

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP040122\  
 Data File : BP009664.D  
 Acq On : 01 Apr 2022 18:38  
 Operator : CG/JU  
 Sample : N2220-01  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 A10015

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-BP031722.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP009664.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         | 5.334       | 393           | 399         | 427          | rBV      | 2779051        | 4960344       | 13.90%          | 3.983%        |
| 2         | 6.922       | 656           | 669         | 689          | rBV      | 2517680        | 5143554       | 14.41%          | 4.130%        |
| 3         | 7.728       | 797           | 806         | 821          | rBV      | 821594         | 1525704       | 4.28%           | 1.225%        |
| 4         | 8.887       | 996           | 1003        | 1013         | rBV      | 1896589        | 3622643       | 10.15%          | 2.909%        |
| 5         | 10.516      | 1272          | 1280        | 1286         | rBV      | 1085878        | 2103041       | 5.89%           | 1.689%        |
| 6         | 12.193      | 1556          | 1565        | 1574         | rVB      | 264099         | 480269        | 1.35%           | 0.386%        |
| 7         | 12.410      | 1596          | 1602        | 1610         | rBV      | 279223         | 472692        | 1.32%           | 0.380%        |
| 8         | 12.998      | 1694          | 1702        | 1722         | rBV      | 5149240        | 8558861       | 23.98%          | 6.873%        |
| 9         | 13.698      | 1815          | 1821        | 1826         | rBV      | 252068         | 441229        | 1.24%           | 0.354%        |
| 10        | 13.751      | 1826          | 1830        | 1841         | rVB      | 211348         | 378809        | 1.06%           | 0.304%        |
| 11        | 14.375      | 1927          | 1936        | 1943         | rBV      | 1690881        | 2801459       | 7.85%           | 2.250%        |
| 12        | 15.175      | 2065          | 2072        | 2076         | rBV3     | 174511         | 365934        | 1.03%           | 0.294%        |
| 13        | 15.428      | 2109          | 2115        | 2120         | rBV2     | 243966         | 406262        | 1.14%           | 0.326%        |
| 14        | 15.881      | 2184          | 2192        | 2202         | rBV      | 3757942        | 6151693       | 17.24%          | 4.940%        |
| 15        | 16.139      | 2227          | 2236        | 2242         | rBV2     | 539125         | 1075996       | 3.02%           | 0.864%        |
| 16        | 16.686      | 2323          | 2329        | 2333         | rBV3     | 184230         | 378714        | 1.06%           | 0.304%        |
| 17        | 17.139      | 2400          | 2406        | 2410         | rBV      | 2128209        | 3217669       | 9.02%           | 2.584%        |
| 18        | 17.186      | 2410          | 2414        | 2420         | rVB      | 1290282        | 1919155       | 5.38%           | 1.541%        |
| 19        | 18.022      | 2551          | 2556        | 2558         | rBV      | 292916         | 393412        | 1.10%           | 0.316%        |
| 20        | 18.063      | 2558          | 2563        | 2571         | rVB2     | 766274         | 1449480       | 4.06%           | 1.164%        |
| 21        | 18.204      | 2582          | 2587        | 2591         | rBV2     | 386353         | 688080        | 1.93%           | 0.553%        |
| 22        | 18.245      | 2591          | 2594        | 2600         | rVB      | 285221         | 418094        | 1.17%           | 0.336%        |
| 23        | 18.916      | 2703          | 2708        | 2711         | rBV      | 262044         | 387280        | 1.09%           | 0.311%        |
| 24        | 19.169      | 2746          | 2751        | 2759         | rBV      | 2172890        | 3136050       | 8.79%           | 2.518%        |
| 25        | 19.522      | 2803          | 2811        | 2814         | rBV2     | 1836944        | 3111709       | 8.72%           | 2.499%        |
| 26        | 19.563      | 2814          | 2818        | 2832         | rVB      | 2107597        | 3809031       | 10.67%          | 3.059%        |
| 27        | 19.727      | 2840          | 2846        | 2851         | rVB      | 9227246        | 11806374      | 33.08%          | 9.481%        |
| 28        | 19.851      | 2861          | 2867        | 2874         | rBV4     | 205577         | 558771        | 1.57%           | 0.449%        |
| 29        | 20.757      | 3017          | 3021        | 3026         | rBV      | 302161         | 435797        | 1.22%           | 0.350%        |
| 30        | 20.986      | 3056          | 3060        | 3066         | rBV3     | 1269048        | 1857828       | 5.21%           | 1.492%        |
| 31        | 21.045      | 3068          | 3070        | 3076         | rVB      | 289096         | 375607        | 1.05%           | 0.302%        |
| 32        | 21.263      | 3100          | 3107        | 3111         | rBV2     | 3226154        | 5966986       | 16.72%          | 4.792%        |
| 33        | 21.298      | 3111          | 3113        | 3118         | rVB      | 1119525        | 1221505       | 3.42%           | 0.981%        |
| 34        | 21.386      | 3121          | 3128        | 3142         | rVV      | 22432444       | 35685722      | 100.00%         | 28.657%       |
| 35        | 22.916      | 3383          | 3388        | 3394         | rBV      | 1117041        | 2576350       | 7.22%           | 2.069%        |
| 36        | 23.427      | 3472          | 3475        | 3484         | rVB      | 575498         | 1053504       | 2.95%           | 0.846%        |

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Sample : N2220-01  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A10015

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-BP031722.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 37 | 23.533 | 3488 | 3493 | 3501 | rVB  | 930215  | 1802871 | 5.05%  | 1.448% |
| 38 | 23.639 | 3504 | 3511 | 3517 | rBV2 | 1800644 | 3788734 | 10.62% | 3.042% |

Sum of corrected areas: 124527213



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Instrument :  
BNA\_P  
ClientSampleId :  
A10015

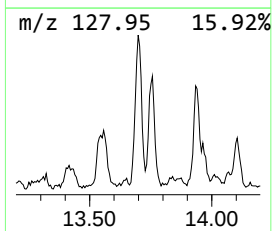
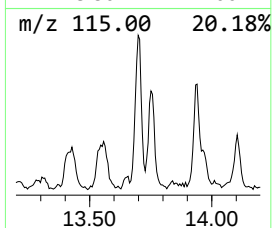
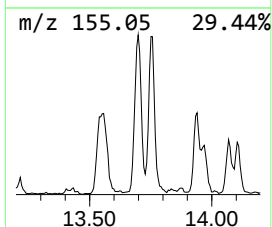
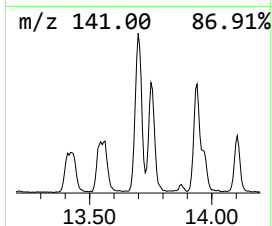
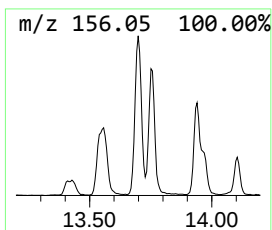
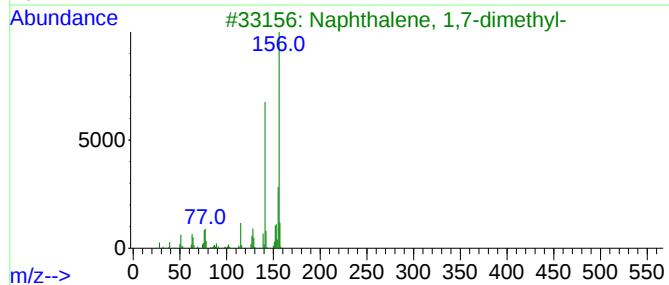
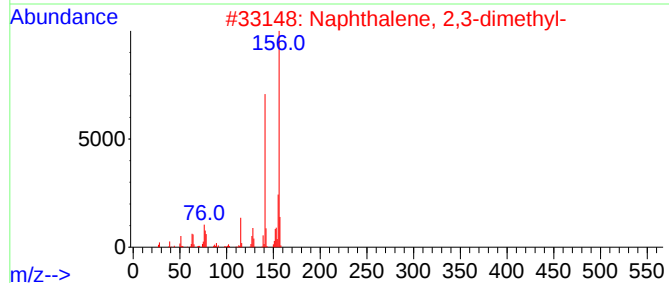
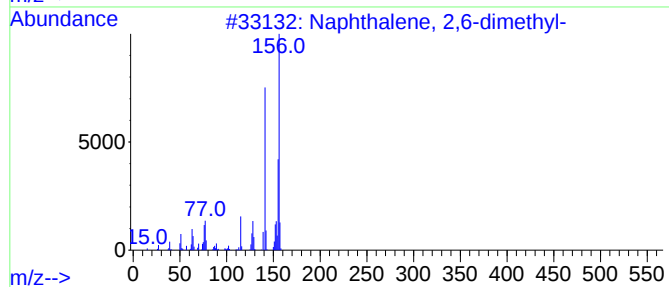
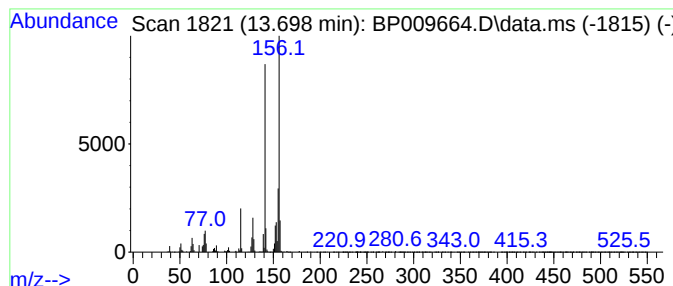
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP031722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 1 Naphthalene, 2,6-dimethyl- Concentration Rank 7

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 13.698 | 3.15 ng | 441229 | Acenaphthene-d10 | 14.375 |

| Hit# | of 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|------|----------------------------|-----|---------|-------------|------|
| 1    |      | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 97   |
| 2    |      | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 97   |
| 3    |      | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 97   |
| 4    |      | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 97   |
| 5    |      | Naphthalene, 1,2-dimethyl- | 156 | C12H12  | 000573-98-8 | 97   |



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Sample : N2220-01  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A10015

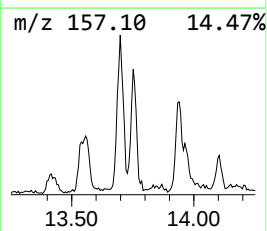
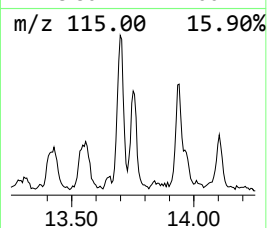
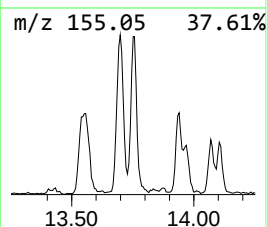
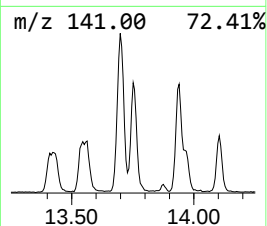
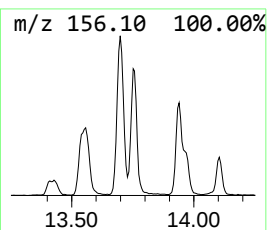
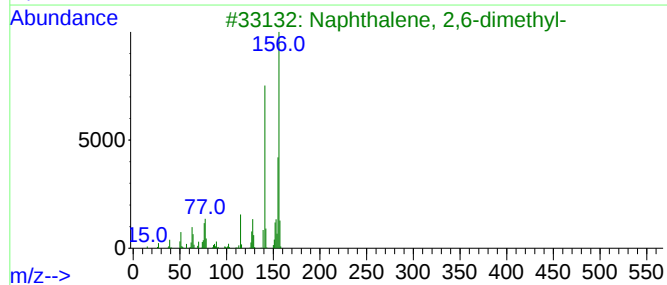
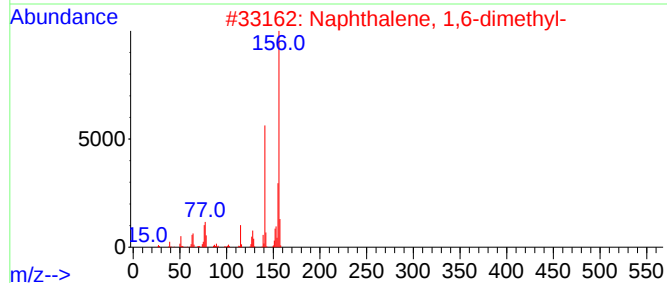
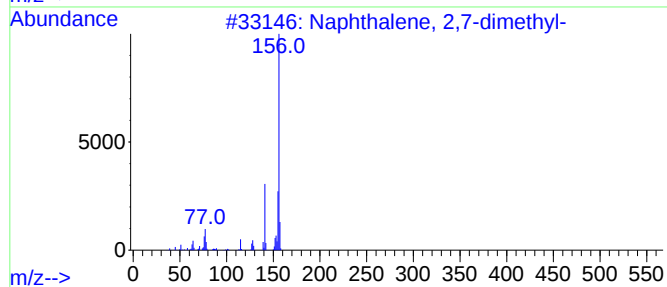
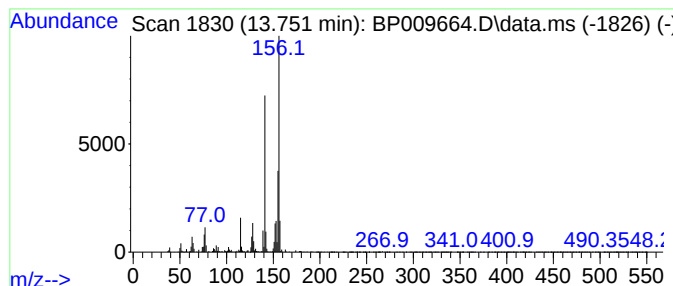
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP031722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 2 Naphthalene, 2,7-dimethyl- Concentration Rank 9

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 13.751 | 2.70 ng | 378809 | Acenaphthene-d10 | 14.375 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 2,7-dimethyl- | 156 | C12H12  | 000582-16-1 | 97   |
| 2    |    |   | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 97   |
| 3    |    |   | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 97   |
| 4    |    |   | Naphthalene, 1,5-dimethyl- | 156 | C12H12  | 000571-61-9 | 97   |
| 5    |    |   | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 97   |



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Sample : N2220-01  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A10015

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP031722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

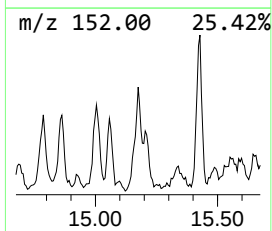
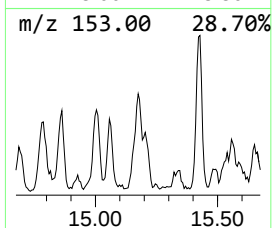
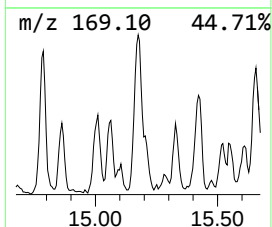
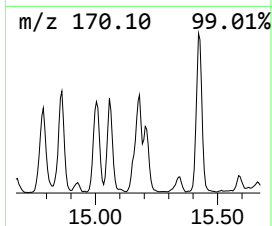
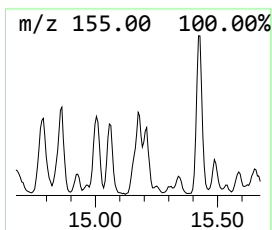
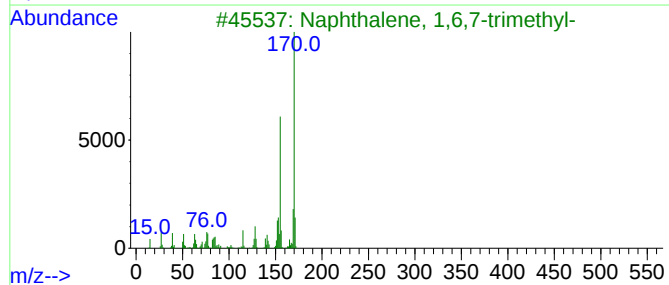
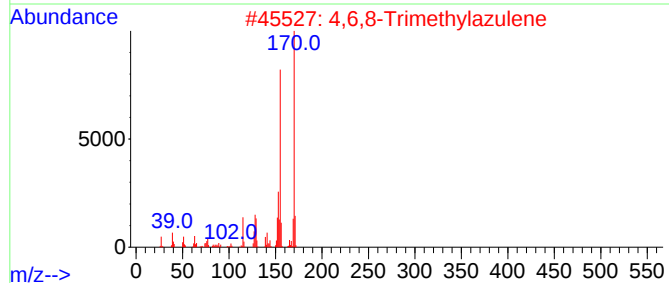
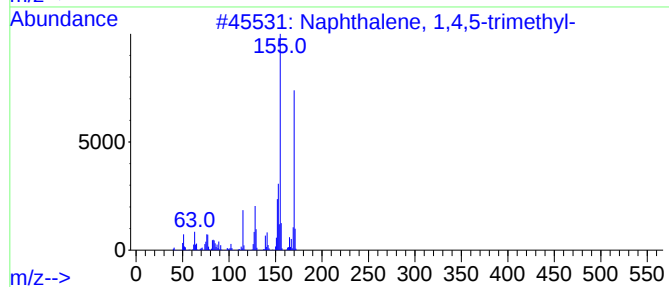
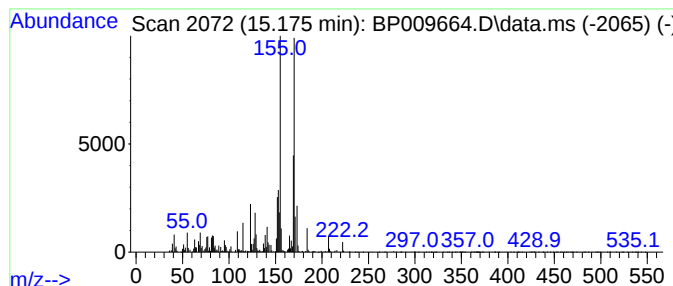
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 Naphthalene, 1,4,5-trimethyl- Concentration Rank 10

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 15.175 | 2.61 ng | 365934 | Acenaphthene-d10 | 14.375 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 1,4,5-trimethyl- | 170 | C13H14  | 002131-41-1 | 92   |
| 2    |    |   | 4,6,8-Trimethylazulene        | 170 | C13H14  | 000941-81-1 | 83   |
| 3    |    |   | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 83   |
| 4    |    |   | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 64   |
| 5    |    |   | Naphthalene, 1,4,6-trimethyl- | 170 | C13H14  | 002131-42-2 | 64   |



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Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A10015

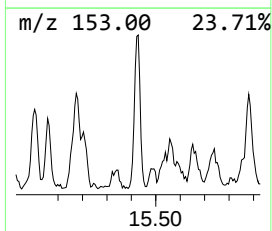
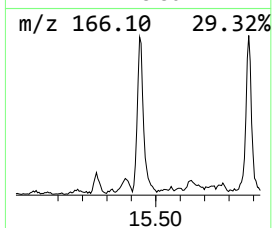
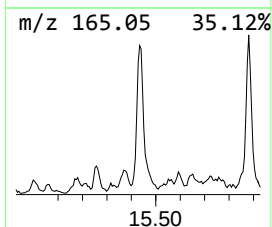
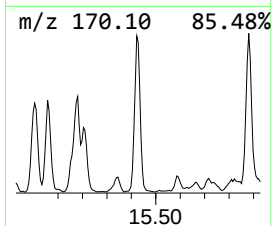
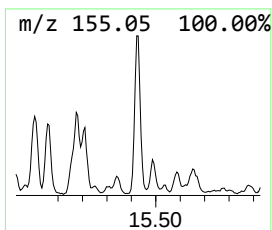
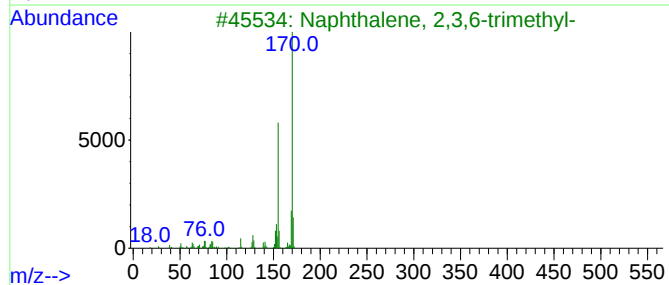
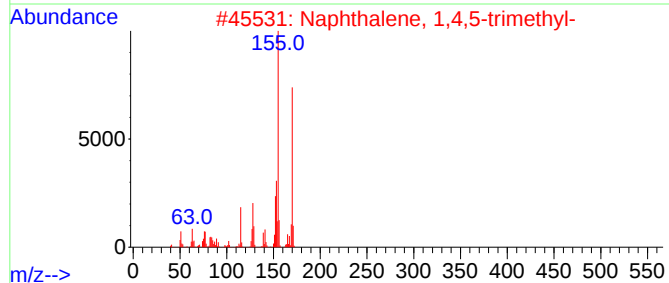
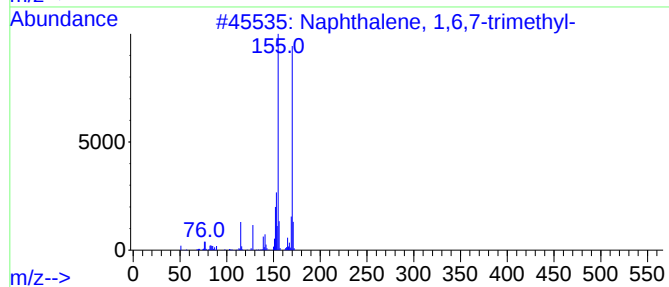
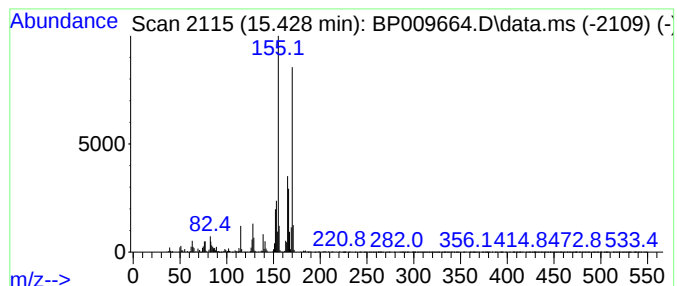
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP031722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 4 Naphthalene, 1,6,7-trimethyl- Concentration Rank 8

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 15.428 | 2.90 ng | 406262 | Acenaphthene-d10 | 14.375 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 97   |
| 2    |    |   | Naphthalene, 1,4,5-trimethyl- | 170 | C13H14  | 002131-41-1 | 97   |
| 3    |    |   | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 90   |
| 4    |    |   | 4,6,8-Trimethylazulene        | 170 | C13H14  | 000941-81-1 | 89   |
| 5    |    |   | Naphthalene, 1,4,6-trimethyl- | 170 | C13H14  | 002131-42-2 | 76   |



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ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A10015

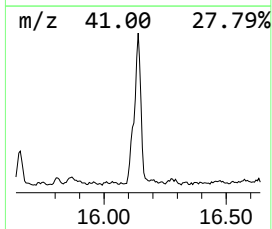
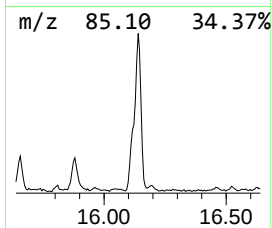
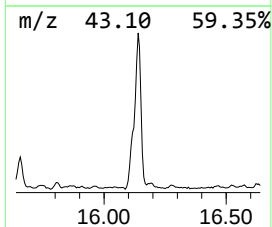
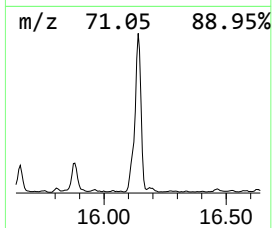
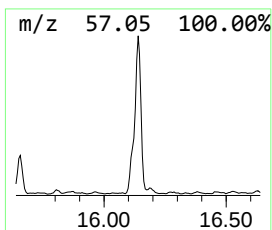
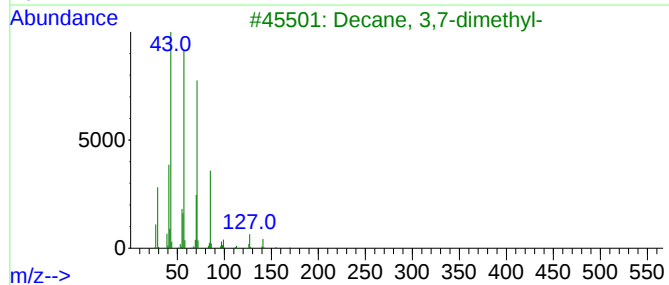
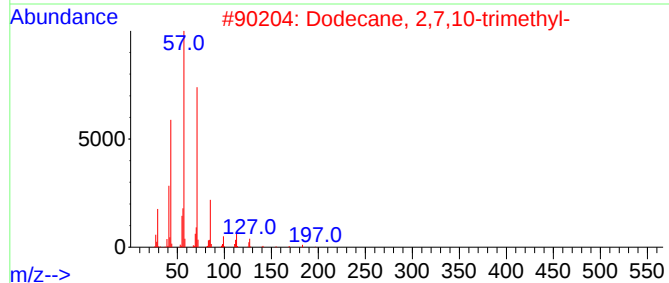
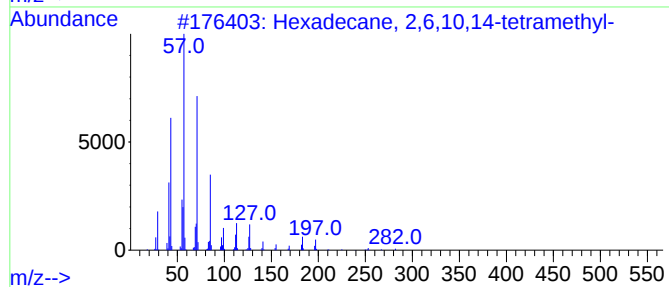
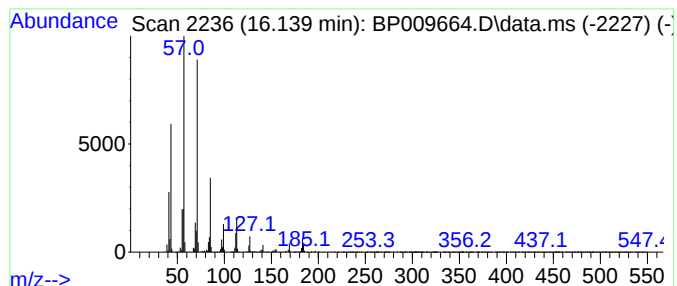
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 5 Hexadecane, 2,6,10,14-tetra... Concentration Rank 3

| R.T.   | EstConc | Area    | Relative to ISTD | R.T.   |
|--------|---------|---------|------------------|--------|
| 16.140 | 6.69 ng | 1076000 | Phenanthrene-d10 | 17.139 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Hexadecane, 2,6,10,14-tetramethyl-  | 282 | C20H42  | 000638-36-8 | 86   |
| 2    |    |   | Dodecane, 2,7,10-trimethyl-         | 212 | C15H32  | 074645-98-0 | 86   |
| 3    |    |   | Decane, 3,7-dimethyl-               | 170 | C12H26  | 017312-54-8 | 83   |
| 4    |    |   | Tridecane, 7-hexyl-                 | 268 | C19H40  | 007225-66-3 | 81   |
| 5    |    |   | Pentadecane, 2,6,10,14-tetramethyl- | 268 | C19H40  | 001921-70-6 | 80   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP040122\  
Data File : BP009664.D  
Acq On : 01 Apr 2022 18:38  
Operator : CG/JU  
Sample : N2220-01  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP031722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

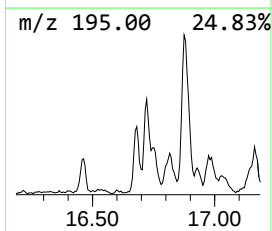
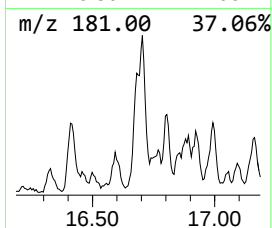
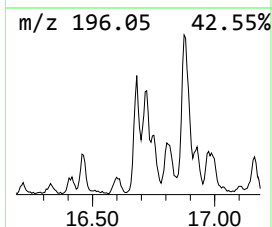
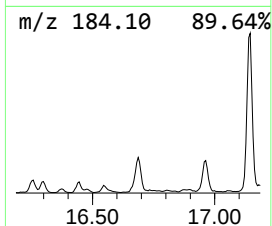
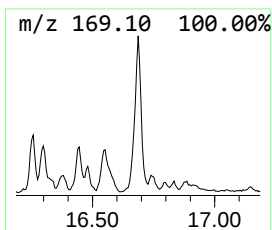
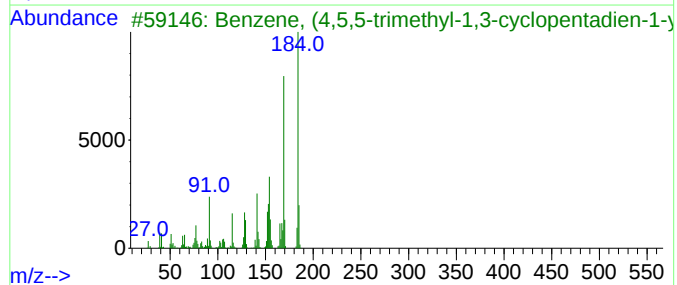
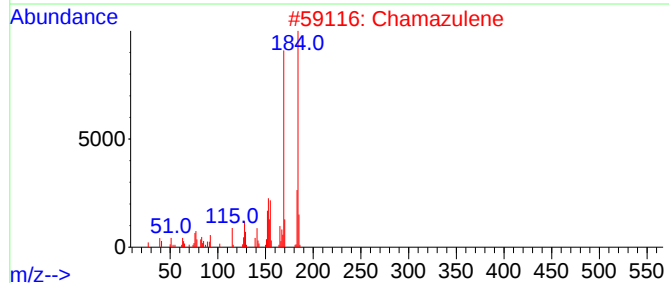
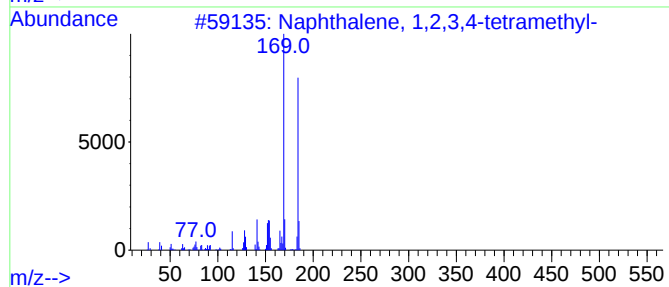
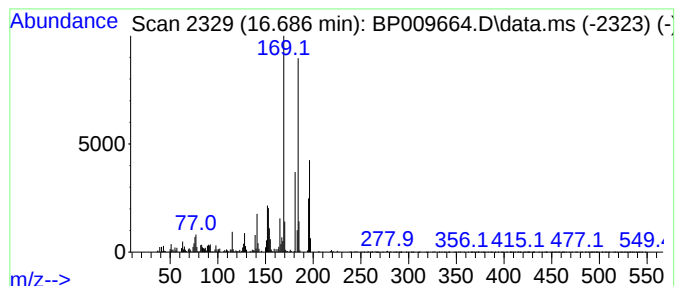
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TIC Integration Parameters: LSCINT.P

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Peak Number 6 Naphthalene, 1,2,3,4-tetram... Concentration Rank 14

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 16.686 | 2.35 ng | 378714 | Phenanthrene-d10 | 17.139 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | Naphthalene, 1,2,3,4-tetramethyl-   | 184 | C14H16  | 003031-15-0  | 91   |
| 2    |    |   | Chamazulene                         | 184 | C14H16  | 000529-05-5  | 83   |
| 3    |    |   | Benzene, (4,5,5-trimethyl-1,3-cy... | 184 | C14H16  | 033930-85-7  | 70   |
| 4    |    |   | 3-Isopropyl- 1-methylnaphthalene    | 184 | C14H16  | 1000383-71-4 | 64   |
| 5    |    |   | Naphthalene, 1-methyl-7-(1-methy... | 184 | C14H16  | 000490-65-3  | 60   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP040122\  
Data File : BP009664.D  
Acq On : 01 Apr 2022 18:38  
Operator : CG/JU  
Sample : N2220-01  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP031722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

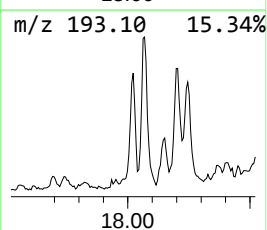
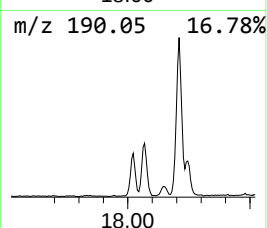
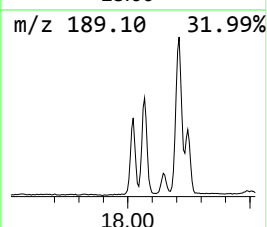
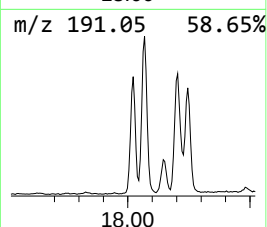
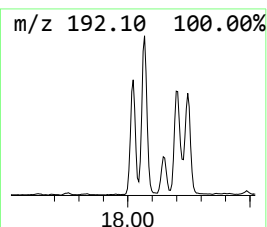
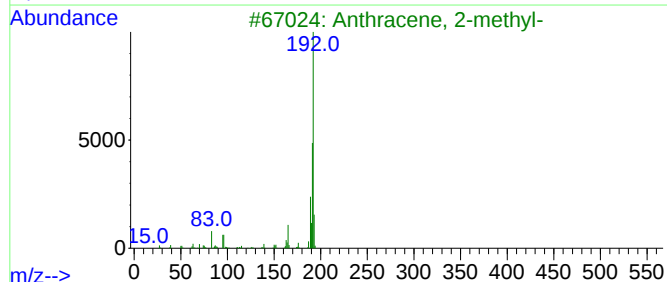
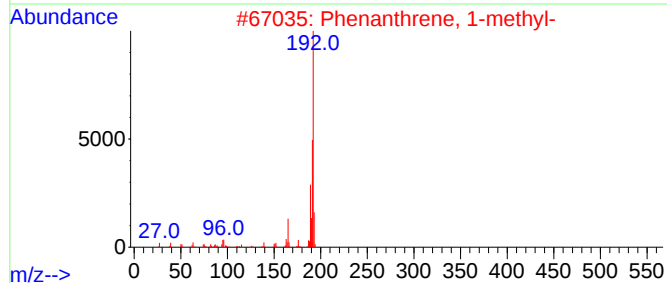
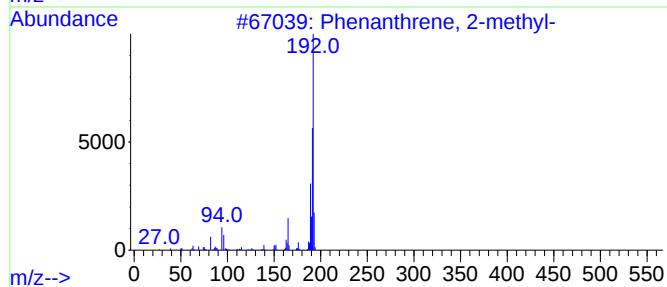
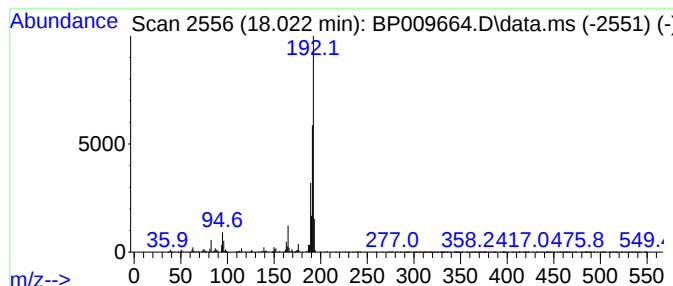
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TIC Integration Parameters: LSCINT.P

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Peak Number 7 Phenanthrene, 2-methyl- Concentration Rank 12

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 18.022 | 2.45 ng | 393412 | Phenanthrene-d10 | 17.139 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Phenanthrene, 2-methyl- | 192 | C15H12  | 002531-84-2 | 96   |
| 2    |    |   | Phenanthrene, 1-methyl- | 192 | C15H12  | 000832-69-9 | 94   |
| 3    |    |   | Anthracene, 2-methyl-   | 192 | C15H12  | 000613-12-7 | 93   |
| 4    |    |   | Phenanthrene, 4-methyl- | 192 | C15H12  | 000832-64-4 | 93   |
| 5    |    |   | Anthracene, 9-methyl-   | 192 | C15H12  | 000779-02-2 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP040122\  
Data File : BP009664.D  
Acq On : 01 Apr 2022 18:38  
Operator : CG/JU  
Sample : N2220-01  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
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Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP031722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

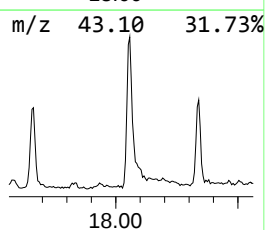
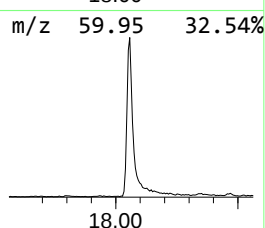
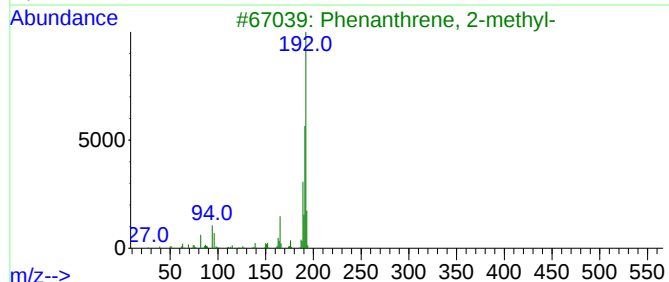
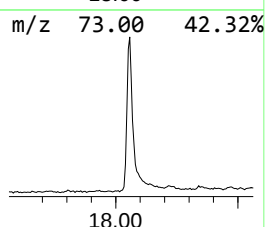
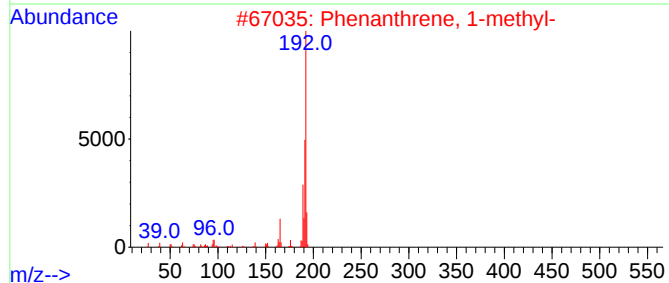
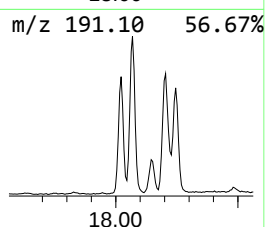
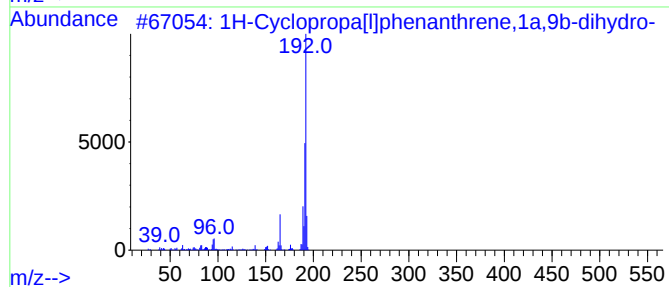
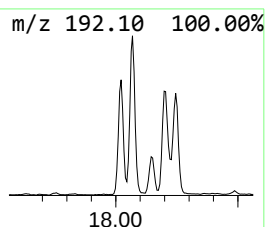
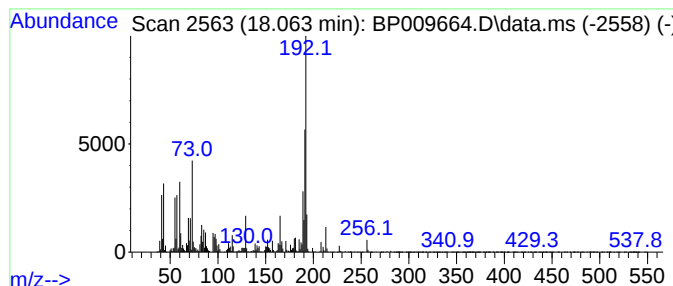
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 1H-Cyclopropa[l]phenanthren... Concentration Rank 2

| R.T.   | EstConc | Area    | Relative to ISTD | R.T.   |
|--------|---------|---------|------------------|--------|
| 18.063 | 9.01 ng | 1449480 | Phenanthrene-d10 | 17.139 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 1H-Cyclopropa[l]phenanthrene,1a,... | 192 | C15H12  | 000949-41-7 | 97   |
| 2    |    |   | Phenanthrene, 1-methyl-             | 192 | C15H12  | 000832-69-9 | 96   |
| 3    |    |   | Phenanthrene, 2-methyl-             | 192 | C15H12  | 002531-84-2 | 96   |
| 4    |    |   | Anthracene, 2-methyl-               | 192 | C15H12  | 000613-12-7 | 96   |
| 5    |    |   | 1H-Indene, 2-phenyl-                | 192 | C15H12  | 004505-48-0 | 92   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP040122\  
Data File : BP009664.D  
Acq On : 01 Apr 2022 18:38  
Operator : CG/JU  
Sample : N2220-01  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A10015

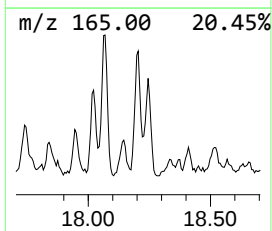
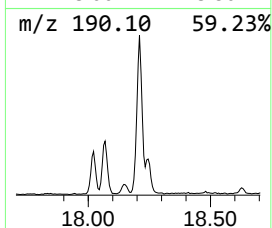
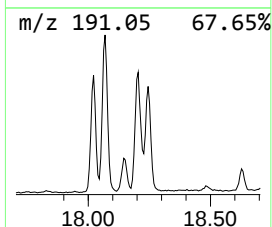
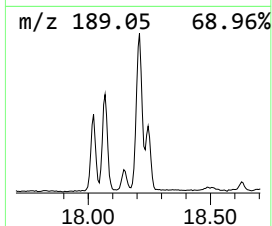
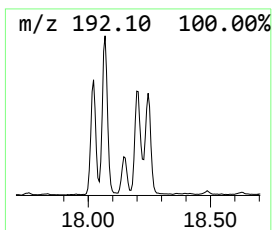
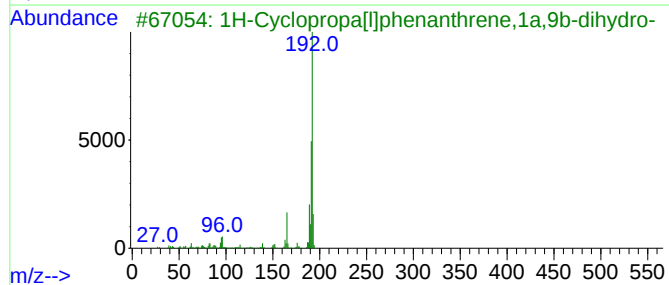
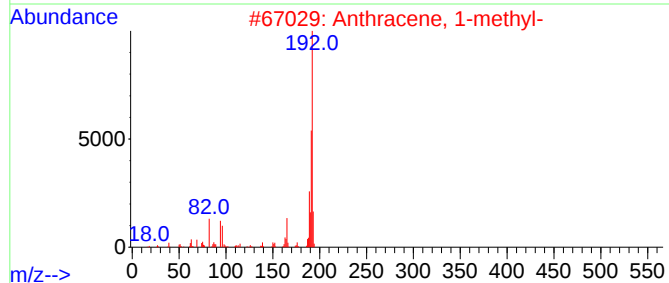
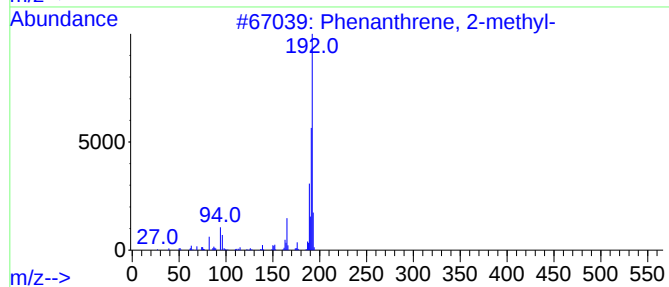
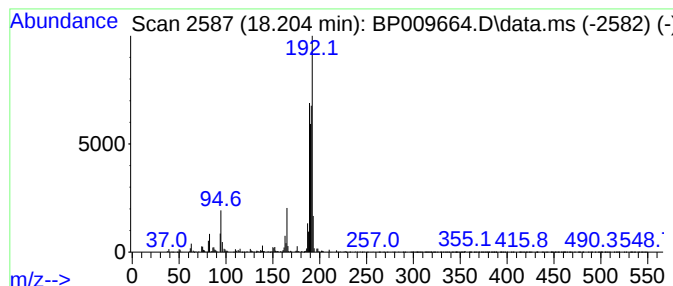
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 9 Anthracene, 1-methyl- Concentration Rank 6

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 18.204 | 4.28 ng | 688080 | Phenanthrene-d10 | 17.139 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Phenanthrene, 2-methyl-             | 192 | C15H12  | 002531-84-2 | 64   |
| 2    |    |   | Anthracene, 1-methyl-               | 192 | C15H12  | 000610-48-0 | 60   |
| 3    |    |   | 1H-Cyclopropa[1]phenanthrene,1a,... | 192 | C15H12  | 000949-41-7 | 60   |
| 4    |    |   | 1H-Indene, 2-phenyl-                | 192 | C15H12  | 004505-48-0 | 60   |
| 5    |    |   | Naphtho[2,3-b]norbornadiene         | 192 | C15H12  | 107426-38-0 | 60   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP040122\  
Data File : BP009664.D  
Acq On : 01 Apr 2022 18:38  
Operator : CG/JU  
Sample : N2220-01  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A10015

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP031722.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

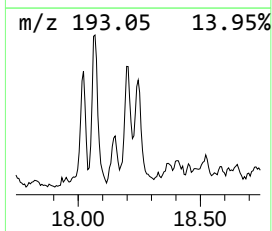
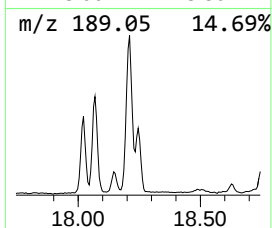
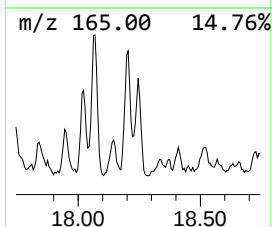
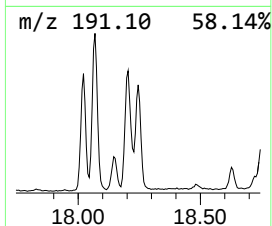
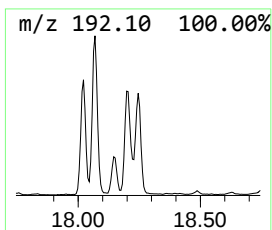
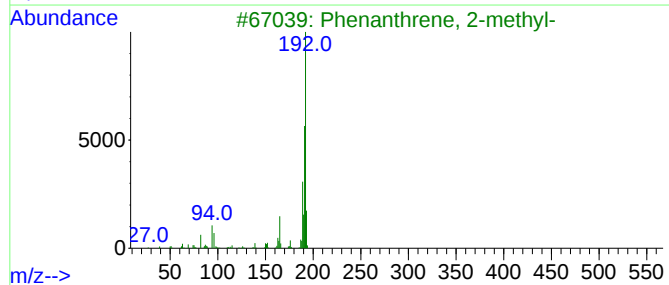
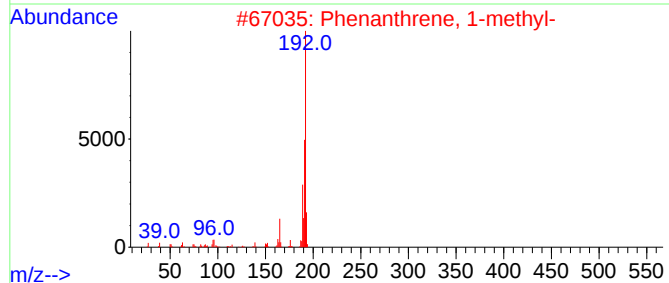
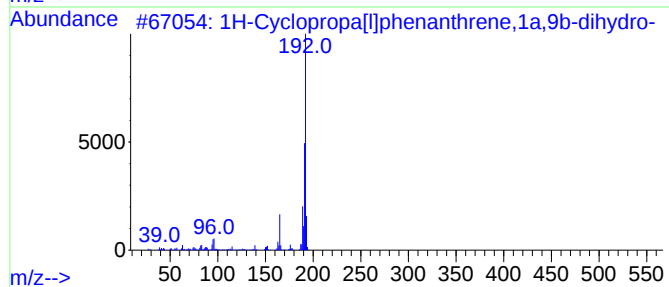
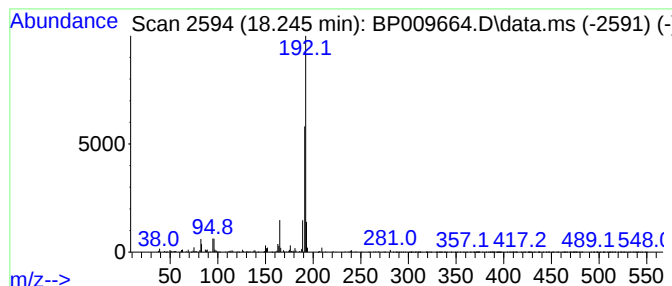
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 Phenanthrene, 1-methyl- Concentration Rank 11

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 18.245 | 2.60 ng | 418094 | Phenanthrene-d10 | 17.139 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------------|-----|---------|-------------|------|
| 1    |      | 1H-Cyclopropa[1]phenanthrene,1a,... | 192 | C15H12  | 000949-41-7 | 93   |
| 2    |      | Phenanthrene, 1-methyl-             | 192 | C15H12  | 000832-69-9 | 93   |
| 3    |      | Phenanthrene, 2-methyl-             | 192 | C15H12  | 002531-84-2 | 93   |
| 4    |      | Anthracene, 1-methyl-               | 192 | C15H12  | 000610-48-0 | 90   |
| 5    |      | Anthracene, 9-methyl-               | 192 | C15H12  | 000779-02-2 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP040122\  
Data File : BP009664.D  
Acq On : 01 Apr 2022 18:38  
Operator : CG/JU  
Sample : N2220-01  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

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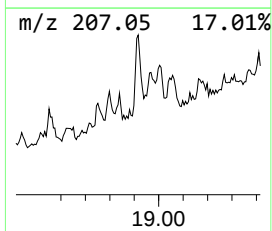
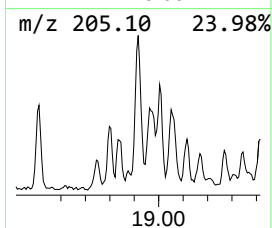
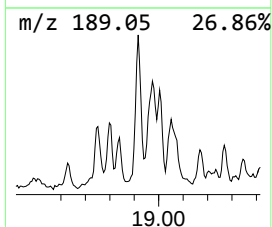
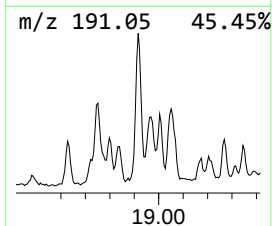
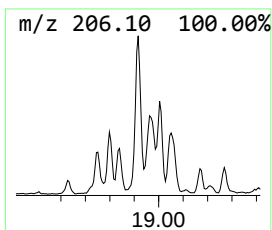
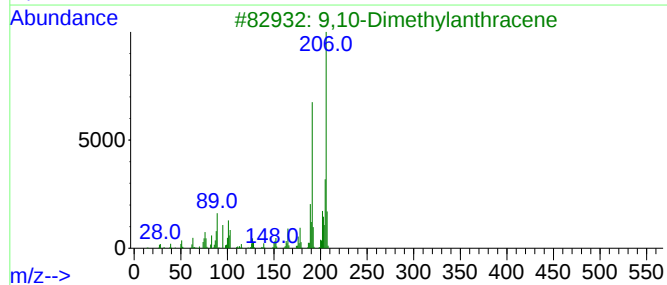
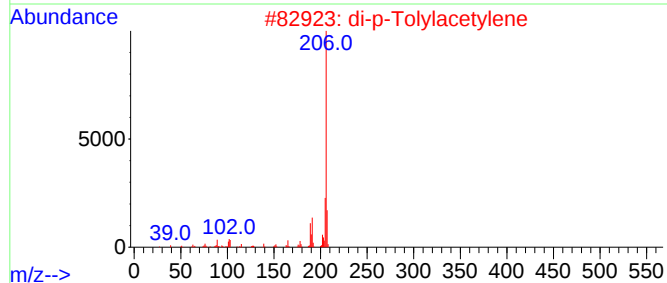
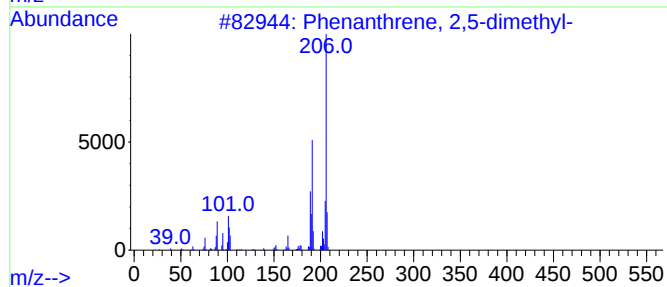
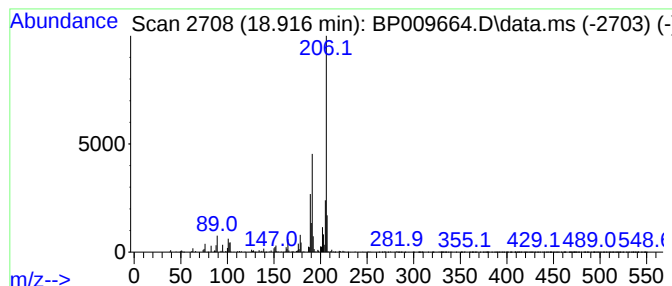
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 11 Phenanthrene, 2,5-dimethyl- Concentration Rank 13

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 18.916 | 2.41 ng | 387280 | Phenanthrene-d10 | 17.139 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Phenanthrene, 2,5-dimethyl- | 206 | C16H14  | 003674-66-6 | 93   |
| 2    |    |   | di-p-Tolylacetylene         | 206 | C16H14  | 002789-88-0 | 91   |
| 3    |    |   | 9,10-Dimethylantracene      | 206 | C16H14  | 000781-43-1 | 91   |
| 4    |    |   | Phenanthrene, 2,3-dimethyl- | 206 | C16H14  | 003674-65-5 | 90   |
| 5    |    |   | Phenanthrene, 1,7-dimethyl- | 206 | C16H14  | 000483-87-4 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP040122\  
Data File : BP009664.D  
Acq On : 01 Apr 2022 18:38  
Operator : CG/JU  
Sample : N2220-01  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A10015

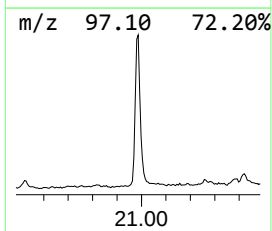
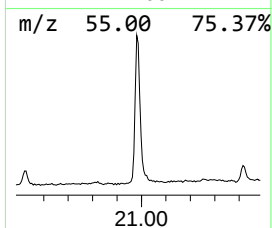
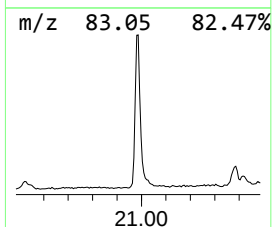
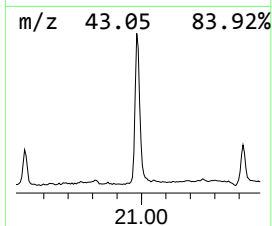
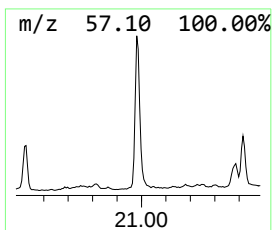
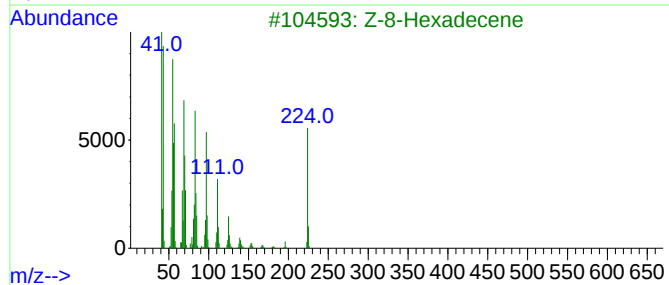
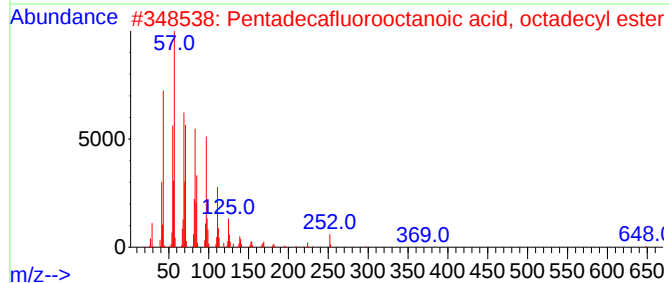
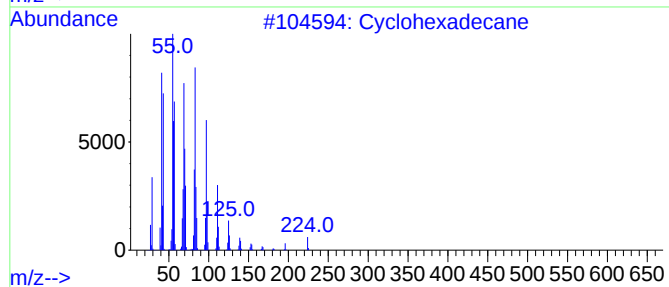
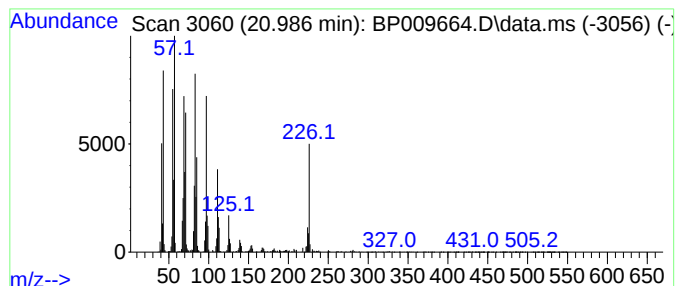
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 12 Cyclohexadecane Concentration Rank 4

| R.T.   | EstConc | Area    | Relative to ISTD | R.T.   |
|--------|---------|---------|------------------|--------|
| 20.986 | 6.23 ng | 1857830 | Chrysene-d12     | 21.263 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-------------|--------------|------|
| 1    |    |   | Cyclohexadecane                     | 224 | C16H32      | 000295-65-8  | 93   |
| 2    |    |   | Pentadecafluorooctanoic acid, oc... | 666 | C26H37F15O2 | 1000406-04-8 | 93   |
| 3    |    |   | Z-8-Hexadecene                      | 224 | C16H32      | 1000130-87-5 | 92   |
| 4    |    |   | Pentadecafluorooctanoic acid, he... | 638 | C24H33F15O2 | 1000406-04-6 | 91   |
| 5    |    |   | Heptadecyl heptafluorobutyrate      | 452 | C21H35F7O2  | 959085-66-6  | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP040122\  
Data File : BP009664.D  
Acq On : 01 Apr 2022 18:38  
Operator : CG/JU  
Sample : N2220-01  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
A10015

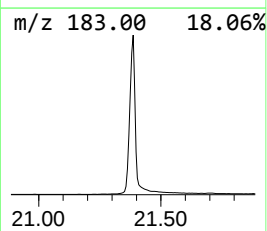
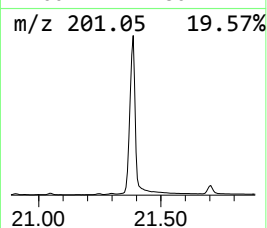
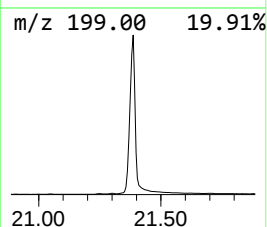
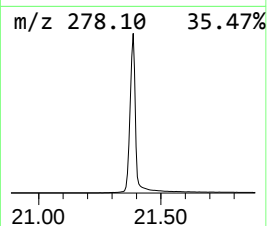
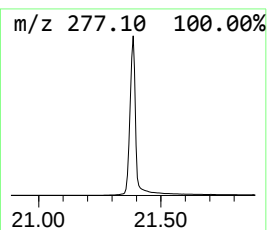
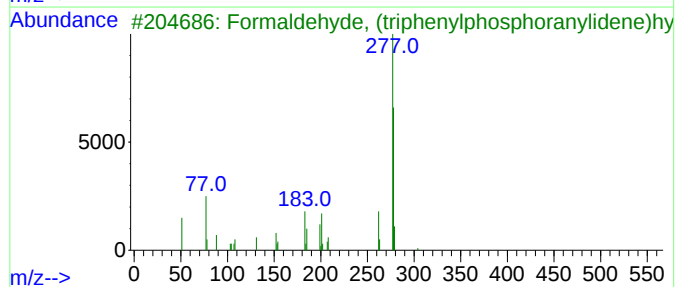
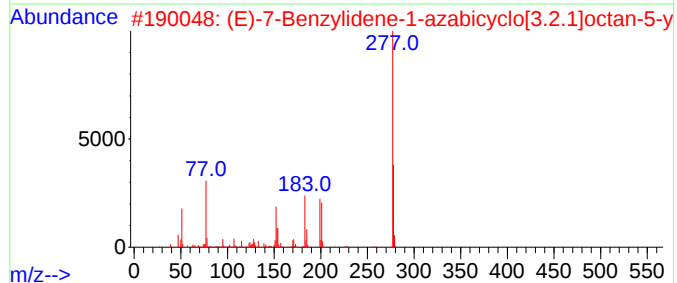
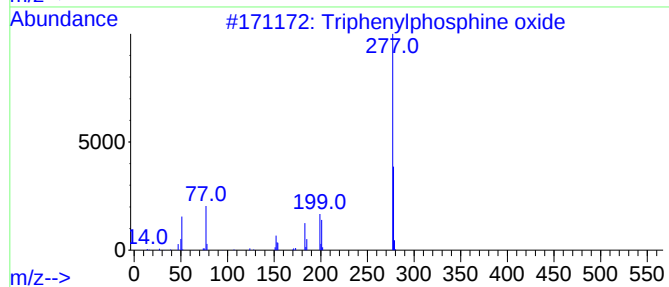
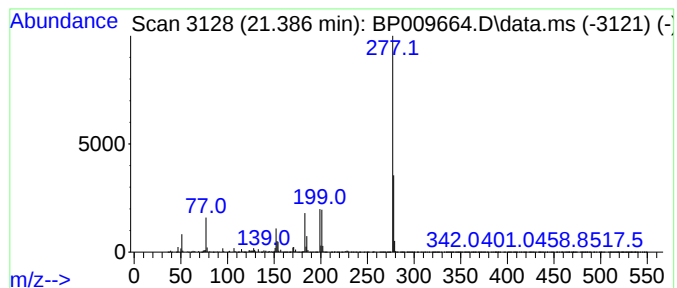
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 13 Triphenylphosphine oxide Concentration Rank 1

| R.T.   | EstConc   | Area     | Relative to ISTD | R.T.   |
|--------|-----------|----------|------------------|--------|
| 21.386 | 119.61 ng | 35685700 | Chrysene-d12     | 21.263 |

| Hit# | of | 5 | Tentative ID                         | MW  | MolForm    | CAS#         | Qual |
|------|----|---|--------------------------------------|-----|------------|--------------|------|
| 1    |    |   | Triphenylphosphine oxide             | 278 | C18H15OP   | 000791-28-6  | 94   |
| 2    |    |   | (E)-7-Benzylidene-1-azabicyclo[3...] | 293 | C15H19NO3S | 1000483-83-4 | 91   |
| 3    |    |   | Formaldehyde, (triphenylphosphor...  | 304 | C19H17N2P  | 015990-54-2  | 37   |
| 4    |    |   | 5-Methyl-8-phenylfuro[3',2':5,6]...  | 277 | C16H11N3O2 | 1000460-07-9 | 25   |
| 5    |    |   | Vanadium, cycloheptatrienyl-(pen...  | 277 | C17H22V    | 1000155-20-5 | 9    |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP040122\  
Data File : BP009664.D  
Acq On : 01 Apr 2022 18:38  
Operator : CG/JU  
Sample : N2220-01  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

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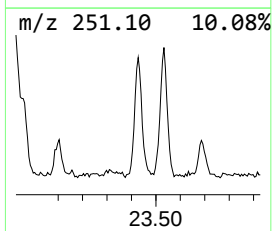
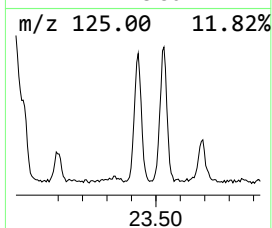
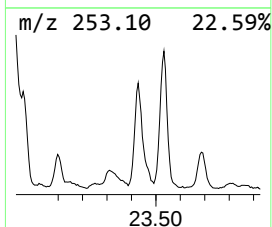
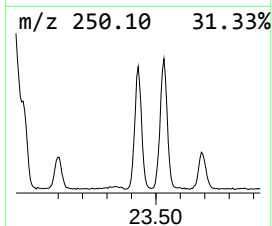
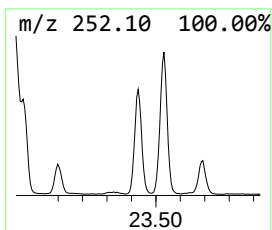
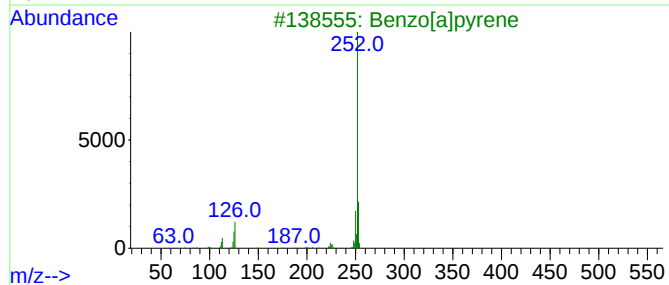
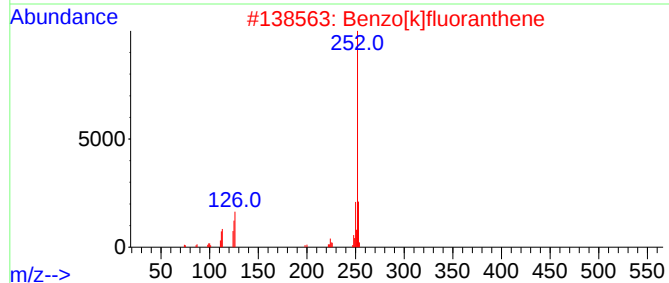
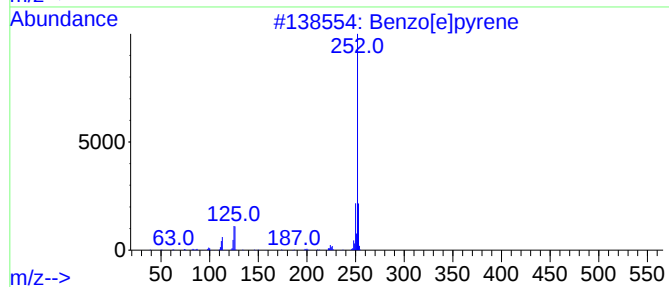
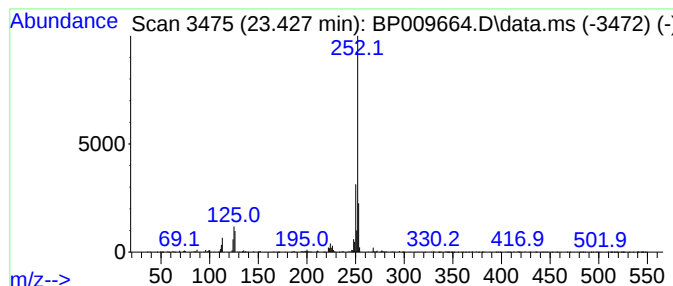
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 14 Benzo[e]pyrene Concentration Rank 5

| R.T.   | EstConc | Area    | Relative to ISTD | R.T.   |
|--------|---------|---------|------------------|--------|
| 23.427 | 5.56 ng | 1053500 | Perylene-d12     | 23.639 |

| Hit# | of | 5 | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------|-----|---------|-------------|------|
| 1    |    |   | Benzo[e]pyrene       | 252 | C20H12  | 000192-97-2 | 97   |
| 2    |    |   | Benzo[k]fluoranthene | 252 | C20H12  | 000207-08-9 | 96   |
| 3    |    |   | Benzo[a]pyrene       | 252 | C20H12  | 000050-32-8 | 95   |
| 4    |    |   | Benzo[j]fluoranthene | 252 | C20H12  | 000205-82-3 | 95   |
| 5    |    |   | Benzo[b]fluoranthene | 252 | C20H12  | 000205-99-2 | 95   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP040122\  
 Data File : BP009664.D  
 Acq On : 01 Apr 2022 18:38  
 Operator : CG/JU  
 Sample : N2220-01  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 A10015

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP031722.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 |
| Naphthalene, 2,... | 13.698 | 3.1     | ng    | 441229   | 3                     | 14.375 | 2801460 | 20.0 |
| Naphthalene, 2,... | 13.751 | 2.7     | ng    | 378809   | 3                     | 14.375 | 2801460 | 20.0 |
| Naphthalene, 1,... | 15.175 | 2.6     | ng    | 365934   | 3                     | 14.375 | 2801460 | 20.0 |
| Naphthalene, 1,... | 15.428 | 2.9     | ng    | 406262   | 3                     | 14.375 | 2801460 | 20.0 |
| Hexadecane, 2,6... | 16.140 | 6.7     | ng    | 1076000  | 4                     | 17.139 | 3217670 | 20.0 |
| Naphthalene, 1,... | 16.686 | 2.4     | ng    | 378714   | 4                     | 17.139 | 3217670 | 20.0 |
| Phenanthrene, 2... | 18.022 | 2.5     | ng    | 393412   | 4                     | 17.139 | 3217670 | 20.0 |
| 1H-Cyclopropa[1... | 18.063 | 9.0     | ng    | 1449480  | 4                     | 17.139 | 3217670 | 20.0 |
| Anthracene, 1-m... | 18.204 | 4.3     | ng    | 688080   | 4                     | 17.139 | 3217670 | 20.0 |
| Phenanthrene, 1... | 18.245 | 2.6     | ng    | 418094   | 4                     | 17.139 | 3217670 | 20.0 |
| Phenanthrene, 2... | 18.916 | 2.4     | ng    | 387280   | 4                     | 17.139 | 3217670 | 20.0 |
| Cyclohexadecane    | 20.986 | 6.2     | ng    | 1857830  | 5                     | 21.263 | 5966990 | 20.0 |
| Triphenylphosph... | 21.386 | 119.6   | ng    | 35685700 | 5                     | 21.263 | 5966990 | 20.0 |
| Benzo[e]pyrene     | 23.427 | 5.6     | ng    | 1053500  | 6                     | 23.639 | 3788730 | 20.0 |