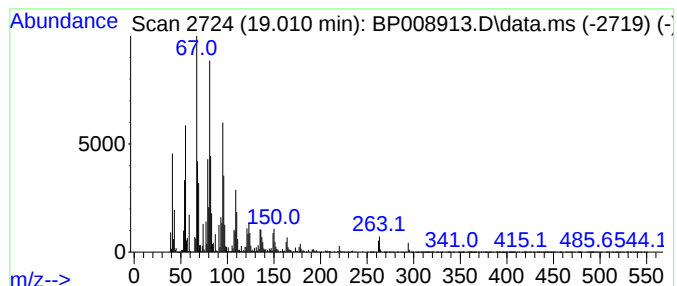
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TIC Library : C:\Database\NIST20.L  
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Peak Number 4 9,12-Octadecadienoic acid, ... Concentration Rank 2

| R.T.      | EstConc                             | Area    | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|---------|------------------|-------------|------|
| 19.010    | 32.56 ng                            | 4493320 | Phenanthrene-d10 | 17.234      |      |
| Hit# of 5 | Tentative ID                        | MW      | MolForm          | CAS#        | Qual |
| 1         | 9,12-Octadecadienoic acid, methy... | 294     | C19H34O2         | 002462-85-3 | 99   |
| 2         | Methyl 9-cis,11-trans-octadecadi... | 294     | C19H34O2         | 013058-52-1 | 99   |
| 3         | 9,12-Octadecadienoic acid (Z,Z)-... | 294     | C19H34O2         | 000112-63-0 | 99   |
| 4         | 9,15-Octadecadienoic acid, methy... | 294     | C19H34O2         | 017309-05-6 | 99   |
| 5         | 10,13-Octadecadienoic acid, meth... | 294     | C19H34O2         | 056554-62-2 | 99   |



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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP011422.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

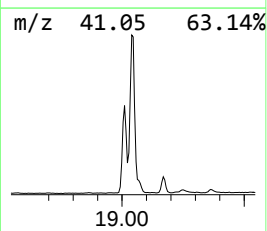
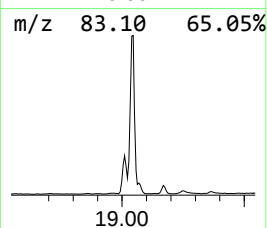
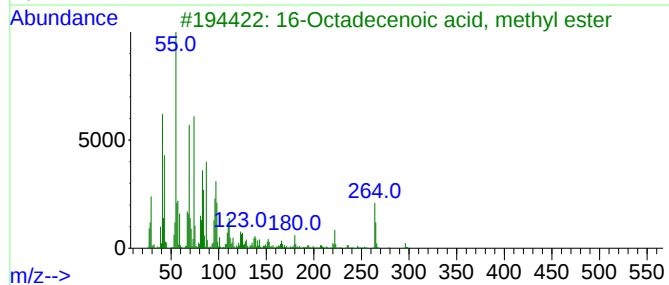
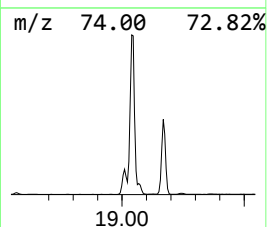
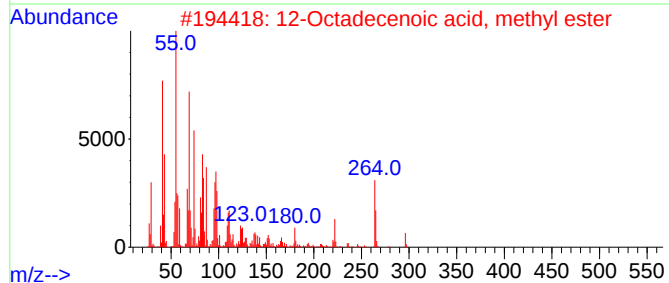
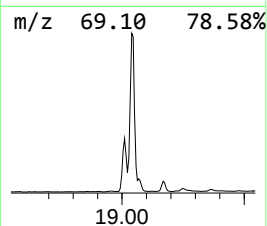
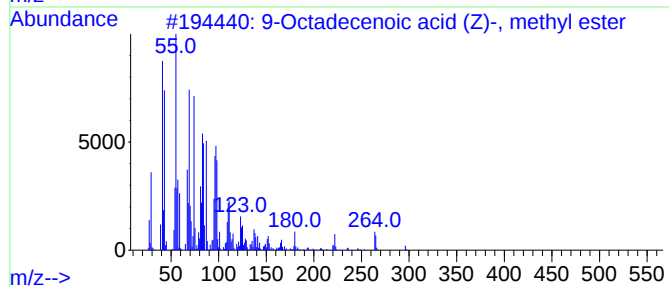
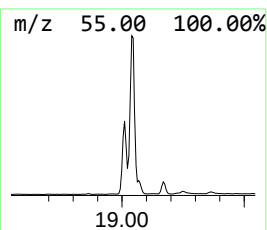
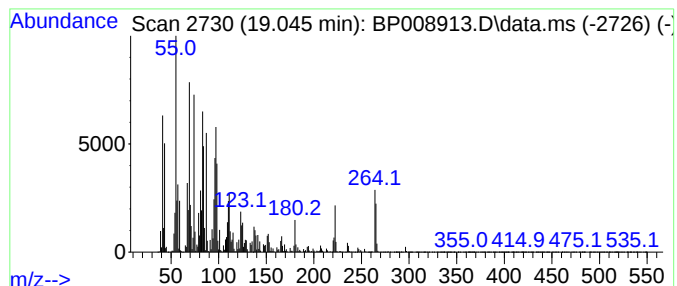
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 9-Octadecenoic acid (Z)-, m... Concentration Rank 1

| R.T.   | EstConc  | Area    | Relative to ISTD | R.T.   |
|--------|----------|---------|------------------|--------|
| 19.045 | 58.58 ng | 8084050 | Phenanthrene-d10 | 17.234 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 9-Octadecenoic acid (Z)-, methyl... | 296 | C19H36O2 | 000112-62-9 | 99   |
| 2    |    |   | 12-Octadecenoic acid, methyl ester  | 296 | C19H36O2 | 056554-46-2 | 99   |
| 3    |    |   | 16-Octadecenoic acid, methyl ester  | 296 | C19H36O2 | 056554-49-5 | 99   |
| 4    |    |   | 10-Octadecenoic acid, methyl ester  | 296 | C19H36O2 | 013481-95-3 | 99   |
| 5    |    |   | 11-Octadecenoic acid, methyl est... | 296 | C19H36O2 | 001937-63-9 | 99   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP012522\  
Data File : BP008913.D  
Acq On : 25 Jan 2022 18:16  
Operator : CG/JU  
Sample : N1233-01  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
RT3876

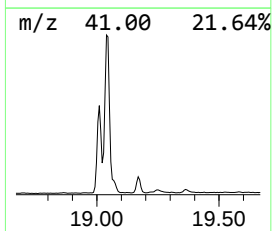
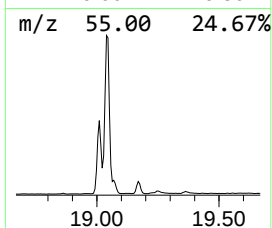
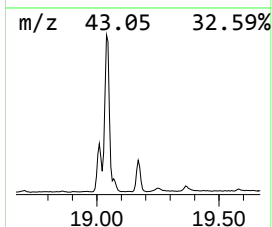
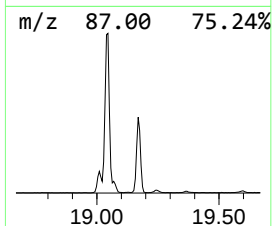
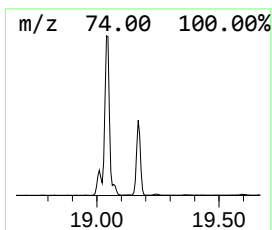
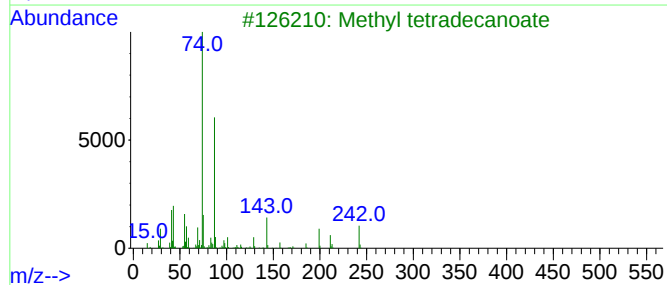
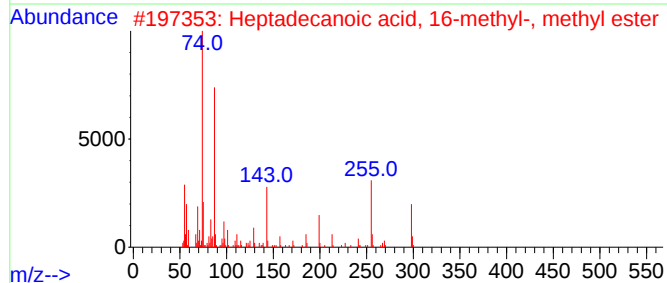
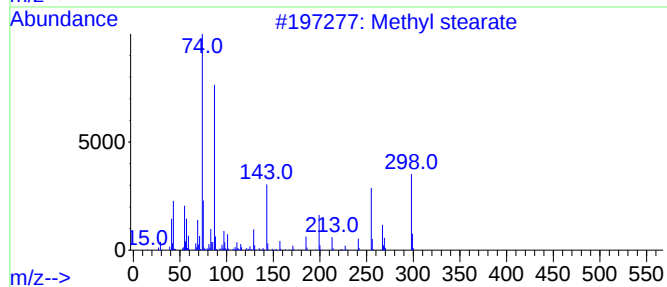
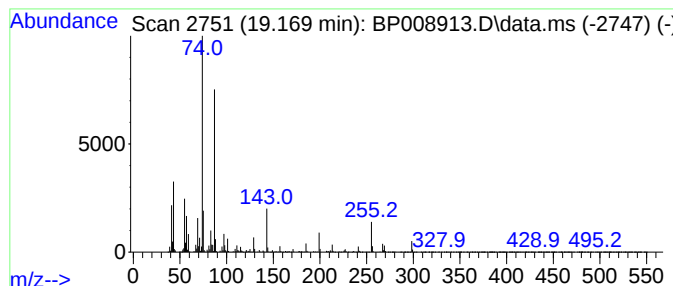
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 Methyl stearate Concentration Rank 7

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 19.169 | 5.93 ng | 818466 | Phenanthrene-d10 | 17.234 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Methyl stearate                     | 298 | C19H38O2 | 000112-61-8 | 98   |
| 2    |    |   | Heptadecanoic acid, 16-methyl-, ... | 298 | C19H38O2 | 005129-61-3 | 93   |
| 3    |    |   | Methyl tetradecanoate               | 242 | C15H30O2 | 000124-10-7 | 90   |
| 4    |    |   | Pentadecanoic acid, methyl ester    | 256 | C16H32O2 | 007132-64-1 | 87   |
| 5    |    |   | Hexadecanoic acid, methyl ester     | 270 | C17H34O2 | 000112-39-0 | 80   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP012522\  
Data File : BP008913.D  
Acq On : 25 Jan 2022 18:16  
Operator : CG/JU  
Sample : N1233-01  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
RT3876

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP011422.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

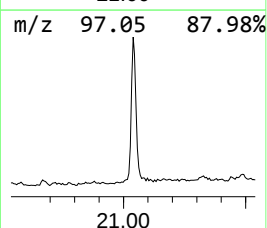
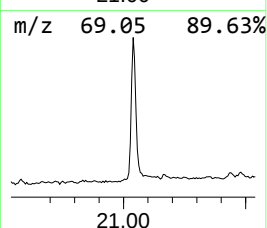
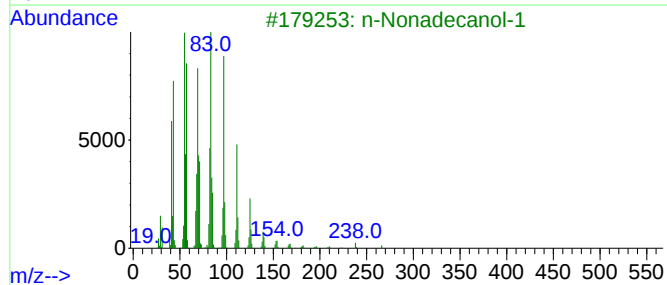
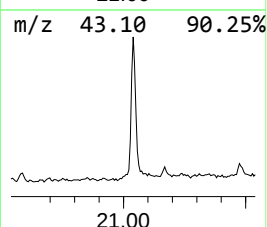
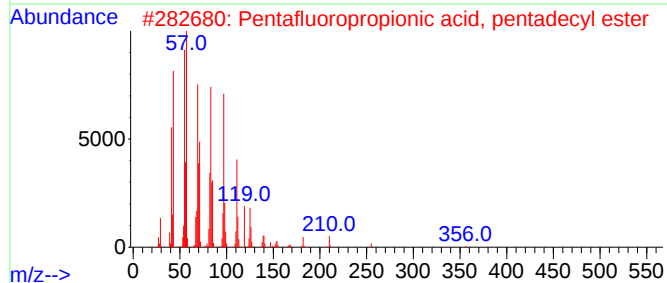
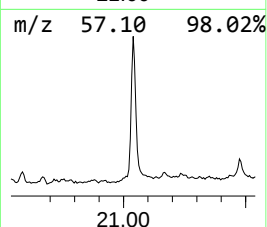
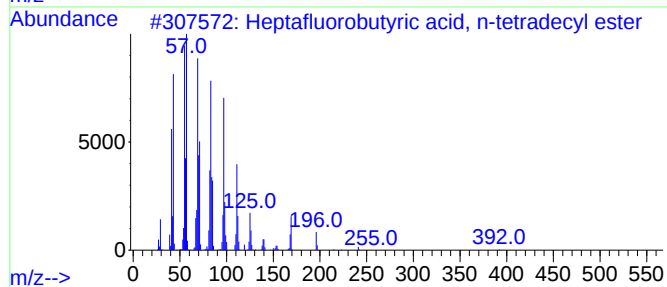
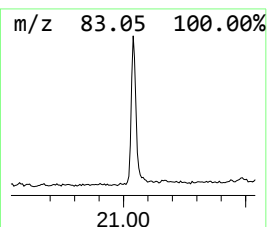
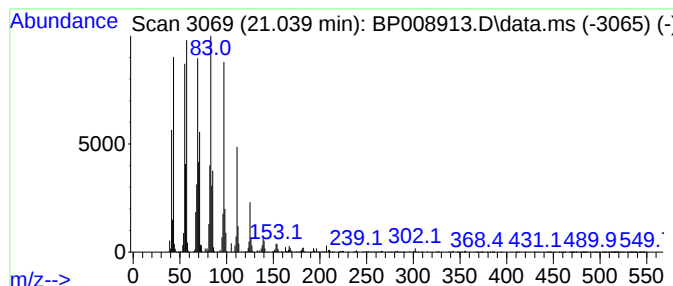
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 Heptafluorobutyric acid, n-... Concentration Rank 6

| R.T.   | EstConc | Area    | Relative to ISTD | R.T.   |
|--------|---------|---------|------------------|--------|
| 21.039 | 7.20 ng | 1294270 | Chrysene-d12     | 21.322 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|------------|-------------|------|
| 1    |    |   | Heptafluorobutyric acid, n-tetra... | 410 | C18H29F7O2 | 007365-36-8 | 94   |
| 2    |    |   | Pentafluoropropionic acid, penta... | 374 | C18H31F5O2 | 959092-08-1 | 94   |
| 3    |    |   | n-Nonadecanol-1                     | 284 | C19H40O    | 001454-84-8 | 94   |
| 4    |    |   | 1-Heneicosanol                      | 312 | C21H44O    | 015594-90-8 | 94   |
| 5    |    |   | Heptafluorobutyric acid, pentade... | 424 | C19H31F7O2 | 959261-23-5 | 94   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP012522\  
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Operator : CG/JU  
Sample : N1233-01  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
RT3876

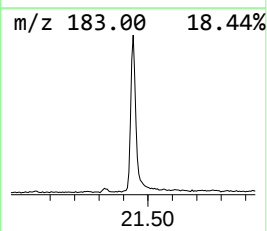
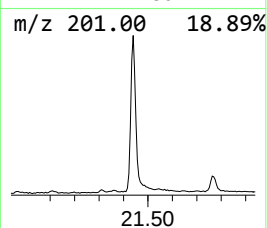
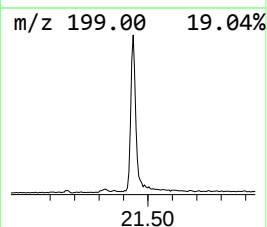
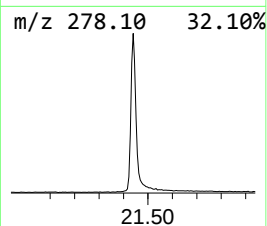
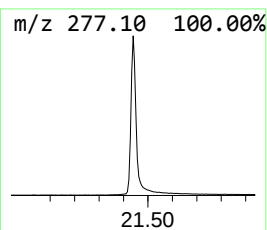
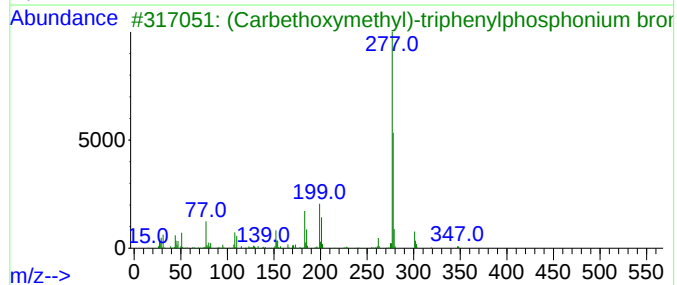
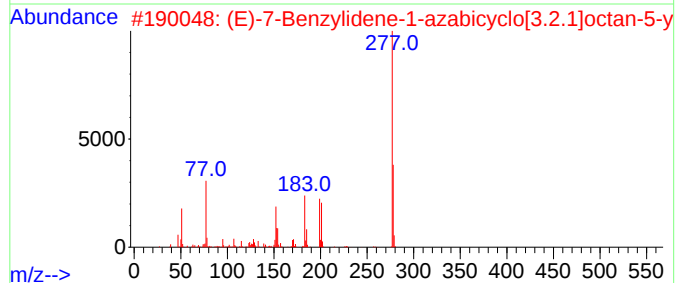
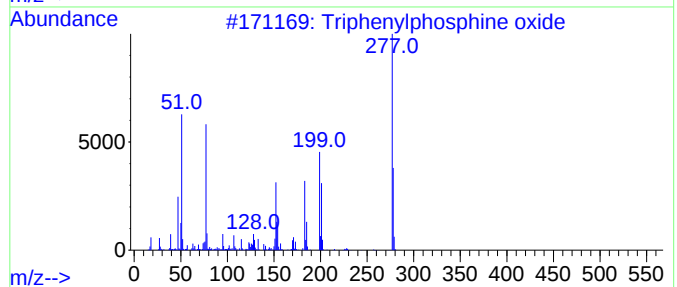
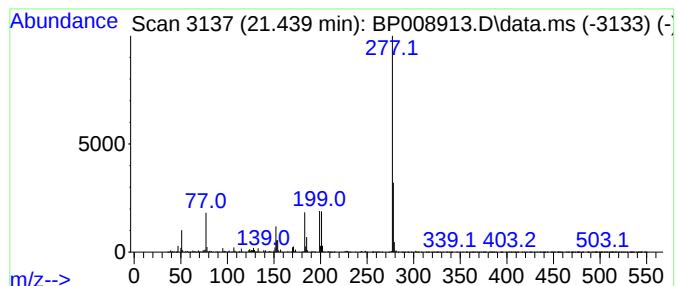
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP011422.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 Triphenylphosphine oxide Concentration Rank 4

| R.T.   | EstConc | Area    | Relative to ISTD | R.T.   |
|--------|---------|---------|------------------|--------|
| 21.439 | 9.22 ng | 1657640 | Chrysene-d12     | 21.322 |

| Hit# | of | 5 | Tentative ID                         | MW  | MolForm     | CAS#         | Qual |
|------|----|---|--------------------------------------|-----|-------------|--------------|------|
| 1    |    |   | Triphenylphosphine oxide             | 278 | C18H15OP    | 000791-28-6  | 95   |
| 2    |    |   | (E)-7-Benzylidene-1-azabicyclo[3...] | 293 | C15H19N03S  | 1000483-83-4 | 91   |
| 3    |    |   | (Carbethoxymethyl)-triphenylphos...  | 428 | C22H22BrO2P | 001530-45-6  | 38   |
| 4    |    |   | Vanadium, cycloheptatrienyl-(pen...  | 277 | C17H22V     | 1000155-20-5 | 33   |
| 5    |    |   | 5-Methyl-8-phenylfuro[3',2':5,6]...  | 277 | C16H11N3O2  | 1000460-07-9 | 28   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP012522\  
Data File : BP008913.D  
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Operator : CG/JU  
Sample : N1233-01  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
RT3876

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP011422.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

| TIC Top Hit name   | RT     | EstConc | Units | Response | --Internal Standard-- |        |         |      |
|--------------------|--------|---------|-------|----------|-----------------------|--------|---------|------|
|                    |        |         |       |          | #                     | RT     | Resp    | Conc |
| Butane, 2-metho... | 3.040  | 9.0     | ng    | 607811   | 1                     | 7.834  | 1345400 | 20.0 |
| Hexadecanoic ac... | 17.910 | 12.8    | ng    | 1770870  | 4                     | 17.234 | 2760180 | 20.0 |
| n-Hexadecanoic ... | 18.128 | 3.9     | ng    | 539381   | 4                     | 17.234 | 2760180 | 20.0 |
| 9,12-Octadecadi... | 19.010 | 32.6    | ng    | 4493320  | 4                     | 17.234 | 2760180 | 20.0 |
| 9-Octadecenoic ... | 19.045 | 58.6    | ng    | 8084050  | 4                     | 17.234 | 2760180 | 20.0 |
| Methyl stearate    | 19.169 | 5.9     | ng    | 818466   | 4                     | 17.234 | 2760180 | 20.0 |
| Heptafluorobuty... | 21.039 | 7.2     | ng    | 1294270  | 5                     | 21.322 | 3595450 | 20.0 |
| Triphenylphosph... | 21.439 | 9.2     | ng    | 1657640  | 5                     | 21.322 | 3595450 | 20.0 |