

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG072722\  
 Data File : BG054423.D  
 Acq On : 27 Jul 2022 19:53  
 Operator : CG/JU  
 Sample : N3906-01  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 NB-7-7-26-22

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\_G\Methods\8270-BG071522.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG054423.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         | 3.118       | 16            | 22          | 36           | rVB      | 72810          | 149278        | 2.56%           | 0.507%        |
| 2         | 5.574       | 433           | 440         | 452          | rBV      | 226343         | 409769        | 7.04%           | 1.391%        |
| 3         | 7.178       | 705           | 713         | 726          | rBV      | 237509         | 440949        | 7.57%           | 1.497%        |
| 4         | 7.989       | 844           | 851         | 860          | rBV      | 63921          | 112917        | 1.94%           | 0.383%        |
| 5         | 9.158       | 1042          | 1050        | 1062         | rBB      | 152535         | 286659        | 4.92%           | 0.973%        |
| 6         | 10.803      | 1322          | 1330        | 1336         | rBV      | 83823          | 154946        | 2.66%           | 0.526%        |
| 7         | 13.247      | 1738          | 1746        | 1755         | rVB      | 499627         | 781636        | 13.42%          | 2.653%        |
| 8         | 14.628      | 1971          | 1981        | 1987         | rBV      | 138812         | 228942        | 3.93%           | 0.777%        |
| 9         | 16.120      | 2227          | 2235        | 2243         | rBV2     | 445096         | 703817        | 12.09%          | 2.389%        |
| 10        | 17.372      | 2442          | 2448        | 2451         | rBV      | 173835         | 280757        | 4.82%           | 0.953%        |
| 11        | 17.419      | 2451          | 2456        | 2462         | rVV      | 367738         | 586220        | 10.07%          | 1.990%        |
| 12        | 17.507      | 2465          | 2471        | 2476         | rVV      | 97130          | 150596        | 2.59%           | 0.511%        |
| 13        | 18.030      | 2554          | 2560        | 2568         | rBV      | 886802         | 1158800       | 19.90%          | 3.933%        |
| 14        | 18.253      | 2593          | 2598        | 2602         | rBV2     | 103212         | 204751        | 3.52%           | 0.695%        |
| 15        | 18.300      | 2602          | 2606        | 2612         | rVB      | 114906         | 182191        | 3.13%           | 0.618%        |
| 16        | 18.376      | 2615          | 2619        | 2624         | rBV      | 37395          | 58741         | 1.01%           | 0.199%        |
| 17        | 18.441      | 2625          | 2630        | 2635         | rBV2     | 123409         | 259336        | 4.45%           | 0.880%        |
| 18        | 18.741      | 2678          | 2681        | 2686         | rBV      | 67635          | 96710         | 1.66%           | 0.328%        |
| 19        | 18.794      | 2686          | 2690        | 2695         | rVB2     | 75607          | 112221        | 1.93%           | 0.381%        |
| 20        | 19.181      | 2748          | 2756        | 2758         | rBV      | 3931482        | 5822361       | 100.00%         | 19.762%       |
| 21        | 19.211      | 2758          | 2761        | 2772         | rVV2     | 2579785        | 3389014       | 58.21%          | 11.503%       |
| 22        | 19.334      | 2774          | 2782        | 2787         | rVB2     | 515384         | 803622        | 13.80%          | 2.728%        |
| 23        | 19.393      | 2788          | 2792        | 2793         | rBV2     | 57194          | 74105         | 1.27%           | 0.252%        |
| 24        | 19.434      | 2793          | 2799        | 2805         | rVB      | 1384745        | 2162603       | 37.14%          | 7.340%        |
| 25        | 19.763      | 2850          | 2855        | 2856         | rBV      | 249319         | 330418        | 5.67%           | 1.121%        |
| 26        | 19.798      | 2856          | 2861        | 2867         | rVB      | 1233626        | 1901558       | 32.66%          | 6.454%        |
| 27        | 19.863      | 2869          | 2872        | 2878         | rVB2     | 77859          | 124435        | 2.14%           | 0.422%        |
| 28        | 19.980      | 2887          | 2892        | 2897         | rBV      | 863117         | 1236629       | 21.24%          | 4.197%        |
| 29        | 20.127      | 2910          | 2917        | 2922         | rBV3     | 122481         | 290427        | 4.99%           | 0.986%        |
| 30        | 20.445      | 2964          | 2971        | 2975         | rBV2     | 155496         | 344668        | 5.92%           | 1.170%        |
| 31        | 20.580      | 2991          | 2994        | 2998         | rBV      | 115609         | 165625        | 2.84%           | 0.562%        |
| 32        | 21.044      | 3069          | 3073        | 3079         | rVB      | 264518         | 401884        | 6.90%           | 1.364%        |
| 33        | 21.267      | 3108          | 3111        | 3115         | rVB      | 119121         | 163665        | 2.81%           | 0.556%        |
| 34        | 21.320      | 3116          | 3120        | 3124         | rVV2     | 128299         | 213610        | 3.67%           | 0.725%        |
| 35        | 21.379      | 3126          | 3130        | 3136         | rVB      | 165070         | 283384        | 4.87%           | 0.962%        |
| 36        | 21.631      | 3166          | 3173        | 3179         | rBV2     | 876667         | 2001905       | 34.38%          | 6.795%        |

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Operator : CG/JU  
Sample : N3906-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
NB-7-7-26-22

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\_G\Methods\8270-BG071522.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |        |         |        |        |
|----|--------|------|------|------|------|--------|---------|--------|--------|
| 37 | 21.696 | 3179 | 3184 | 3196 | rVB  | 682034 | 1357682 | 23.32% | 4.608% |
| 38 | 23.852 | 3543 | 3551 | 3557 | rBV3 | 381214 | 1262053 | 21.68% | 4.284% |
| 39 | 24.716 | 3692 | 3698 | 3712 | rVB  | 267972 | 773275  | 13.28% | 2.625% |

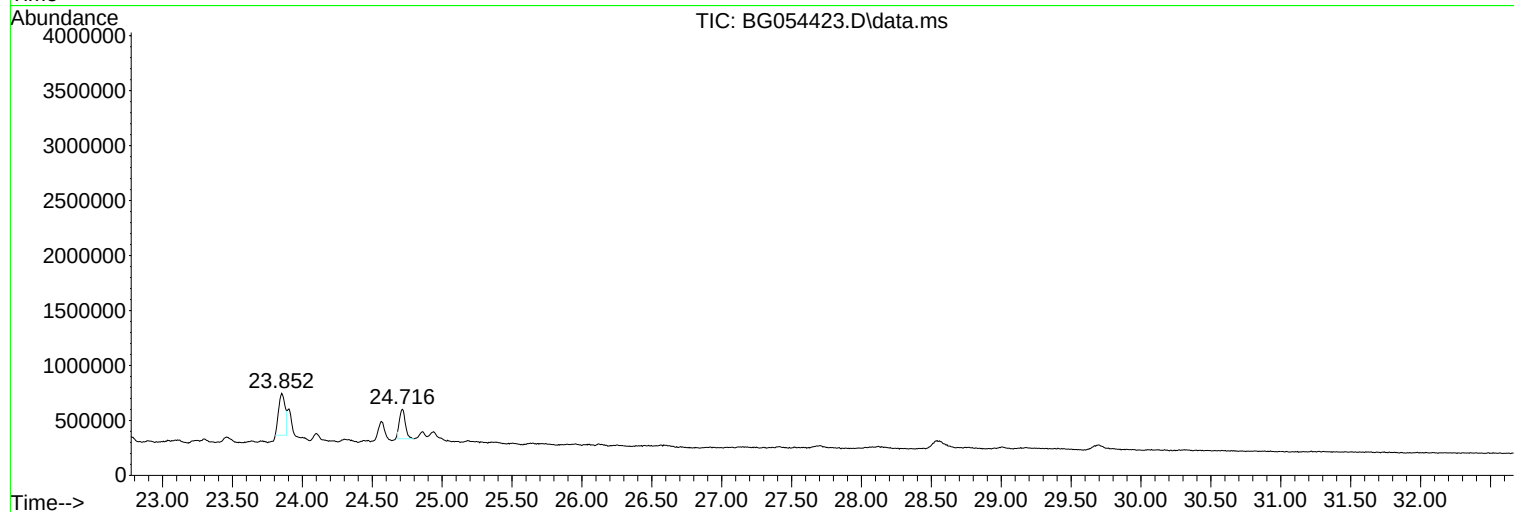
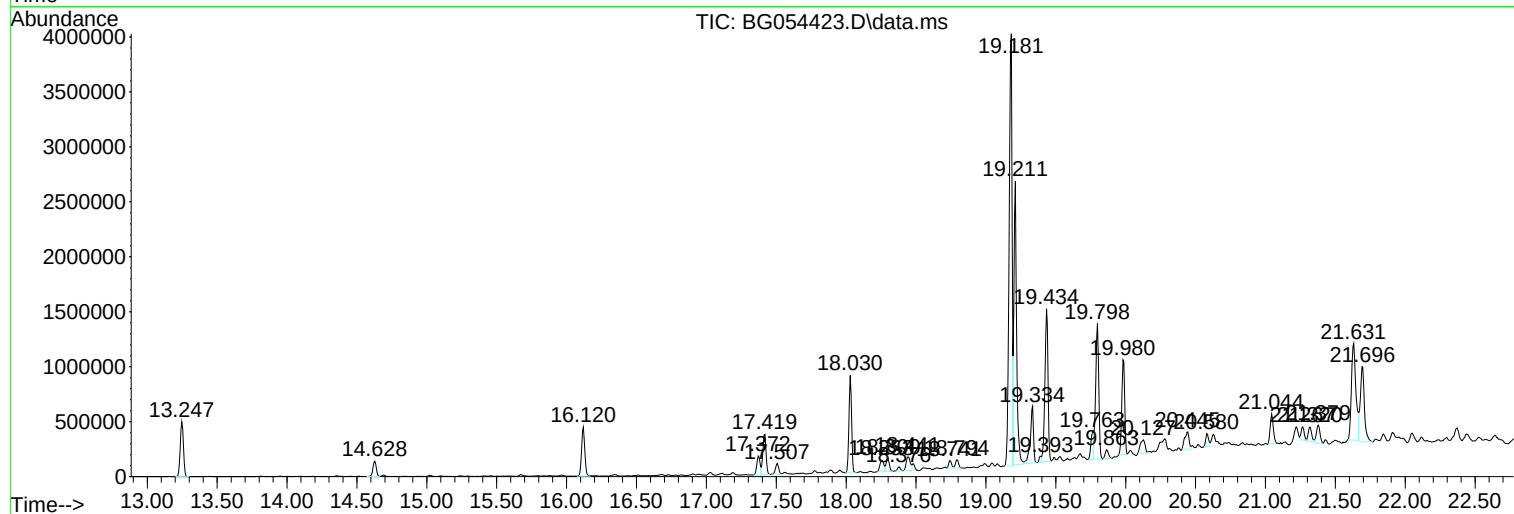
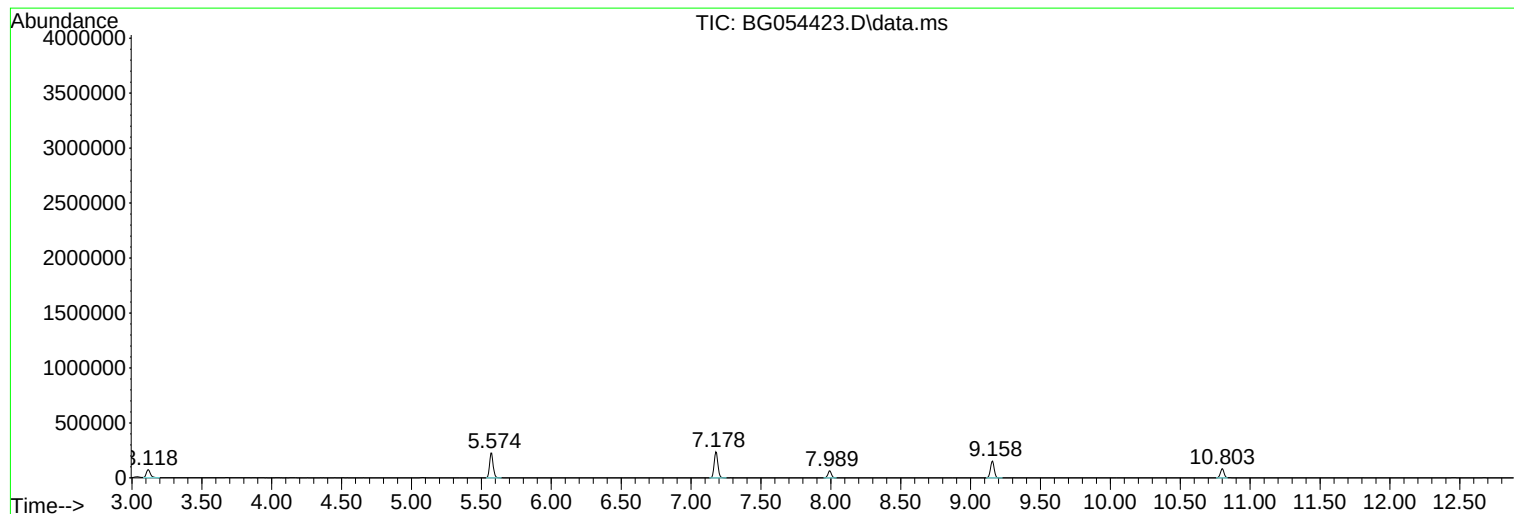
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Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG072722\  
Data File : BG054423.D  
Acq On : 27 Jul 2022 19:53  
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Sample : N3906-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
NB-7-7-26-22

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG071522.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\_G\Data\BG072722\  
Data File : BG054423.D  
Acq On : 27 Jul 2022 19:53  
Operator : CG/JU  
Sample : N3906-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
NB-7-7-26-22

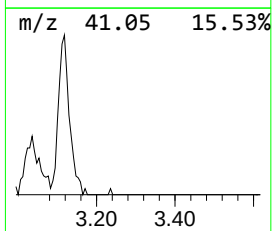
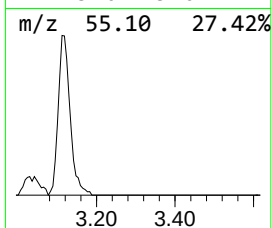
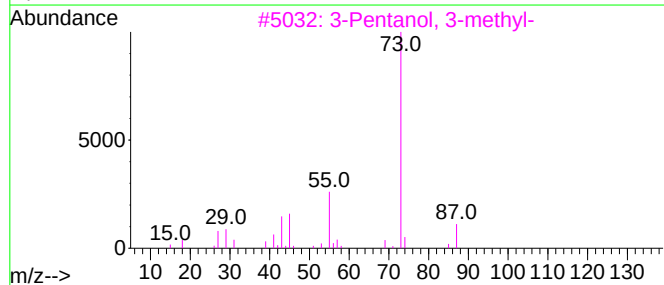
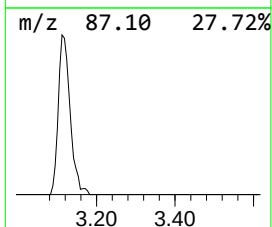
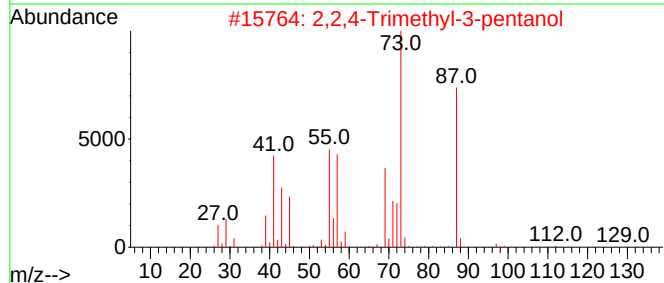
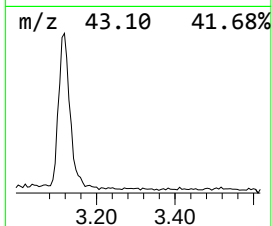
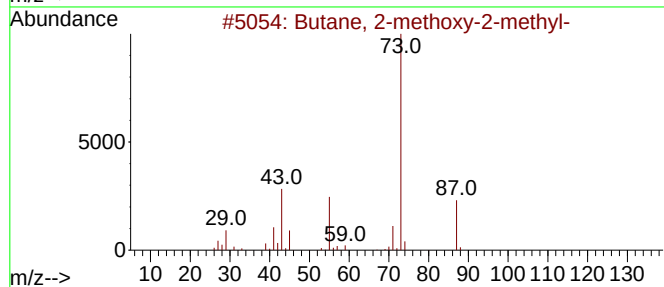
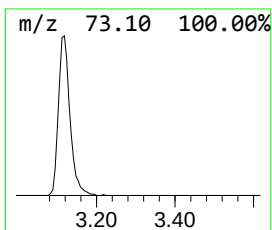
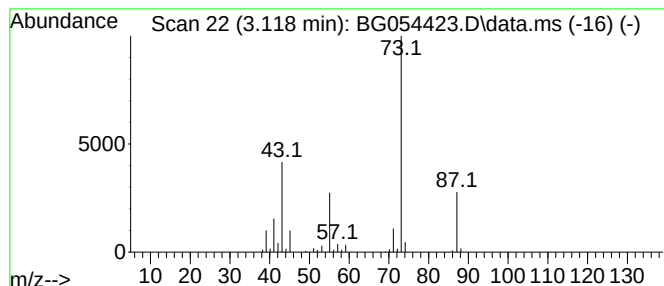
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG071522.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

| R.T.  | EstConc  | Area   | Relative to ISTD       | R.T.  |
|-------|----------|--------|------------------------|-------|
| 3.118 | 26.44 ng | 149278 | 1,4-Dichlorobenzene-d4 | 7.989 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Butane, 2-methoxy-2-methyl- | 102 | C6H14O  | 000994-05-8 | 83   |
| 2    |    |   | 2,2,4-Trimethyl-3-pentanol  | 130 | C8H18O  | 005162-48-1 | 40   |
| 3    |    |   | 3-Pentanol, 3-methyl-       | 102 | C6H14O  | 000077-74-7 | 25   |
| 4    |    |   | N-Ethylformamide            | 73  | C3H7NO  | 000627-45-2 | 9    |
| 5    |    |   | Silane, tetramethyl-        | 88  | C4H12Si | 000075-76-3 | 9    |



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Sample : N3906-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
NB-7-7-26-22

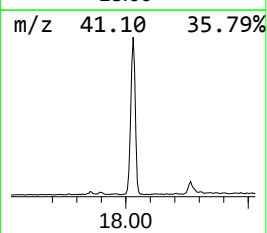
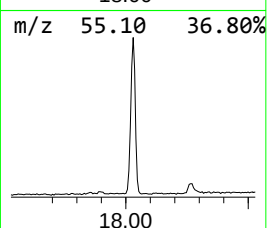
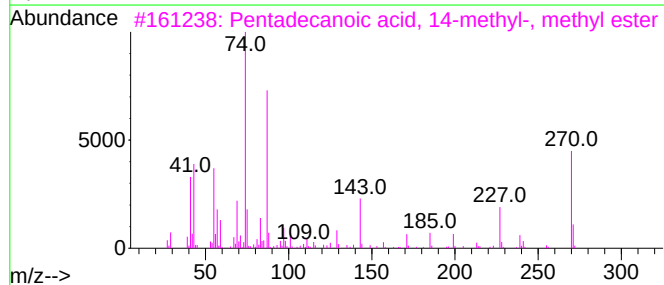
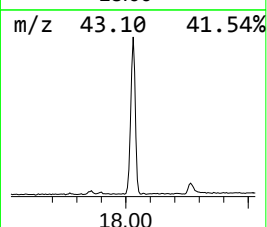
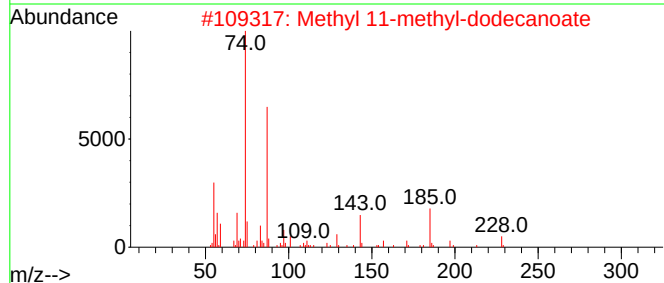
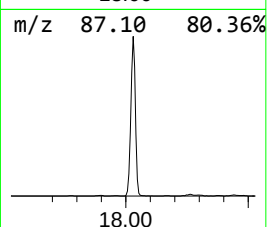
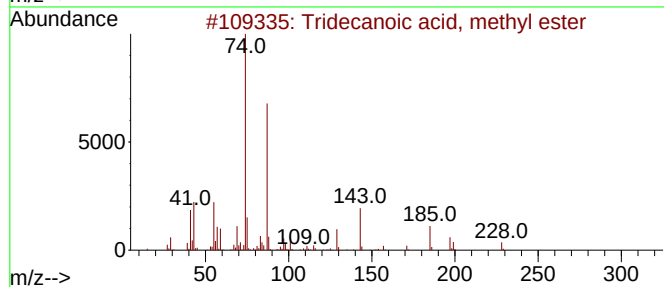
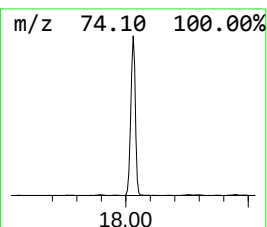
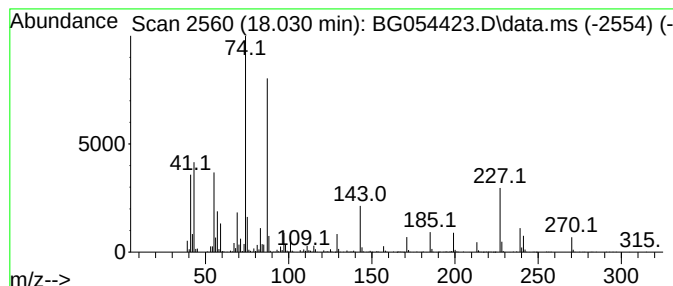
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG071522.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 2 Tridecanoic acid, methyl ester Concentration Rank 3

| R.T.      | EstConc                             | Area    | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|---------|------------------|--------------|------|
| 18.030    | 82.55 ng                            | 1158800 | Phenanthrene-d10 | 17.372       |      |
| Hit# of 5 | Tentative ID                        | MW      | MolForm          | CAS#         | Qual |
| 1         | Tridecanoic acid, methyl ester      | 228     | C14H28O2         | 001731-88-0  | 96   |
| 2         | Methyl 11-methyl-dodecanoate        | 228     | C14H28O2         | 1000336-45-1 | 93   |
| 3         | Pentadecanoic acid, 14-methyl-, ... | 270     | C17H34O2         | 005129-60-2  | 86   |
| 4         | Methyl tetradecanoate               | 242     | C15H30O2         | 000124-10-7  | 81   |
| 5         | Hexadecanoic acid, 2-methyl-        | 270     | C17H34O2         | 027147-71-3  | 81   |



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Sample : N3906-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
NB-7-7-26-22

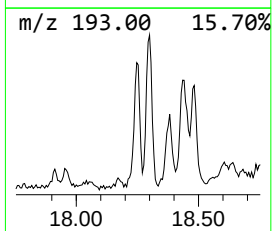
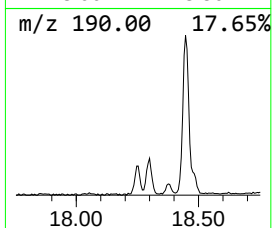
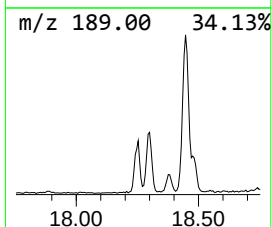
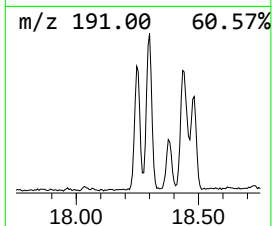
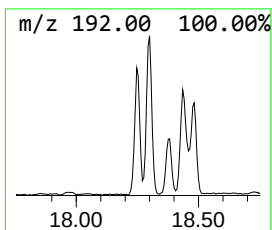
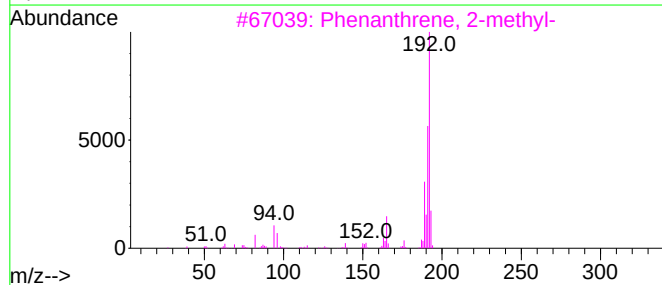
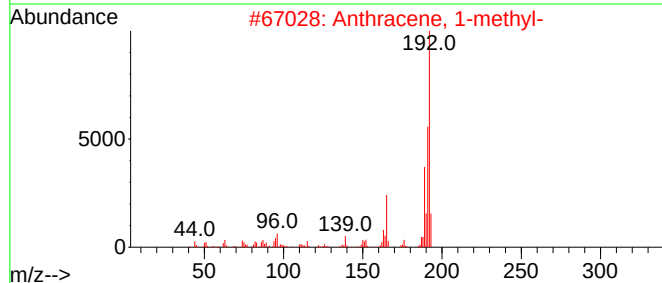
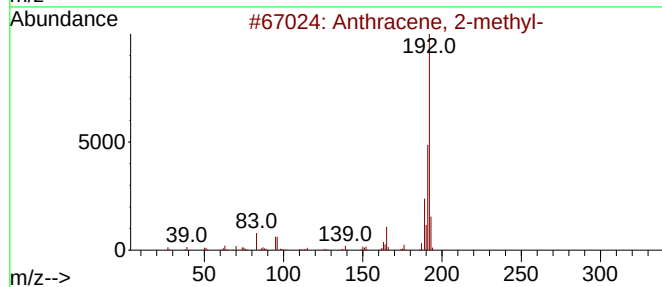
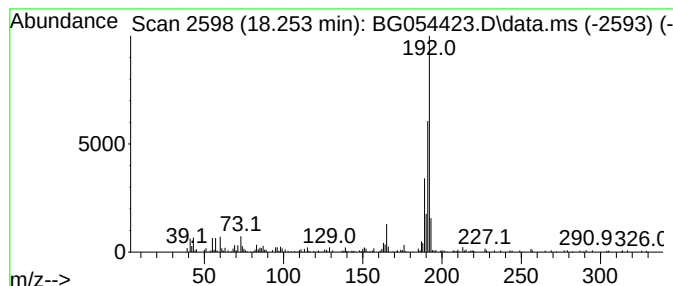
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG071522.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 3 Anthracene, 2-methyl- Concentration Rank 7

| R.T.   | EstConc  | Area   | Relative to ISTD | R.T.   |
|--------|----------|--------|------------------|--------|
| 18.253 | 14.59 ng | 204751 | Phenanthrene-d10 | 17.372 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Anthracene, 2-methyl-   | 192 | C15H12  | 000613-12-7 | 95   |
| 2    |    |   | Anthracene, 1-methyl-   | 192 | C15H12  | 000610-48-0 | 93   |
| 3    |    |   | Phenanthrene, 2-methyl- | 192 | C15H12  | 002531-84-2 | 91   |
| 4    |    |   | Phenanthrene, 1-methyl- | 192 | C15H12  | 000832-69-9 | 91   |
| 5    |    |   | Anthracene, 9-methyl-   | 192 | C15H12  | 000779-02-2 | 91   |



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Sample : N3906-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
NB-7-7-26-22

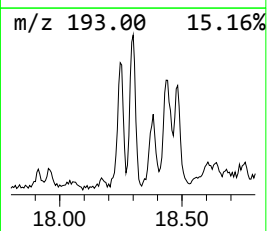
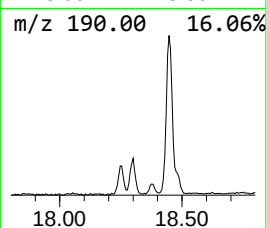
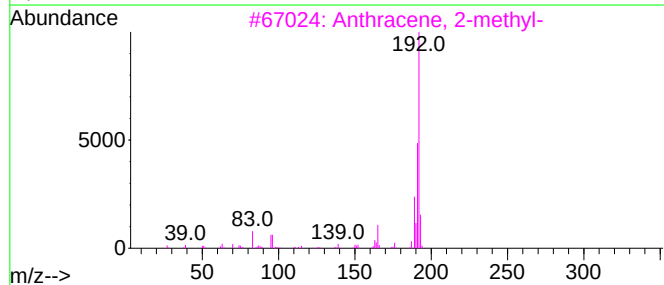
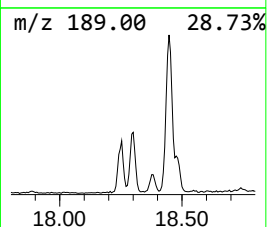
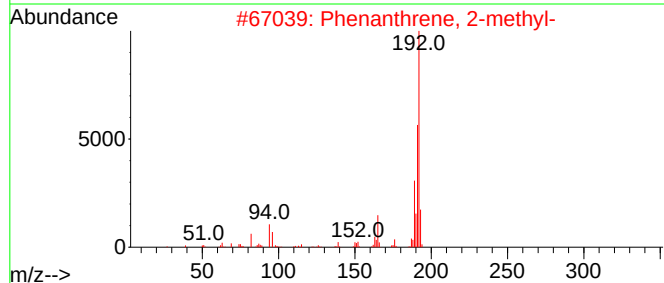
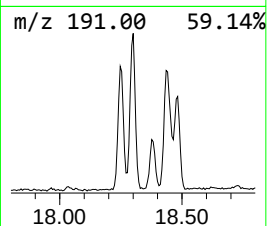
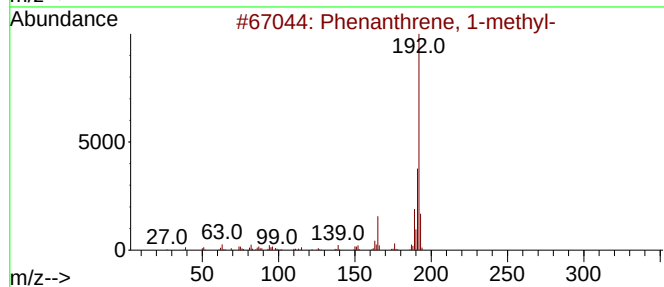
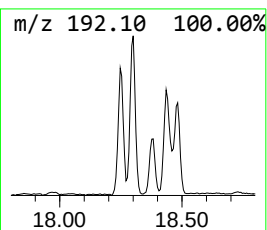
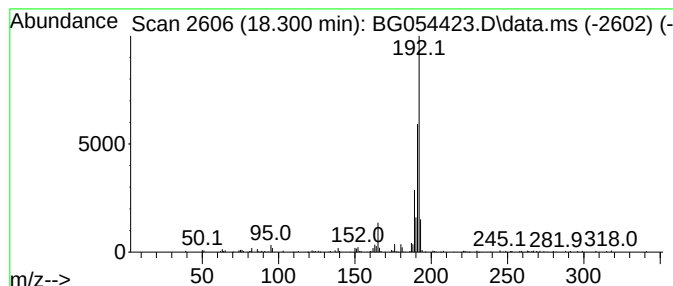
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 4 Phenanthrene, 1-methyl- Concentration Rank 8

| R.T.   | EstConc  | Area   | Relative to ISTD | R.T.   |
|--------|----------|--------|------------------|--------|
| 18.300 | 12.98 ng | 182191 | Phenanthrene-d10 | 17.372 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Phenanthrene, 1-methyl- | 192 | C15H12  | 000832-69-9 | 95   |
| 2    |    |   | Phenanthrene, 2-methyl- | 192 | C15H12  | 002531-84-2 | 93   |
| 3    |    |   | Anthracene, 2-methyl-   | 192 | C15H12  | 000613-12-7 | 93   |
| 4    |    |   | Anthracene, 9-methyl-   | 192 | C15H12  | 000779-02-2 | 91   |
| 5    |    |   | Anthracene, 1-methyl-   | 192 | C15H12  | 000610-48-0 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG072722\  
Data File : BG054423.D  
Acq On : 27 Jul 2022 19:53  
Operator : CG/JU  
Sample : N3906-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
NB-7-7-26-22

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

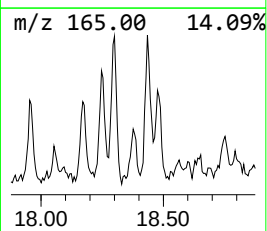
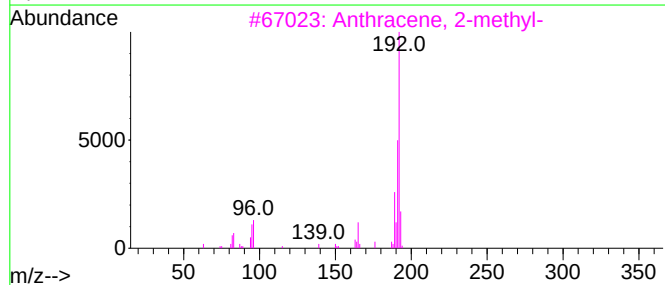
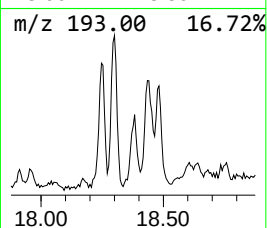
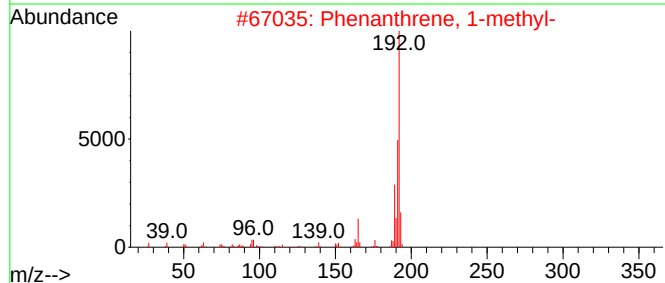
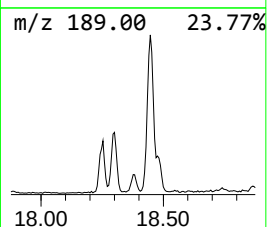
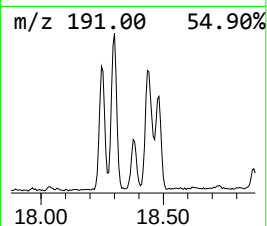
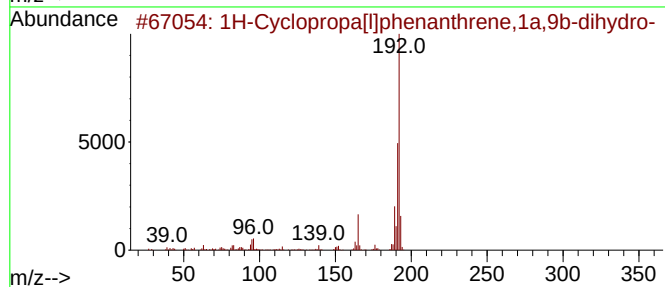
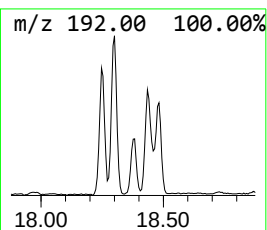
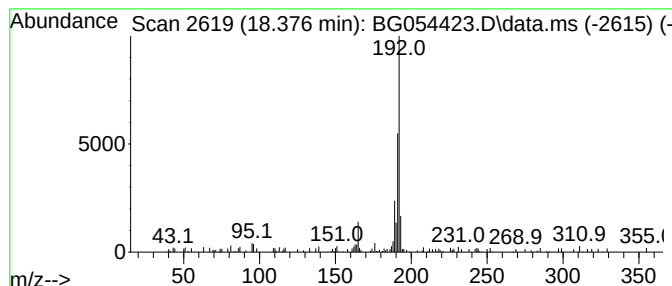
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TIC Integration Parameters: LSCINT.P

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Peak Number 5 1H-Cyclopropa[l]phenanthren... Concentration Rank 11

| R.T.   | EstConc | Area  | Relative to ISTD | R.T.   |
|--------|---------|-------|------------------|--------|
| 18.376 | 4.18 ng | 58741 | Phenanthrene-d10 | 17.372 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 1H-Cyclopropa[l]phenanthrene,1a,... | 192 | C15H12  | 000949-41-7 | 96   |
| 2    |    |   | Phenanthrene, 1-methyl-             | 192 | C15H12  | 000832-69-9 | 96   |
| 3    |    |   | Anthracene, 2-methyl-               | 192 | C15H12  | 000613-12-7 | 94   |
| 4    |    |   | 1H-Indene, 2-phenyl-                | 192 | C15H12  | 004505-48-0 | 93   |
| 5    |    |   | Phenanthrene, 2-methyl-             | 192 | C15H12  | 002531-84-2 | 92   |





Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG072722\  
Data File : BG054423.D  
Acq On : 27 Jul 2022 19:53  
Operator : CG/JU  
Sample : N3906-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
NB-7-7-26-22

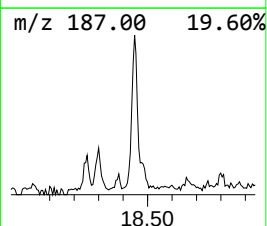
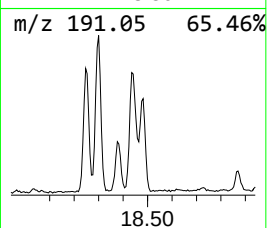
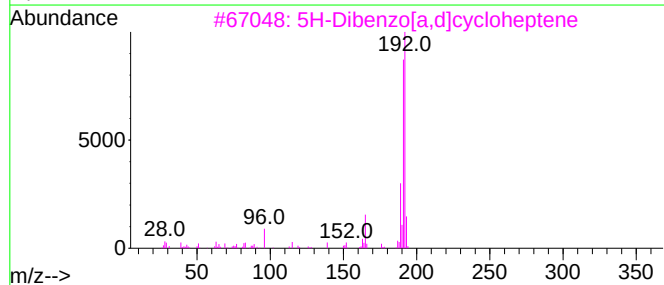
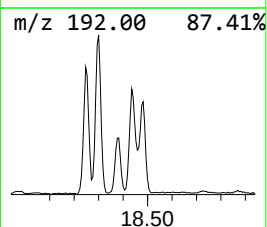
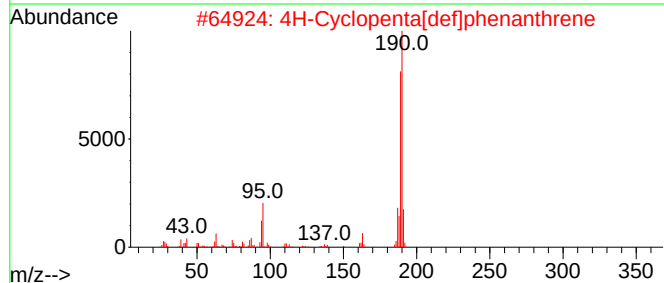
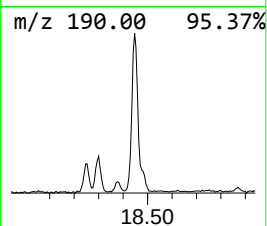
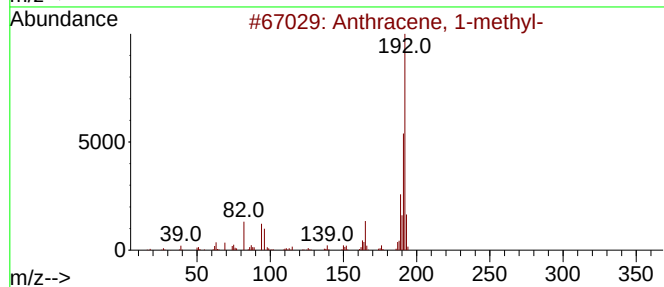
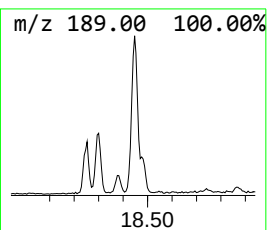
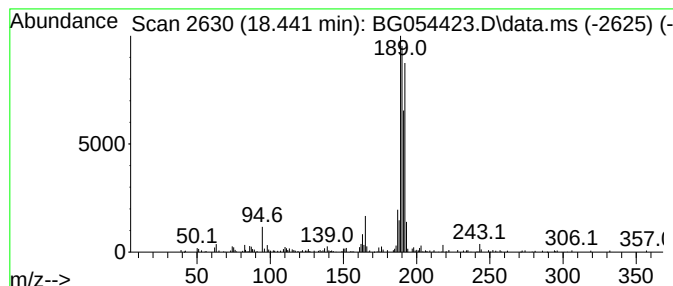
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 6 Anthracene, 1-methyl- Concentration Rank 6

| R.T.   | EstConc  | Area   | Relative to ISTD | R.T.   |
|--------|----------|--------|------------------|--------|
| 18.441 | 18.47 ng | 259336 | Phenanthrene-d10 | 17.372 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Anthracene, 1-methyl-          | 192 | C15H12  | 000610-48-0 | 60   |
| 2    |    |   | 4H-Cyclopenta[def]phenanthrene | 190 | C15H10  | 000203-64-5 | 55   |
| 3    |    |   | 5H-Dibenzo[a,d]cycloheptene    | 192 | C15H12  | 000256-81-5 | 42   |
| 4    |    |   | Phenanthrene, 4-methyl-        | 192 | C15H12  | 000832-64-4 | 42   |
| 5    |    |   | 1H-Indene, 2-phenyl-           | 192 | C15H12  | 004505-48-0 | 42   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG072722\  
Data File : BG054423.D  
Acq On : 27 Jul 2022 19:53  
Operator : CG/JU  
Sample : N3906-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
NB-7-7-26-22

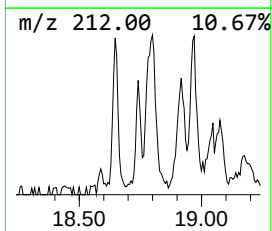
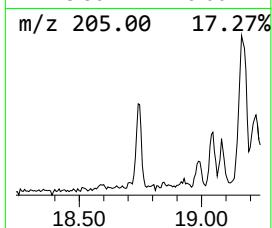
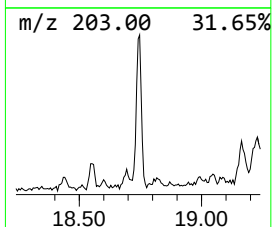
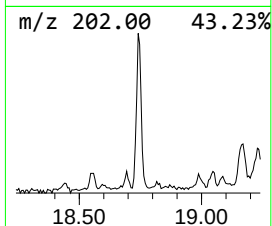
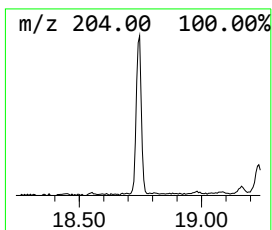
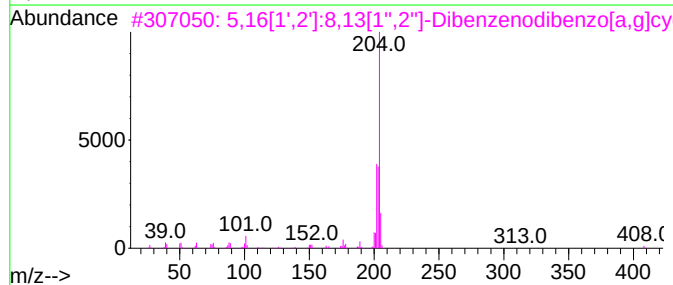
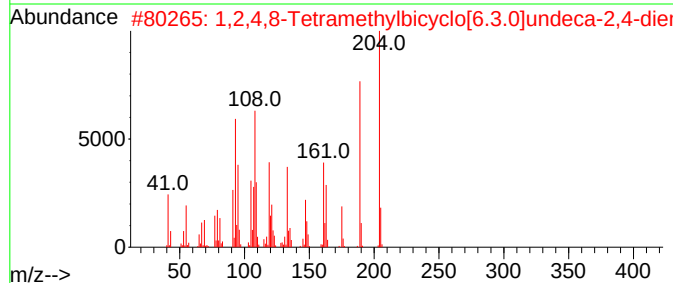
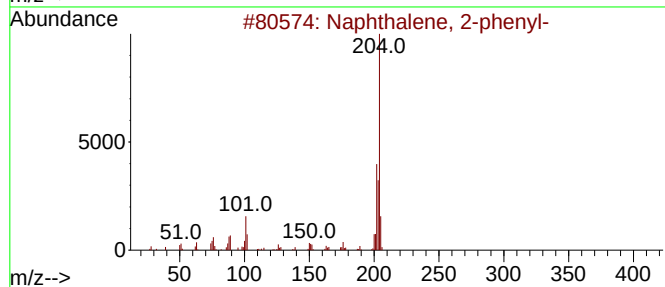
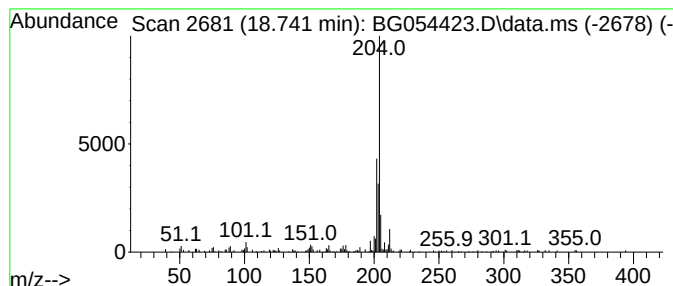
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 7 Naphthalene, 2-phenyl- Concentration Rank 10

| R.T.   | EstConc | Area  | Relative to ISTD | R.T.   |
|--------|---------|-------|------------------|--------|
| 18.741 | 6.89 ng | 96710 | Phenanthrene-d10 | 17.372 |

| Hit# | of | 5 | Tentative ID                         | MW  | MolForm  | CAS#         | Qual |
|------|----|---|--------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Naphthalene, 2-phenyl-               | 204 | C16H12   | 000612-94-2  | 90   |
| 2    |    |   | 1,2,4,8-Tetramethylbicyclo[6.3.0...] | 204 | C15H24   | 137235-51-9  | 86   |
| 3    |    |   | 5,16[1',2']:8,13[1'',2'']-Dibenz...  | 408 | C32H24   | 005672-97-9  | 62   |
| 4    |    |   | Tricyclo[8.2.2.2(4,7)]hexadeca-2...  | 204 | C16H12   | 006572-60-7  | 60   |
| 5    |    |   | Anthracene, 9,10-bis(0-anisilyoxy... | 450 | C30H26O4 | 1000154-05-7 | 56   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG072722\  
Data File : BG054423.D  
Acq On : 27 Jul 2022 19:53  
Operator : CG/JU  
Sample : N3906-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
NB-7-7-26-22

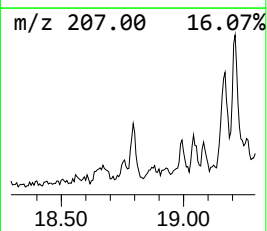
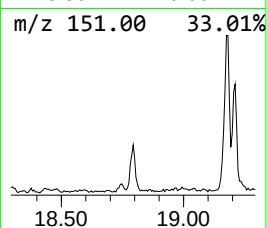
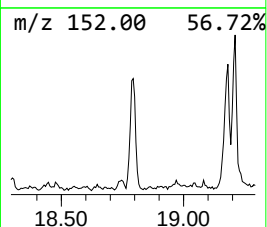
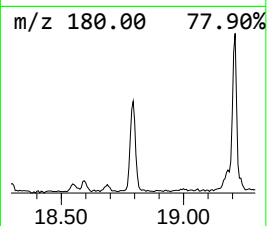
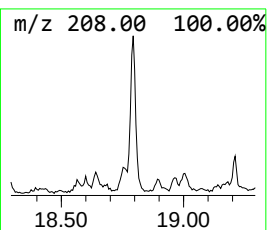
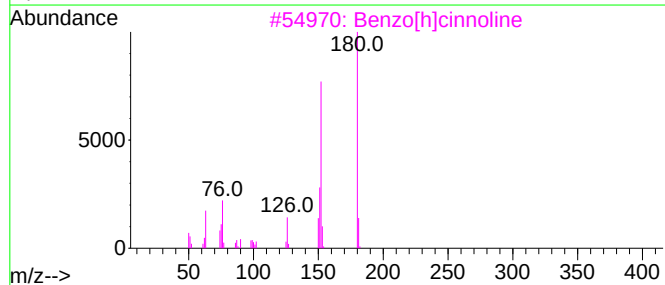
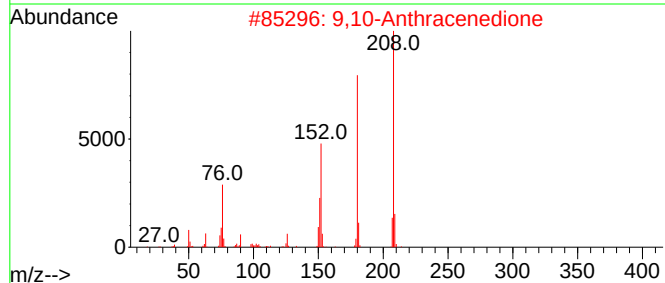
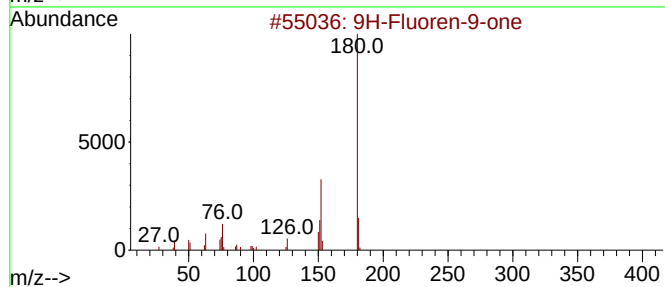
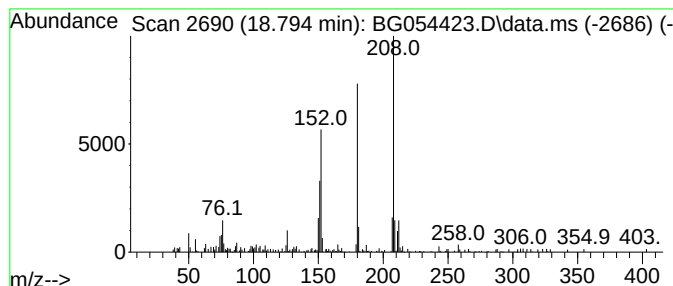
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 8 9H-Fluoren-9-one Concentration Rank 9

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 18.794 | 7.99 ng | 112221 | Phenanthrene-d10 | 17.372 |

| Hit# | of | 5 | Tentative ID                      | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-----------------------------------|-----|----------|-------------|------|
| 1    |    |   | 9H-Fluoren-9-one                  | 180 | C13H8O   | 000486-25-9 | 94   |
| 2    |    |   | 9,10-Anthracenedione              | 208 | C14H8O2  | 000084-65-1 | 93   |
| 3    |    |   | Benzo[h]cinnoline                 | 180 | C12H8N2  | 000230-31-9 | 60   |
| 4    |    |   | Benzo[c]cinnoline                 | 180 | C12H8N2  | 000230-17-1 | 50   |
| 5    |    |   | Benzaldehyde, 3-butoxy-4-methoxy- | 208 | C12H16O3 | 034127-96-3 | 43   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG072722\  
Data File : BG054423.D  
Acq On : 27 Jul 2022 19:53  
Operator : CG/JU  
Sample : N3906-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
NB-7-7-26-22

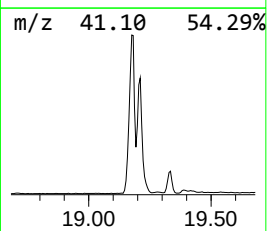
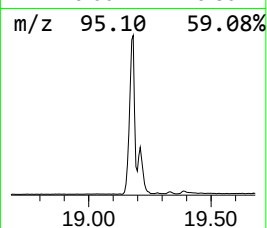
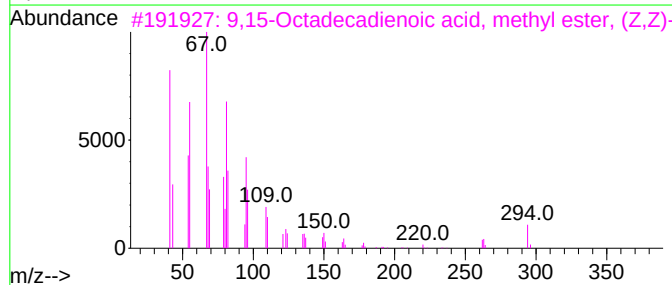
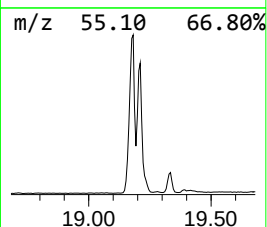
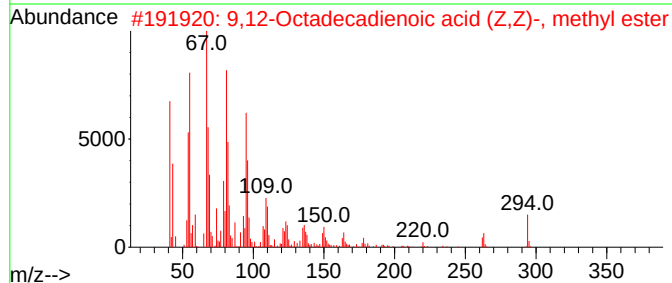
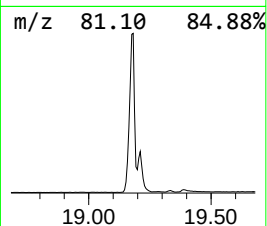
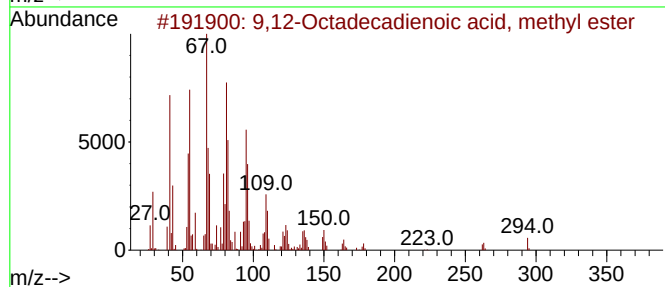
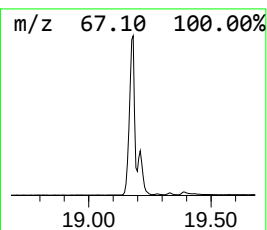
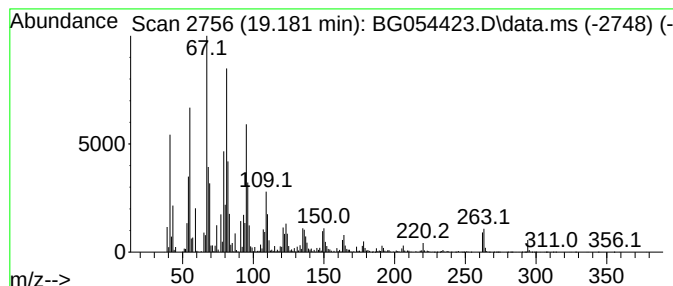
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

| R.T.   | EstConc   | Area                                | Relative to ISTD |          | R.T.         |      |
|--------|-----------|-------------------------------------|------------------|----------|--------------|------|
| 19.181 | 414.76 ng | 5822360                             | Phenanthrene-d10 |          | 17.372       |      |
| Hit#   | of 5      | Tentative ID                        | MW               | MolForm  | CAS#         | Qual |
| 1      |           | 9,12-Octadecadienoic acid, methy... | 294              | C19H34O2 | 002462-85-3  | 99   |
| 2      |           | 9,12-Octadecadienoic acid (Z,Z)-... | 294              | C19H34O2 | 000112-63-0  | 99   |
| 3      |           | 9,15-Octadecadienoic acid, methy... | 294              | C19H34O2 | 017309-05-6  | 99   |
| 4      |           | Methyl 10-trans,12-cis-octadecad... | 294              | C19H34O2 | 1000336-44-2 | 94   |
| 5      |           | Methyl 9-cis,11-trans-octadecadi... | 294              | C19H34O2 | 013058-52-1  | 94   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG072722\  
Data File : BG054423.D  
Acq On : 27 Jul 2022 19:53  
Operator : CG/JU  
Sample : N3906-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
NB-7-7-26-22

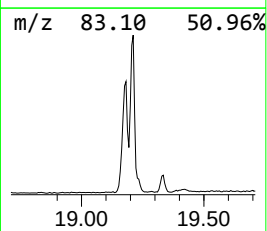
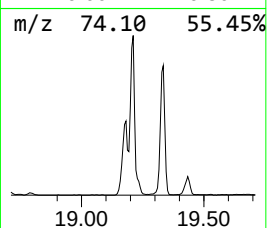
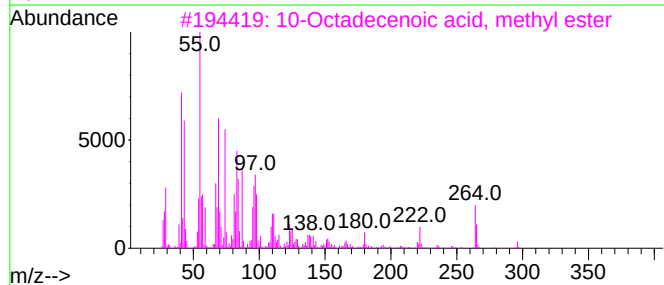
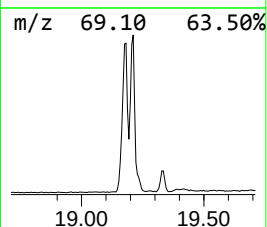
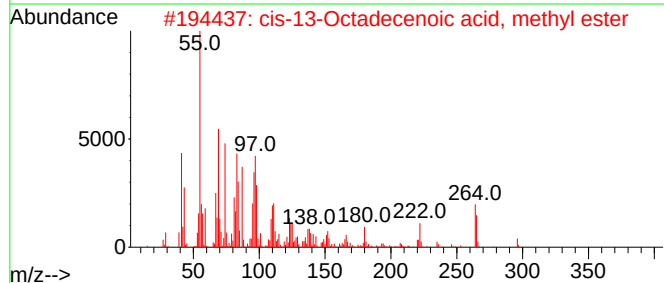
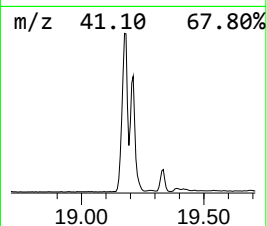
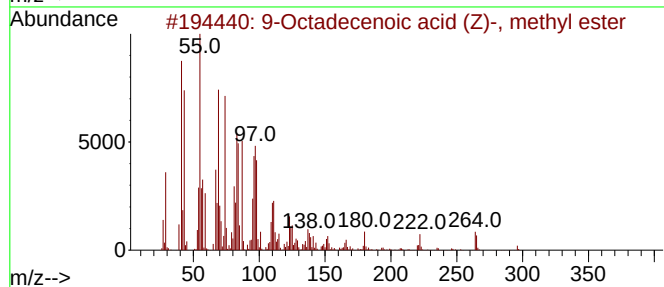
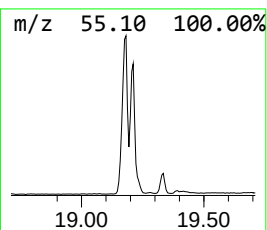
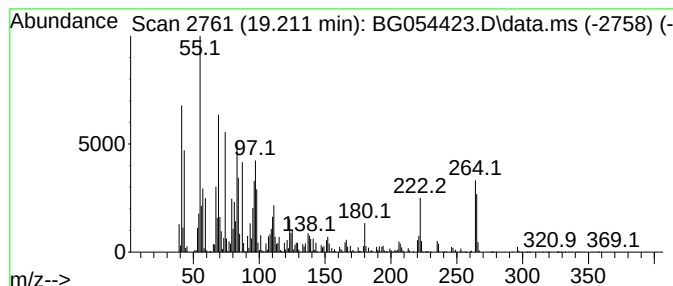
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 10 9-Octadecenoic acid (Z)-, m... Concentration Rank 2

| R.T.      | EstConc                             | Area    | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|---------|------------------|--------------|------|
| 19.211    | 241.42 ng                           | 3389010 | Phenanthrene-d10 | 17.372       |      |
| Hit# of 5 | Tentative ID                        | MW      | MolForm          | CAS#         | Qual |
| 1         | 9-Octadecenoic acid (Z)-, methyl... | 296     | C19H36O2         | 000112-62-9  | 99   |
| 2         | cis-13-Octadecenoic acid, methyl... | 296     | C19H36O2         | 1010333-58-3 | 98   |
| 3         | 10-Octadecenoic acid, methyl ester  | 296     | C19H36O2         | 013481-95-3  | 98   |
| 4         | 11-Octadecenoic acid, methyl ester  | 296     | C19H36O2         | 052380-33-3  | 96   |
| 5         | trans-13-Octadecenoic acid, meth... | 296     | C19H36O2         | 1000333-61-3 | 96   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG072722\  
Data File : BG054423.D  
Acq On : 27 Jul 2022 19:53  
Operator : CG/JU  
Sample : N3906-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
NB-7-7-26-22

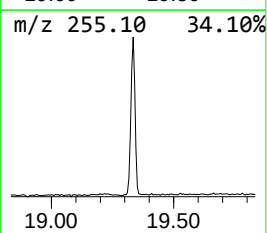
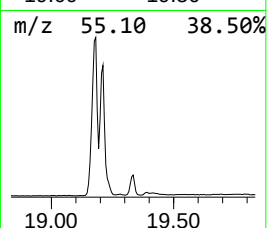
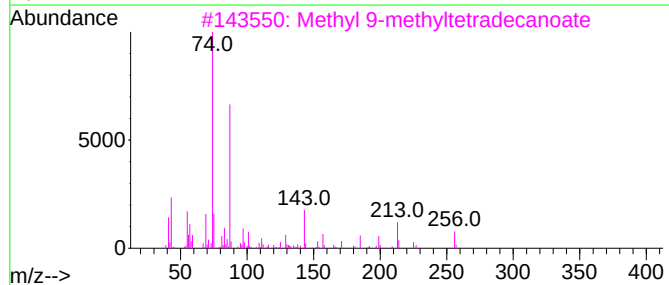
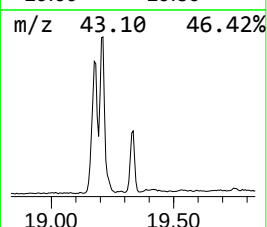
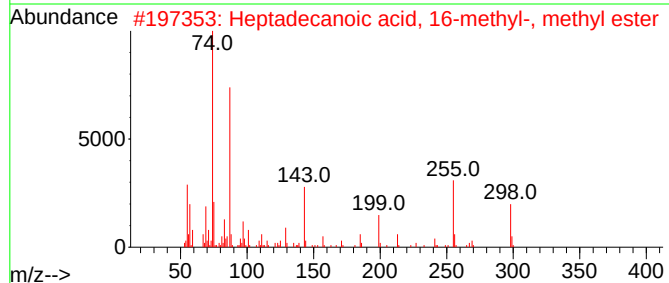
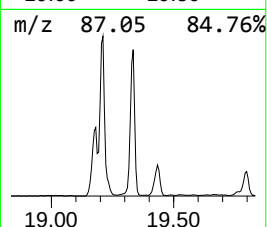
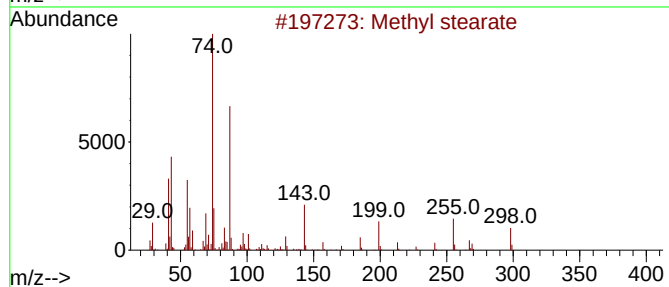
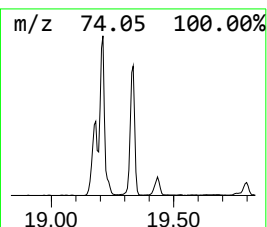
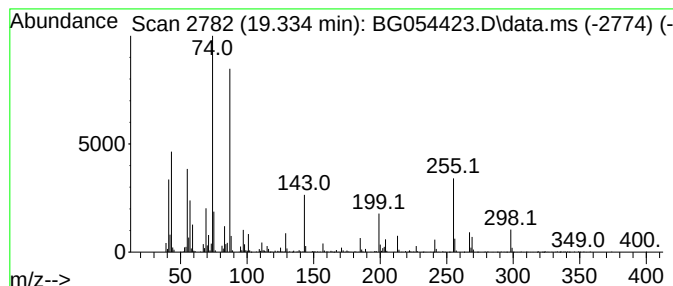
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 11 Methyl stearate Concentration Rank 4

| R.T.   | EstConc  | Area   | Relative to ISTD | R.T.   |
|--------|----------|--------|------------------|--------|
| 19.334 | 57.25 ng | 803622 | Phenanthrene-d10 | 17.372 |

| 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 9-methyltetradecanoate       | 256 | C16H32O2 | 213617-69-7 | 81   |
| 4    |    |   | Pentadecanoic acid, methyl ester    | 256 | C16H32O2 | 007132-64-1 | 76   |
| 5    |    |   | Nonadecanoic acid, methyl ester     | 312 | C20H40O2 | 001731-94-8 | 74   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG072722\  
Data File : BG054423.D  
Acq On : 27 Jul 2022 19:53  
Operator : CG/JU  
Sample : N3906-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
NB-7-7-26-22

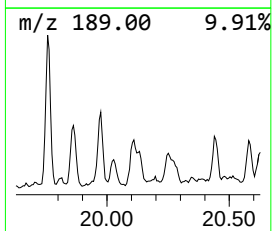
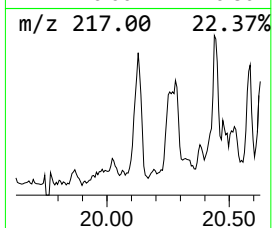
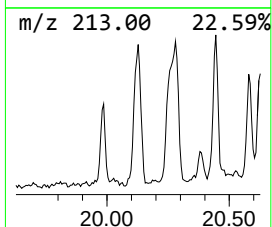
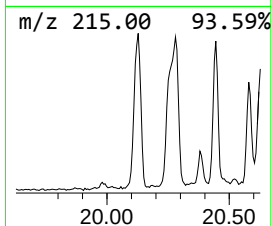
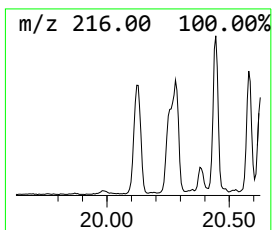
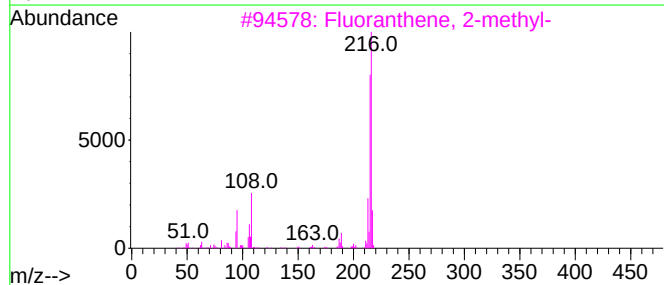
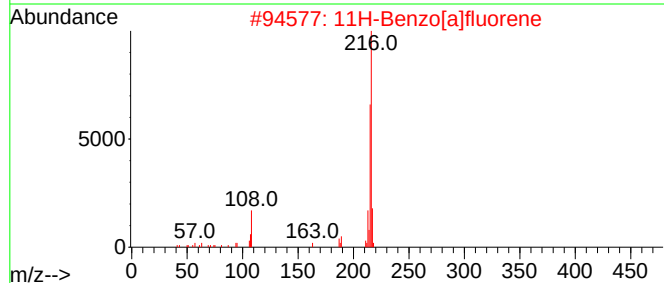
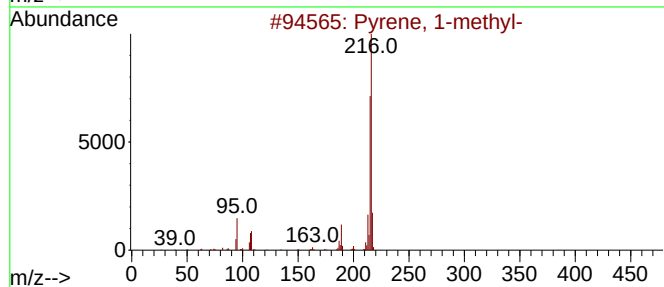
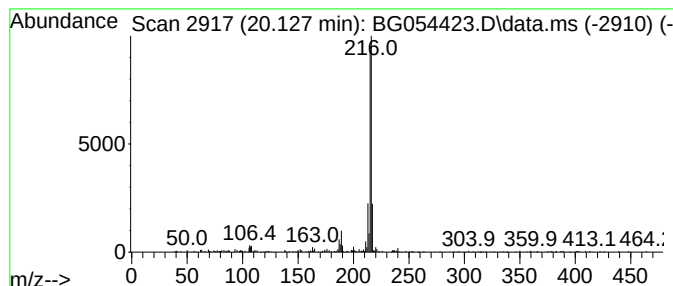
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 12 Pyrene, 1-methyl- Concentration Rank 14

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 20.127 | 2.90 ng | 290427 | Chrysene-d12     | 21.643 |

| Hit# | of 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------|-----|---------|-------------|------|
| 1    |      | Pyrene, 1-methyl-       | 216 | C17H12  | 002381-21-7 | 93   |
| 2    |      | 11H-Benzo[a]fluorene    | 216 | C17H12  | 000238-84-6 | 87   |
| 3    |      | Fluoranthene, 2-methyl- | 216 | C17H12  | 033543-31-6 | 87   |
| 4    |      | 11H-Benzo[b]fluorene    | 216 | C17H12  | 000243-17-4 | 87   |
| 5    |      | 7H-Benzo[c]fluorene     | 216 | C17H12  | 000205-12-9 | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG072722\  
Data File : BG054423.D  
Acq On : 27 Jul 2022 19:53  
Operator : CG/JU  
Sample : N3906-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
NB-7-7-26-22

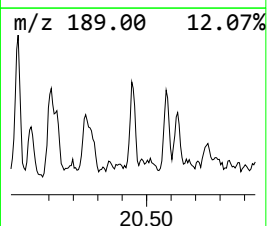
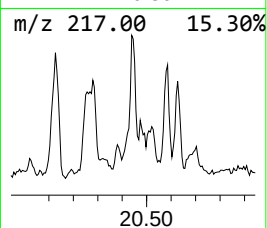
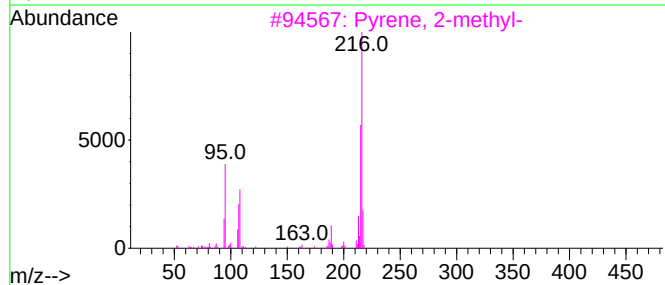
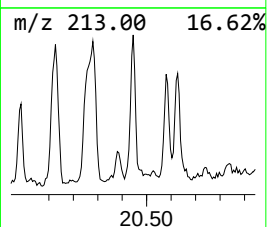
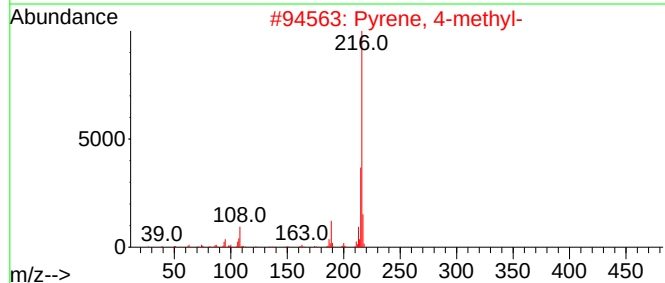
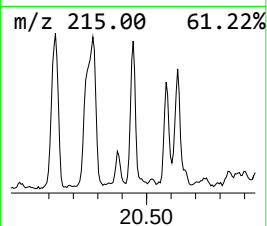
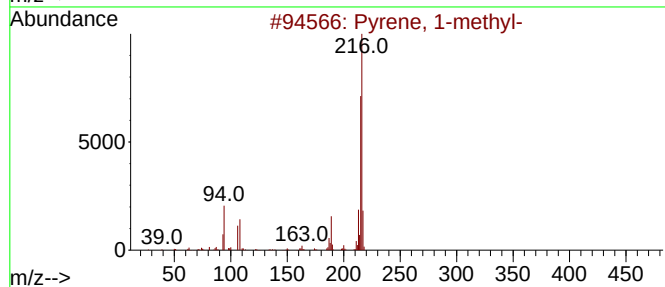
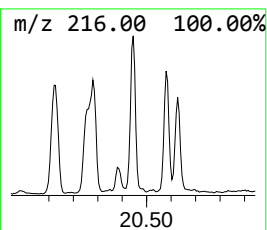
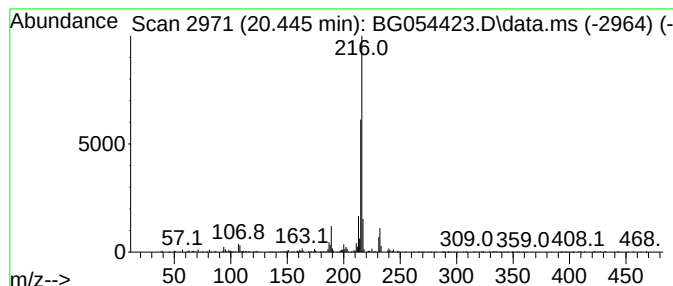
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 13 Pyrene, 4-methyl- Concentration Rank 13

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 20.445 | 3.44 ng | 344668 | Chrysene-d12     | 21.643 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Pyrene, 1-methyl-       | 216 | C17H12  | 002381-21-7 | 94   |
| 2    |    |   | Pyrene, 4-methyl-       | 216 | C17H12  | 003353-12-6 | 94   |
| 3    |    |   | Pyrene, 2-methyl-       | 216 | C17H12  | 003442-78-2 | 93   |
| 4    |    |   | Fluoranthene, 2-methyl- | 216 | C17H12  | 033543-31-6 | 91   |
| 5    |    |   | 11H-Benzo[a]fluorene    | 216 | C17H12  | 000238-84-6 | 86   |





Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG072722\  
Data File : BG054423.D  
Acq On : 27 Jul 2022 19:53  
Operator : CG/JU  
Sample : N3906-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
NB-7-7-26-22

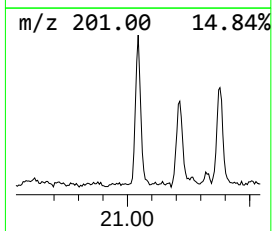
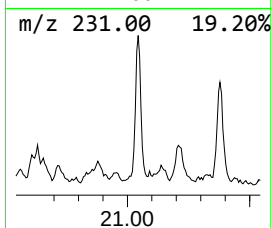
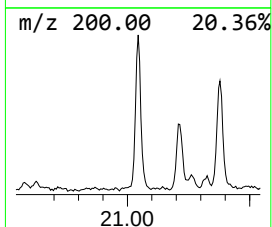
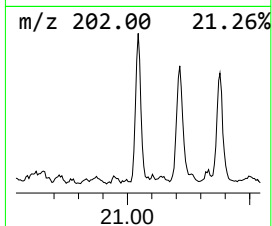
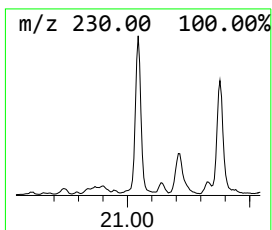
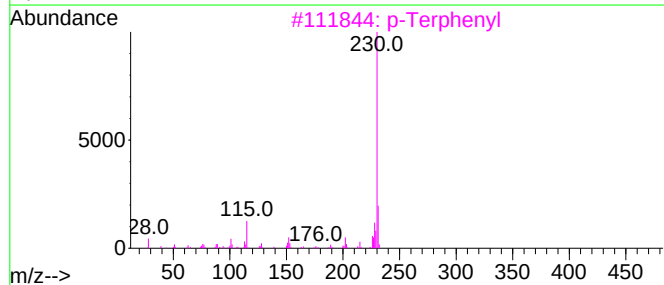
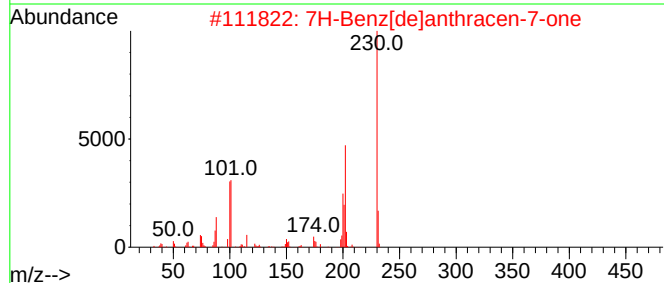
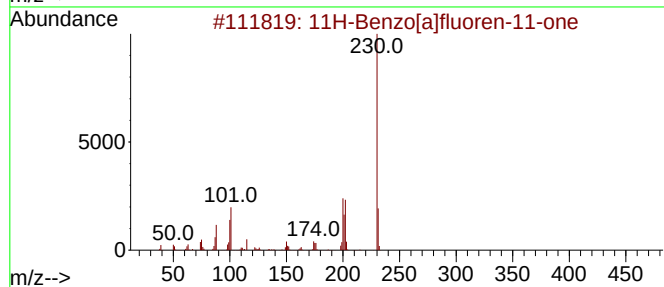
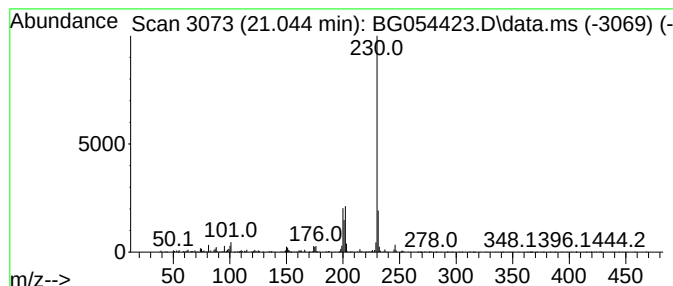
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 14 11H-Benzo[a]fluoren-11-one Concentration Rank 12

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 21.044 | 4.02 ng | 401884 | Chrysene-d12     | 21.643 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | 11H-Benzo[a]fluoren-11-one          | 230 | C17H10O    | 000479-79-8  | 95   |
| 2    |    |   | 7H-Benz[de]anthracen-7-one          | 230 | C17H10O    | 000082-05-3  | 87   |
| 3    |    |   | p-Terphenyl                         | 230 | C18H14     | 000092-94-4  | 50   |
| 4    |    |   | 3H-pyrazol-3-one, 2-(1,3-dihydro... | 230 | C12H10N2O3 | 1000396-36-9 | 50   |
| 5    |    |   | m-Terphenyl                         | 230 | C18H14     | 000092-06-8  | 50   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG072722\  
Data File : BG054423.D  
Acq On : 27 Jul 2022 19:53  
Operator : CG/JU  
Sample : N3906-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
NB-7-7-26-22

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG071522.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

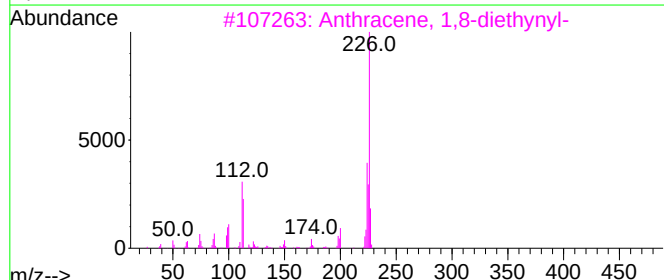
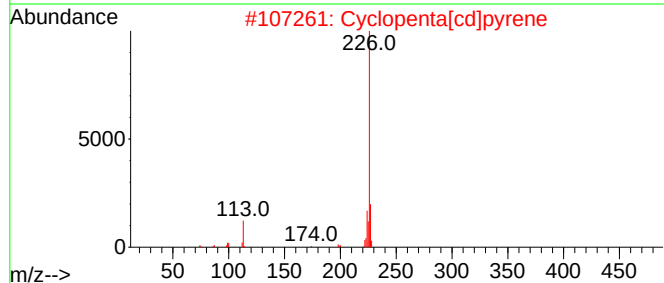
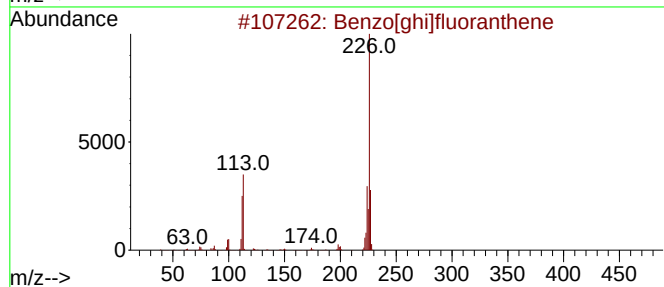
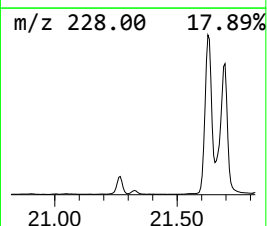
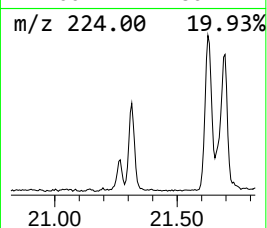
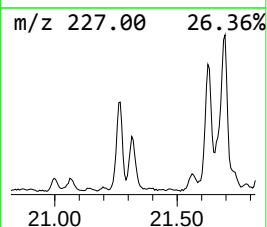
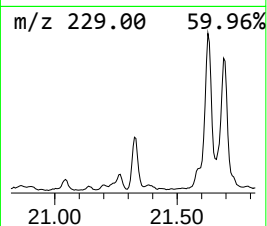
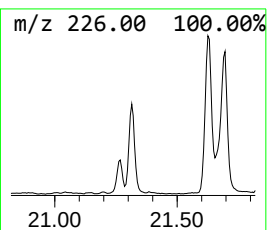
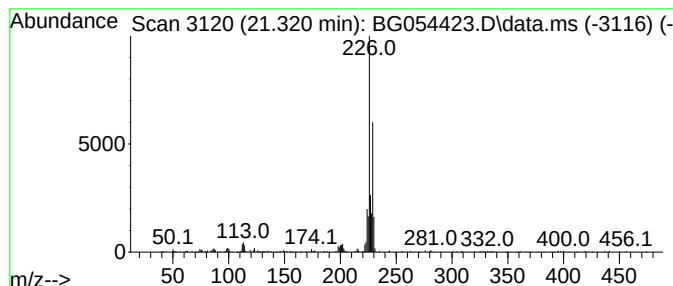
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 15 unknown21.320 Concentration Rank 16

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 21.320 | 2.13 ng | 213610 | Chrysene-d12     | 21.643 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Benzo[ghi]fluoranthene              | 226 | C18H10    | 000203-12-3  | 43   |
| 2    |    |   | Cyclopenta[cd]pyrene                | 226 | C18H10    | 027208-37-3  | 43   |
| 3    |    |   | Anthracene, 1,8-diethynyl-          | 226 | C18H10    | 078053-58-4  | 38   |
| 4    |    |   | 6-Bromo-3,4-dihydro-2(1H)-quinaz... | 226 | C8H7BrN2O | 1246765-38-7 | 37   |
| 5    |    |   | N,N-Dimethylindooaniline            | 226 | C14H14N2O | 002150-58-5  | 37   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG072722\  
Data File : BG054423.D  
Acq On : 27 Jul 2022 19:53  
Operator : CG/JU  
Sample : N3906-01  
Misc :  
ALS Vial : 16 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
NB-7-7-26-22

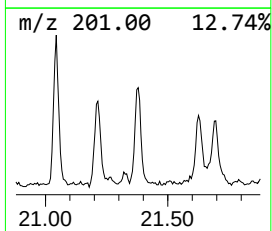
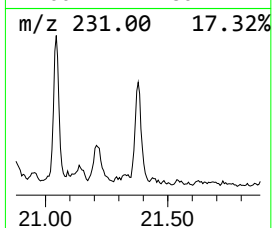
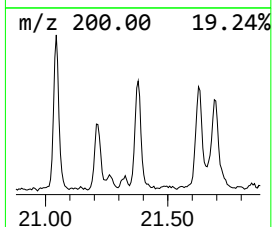
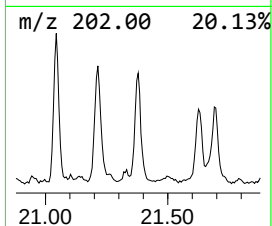
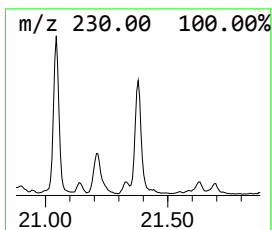
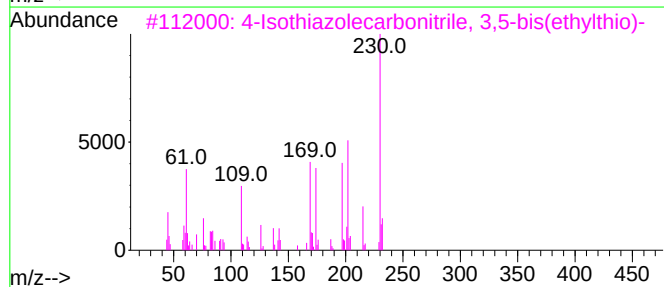
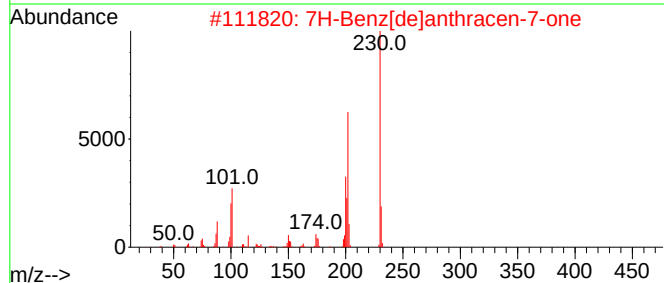
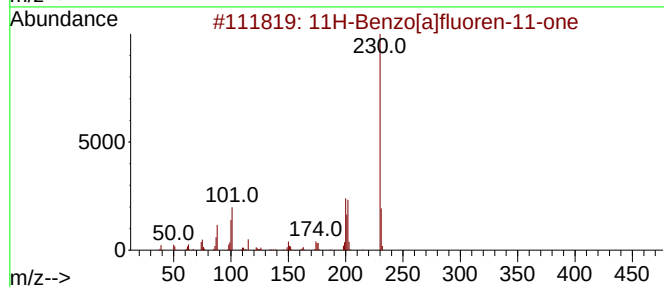
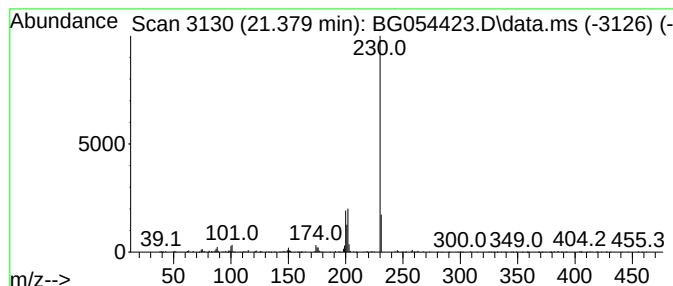
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 16 7H-Benz[de]anthracen-7-one Concentration Rank 15

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|-------------|------|
| 21.379    | 2.83 ng                             | 283384 | Chrysene-d12     | 21.643      |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual |
| 1         | 11H-Benzo[a]fluoren-11-one          | 230    | C17H10O          | 000479-79-8 | 94   |
| 2         | 7H-Benz[de]anthracen-7-one          | 230    | C17H10O          | 000082-05-3 | 83   |
| 3         | 4-Isothiazolecarbonitrile, 3,5-b... | 230    | C8H10N2S3        | 024135-13-5 | 59   |
| 4         | p-Terphenyl                         | 230    | C18H14           | 000092-94-4 | 50   |
| 5         | m-Terphenyl                         | 230    | C18H14           | 000092-06-8 | 50   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG072722\  
 Data File : BG054423.D  
 Acq On : 27 Jul 2022 19:53  
 Operator : CG/JU  
 Sample : N3906-01  
 Misc :  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 NB-7-7-26-22

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG071522.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 |
| Butane, 2-metho... | 3.118  | 26.4    | ng    | 149278   | 1                     | 7.989  | 112917  | 20.0 |
| Tridecanoic aci... | 18.030 | 82.5    | ng    | 1158800  | 4                     | 17.372 | 280757  | 20.0 |
| Anthracene, 2-m... | 18.253 | 14.6    | ng    | 204751   | 4                     | 17.372 | 280757  | 20.0 |
| Phenanthrene, 1... | 18.300 | 13.0    | ng    | 182191   | 4                     | 17.372 | 280757  | 20.0 |
| 1H-Cyclopropa[1... | 18.376 | 4.2     | ng    | 58741    | 4                     | 17.372 | 280757  | 20.0 |
| Anthracene, 1-m... | 18.441 | 18.5    | ng    | 259336   | 4                     | 17.372 | 280757  | 20.0 |
| Naphthalene, 2-... | 18.741 | 6.9     | ng    | 96710    | 4                     | 17.372 | 280757  | 20.0 |
| 9H-Fluoren-9-one   | 18.794 | 8.0     | ng    | 112221   | 4                     | 17.372 | 280757  | 20.0 |
| 9,12-Octadecadi... | 19.181 | 414.8   | ng    | 5822360  | 4                     | 17.372 | 280757  | 20.0 |
| 9-Octadecenoic ... | 19.211 | 241.4   | ng    | 3389010  | 4                     | 17.372 | 280757  | 20.0 |
| Methyl stearate    | 19.334 | 57.3    | ng    | 803622   | 4                     | 17.372 | 280757  | 20.0 |
| Pyrene, 1-methyl-  | 20.127 | 2.9     | ng    | 290427   | 5                     | 21.643 | 2001910 | 20.0 |
| Pyrene, 4-methyl-  | 20.445 | 3.4     | ng    | 344668   | 5                     | 21.643 | 2001910 | 20.0 |
| 11H-Benzo[a]flu... | 21.044 | 4.0     | ng    | 401884   | 5                     | 21.643 | 2001910 | 20.0 |
| unknown21.320      | 21.320 | 2.1     | ng    | 213610   | 5                     | 21.643 | 2001910 | 20.0 |
| 7H-Benz[de]anth... | 21.379 | 2.8     | ng    | 283384   | 5                     | 21.643 | 2001910 | 20.0 |