

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092721\  
 Data File : BP007219.D  
 Acq On : 27 Sep 2021 22:57  
 Operator : CG/JU  
 Sample : M3955-01  
 Misc :  
 ALS Vial : 19 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SS01-NA

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-BP092121.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP007219.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         | 5.463       | 397           | 404         | 411          | rBV      | 1958821        | 3073742       | 2.84%           | 0.654%        |
| 2         | 7.063       | 667           | 676         | 688          | rBV      | 2242010        | 3736243       | 3.46%           | 0.795%        |
| 3         | 7.881       | 808           | 815         | 822          | rBV      | 753038         | 1205100       | 1.12%           | 0.257%        |
| 4         | 9.046       | 1006          | 1013        | 1035         | rVB      | 1544514        | 2777412       | 2.57%           | 0.591%        |
| 5         | 10.687      | 1285          | 1292        | 1298         | rBV      | 1135026        | 1886541       | 1.75%           | 0.402%        |
| 6         | 12.363      | 1565          | 1577        | 1585         | rBV      | 958272         | 2015167       | 1.86%           | 0.429%        |
| 7         | 13.145      | 1703          | 1710        | 1717         | rVB      | 5500647        | 8146254       | 7.54%           | 1.734%        |
| 8         | 14.522      | 1936          | 1944        | 1948         | rVV      | 2094114        | 3274153       | 3.03%           | 0.697%        |
| 9         | 15.857      | 2166          | 2171        | 2175         | rBV      | 1133869        | 1577133       | 1.46%           | 0.336%        |
| 10        | 16.016      | 2192          | 2198        | 2202         | rVV      | 5129145        | 7668603       | 7.10%           | 1.632%        |
| 11        | 16.063      | 2202          | 2206        | 2217         | rVB      | 2896992        | 4819481       | 4.46%           | 1.026%        |
| 12        | 16.233      | 2229          | 2235        | 2249         | rVB      | 2929019        | 5015083       | 4.64%           | 1.067%        |
| 13        | 16.392      | 2257          | 2262        | 2269         | rBV2     | 1206237        | 2135374       | 1.98%           | 0.455%        |
| 14        | 16.598      | 2292          | 2297        | 2305         | rVB      | 838844         | 1386352       | 1.28%           | 0.295%        |
| 15        | 16.739      | 2305          | 2321        | 2328         | rBV      | 8044784        | 23687336      | 21.92%          | 5.042%        |
| 16        | 16.892      | 2342          | 2347        | 2358         | rVB      | 4040401        | 6247238       | 5.78%           | 1.330%        |
| 17        | 17.075      | 2374          | 2378        | 2383         | rVB4     | 763742         | 1216207       | 1.13%           | 0.259%        |
| 18        | 17.210      | 2393          | 2401        | 2406         | rBV2     | 4379755        | 8513517       | 7.88%           | 1.812%        |
| 19        | 17.269      | 2406          | 2411        | 2415         | rBV2     | 5977527        | 8601015       | 7.96%           | 1.831%        |
| 20        | 17.322      | 2415          | 2420        | 2424         | rVV2     | 12297812       | 19032616      | 17.61%          | 4.051%        |
| 21        | 17.363      | 2424          | 2427        | 2433         | rVB2     | 6429202        | 9546267       | 8.83%           | 2.032%        |
| 22        | 17.498      | 2441          | 2450        | 2454         | rBV      | 8785118        | 17490762      | 16.19%          | 3.723%        |
| 23        | 17.533      | 2454          | 2456        | 2468         | rVV3     | 1576122        | 3060037       | 2.83%           | 0.651%        |
| 24        | 17.645      | 2469          | 2475        | 2476         | rVV      | 6315381        | 8788100       | 8.13%           | 1.871%        |
| 25        | 17.669      | 2476          | 2479        | 2484         | rVB      | 9272355        | 13054464      | 12.08%          | 2.779%        |
| 26        | 17.939      | 2520          | 2525        | 2531         | rVV5     | 1268980        | 2374766       | 2.20%           | 0.505%        |
| 27        | 18.110      | 2542          | 2554        | 2566         | rVV2     | 9682556        | 52152905      | 48.26%          | 11.101%       |
| 28        | 18.322      | 2567          | 2590        | 2593         | rVV3     | 16899295       | 108060585     | 100.00%         | 23.001%       |
| 29        | 18.369      | 2595          | 2598        | 2603         | rVB      | 2799052        | 3710101       | 3.43%           | 0.790%        |
| 30        | 18.510      | 2618          | 2622        | 2631         | rVB      | 1789796        | 3141135       | 2.91%           | 0.669%        |
| 31        | 18.622      | 2637          | 2641        | 2646         | rBV2     | 2091395        | 2874669       | 2.66%           | 0.612%        |
| 32        | 18.680      | 2647          | 2651        | 2656         | rVB3     | 1283176        | 1794858       | 1.66%           | 0.382%        |
| 33        | 18.739      | 2657          | 2661        | 2663         | rBV      | 1569767        | 2031577       | 1.88%           | 0.432%        |
| 34        | 18.774      | 2664          | 2667        | 2673         | rVB2     | 1395485        | 2246138       | 2.08%           | 0.478%        |
| 35        | 18.845      | 2674          | 2679        | 2681         | rBV      | 3098487        | 4171904       | 3.86%           | 0.888%        |
| 36        | 18.998      | 2701          | 2705        | 2711         | rBV      | 4148533        | 5715749       | 5.29%           | 1.217%        |

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 Sample : M3955-01  
 Misc :  
 ALS Vial : 19 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SS01-NA

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-BP092121.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |          |        |        |
|----|--------|------|------|------|------|---------|----------|--------|--------|
| 37 | 19.057 | 2712 | 2715 | 2718 | rVB  | 1205220 | 1390009  | 1.29%  | 0.296% |
| 38 | 19.151 | 2726 | 2731 | 2735 | rBV  | 5295468 | 7541625  | 6.98%  | 1.605% |
| 39 | 19.286 | 2749 | 2754 | 2756 | rBV2 | 4632562 | 7519890  | 6.96%  | 1.601% |
| 40 | 19.339 | 2760 | 2763 | 2765 | rVV2 | 2976347 | 4379941  | 4.05%  | 0.932% |
| 41 | 19.439 | 2775 | 2780 | 2792 | rVB2 | 5834650 | 9472874  | 8.77%  | 2.016% |
| 42 | 19.580 | 2801 | 2804 | 2807 | rBV2 | 1429713 | 1636114  | 1.51%  | 0.348% |
| 43 | 19.839 | 2843 | 2848 | 2854 | rVB  | 9032005 | 10708684 | 9.91%  | 2.279% |
| 44 | 20.110 | 2891 | 2894 | 2896 | rBV2 | 1437796 | 1947294  | 1.80%  | 0.414% |
| 45 | 20.133 | 2896 | 2898 | 2902 | rVB  | 2482533 | 2547312  | 2.36%  | 0.542% |
| 46 | 20.674 | 2988 | 2990 | 2997 | rVB  | 2615891 | 3246797  | 3.00%  | 0.691% |
| 47 | 21.068 | 3054 | 3057 | 3061 | rVB  | 2608462 | 2984740  | 2.76%  | 0.635% |
| 48 | 23.045 | 3389 | 3393 | 3396 | rBV2 | 2982017 | 4776886  | 4.42%  | 1.017% |
| 49 | 23.092 | 3397 | 3401 | 3408 | rVB2 | 5416683 | 9404883  | 8.70%  | 2.002% |
| 50 | 24.421 | 3622 | 3627 | 3634 | rVB  | 3728033 | 8069815  | 7.47%  | 1.718% |
| 51 | 25.256 | 3761 | 3769 | 3783 | rVB5 | 7077763 | 20188478 | 18.68% | 4.297% |
| 52 | 26.586 | 3986 | 3995 | 4004 | rVB4 | 5810249 | 17765513 | 16.44% | 3.781% |

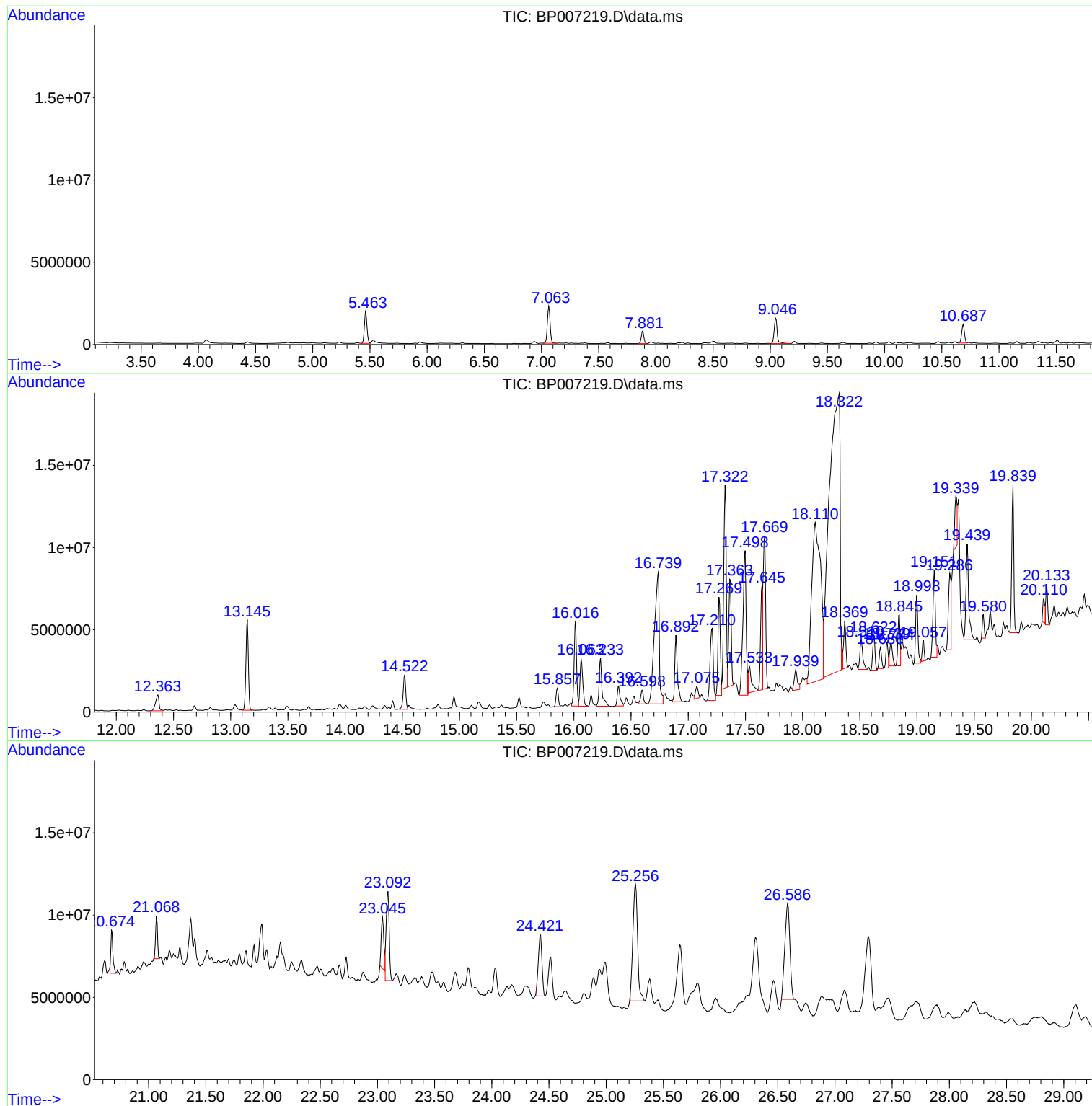
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Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092721\  
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Sample : M3955-01  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SS01-NA

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP092121.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092721\  
Data File : BP007219.D  
Acq On : 27 Sep 2021 22:57  
Operator : CG/JU  
Sample : M3955-01  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SS01-NA

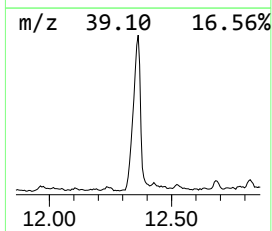
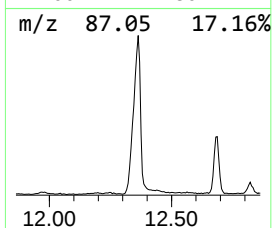
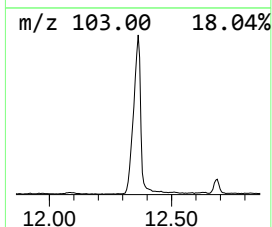
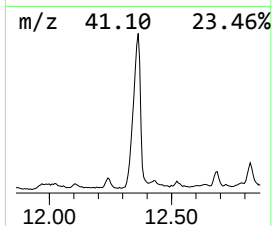
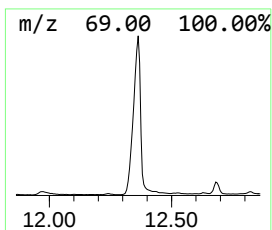
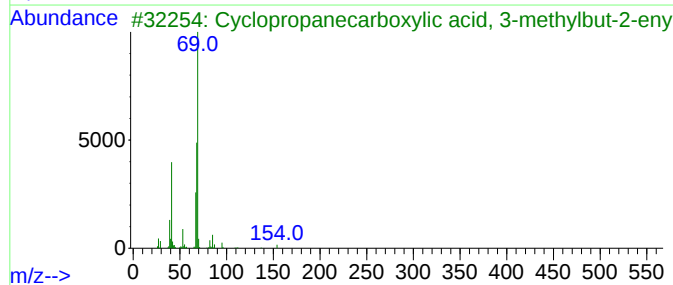
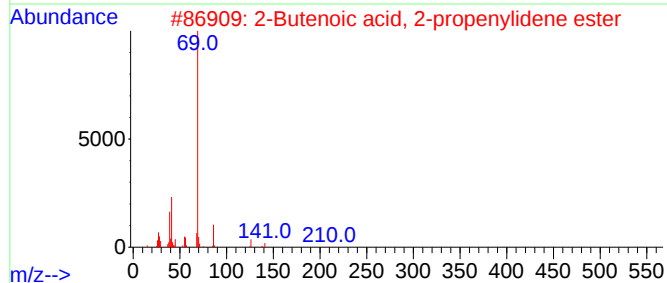
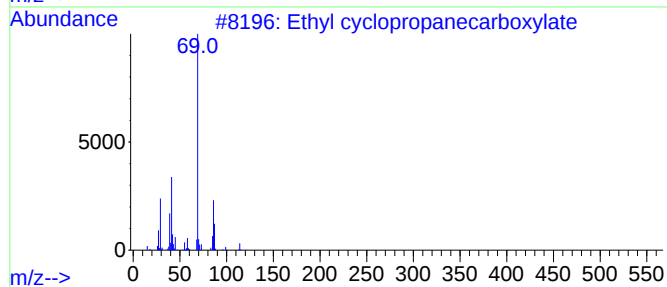
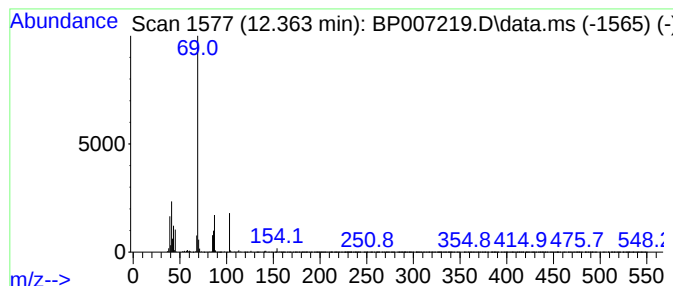
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP092121.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 1 Ethyl cyclopropanecarboxylate Concentration Rank 11

| R.T.   | EstConc  | Area    | Relative to ISTD | R.T.   |
|--------|----------|---------|------------------|--------|
| 12.363 | 21.36 ng | 2015170 | Naphthalene-d8   | 10.687 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Ethyl cyclopropanecarboxylate       | 114 | C6H10O2  | 004606-07-9  | 50   |
| 2    |    |   | 2-Butenoic acid, 2-propenylidene... | 210 | C11H14O4 | 055030-70-1  | 42   |
| 3    |    |   | Cyclopropanecarboxylic acid, 3-m... | 154 | C9H14O2  | 1000299-37-4 | 9    |
| 4    |    |   | 2-Butenoic acid, 2-propenyl ester   | 126 | C7H10O2  | 020474-93-5  | 9    |
| 5    |    |   | Cyclopropanecarboxylic acid, 2-m... | 170 | C10H18O2 | 1000354-65-8 | 9    |



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ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SS01-NA

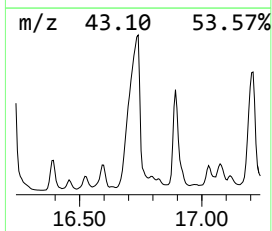
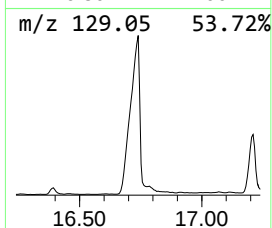
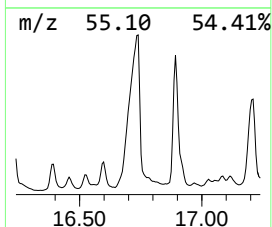
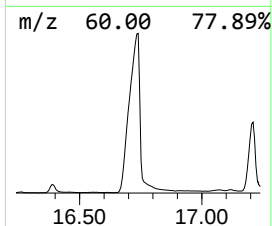
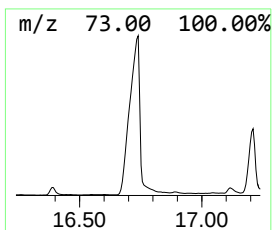
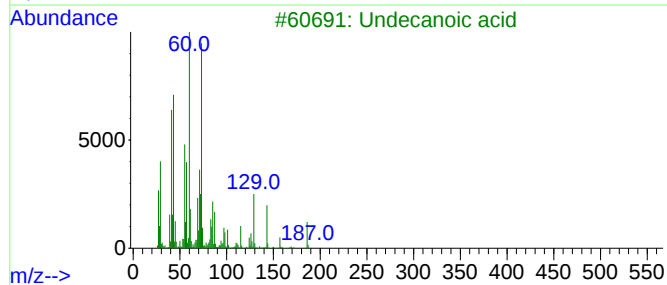
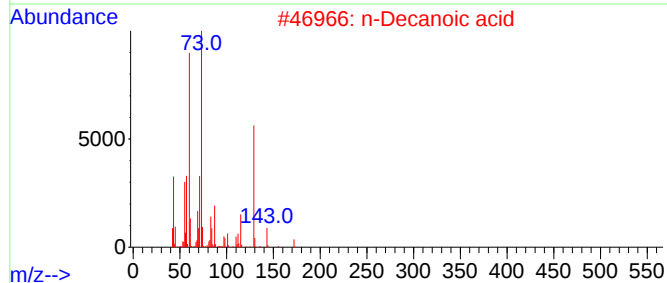
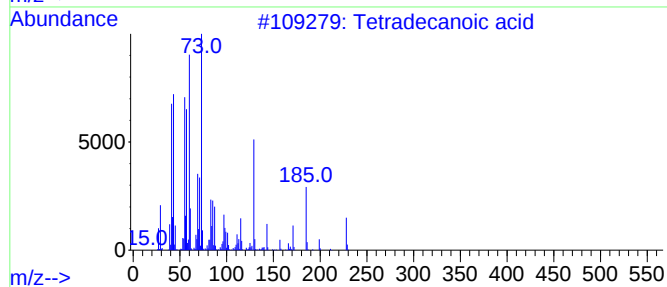
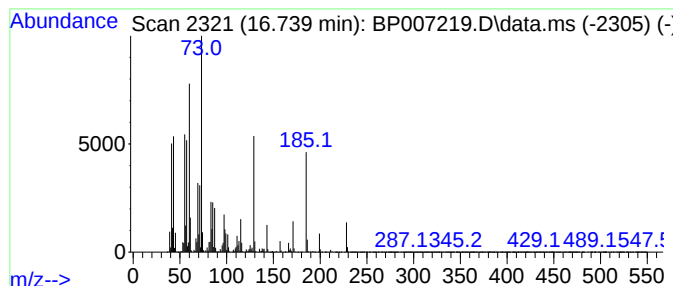
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP092121.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 2 Tetradecanoic acid Concentration Rank 4

| R.T.   | EstConc  | Area     | Relative to ISTD | R.T.   |
|--------|----------|----------|------------------|--------|
| 16.739 | 55.08 ng | 23687300 | Phenanthrene-d10 | 17.275 |

| Hit# | of | 5 | Tentative ID       | MW  | MolForm  | CAS#        | Qual |
|------|----|---|--------------------|-----|----------|-------------|------|
| 1    |    |   | Tetradecanoic acid | 228 | C14H28O2 | 000544-63-8 | 99   |
| 2    |    |   | n-Decanoic acid    | 172 | C10H20O2 | 000334-48-5 | 64   |
| 3    |    |   | Undecanoic acid    | 186 | C11H22O2 | 000112-37-8 | 58   |
| 4    |    |   | Dodecanoic acid    | 200 | C12H24O2 | 000143-07-7 | 58   |
| 5    |    |   | Tridecanoic acid   | 214 | C13H26O2 | 000638-53-9 | 52   |



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Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SS01-NA

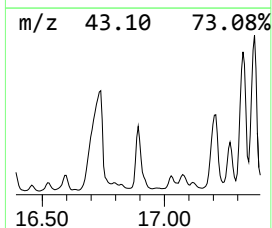
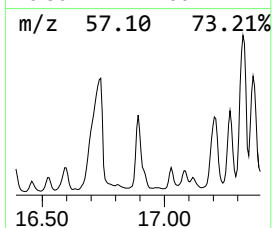
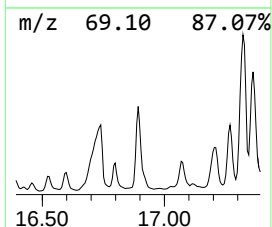
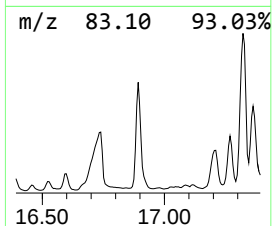
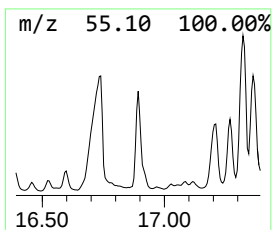
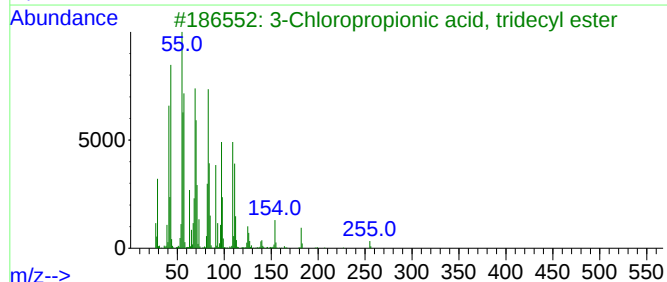
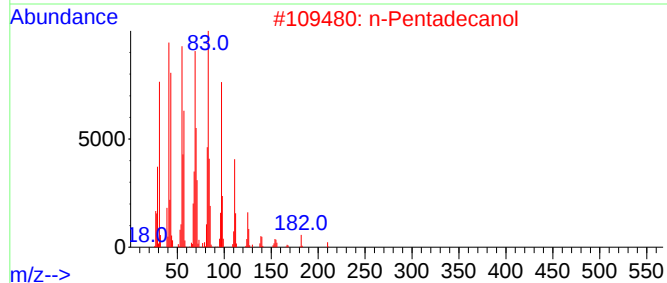
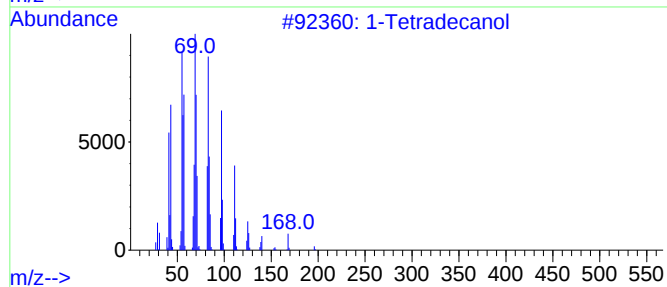
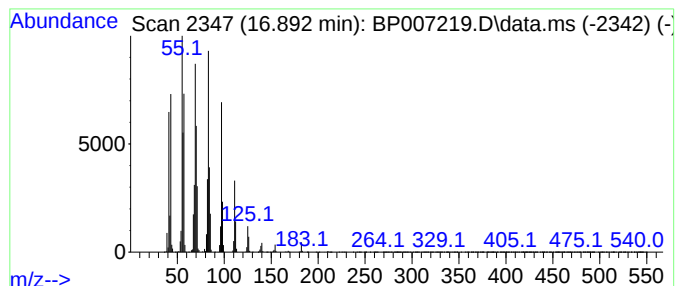
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP092121.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 3 1-Tetradecanol Concentration Rank 18

| R.T.   | EstConc  | Area    | Relative to ISTD | R.T.   |
|--------|----------|---------|------------------|--------|
| 16.892 | 14.53 ng | 6247240 | Phenanthrene-d10 | 17.275 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | 1-Tetradecanol                      | 214 | C14H30O    | 000112-72-1  | 94   |
| 2    |    |   | n-Pentadecanol                      | 228 | C15H32O    | 000629-76-5  | 94   |
| 3    |    |   | 3-Chloropropionic acid, tridecyl... | 290 | C16H31ClO2 | 1000281-77-4 | 91   |
| 4    |    |   | n-Tridecan-1-ol                     | 200 | C13H28O    | 000112-70-9  | 91   |
| 5    |    |   | 1-Hexadecanol                       | 242 | C16H34O    | 036653-82-4  | 91   |



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Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SS01-NA

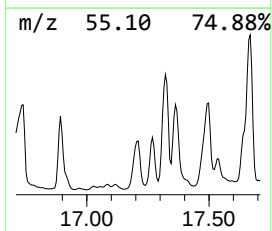
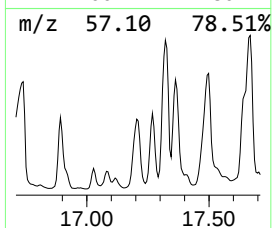
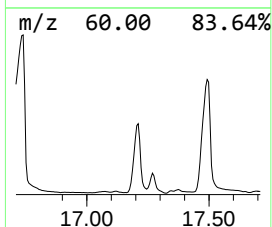
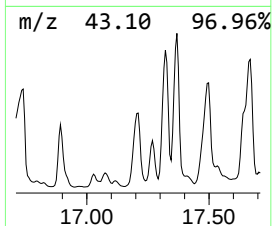
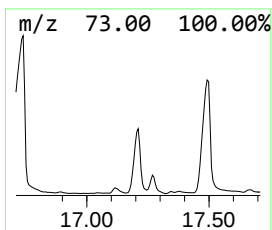
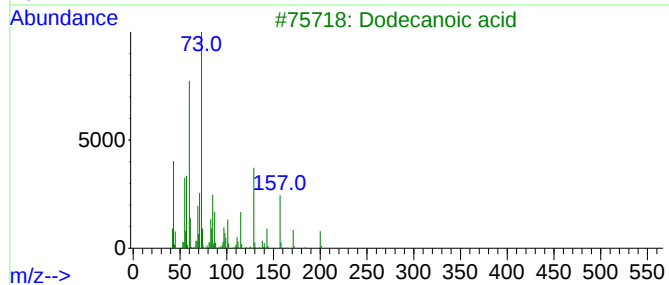
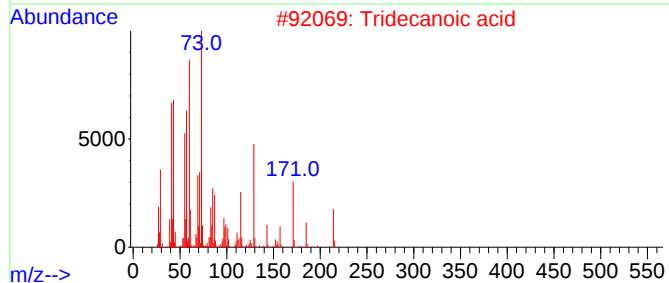
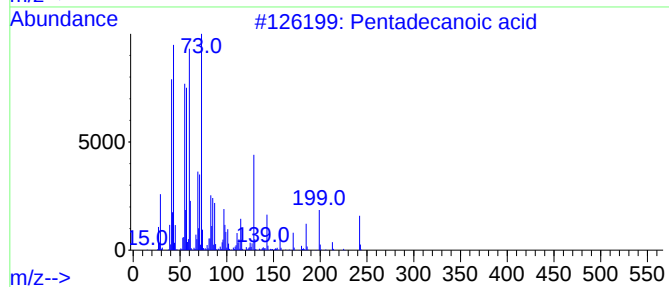
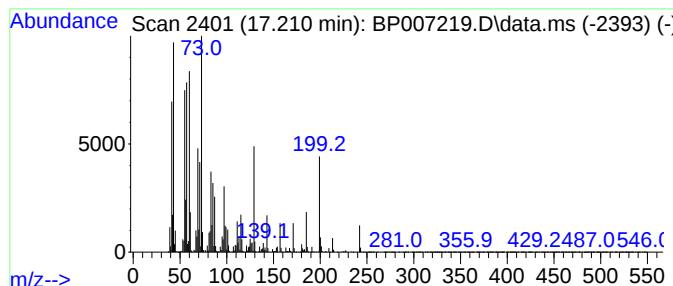
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 4 Pentadecanoic acid Concentration Rank 15

| R.T.   | EstConc  | Area    | Relative to ISTD | R.T.   |
|--------|----------|---------|------------------|--------|
| 17.210 | 19.80 ng | 8513520 | Phenanthrene-d10 | 17.275 |

| Hit# | of | 5 | Tentative ID       | MW  | MolForm  | CAS#        | Qual |
|------|----|---|--------------------|-----|----------|-------------|------|
| 1    |    |   | Pentadecanoic acid | 242 | C15H30O2 | 001002-84-2 | 97   |
| 2    |    |   | Tridecanoic acid   | 214 | C13H26O2 | 000638-53-9 | 72   |
| 3    |    |   | Dodecanoic acid    | 200 | C12H24O2 | 000143-07-7 | 70   |
| 4    |    |   | Tetradecanoic acid | 228 | C14H28O2 | 000544-63-8 | 64   |
| 5    |    |   | Undecanoic acid    | 186 | C11H22O2 | 000112-37-8 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092721\  
Data File : BP007219.D  
Acq On : 27 Sep 2021 22:57  
Operator : CG/JU  
Sample : M3955-01  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SS01-NA

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP092121.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

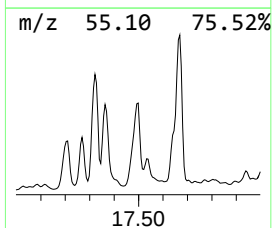
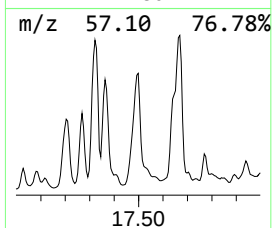
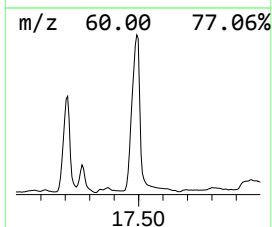
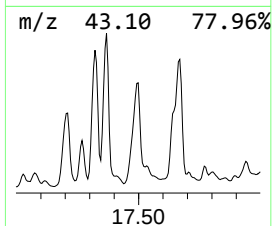
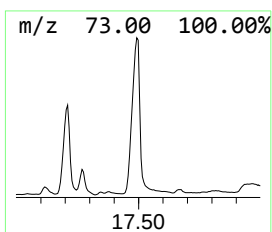
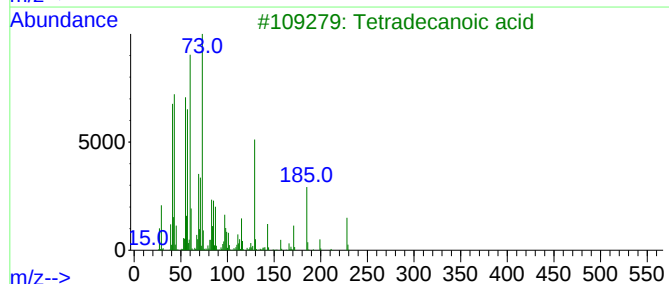
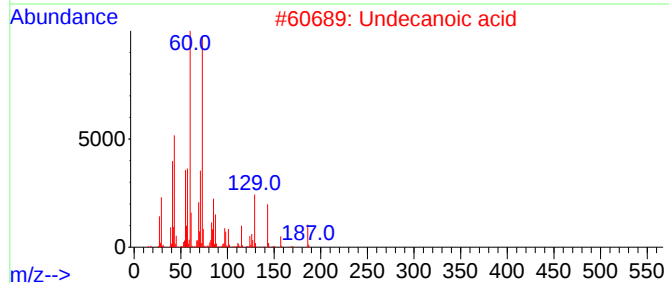
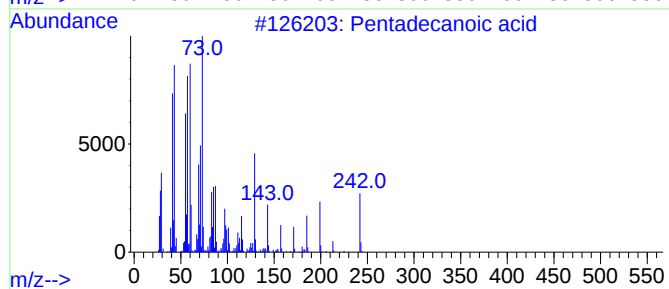
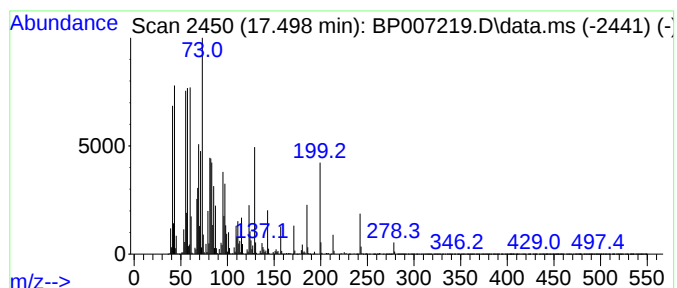
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Peak Number 5 Undecanoic acid Concentration Rank 7

| R.T.   | EstConc  | Area     | Relative to ISTD | R.T.   |
|--------|----------|----------|------------------|--------|
| 17.498 | 40.67 ng | 17490800 | Phenanthrene-d10 | 17.275 |

| Hit# | of | 5 | Tentative ID       | MW  | MolForm  | CAS#        | Qual |
|------|----|---|--------------------|-----|----------|-------------|------|
| 1    |    |   | Pentadecanoic acid | 242 | C15H30O2 | 001002-84-2 | 98   |
| 2    |    |   | Undecanoic acid    | 186 | C11H22O2 | 000112-37-8 | 55   |
| 3    |    |   | Tetradecanoic acid | 228 | C14H28O2 | 000544-63-8 | 52   |
| 4    |    |   | n-Decanoic acid    | 172 | C10H20O2 | 000334-48-5 | 50   |
| 5    |    |   | Tridecanoic acid   | 214 | C13H26O2 | 000638-53-9 | 49   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092721\  
Data File : BP007219.D  
Acq On : 27 Sep 2021 22:57  
Operator : CG/JU  
Sample : M3955-01  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SS01-NA

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP092121.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

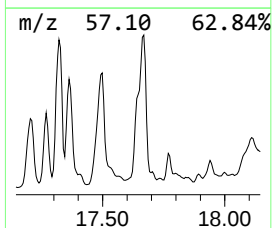
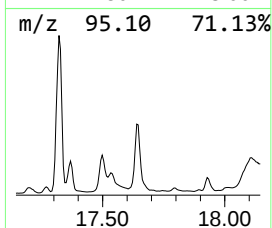
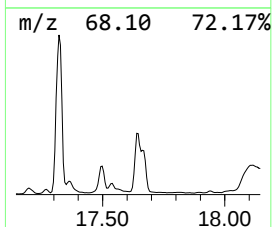
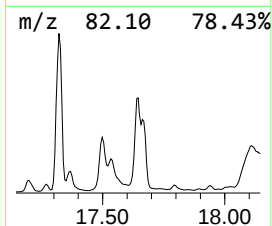
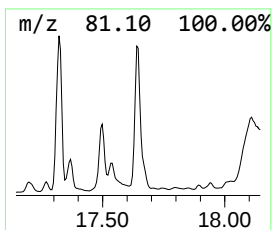
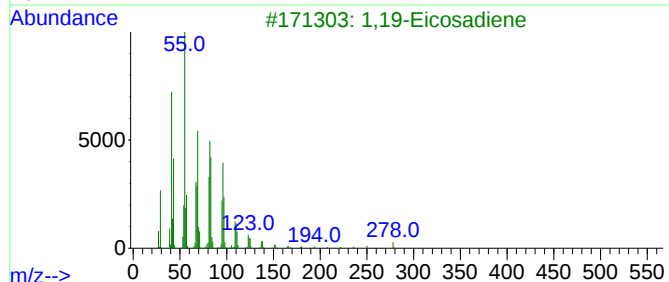
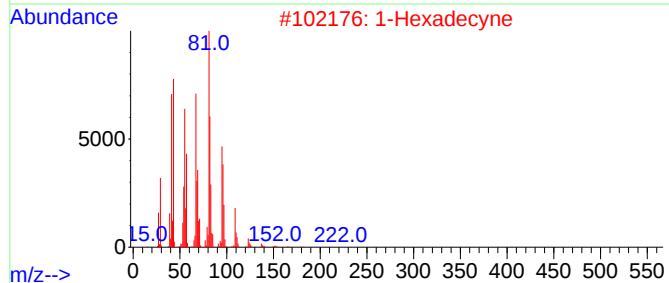
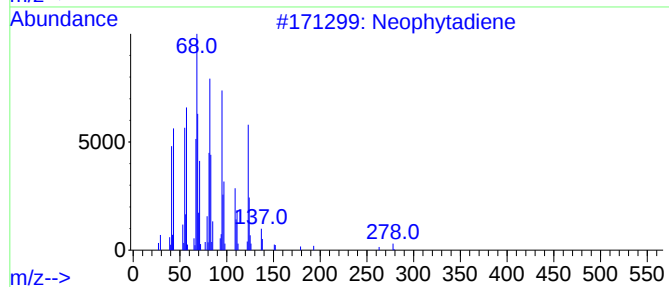
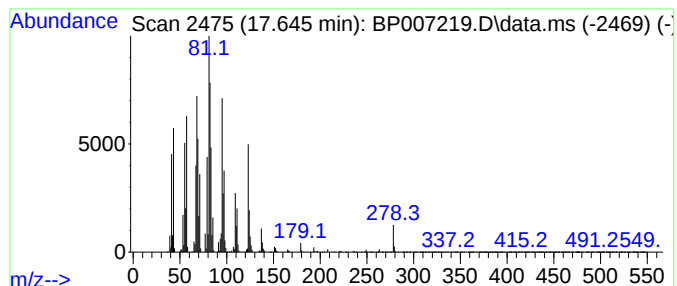
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 Neophytadiene Concentration Rank 12

| R.T.   | EstConc  | Area    | Relative to ISTD | R.T.   |
|--------|----------|---------|------------------|--------|
| 17.645 | 20.43 ng | 8788100 | Phenanthrene-d10 | 17.275 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------|-----|---------|--------------|------|
| 1    |    |   | Neophytadiene                 | 278 | C20H38  | 000504-96-1  | 70   |
| 2    |    |   | 1-Hexadecyne                  | 222 | C16H30  | 000629-74-3  | 53   |
| 3    |    |   | 1,19-Eicosadiene              | 278 | C20H38  | 014811-95-1  | 43   |
| 4    |    |   | Bicyclo[10.8.0]eicosane, cis- | 278 | C20H38  | 1000155-82-2 | 43   |
| 5    |    |   | Cyclohexanol, 1-ethynyl-      | 124 | C8H12O  | 000078-27-3  | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092721\  
Data File : BP007219.D  
Acq On : 27 Sep 2021 22:57  
Operator : CG/JU  
Sample : M3955-01  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SS01-NA

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP092121.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

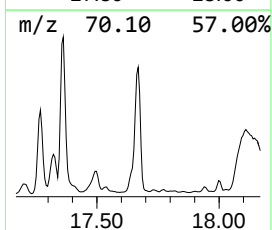
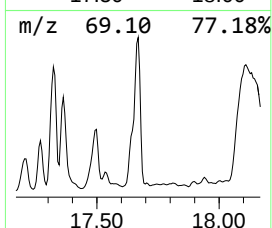
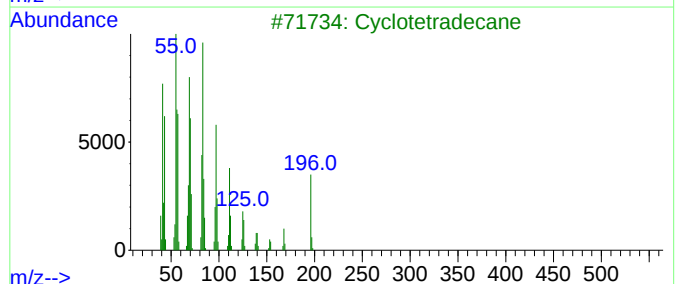
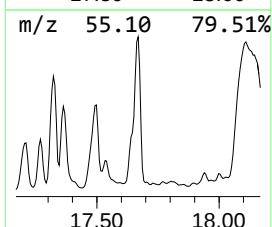
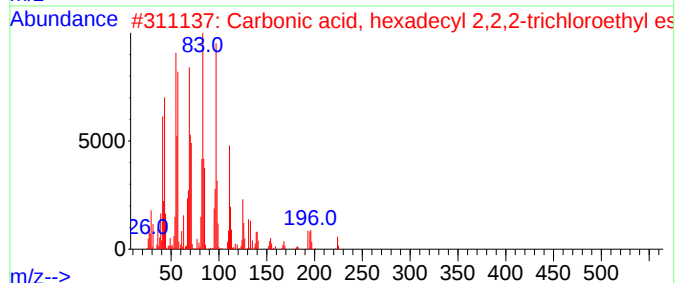
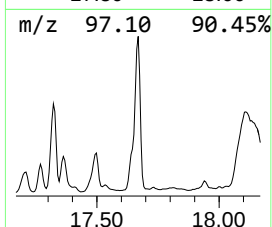
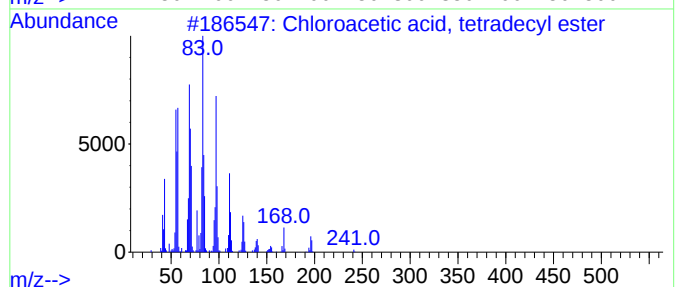
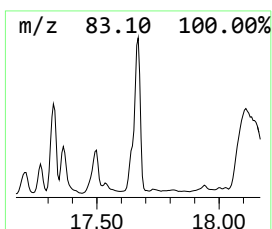
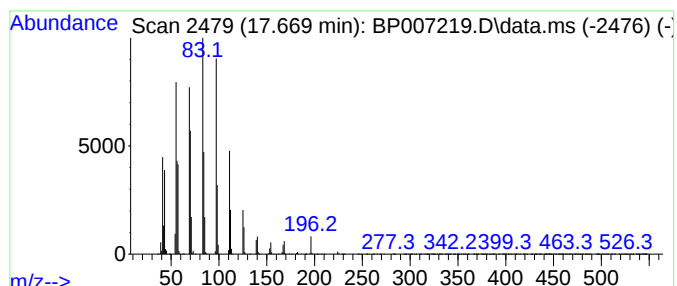
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 Chloroacetic acid, tetradec... Concentration Rank 8

| R.T.   | EstConc  | Area     | Relative to ISTD | R.T.   |
|--------|----------|----------|------------------|--------|
| 17.669 | 30.36 ng | 13054500 | Phenanthrene-d10 | 17.275 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-------------|--------------|------|
| 1    |    |   | Chloroacetic acid, tetradecyl ester | 290 | C16H31ClO2  | 018277-86-6  | 91   |
| 2    |    |   | Carbonic acid, hexadecyl 2,2,2-t... | 416 | C19H35Cl3O3 | 1000314-56-2 | 90   |
| 3    |    |   | Cyclotetradecane                    | 196 | C14H28      | 000295-17-0  | 83   |
| 4    |    |   | 1-Dodecanol                         | 186 | C12H26O     | 000112-53-8  | 72   |
| 5    |    |   | Cyclohexane, 1,1,2-trimethyl-       | 126 | C9H18       | 007094-26-0  | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092721\  
Data File : BP007219.D  
Acq On : 27 Sep 2021 22:57  
Operator : CG/JU  
Sample : M3955-01  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

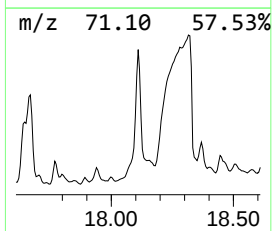
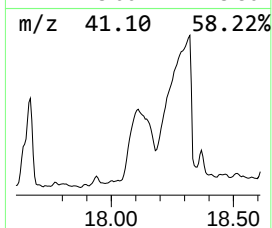
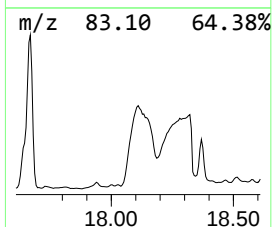
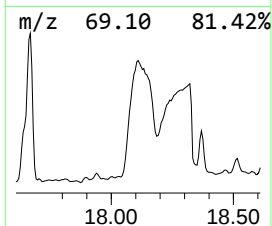
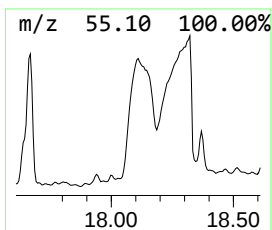
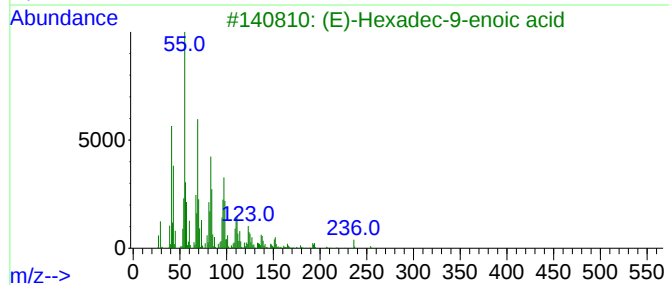
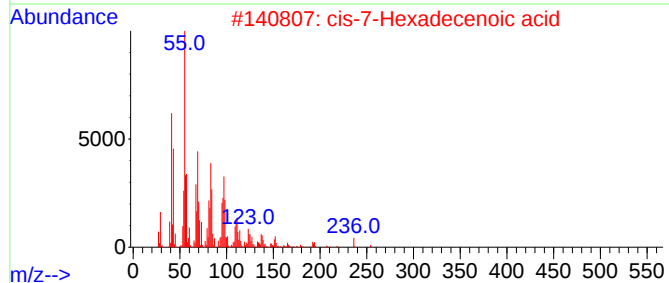
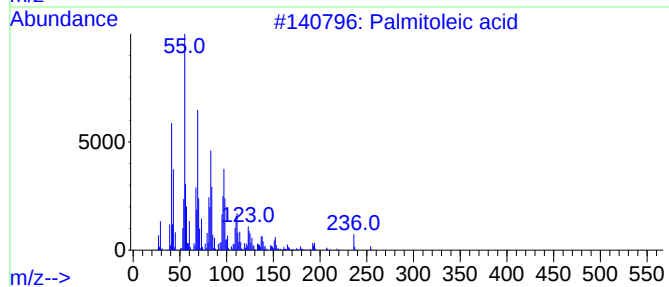
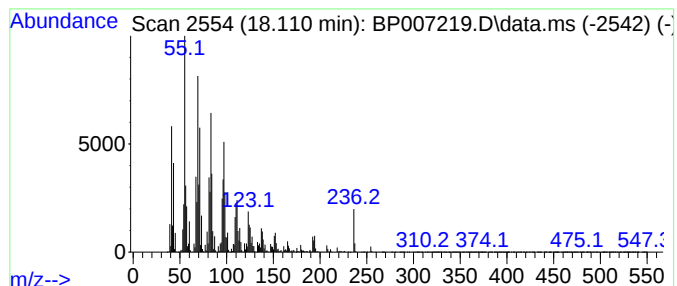
Instrument :  
BNA\_P  
ClientSampleId :  
SS01-NA

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP092121.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 Palmitoleic acid Concentration Rank 2

| R.T.      | EstConc                             | Area     | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|----------|------------------|-------------|------|
| 18.110    | 121.27 ng                           | 52152900 | Phenanthrene-d10 | 17.275      |      |
| Hit# of 5 | Tentative ID                        | MW       | MolForm          | CAS#        | Qual |
| 1         | Palmitoleic acid                    | 254      | C16H30O2         | 000373-49-9 | 99   |
| 2         | cis-7-Hexadecenoic acid             | 254      | C16H30O2         | 002416-19-5 | 92   |
| 3         | (E)-Hexadec-9-enoic acid            | 254      | C16H30O2         | 010030-73-6 | 92   |
| 4         | 6-Pentadecenoic acid, 13-methyl-... | 254      | C16H30O2         | 682751-34-4 | 91   |
| 5         | 4-n-Hexylthiane, S,S-dioxide        | 218      | C11H22O2S        | 070928-52-8 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092721\  
Data File : BP007219.D  
Acq On : 27 Sep 2021 22:57  
Operator : CG/JU  
Sample : M3955-01  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SS01-NA

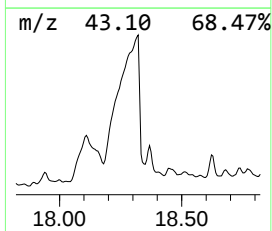
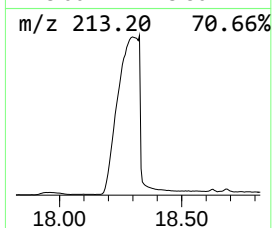
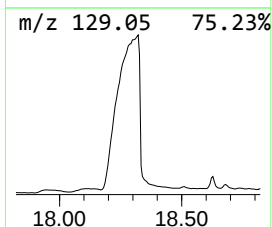
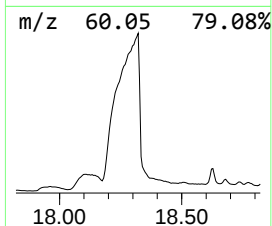
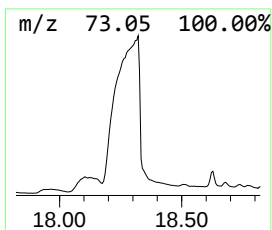
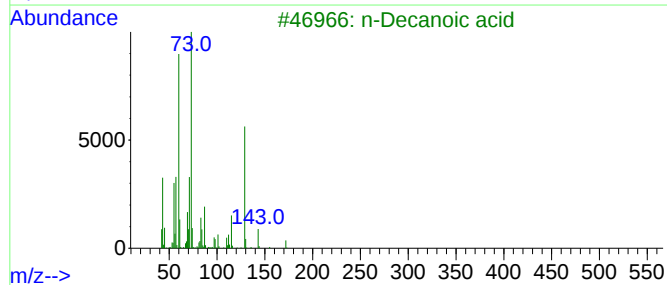
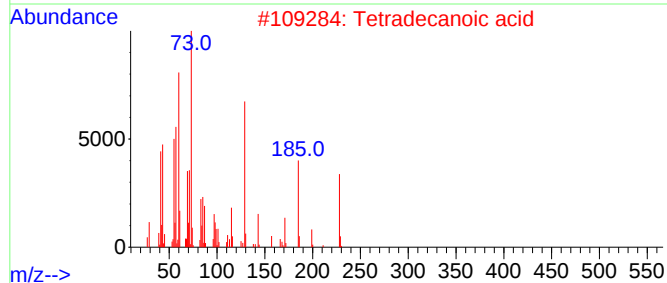
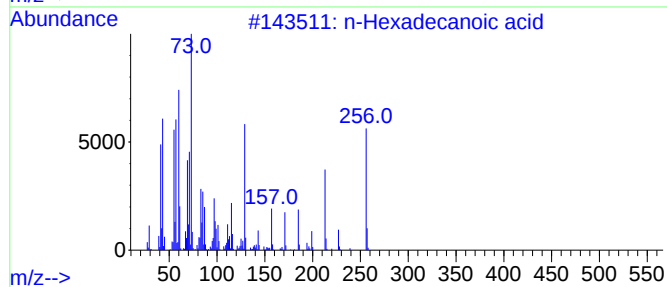
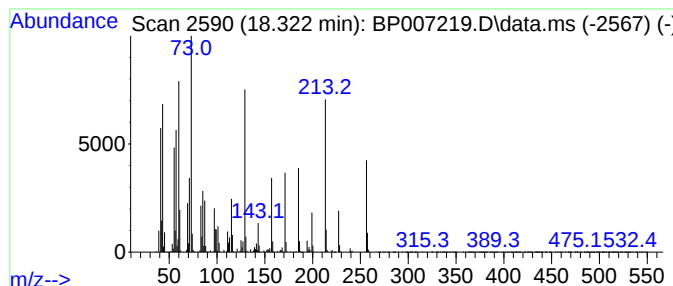
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 9 n-Hexadecanoic acid Concentration Rank 1

| R.T.   | EstConc   | Area      | Relative to ISTD | R.T.   |
|--------|-----------|-----------|------------------|--------|
| 18.322 | 251.27 ng | 108061000 | Phenanthrene-d10 | 17.275 |

| Hit# | of | 5 | Tentative ID        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|---------------------|-----|----------|-------------|------|
| 1    |    |   | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 96   |
| 2    |    |   | Tetradecanoic acid  | 228 | C14H28O2 | 000544-63-8 | 53   |
| 3    |    |   | n-Decanoic acid     | 172 | C10H20O2 | 000334-48-5 | 46   |
| 4    |    |   | Tridecanoic acid    | 214 | C13H26O2 | 000638-53-9 | 44   |
| 5    |    |   | Pentadecanoic acid  | 242 | C15H30O2 | 001002-84-2 | 25   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092721\  
Data File : BP007219.D  
Acq On : 27 Sep 2021 22:57  
Operator : CG/JU  
Sample : M3955-01  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

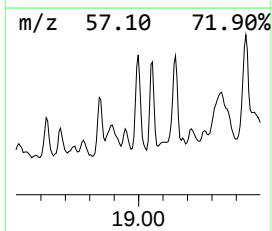
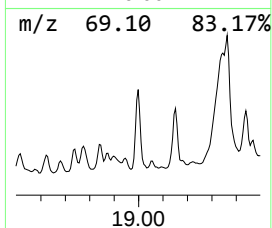
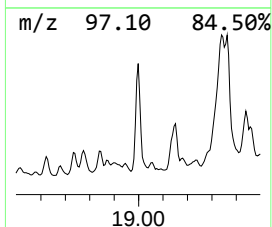
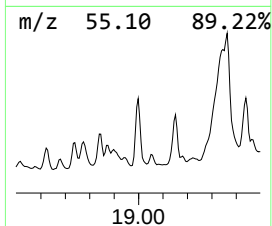
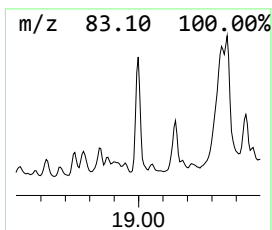
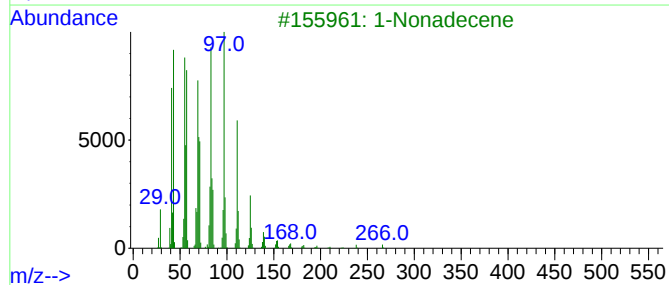
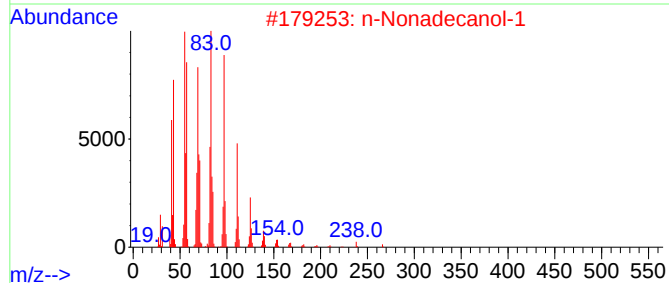
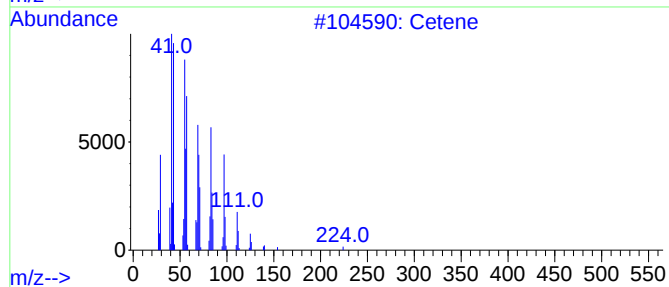
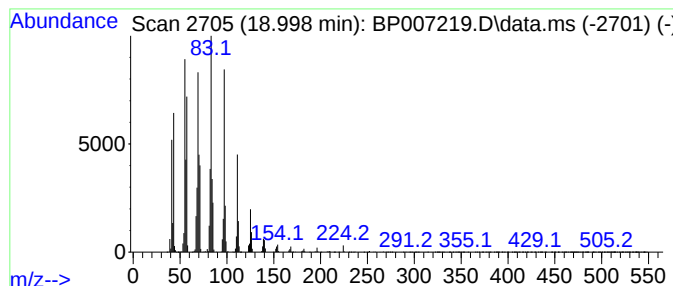
Instrument :  
BNA\_P  
ClientSampleId :  
SS01-NA

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP092121.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 10 Cetene Concentration Rank 19

| R.T.      | EstConc          | Area    | Relative to ISTD |         | R.T.        |      |
|-----------|------------------|---------|------------------|---------|-------------|------|
| 18.998    | 13.29 ng         | 5715750 | Phenanthrene-d10 |         | 17.275      |      |
| Hit# of 5 | Tentative ID     |         | MW               | MolForm | CAS#        | Qual |
| 1         | Cetene           |         | 224              | C16H32  | 000629-73-2 | 98   |
| 2         | n-Nonadecanol-1  |         | 284              | C19H40O | 001454-84-8 | 95   |
| 3         | 1-Nonadecene     |         | 266              | C19H38  | 018435-45-5 | 95   |
| 4         | n-Heptadecanol-1 |         | 256              | C17H36O | 001454-85-9 | 94   |
| 5         | 1-Heneicosanol   |         | 312              | C21H44O | 015594-90-8 | 94   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092721\  
Data File : BP007219.D  
Acq On : 27 Sep 2021 22:57  
Operator : CG/JU  
Sample : M3955-01  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SS01-NA

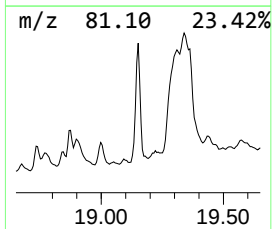
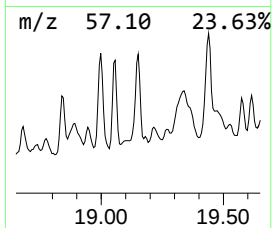
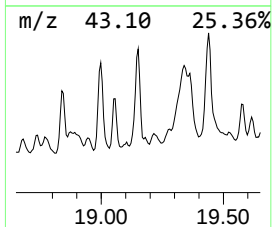
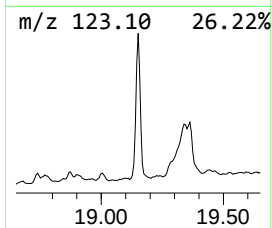
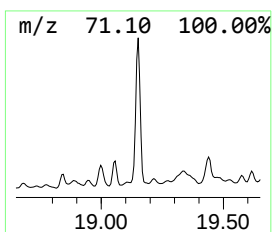
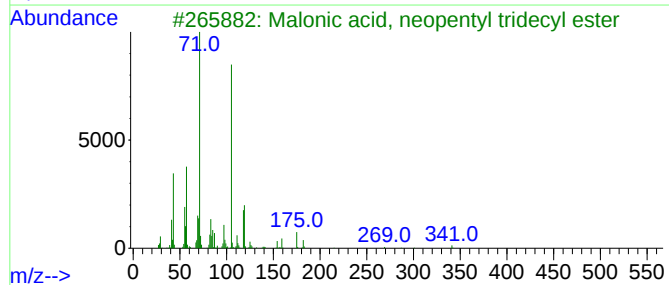
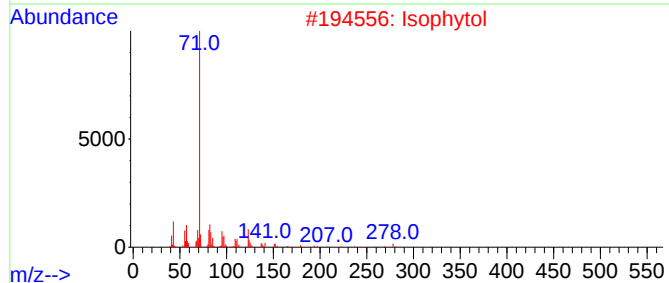
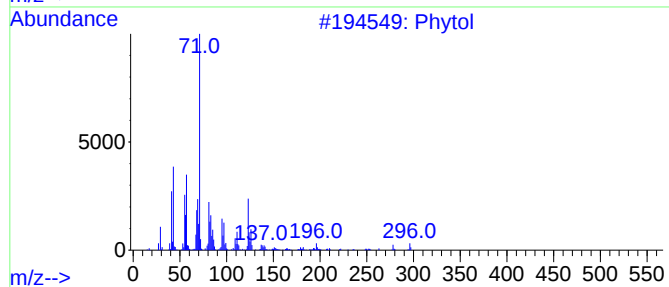
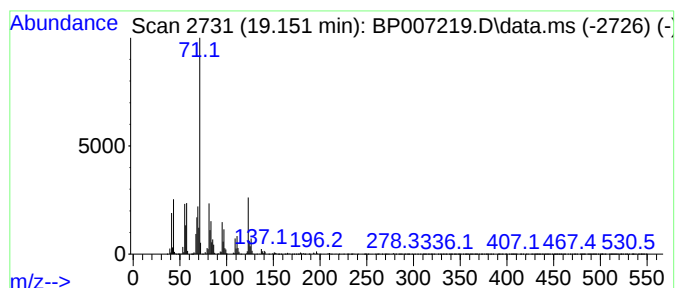
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 11 Phytol Concentration Rank 16

| R.T.      | EstConc                             | Area    | Relative to ISTD |          | R.T.         |      |
|-----------|-------------------------------------|---------|------------------|----------|--------------|------|
| 19.151    | 17.54 ng                            | 7541630 | Phenanthrene-d10 |          | 17.275       |      |
| Hit# of 5 | Tentative ID                        |         | MW               | MolForm  | CAS#         | Qual |
| 1         | Phytol                              |         | 296              | C20H40O  | 000150-86-7  | 98   |
| 2         | Isophytol                           |         | 296              | C20H40O  | 000505-32-8  | 87   |
| 3         | Malonic acid, neopentyl tridecyl... |         | 356              | C21H40O4 | 1000363-93-6 | 53   |
| 4         | Butyric acid, 3-tetradecyl ester    |         | 284              | C18H36O2 | 1000280-57-1 | 53   |
| 5         | Butyric acid, 3-pentadecyl ester    |         | 298              | C19H38O2 | 1000280-57-3 | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092721\  
Data File : BP007219.D  
Acq On : 27 Sep 2021 22:57  
Operator : CG/JU  
Sample : M3955-01  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

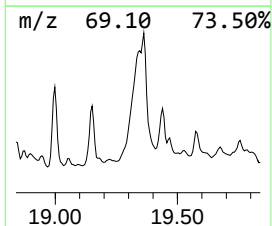
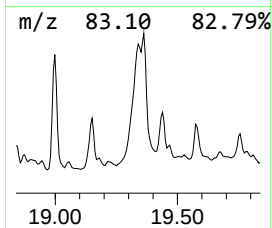
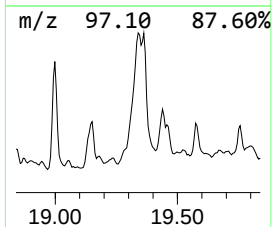
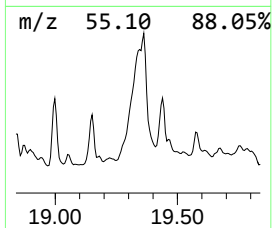
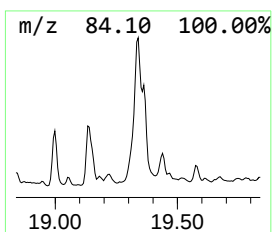
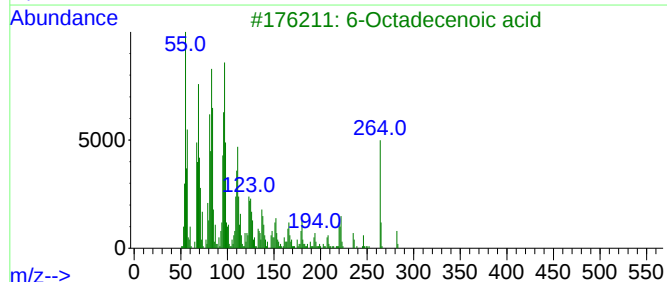
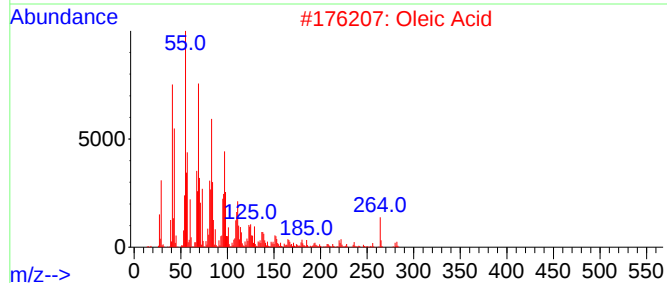
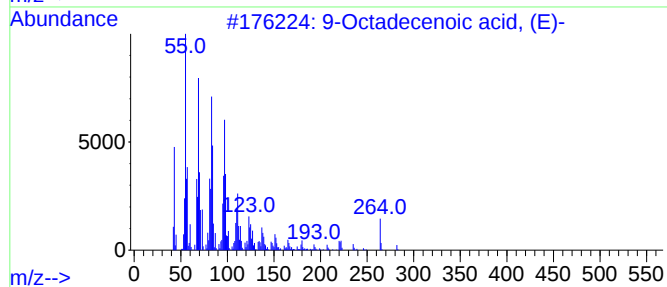
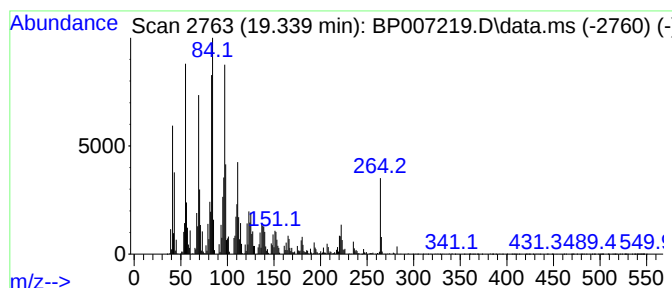
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 12 9-Octadecenoic acid, (E)- Concentration Rank 9

| R.T.      | EstConc                    | Area    | Relative to ISTD |          | R.T.         |      |
|-----------|----------------------------|---------|------------------|----------|--------------|------|
| 19.339    | 29.35 ng                   | 4379940 | Chrysene-d12     |          | 21.368       |      |
| Hit# of 5 | Tentative ID               |         | MW               | MolForm  | CAS#         | Qual |
| 1         | 9-Octadecenoic acid, (E)-  |         | 282              | C18H34O2 | 000112-79-8  | 98   |
| 2         | Oleic Acid                 |         | 282              | C18H34O2 | 000112-80-1  | 93   |
| 3         | 6-Octadecenoic acid        |         | 282              | C18H34O2 | 1000336-66-8 | 86   |
| 4         | cis-Vaccenic acid          |         | 282              | C18H34O2 | 000506-17-2  | 83   |
| 5         | trans-13-Octadecenoic acid |         | 282              | C18H34O2 | 000693-71-0  | 74   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092721\  
Data File : BP007219.D  
Acq On : 27 Sep 2021 22:57  
Operator : CG/JU  
Sample : M3955-01  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

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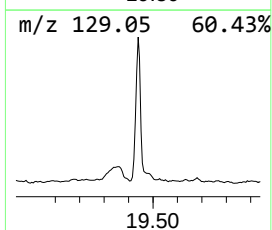
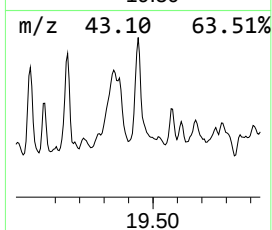
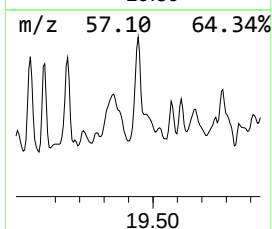
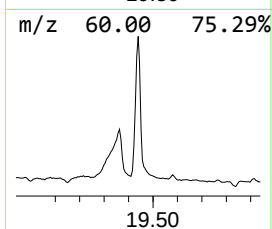
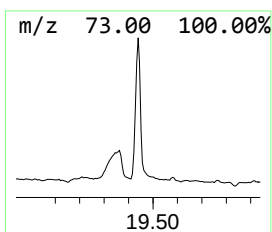
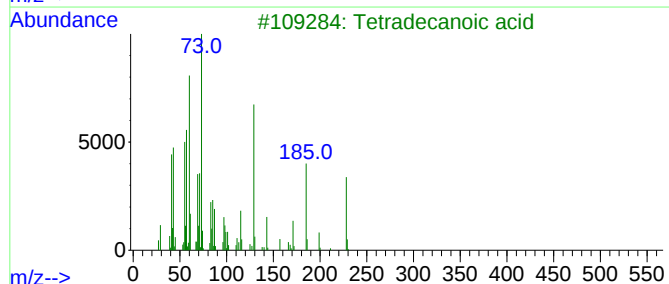
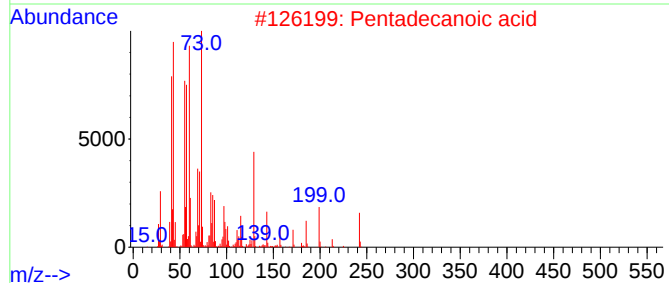
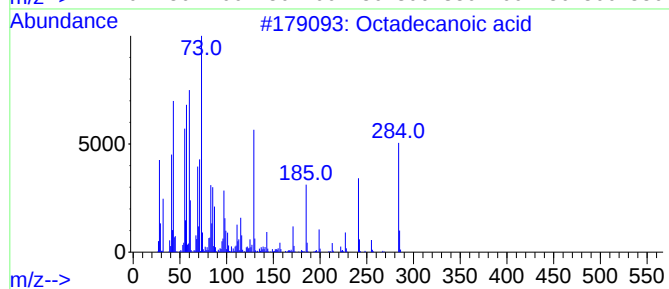
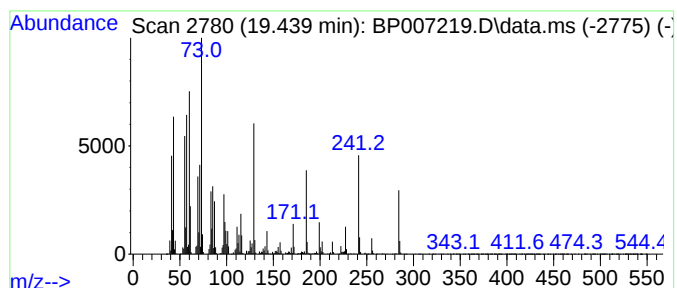
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 13 Octadecanoic acid Concentration Rank 3

| R.T.   | EstConc  | Area    | Relative to ISTD | R.T.   |
|--------|----------|---------|------------------|--------|
| 19.439 | 63.48 ng | 9472870 | Chrysene-d12     | 21.368 |

| Hit# | of | 5 | Tentative ID        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|---------------------|-----|----------|-------------|------|
| 1    |    |   | Octadecanoic acid   | 284 | C18H36O2 | 000057-11-4 | 99   |
| 2    |    |   | Pentadecanoic acid  | 242 | C15H30O2 | 001002-84-2 | 93   |
| 3    |    |   | Tetradecanoic acid  | 228 | C14H28O2 | 000544-63-8 | 81   |
| 4    |    |   | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 58   |
| 5    |    |   | Undecanoic acid     | 186 | C11H22O2 | 000112-37-8 | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092721\  
Data File : BP007219.D  
Acq On : 27 Sep 2021 22:57  
Operator : CG/JU  
Sample : M3955-01  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
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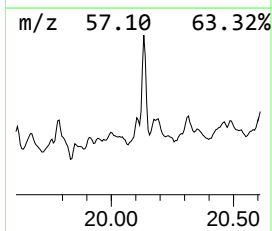
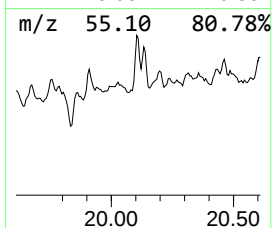
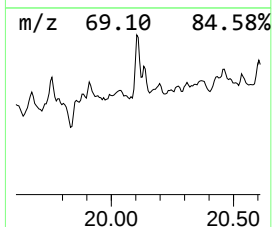
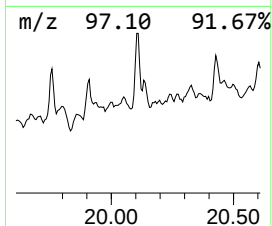
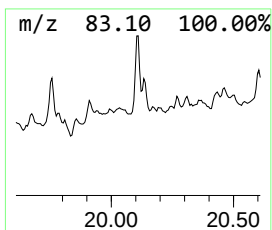
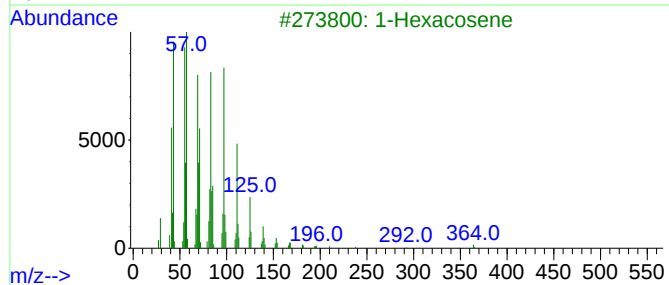
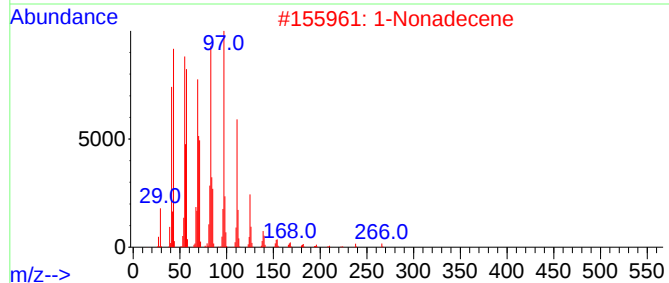
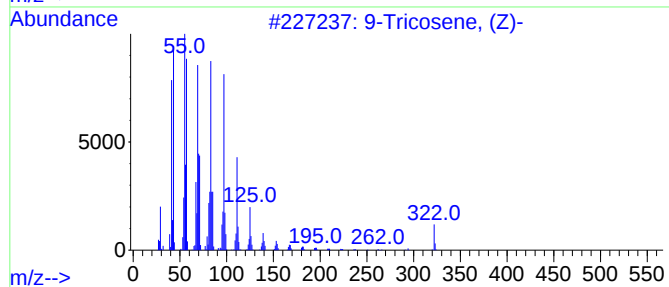
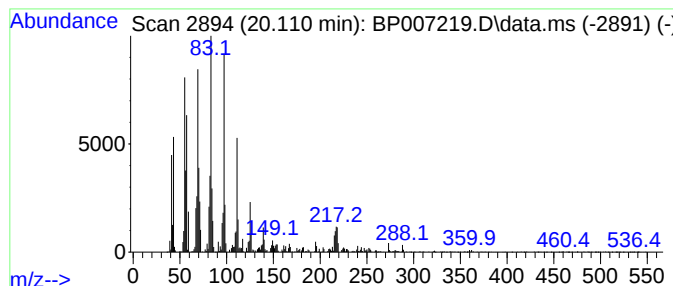
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TIC Integration Parameters: LSCINT.P

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Peak Number 14 9-Tricosene, (Z)- Concentration Rank 20

| R.T.   | EstConc  | Area    | Relative to ISTD | R.T.   |
|--------|----------|---------|------------------|--------|
| 20.110 | 13.05 ng | 1947290 | Chrysene-d12     | 21.368 |

| Hit# | of 5                                | Tentative ID | MW          | MolForm      | CAS# | Qual |
|------|-------------------------------------|--------------|-------------|--------------|------|------|
| 1    | 9-Tricosene, (Z)-                   | 322          | C23H46      | 027519-02-4  | 91   |      |
| 2    | 1-Nonadecene                        | 266          | C19H38      | 018435-45-5  | 91   |      |
| 3    | 1-Hexacosene                        | 364          | C26H52      | 018835-33-1  | 91   |      |
| 4    | Behenic alcohol                     | 326          | C22H46O     | 000661-19-8  | 91   |      |
| 5    | Pentadecafluorooctanoic acid, oc... | 666          | C26H37F15O2 | 1000406-04-8 | 90   |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092721\  
Data File : BP007219.D  
Acq On : 27 Sep 2021 22:57  
Operator : CG/JU  
Sample : M3955-01  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

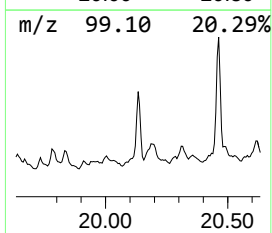
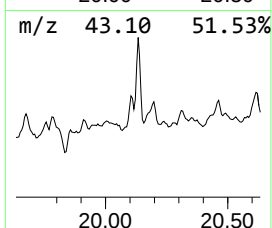
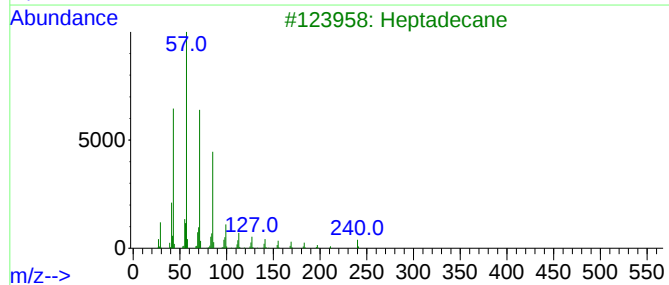
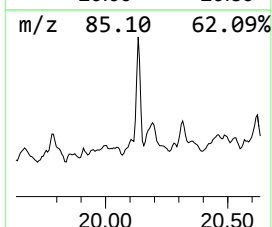
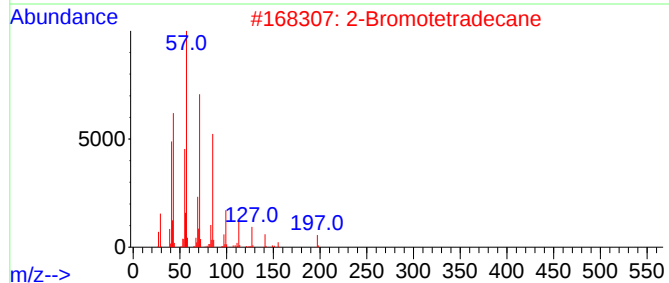
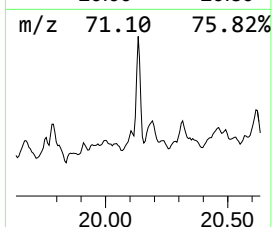
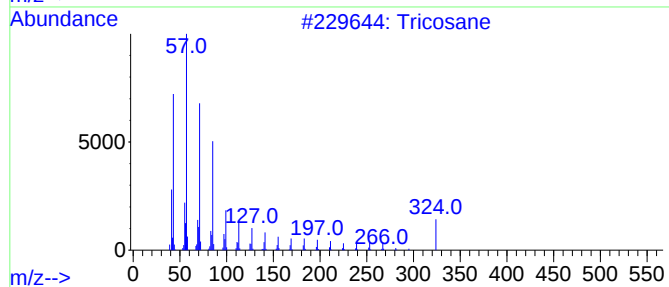
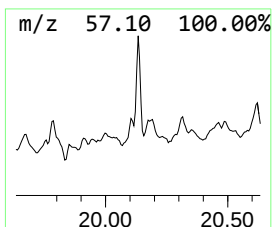
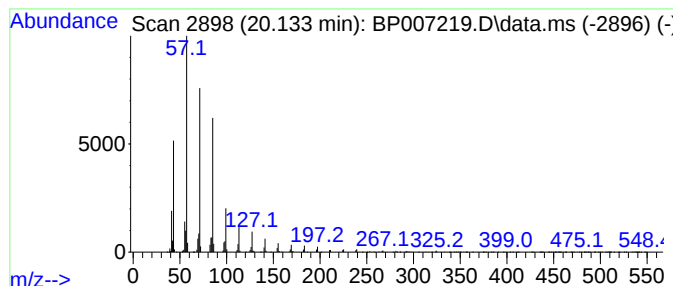
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 15 Tricosane Concentration Rank 17

| R.T.      | EstConc            | Area    | Relative to ISTD |          | R.T.        |      |
|-----------|--------------------|---------|------------------|----------|-------------|------|
| 20.133    | 17.07 ng           | 2547310 | Chrysene-d12     |          | 21.368      |      |
| Hit# of 5 | Tentative ID       |         | MW               | MolForm  | CAS#        | Qual |
| 1         | Tricosane          |         | 324              | C23H48   | 000638-67-5 | 94   |
| 2         | 2-Bromotetradecane |         | 276              | C14H29Br | 074036-95-6 | 90   |
| 3         | Heptadecane        |         | 240              | C17H36   | 000629-78-7 | 86   |
| 4         | Octadecane         |         | 254              | C18H38   | 000593-45-3 | 72   |
| 5         | Docosane           |         | 310              | C22H46   | 000629-97-0 | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092721\  
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Acq On : 27 Sep 2021 22:57  
Operator : CG/JU  
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ALS Vial : 19 Sample Multiplier: 1

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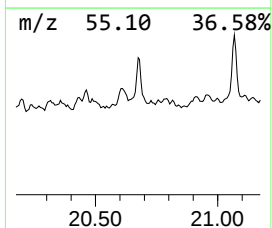
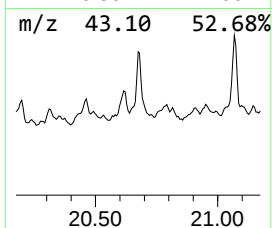
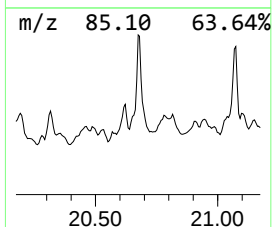
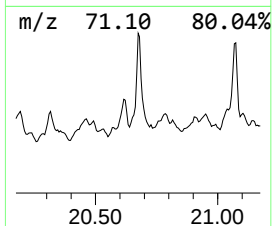
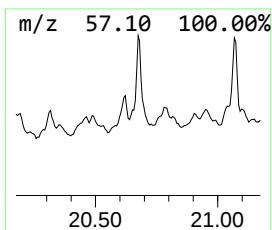
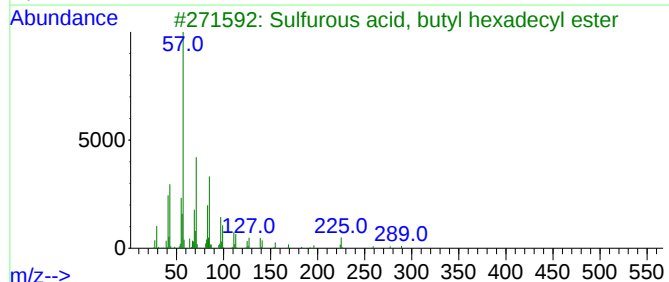
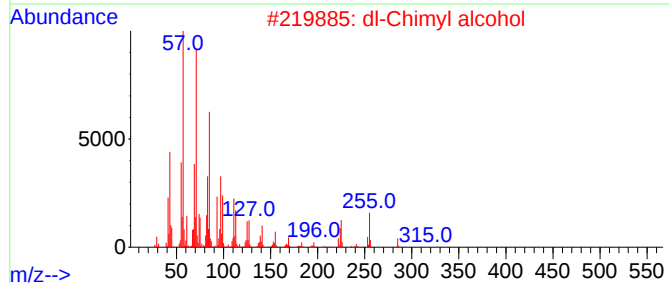
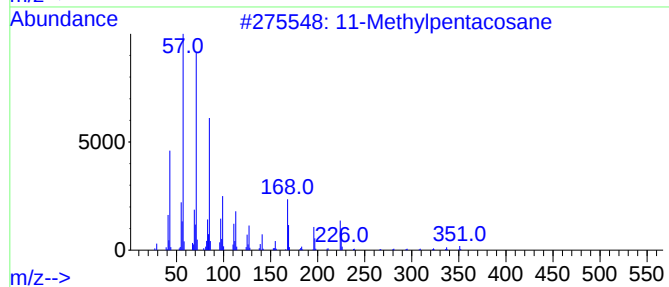
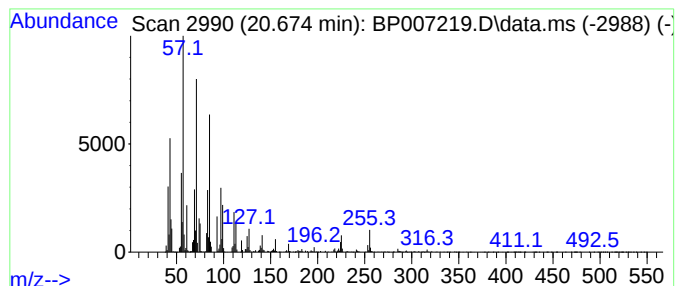
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 16 11-Methylpentacosane Concentration Rank 10

| R.T.   | EstConc  | Area    | Relative to ISTD | R.T.   |
|--------|----------|---------|------------------|--------|
| 20.674 | 21.76 ng | 3246800 | Chrysene-d12     | 21.368 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | 11-Methylpentacosane                | 366 | C26H54    | 015689-71-1  | 90   |
| 2    |    |   | dl-Chimyl alcohol                   | 316 | C19H40O3  | 006145-69-3  | 87   |
| 3    |    |   | Sulfurous acid, butyl hexadecyl ... | 362 | C20H42O3S | 1000309-18-3 | 83   |
| 4    |    |   | Heptadecane, 2-methyl-              | 254 | C18H38    | 001560-89-0  | 70   |
| 5    |    |   | Heneicosane                         | 296 | C21H44    | 000629-94-7  | 68   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092721\  
Data File : BP007219.D  
Acq On : 27 Sep 2021 22:57  
Operator : CG/JU  
Sample : M3955-01  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SS01-NA

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP092121.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

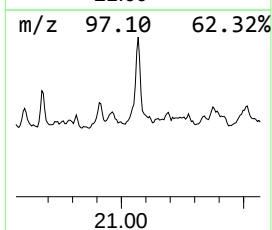
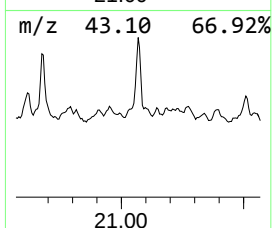
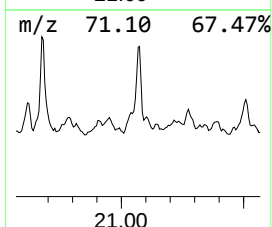
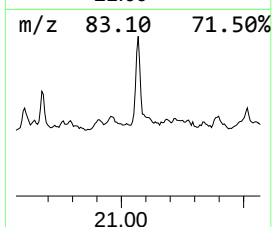
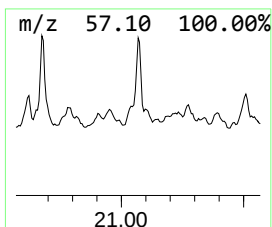
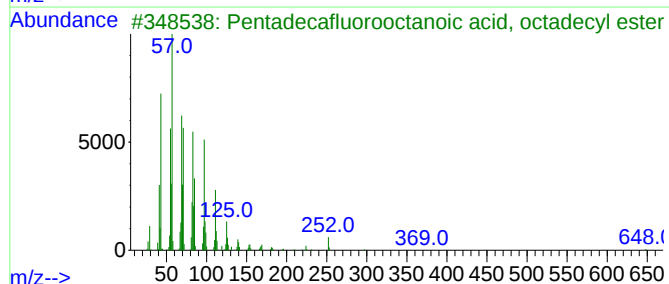
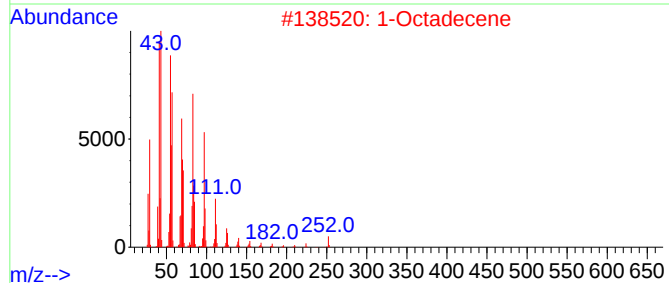
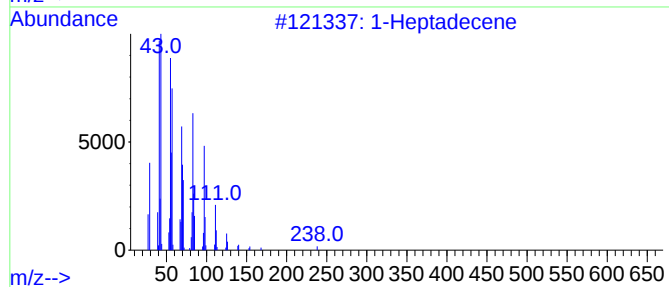
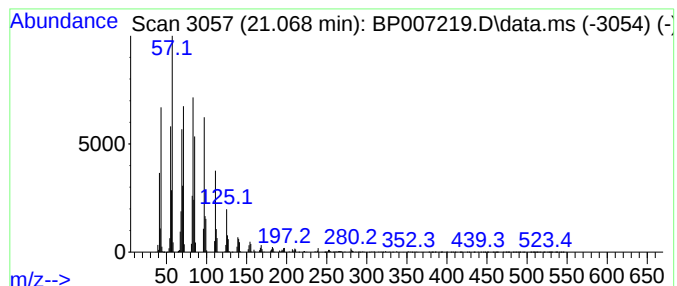
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Peak Number 17 1-Heptadecene Concentration Rank 13

| R.T.   | EstConc  | Area    | Relative to ISTD | R.T.   |
|--------|----------|---------|------------------|--------|
| 21.068 | 20.00 ng | 2984740 | Chrysene-d12     | 21.368 |

| Hit# | of 5                                | Tentative ID | MW  | MolForm     | CAS#         | Qual |
|------|-------------------------------------|--------------|-----|-------------|--------------|------|
| 1    | 1-Heptadecene                       |              | 238 | C17H34      | 006765-39-5  | 94   |
| 2    | 1-Octadecene                        |              | 252 | C18H36      | 000112-88-9  | 93   |
| 3    | Pentadecafluorooctanoic acid, oc... |              | 666 | C26H37F15O2 | 1000406-04-8 | 91   |
| 4    | Octacosyl trifluoroacetate          |              | 506 | C30H57F3O2  | 1000351-74-9 | 91   |
| 5    | 1-Decanol, 2-hexyl-                 |              | 242 | C16H34O     | 002425-77-6  | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092721\  
Data File : BP007219.D  
Acq On : 27 Sep 2021 22:57  
Operator : CG/JU  
Sample : M3955-01  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SS01-NA

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP092121.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

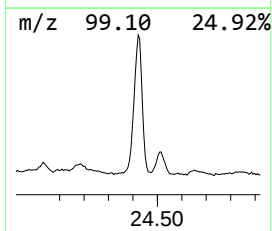
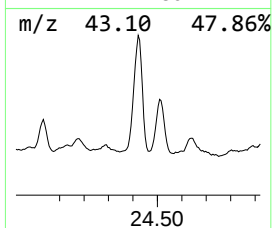
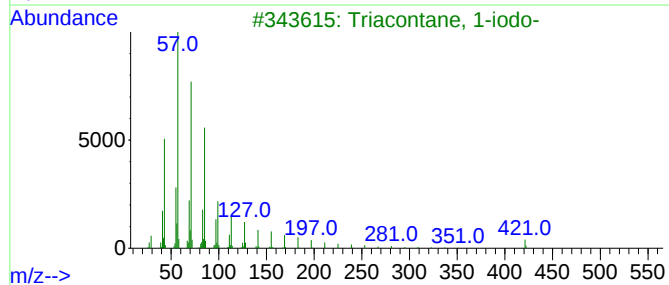
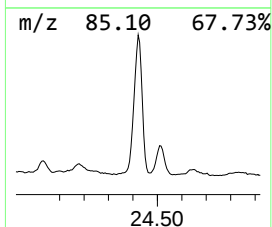
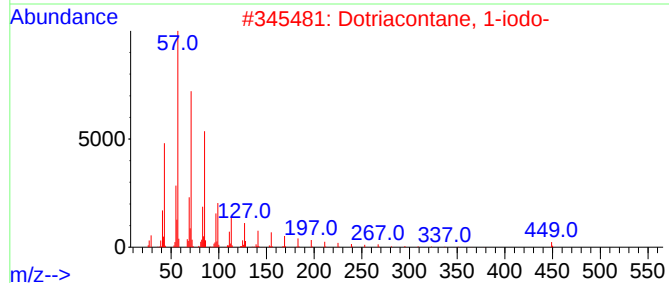
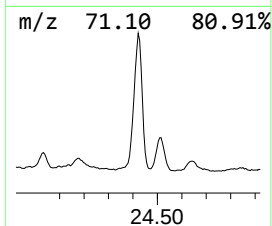
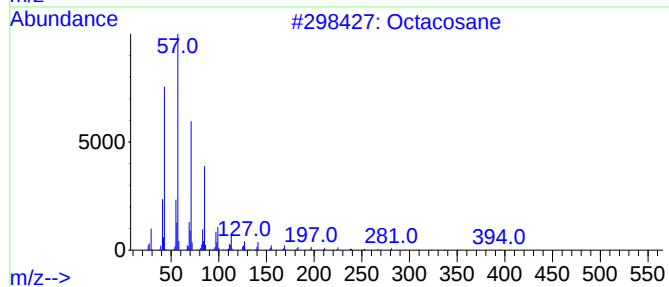
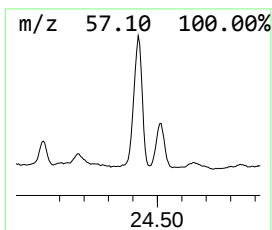
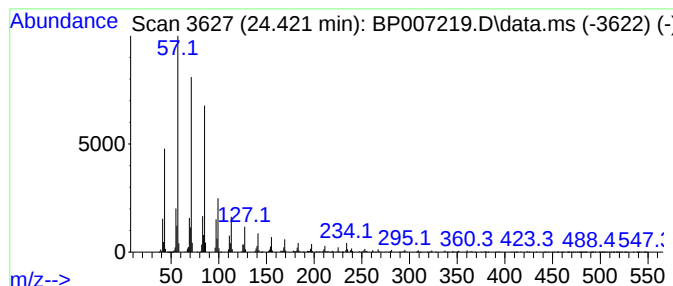
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Peak Number 18 Octacosane Concentration Rank 14

| R.T.   | EstConc  | Area    | Relative to ISTD | R.T.   |
|--------|----------|---------|------------------|--------|
| 24.421 | 20.00 ng | 8069820 | Perylene-d12     | 23.792 |

| Hit# | of 5 | Tentative ID           | MW  | MolForm | CAS#         | Qual |
|------|------|------------------------|-----|---------|--------------|------|
| 1    |      | Octacosane             | 394 | C28H58  | 000630-02-4  | 96   |
| 2    |      | Dotriacontane, 1-iodo- | 576 | C32H65I | 1000406-32-4 | 95   |
| 3    |      | Triacontane, 1-iodo-   | 548 | C30H61I | 1000406-32-3 | 94   |
| 4    |      | Octadecane             | 254 | C18H38  | 000593-45-3  | 94   |
| 5    |      | Octacosane, 1-iodo-    | 520 | C28H57I | 1000406-32-2 | 94   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092721\  
Data File : BP007219.D  
Acq On : 27 Sep 2021 22:57  
Operator : CG/JU  
Sample : M3955-01  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SS01-NA

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP092121.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

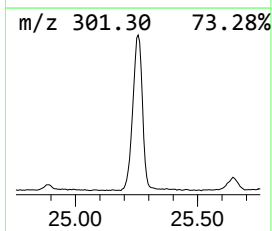
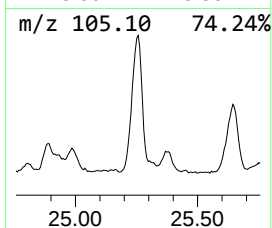
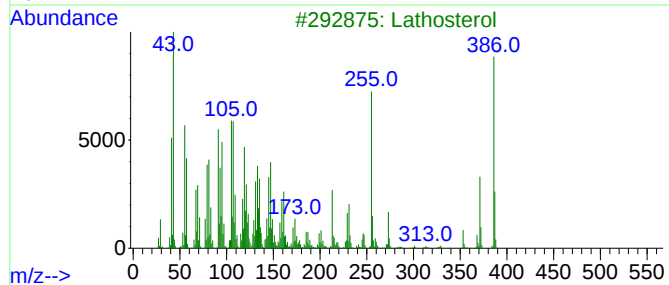
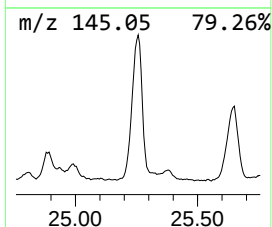
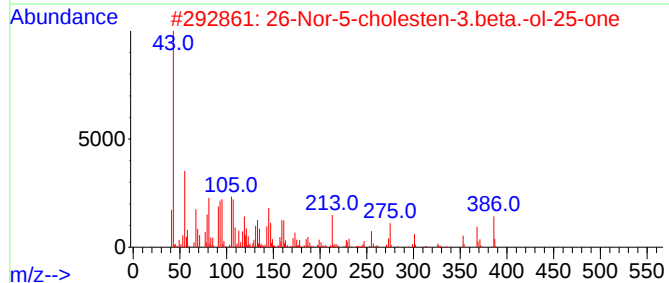
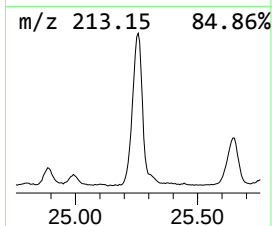
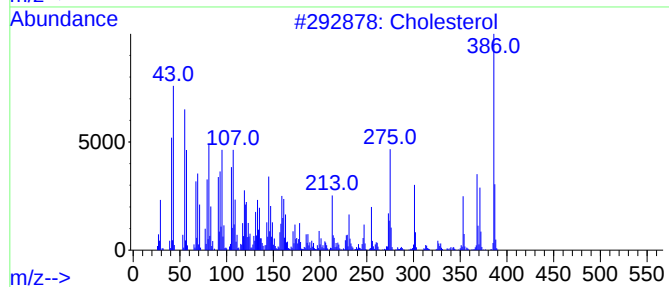
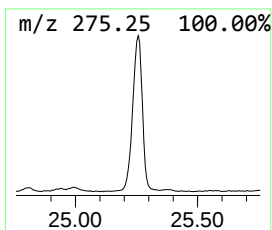
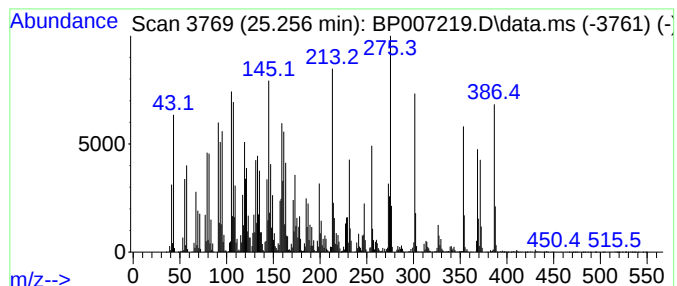
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 Cholesterol Concentration Rank 5

| R.T.   | EstConc  | Area     | Relative to ISTD | R.T.   |
|--------|----------|----------|------------------|--------|
| 25.256 | 50.03 ng | 20188500 | Perylene-d12     | 23.792 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Cholesterol                         | 386 | C27H46O   | 000057-88-5  | 99   |
| 2    |    |   | 26-Nor-5-cholesten-3.beta.-ol-25... | 386 | C26H42O2  | 007494-34-0  | 99   |
| 3    |    |   | Lathosterol                         | 386 | C27H46O   | 000080-99-9  | 56   |
| 4    |    |   | benzeneacetonitrile, .alpha.-[(4... | 275 | C16H9N3O2 | 1000401-81-2 | 46   |
| 5    |    |   | Cholest-8-en-3-ol, (3.beta.)-       | 386 | C27H46O   | 007199-91-9  | 46   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092721\  
Data File : BP007219.D  
Acq On : 27 Sep 2021 22:57  
Operator : CG/JU  
Sample : M3955-01  
Misc :  
ALS Vial : 19 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
SS01-NA

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP092121.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

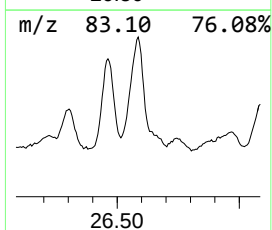
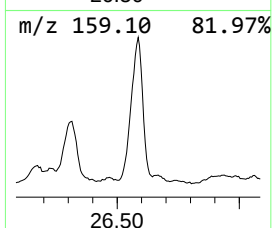
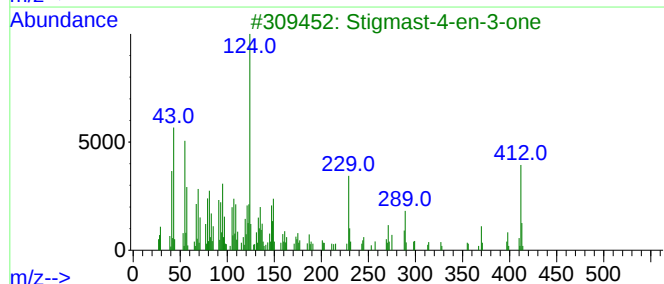
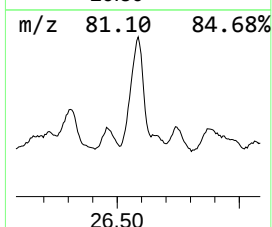
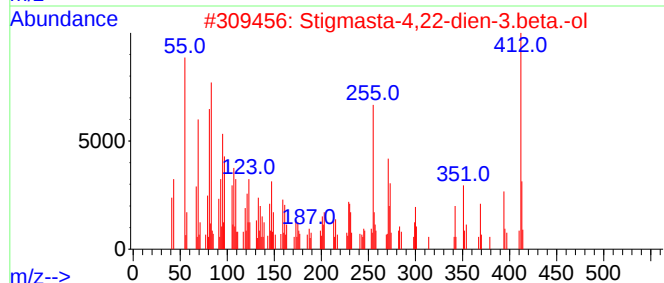
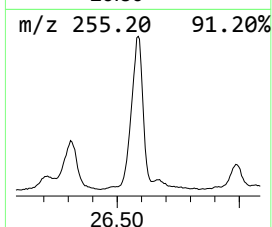
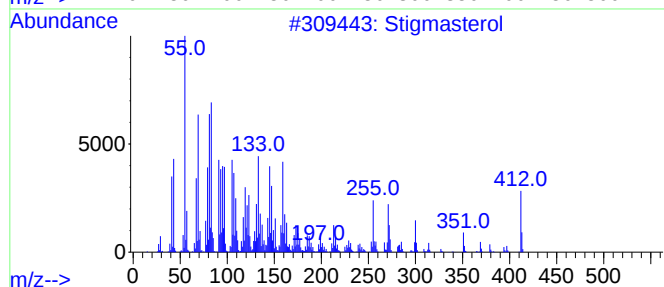
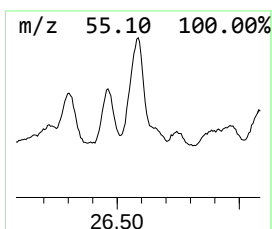
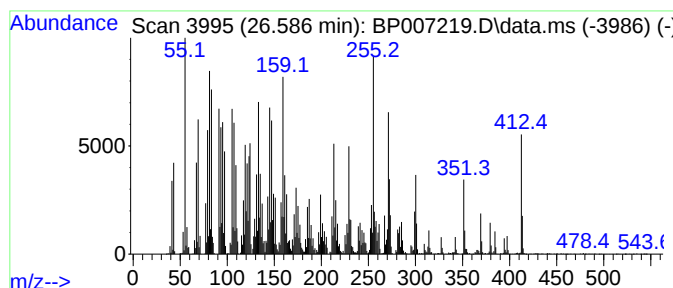
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 Stigmasterol Concentration Rank 6

| R.T.   | EstConc  | Area     | Relative to ISTD | R.T.   |
|--------|----------|----------|------------------|--------|
| 26.586 | 44.03 ng | 17765500 | Perylene-d12     | 23.792 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | Stigmasterol                        | 412 | C29H48O | 000083-48-7  | 99   |
| 2    |    |   | Stigmasta-4,22-dien-3.beta.-ol      | 412 | C29H48O | 057815-94-8  | 92   |
| 3    |    |   | Stigmast-4-en-3-one                 | 412 | C29H48O | 001058-61-3  | 46   |
| 4    |    |   | (4S,14R)-4,10,13,14-tetramethyl-... | 412 | C29H48O | 1000456-24-3 | 44   |
| 5    |    |   | Chondrillasterol                    | 412 | C29H48O | 000481-17-4  | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP092721\  
 Data File : BP007219.D  
 Acq On : 27 Sep 2021 22:57  
 Operator : CG/JU  
 Sample : M3955-01  
 Misc :  
 ALS Vial : 19 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 SS01-NA

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP092121.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST0.L  
 TIC Integration Parameters: LSCINT.P

| TIC Top Hit name   | RT     | EstConc | Units | Response  | --Internal Standard-- |        |         |      |
|--------------------|--------|---------|-------|-----------|-----------------------|--------|---------|------|
|                    |        |         |       |           | #                     | RT     | Resp    | Conc |
| Ethyl cycloprop... | 12.363 | 21.4    | ng    | 2015170   | 2                     | 10.687 | 1886540 | 20.0 |
| Tetradecanoic acid | 16.739 | 55.1    | ng    | 23687300  | 4                     | 17.275 | 8601020 | 20.0 |
| 1-Tetradecanol     | 16.892 | 14.5    | ng    | 6247240   | 4                     | 17.275 | 8601020 | 20.0 |
| Pentadecanoic acid | 17.210 | 19.8    | ng    | 8513520   | 4                     | 17.275 | 8601020 | 20.0 |
| Undecanoic acid    | 17.498 | 40.7    | ng    | 17490800  | 4                     | 17.275 | 8601020 | 20.0 |
| Neophytadiene      | 17.645 | 20.4    | ng    | 8788100   | 4                     | 17.275 | 8601020 | 20.0 |
| Chloroacetic ac... | 17.669 | 30.4    | ng    | 13054500  | 4                     | 17.275 | 8601020 | 20.0 |
| Palmitoleic acid   | 18.110 | 121.3   | ng    | 52152900  | 4                     | 17.275 | 8601020 | 20.0 |
| n-Hexadecanoic ... | 18.322 | 251.3   | ng    | 108061000 | 4                     | 17.275 | 8601020 | 20.0 |
| Cetene             | 18.998 | 13.3    | ng    | 5715750   | 4                     | 17.275 | 8601020 | 20.0 |
| Phytol             | 19.151 | 17.5    | ng    | 7541630   | 4                     | 17.275 | 8601020 | 20.0 |
| 9-Octadecenoic ... | 19.339 | 29.4    | ng    | 4379940   | 5                     | 21.368 | 2984740 | 20.0 |
| Octadecanoic acid  | 19.439 | 63.5    | ng    | 9472870   | 5                     | 21.368 | 2984740 | 20.0 |
| 9-Tricosene, (Z)-  | 20.110 | 13.1    | ng    | 1947290   | 5                     | 21.368 | 2984740 | 20.0 |
| Tricosane          | 20.133 | 17.1    | ng    | 2547310   | 5                     | 21.368 | 2984740 | 20.0 |
| 11-Methylpentac... | 20.674 | 21.8    | ng    | 3246800   | 5                     | 21.368 | 2984740 | 20.0 |
| 1-Heptadecene      | 21.068 | 20.0    | ng    | 2984740   | 5                     | 21.368 | 2984740 | 20.0 |
| Octacosane         | 24.421 | 20.0    | ng    | 8069820   | 6                     | 23.792 | 8069820 | 20.0 |
| Cholesterol        | 25.256 | 50.0    | ng    | 20188500  | 6                     | 23.792 | 8069820 | 20.0 |
| Stigmasterol       | 26.586 | 44.0    | ng    | 17765500  | 6                     | 23.792 | 8069820 | 20.0 |