

Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP081125\  
 Data File : BP025355.D  
 Acq On : 11 Aug 2025 19:34  
 Operator : CG/JU  
 Sample : Q2795-03  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_P  
 ClientSampleId :  
 COMP-3

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-BP080525.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP025355.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         | 2.967       | 3             | 5           | 13           | rVB2     | 213472         | 362934        | 3.61%           | 0.457%        |
| 2         | 3.049       | 13            | 19          | 36           | rVB      | 2065034        | 2844445       | 28.32%          | 3.579%        |
| 3         | 4.949       | 335           | 342         | 349          | rBV      | 445073         | 621788        | 6.19%           | 0.782%        |
| 4         | 5.414       | 413           | 421         | 431          | rBV      | 4292614        | 6298344       | 62.71%          | 7.925%        |
| 5         | 6.972       | 678           | 686         | 694          | rBV      | 3908334        | 6252350       | 62.25%          | 7.867%        |
| 6         | 7.796       | 818           | 826         | 834          | rBV      | 1085143        | 1743931       | 17.36%          | 2.194%        |
| 7         | 8.937       | 1011          | 1020        | 1029         | rBV      | 2700737        | 4363296       | 43.44%          | 5.490%        |
| 8         | 10.572      | 1288          | 1298        | 1305         | rBV      | 1272940        | 2292589       | 22.83%          | 2.885%        |
| 9         | 13.037      | 1709          | 1717        | 1727         | rBV      | 5622676        | 8761964       | 87.24%          | 11.025%       |
| 10        | 14.419      | 1943          | 1952        | 1959         | rBV      | 1570402        | 2558441       | 25.47%          | 3.219%        |
| 11        | 14.490      | 1959          | 1964        | 1971         | rVB      | 102907         | 173273        | 1.73%           | 0.218%        |
| 12        | 14.825      | 2015          | 2021        | 2030         | rBV      | 65109          | 125193        | 1.25%           | 0.158%        |
| 13        | 15.484      | 2126          | 2133        | 2140         | rBV2     | 102664         | 161961        | 1.61%           | 0.204%        |
| 14        | 15.795      | 2179          | 2186        | 2196         | rVB      | 554812         | 872446        | 8.69%           | 1.098%        |
| 15        | 15.931      | 2202          | 2209        | 2218         | rBV      | 3762493        | 5773543       | 57.48%          | 7.265%        |
| 16        | 17.219      | 2422          | 2428        | 2432         | rBV2     | 1734425        | 2573735       | 25.63%          | 3.238%        |
| 17        | 17.260      | 2432          | 2435        | 2441         | rVV      | 997978         | 1385620       | 13.80%          | 1.744%        |
| 18        | 17.354      | 2441          | 2451        | 2457         | rVV      | 245740         | 446284        | 4.44%           | 0.562%        |
| 19        | 17.625      | 2489          | 2497        | 2503         | rBV2     | 107106         | 206795        | 2.06%           | 0.260%        |
| 20        | 17.931      | 2544          | 2549        | 2554         | rBV      | 172871         | 221874        | 2.21%           | 0.279%        |
| 21        | 18.119      | 2577          | 2581        | 2584         | rBV      | 121933         | 166131        | 1.65%           | 0.209%        |
| 22        | 18.166      | 2584          | 2589        | 2597         | rVV2     | 1323454        | 1983629       | 19.75%          | 2.496%        |
| 23        | 18.260      | 2602          | 2605        | 2610         | rVV      | 111273         | 171259        | 1.71%           | 0.215%        |
| 24        | 18.325      | 2611          | 2616        | 2620         | rVV3     | 201325         | 371156        | 3.70%           | 0.467%        |
| 25        | 18.360      | 2620          | 2622        | 2629         | rVB      | 110546         | 153824        | 1.53%           | 0.194%        |
| 26        | 18.642      | 2664          | 2670        | 2673         | rBV      | 96119          | 145385        | 1.45%           | 0.183%        |
| 27        | 19.060      | 2734          | 2741        | 2748         | rBV3     | 205092         | 367911        | 3.66%           | 0.463%        |
| 28        | 19.348      | 2784          | 2790        | 2799         | rBV      | 1310209        | 1830793       | 18.23%          | 2.304%        |
| 29        | 19.489      | 2810          | 2814        | 2820         | rBV      | 599313         | 735806        | 7.33%           | 0.926%        |
| 30        | 19.630      | 2834          | 2838        | 2843         | rVB      | 136071         | 191574        | 1.91%           | 0.241%        |
| 31        | 19.689      | 2843          | 2848        | 2850         | rBV2     | 256015         | 400775        | 3.99%           | 0.504%        |
| 32        | 19.719      | 2850          | 2853        | 2860         | rVB      | 973626         | 1333029       | 13.27%          | 1.677%        |
| 33        | 19.936      | 2882          | 2890        | 2901         | rVB      | 7748731        | 10043794      | 100.00%         | 12.638%       |
| 34        | 20.054      | 2905          | 2910        | 2915         | rVV4     | 83690          | 190038        | 1.89%           | 0.239%        |
| 35        | 20.242      | 2939          | 2942        | 2947         | rVB      | 272063         | 348494        | 3.47%           | 0.439%        |
| 36        | 20.336      | 2955          | 2958        | 2962         | rBV2     | 135232         | 146060        | 1.45%           | 0.184%        |

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Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
COMP-3

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-BP080525.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 37 | 20.395 | 2965 | 2968 | 2972 | rVB3 | 150467  | 186159  | 1.85%  | 0.234% |
| 38 | 21.330 | 3123 | 3127 | 3142 | rVB4 | 807666  | 1563969 | 15.57% | 1.968% |
| 39 | 21.660 | 3175 | 3183 | 3188 | rBV3 | 1938691 | 4383504 | 43.64% | 5.516% |
| 40 | 21.707 | 3188 | 3191 | 3202 | rVB  | 575264  | 937102  | 9.33%  | 1.179% |
| 41 | 23.295 | 3455 | 3461 | 3471 | rBV  | 744771  | 1625121 | 16.18% | 2.045% |
| 42 | 23.960 | 3569 | 3574 | 3582 | rBV2 | 233722  | 587524  | 5.85%  | 0.739% |
| 43 | 25.042 | 3749 | 3758 | 3767 | rBV  | 1410589 | 3569333 | 35.54% | 4.491% |

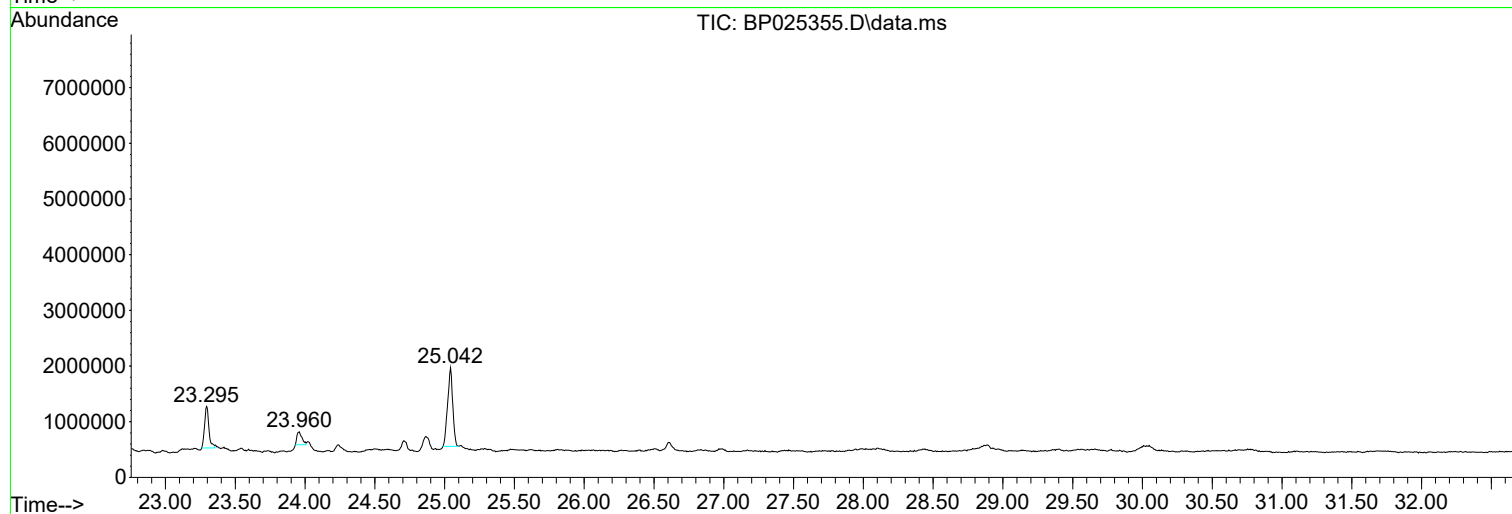
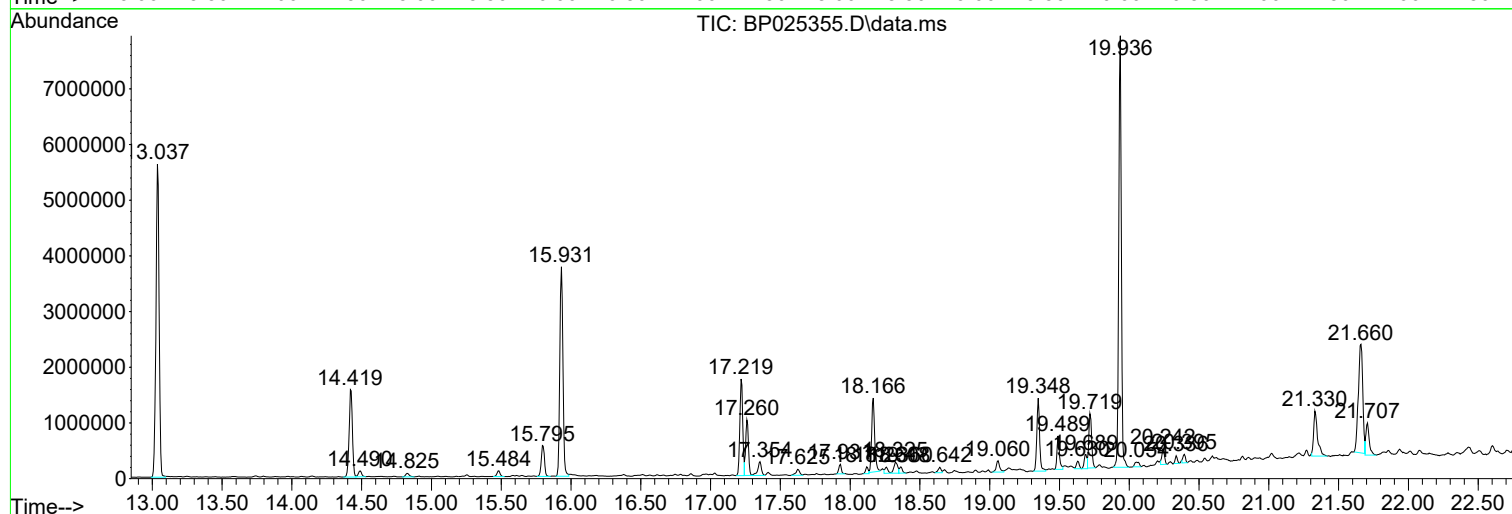
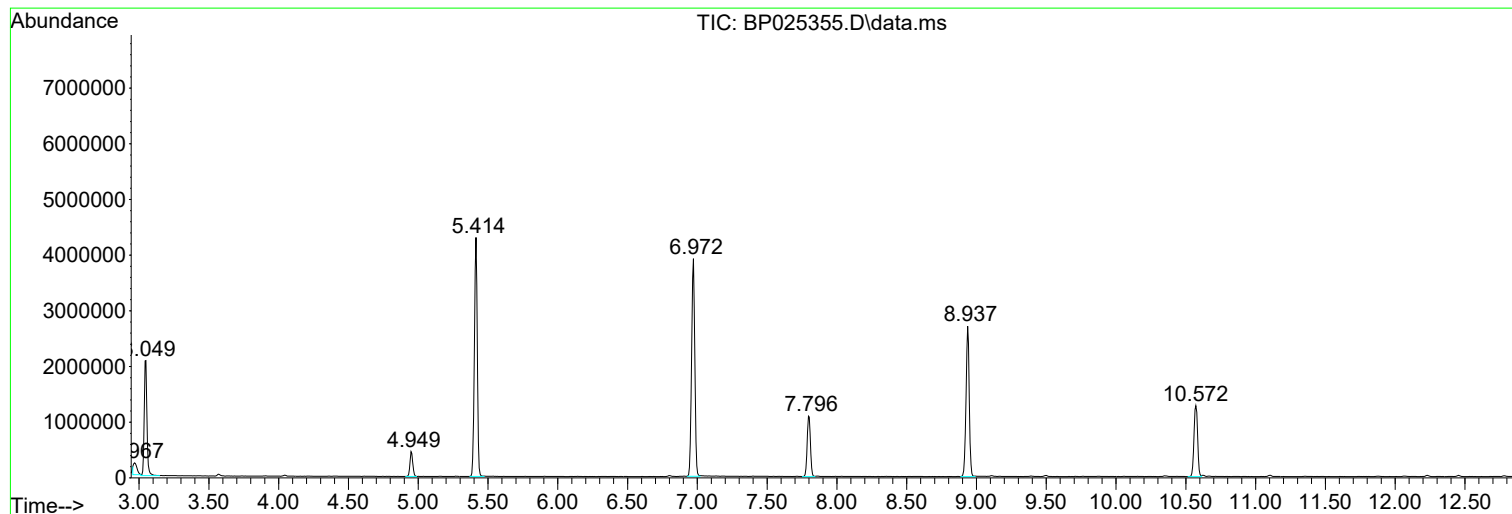
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Data File : BP025355.D  
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Sample : Q2795-03  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
COMP-3

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP080525.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP081125\  
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Operator : CG/JU  
Sample : Q2795-03  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
COMP-3

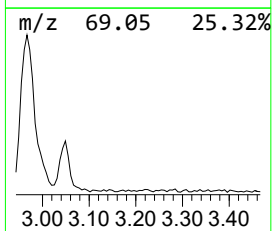
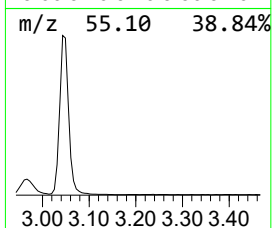
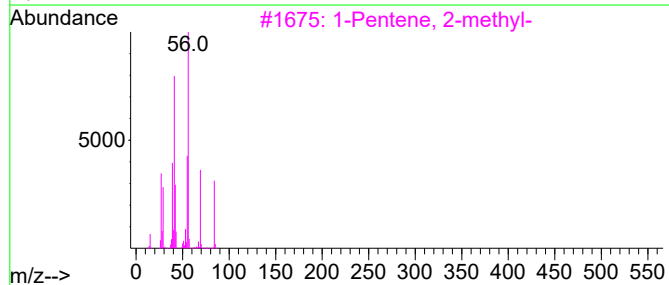
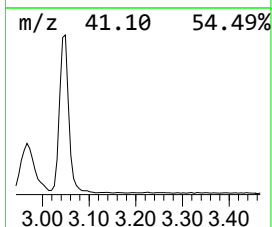
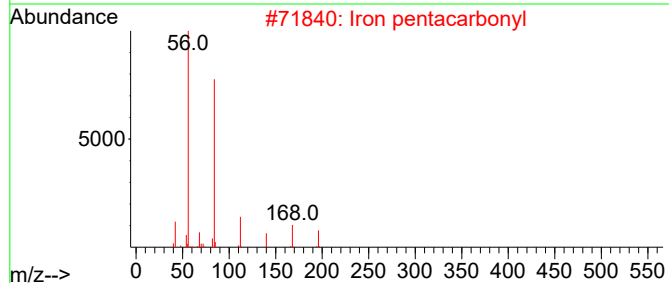
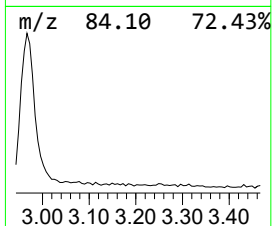
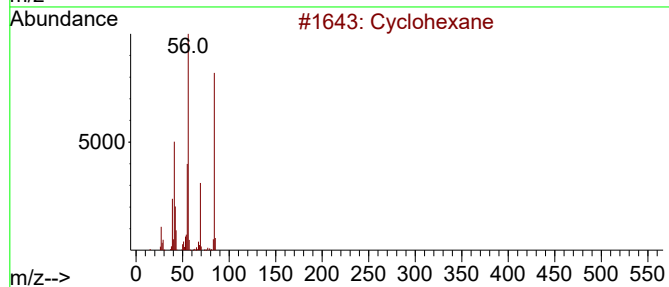
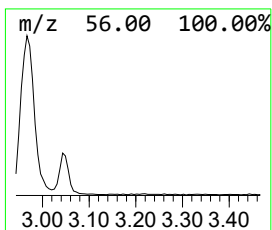
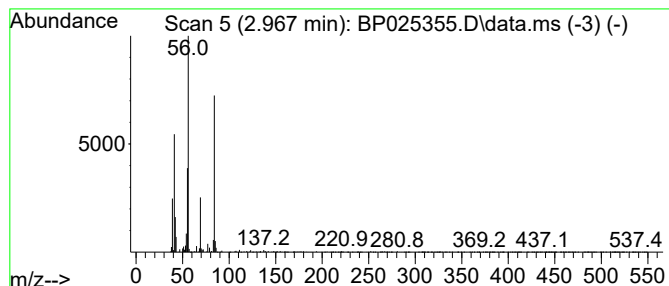
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP080525.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 1 Cyclohexane Concentration Rank 7

| R.T.  | EstConc | Area   | Relative to ISTD       | R.T.  |
|-------|---------|--------|------------------------|-------|
| 2.967 | 4.16 ng | 362934 | 1,4-Dichlorobenzene-d4 | 7.802 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane                         | 84  | C6H12   | 000110-82-7 | 87   |
| 2    |    |   | Iron pentacarbonyl                  | 196 | C5Fe05  | 013463-40-6 | 72   |
| 3    |    |   | 1-Pentene, 2-methyl-                | 84  | C6H12   | 000763-29-1 | 47   |
| 4    |    |   | 1,4-Dioxaspiro[2.4]heptan-5-one,... | 142 | C7H10O3 | 126091-79-0 | 47   |
| 5    |    |   | Cyclopentanone, 2-ethyl-            | 112 | C7H12O  | 004971-18-0 | 40   |



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Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
COMP-3

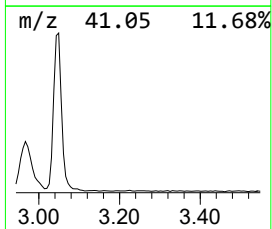
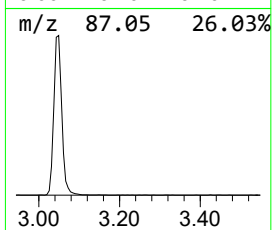
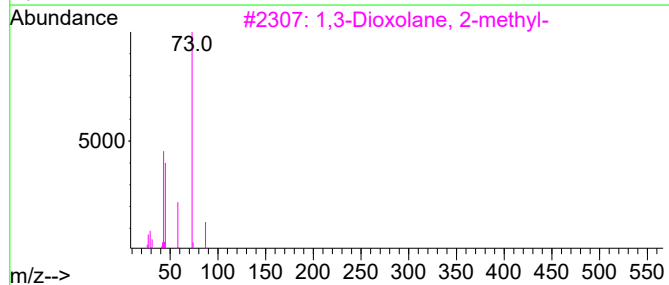
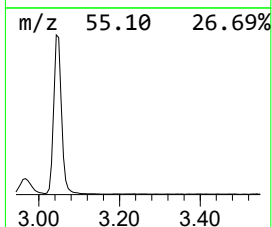
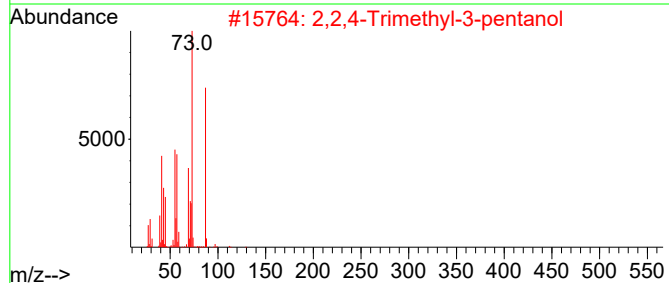
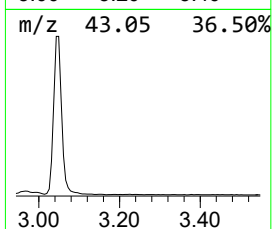
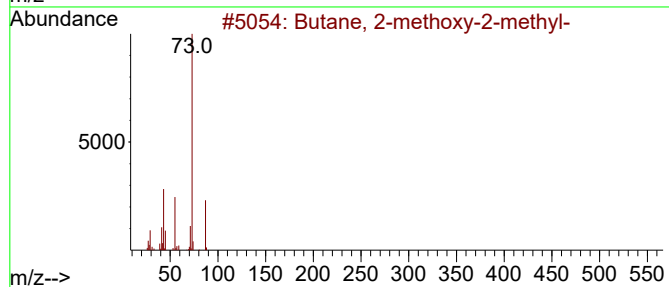
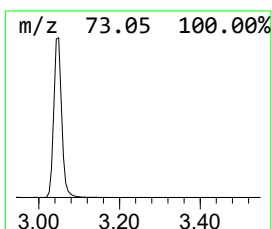
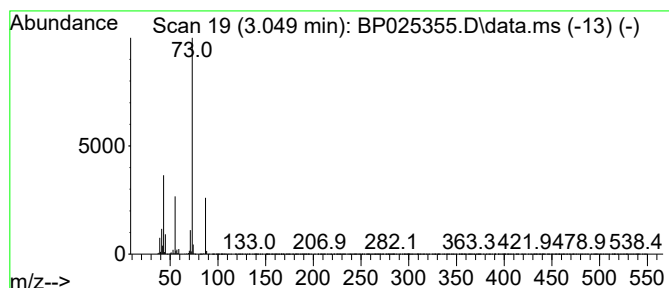
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP080525.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

| R.T.  | EstConc  | Area    | Relative to ISTD       | R.T.  |
|-------|----------|---------|------------------------|-------|
| 3.049 | 32.62 ng | 2844450 | 1,4-Dichlorobenzene-d4 | 7.802 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Butane, 2-methoxy-2-methyl- | 102 | C6H14O  | 000994-05-8 | 78   |
| 2    |    |   | 2,2,4-Trimethyl-3-pentanol  | 130 | C8H18O  | 005162-48-1 | 40   |
| 3    |    |   | 1,3-Dioxolane, 2-methyl-    | 88  | C4H8O2  | 000497-26-7 | 25   |
| 4    |    |   | Pentane, 3-methoxy-         | 102 | C6H14O  | 036839-67-5 | 17   |
| 5    |    |   | Silane, tetramethyl-        | 88  | C4H12Si | 000075-76-3 | 12   |



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ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
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COMP-3

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP080525.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

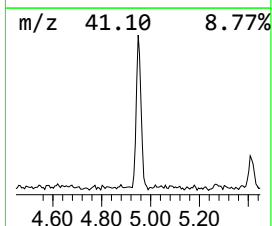
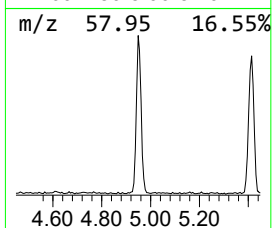
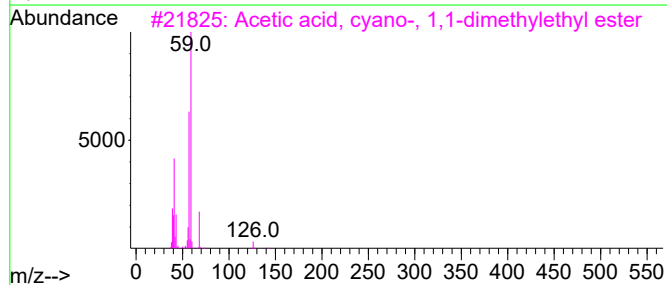
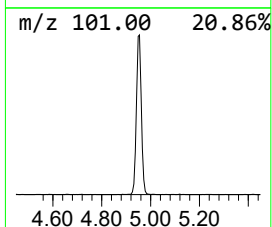
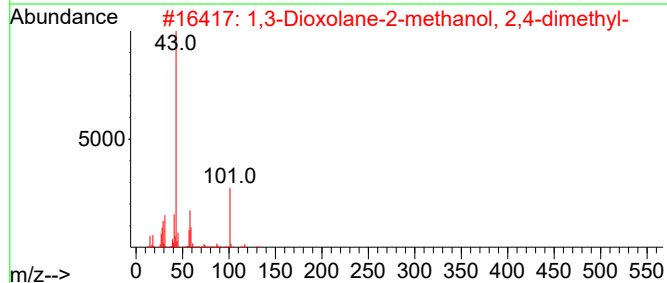
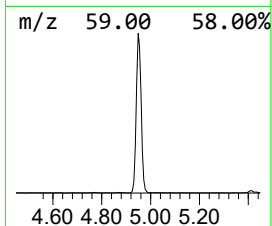
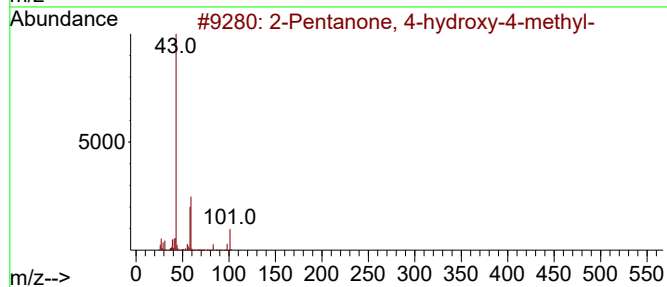
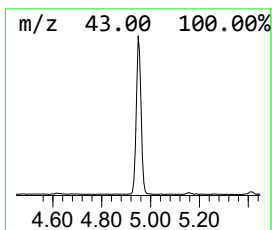
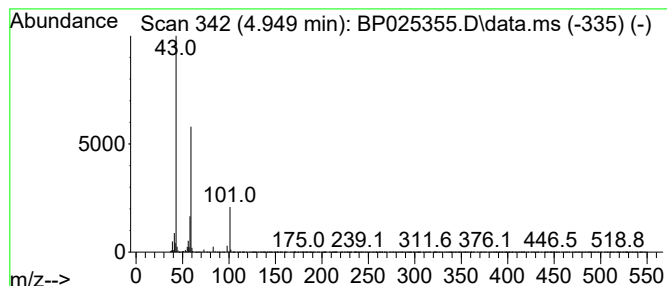
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

| R.T.  | EstConc | Area   | Relative to ISTD       | R.T.  |
|-------|---------|--------|------------------------|-------|
| 4.949 | 7.13 ng | 621788 | 1,4-Dichlorobenzene-d4 | 7.802 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2  | 000123-42-2 | 64   |
| 2    |    |   | 1,3-Dioxolane-2-methanol, 2,4-di... | 132 | C6H12O3  | 053951-43-2 | 28   |
| 3    |    |   | Acetic acid, cyano-, 1,1-dimethy... | 141 | C7H11NO2 | 001116-98-9 | 23   |
| 4    |    |   | 1-Propen-2-ol, acetate              | 100 | C5H8O2   | 000108-22-5 | 10   |
| 5    |    |   | Morpholine, 4-methyl-               | 101 | C5H11NO  | 000109-02-4 | 9    |



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COMP-3

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP080525.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

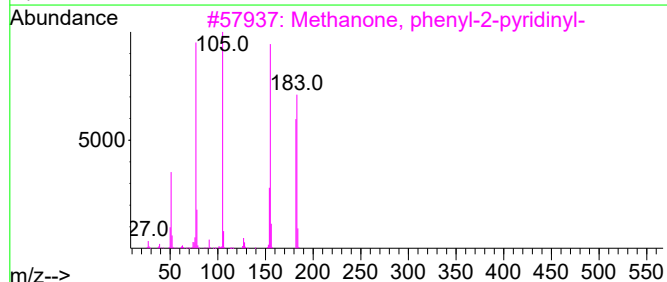
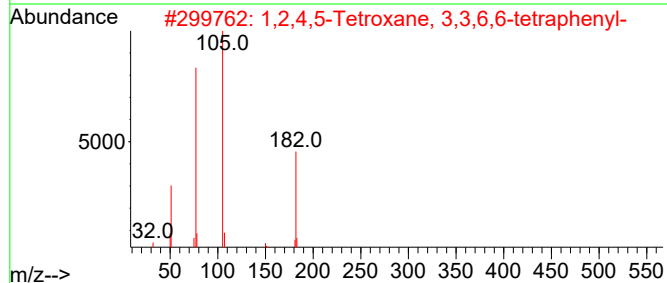
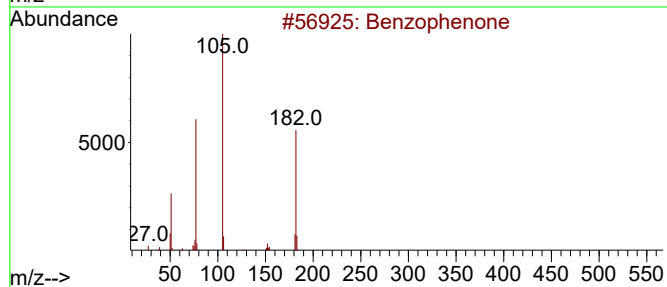
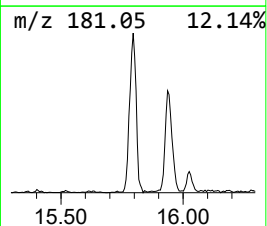
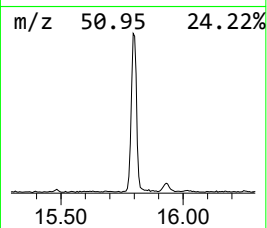
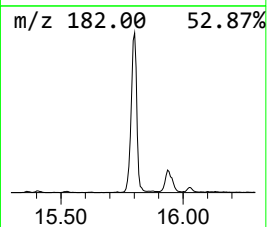
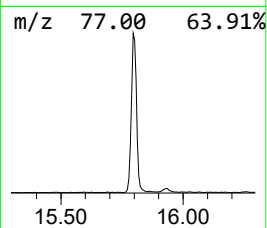
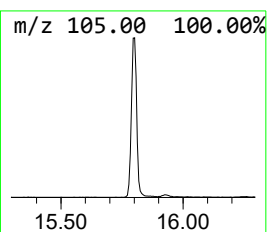
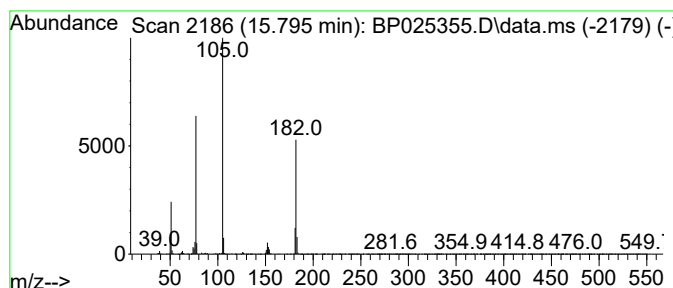
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TIC Integration Parameters: LSCINT.P

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Peak Number 4 Benzophenone Concentration Rank 6

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 15.795 | 6.82 ng | 872446 | Acenaphthene-d10 | 14.419 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Benzophenone                        | 182 | C13H10O  | 000119-61-9 | 96   |
| 2    |    |   | 1,2,4,5-Tetroxane, 3,3,6,6-tetra... | 396 | C26H20O4 | 016204-36-7 | 64   |
| 3    |    |   | Methanone, phenyl-2-pyridinyl-      | 183 | C12H9NO  | 000091-02-1 | 56   |
| 4    |    |   | 1,2,4-Trioxolane, 3,3,5-triphenyl-  | 304 | C20H16O3 | 023246-12-0 | 45   |
| 5    |    |   | Azobenzene                          | 182 | C12H10N2 | 000103-33-3 | 42   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP081125\  
Data File : BP025355.D  
Acq On : 11 Aug 2025 19:34  
Operator : CG/JU  
Sample : Q2795-03  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
COMP-3

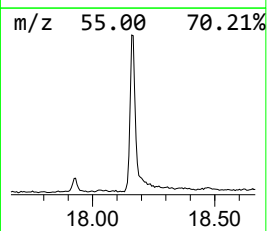
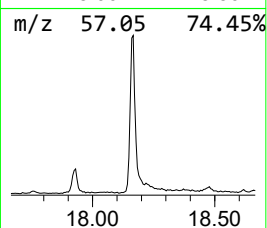
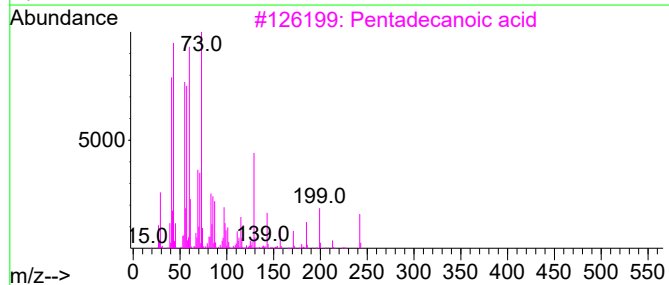
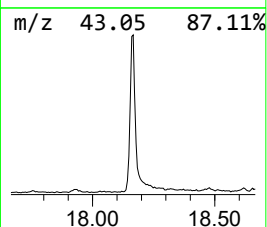
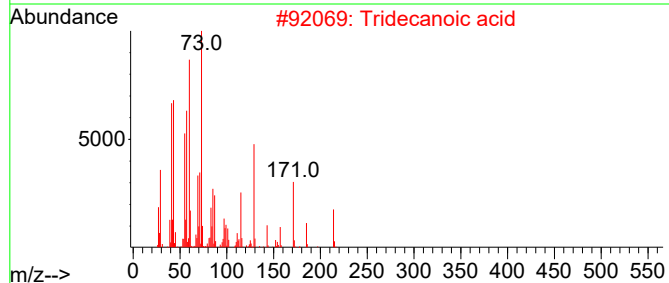
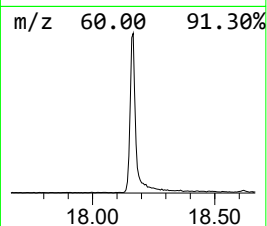
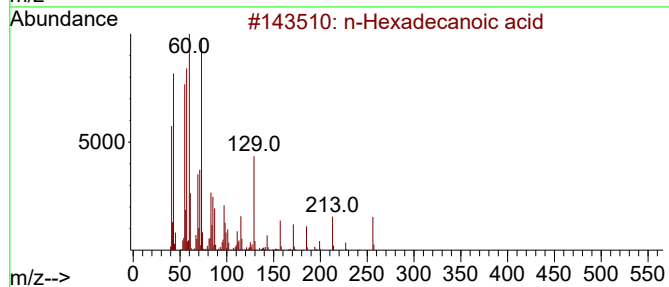
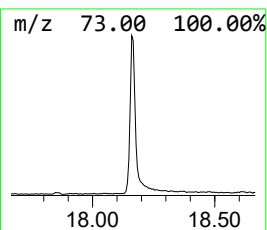
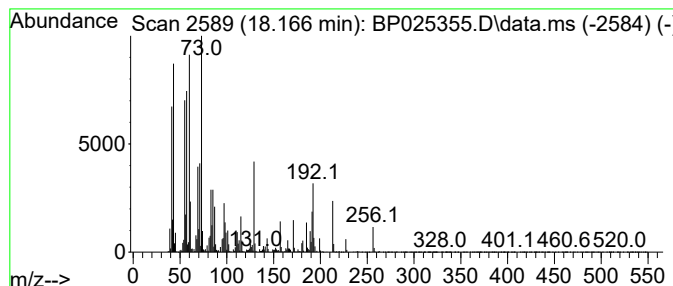
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 5 n-Hexadecanoic acid Concentration Rank 2

| R.T.      | EstConc             | Area    | Relative to ISTD |          | R.T.        |      |
|-----------|---------------------|---------|------------------|----------|-------------|------|
| 18.166    | 15.41 ng            | 1983630 | Phenanthrene-d10 |          | 17.219      |      |
| Hit# of 5 | Tentative ID        |         | MW               | MolForm  | CAS#        | Qual |
| 1         | n-Hexadecanoic acid |         | 256              | C16H32O2 | 000057-10-3 | 99   |
| 2         | Tridecanoic acid    |         | 214              | C13H26O2 | 000638-53-9 | 94   |
| 3         | Pentadecanoic acid  |         | 242              | C15H30O2 | 001002-84-2 | 91   |
| 4         | Tetradecanoic acid  |         | 228              | C14H28O2 | 000544-63-8 | 90   |
| 5         | n-Decanoic acid     |         | 172              | C10H20O2 | 000334-48-5 | 70   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP081125\  
Data File : BP025355.D  
Acq On : 11 Aug 2025 19:34  
Operator : CG/JU  
Sample : Q2795-03  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
COMP-3

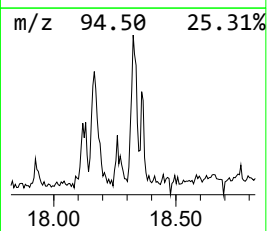
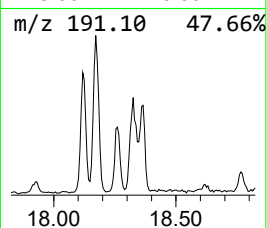
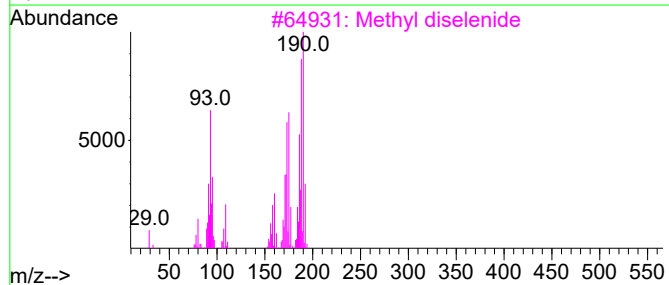
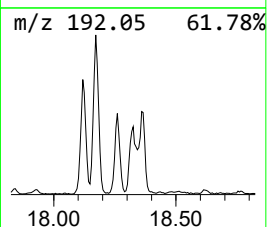
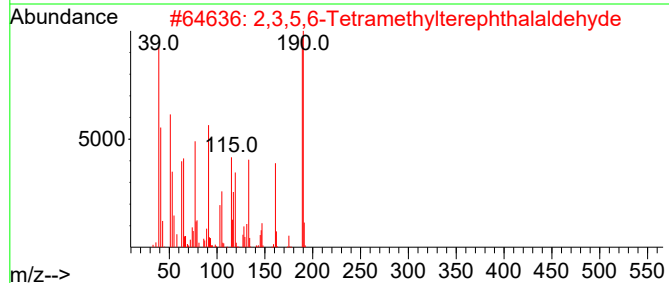
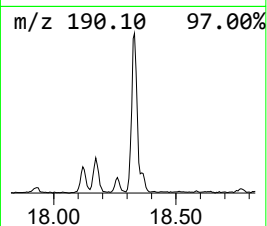
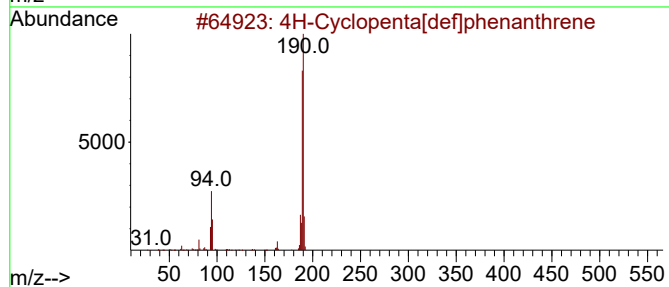
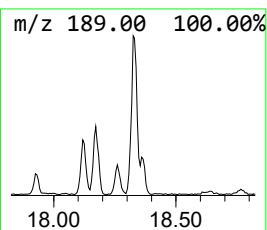
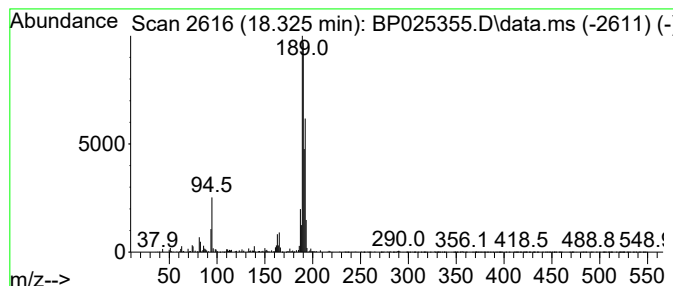
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 6 4H-Cyclopenta[def]phenanthrene Concentration Rank 9

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 18.325 | 2.88 ng | 371156 | Phenanthrene-d10 | 17.219 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 4H-Cyclopenta[def]phenanthrene      | 190 | C15H10   | 000203-64-5 | 92   |
| 2    |    |   | 2,3,5,6-Tetramethylterephthalald... | 190 | C12H14O2 | 007072-01-7 | 43   |
| 3    |    |   | Methyl diselenide                   | 190 | C2H6Se2  | 007101-31-7 | 27   |
| 4    |    |   | 5H-Dibenzo[a,d]cycloheptene         | 192 | C15H12   | 000256-81-5 | 25   |
| 5    |    |   | Phenanthrene, 2-methyl-             | 192 | C15H12   | 002531-84-2 | 25   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP081125\  
Data File : BP025355.D  
Acq On : 11 Aug 2025 19:34  
Operator : CG/JU  
Sample : Q2795-03  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
COMP-3

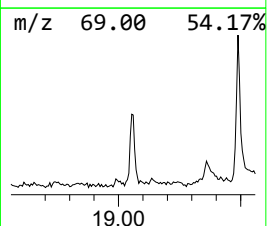
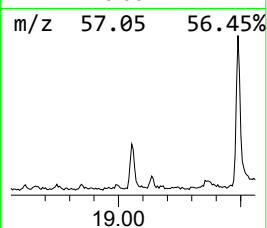
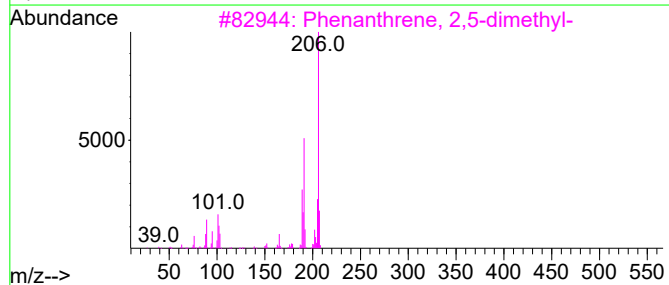
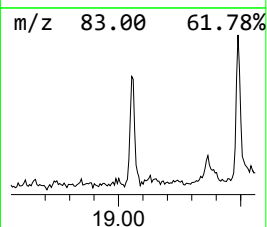
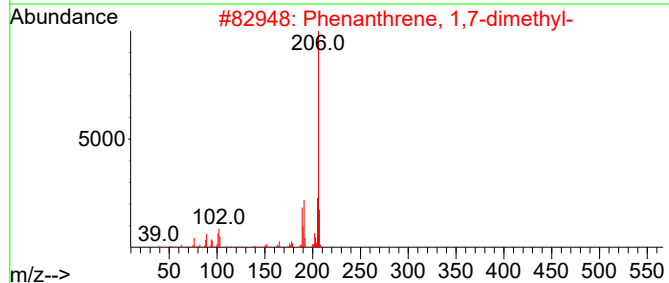
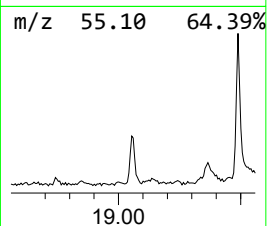
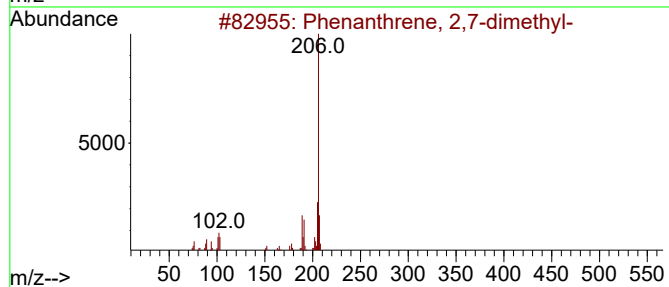
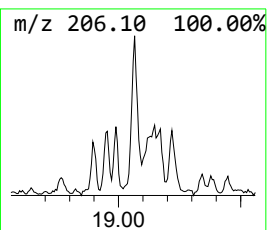
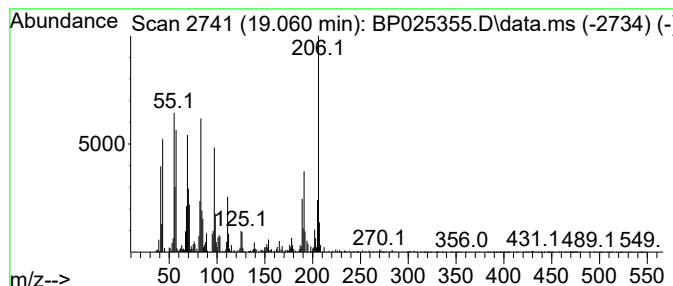
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 7 Phenanthrene, 2,7-dimethyl- Concentration Rank 10

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 19.060 | 2.86 ng | 367911 | Phenanthrene-d10 | 17.219 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Phenanthrene, 2,7-dimethyl- | 206 | C16H14  | 001576-69-8 | 94   |
| 2    |    |   | Phenanthrene, 1,7-dimethyl- | 206 | C16H14  | 000483-87-4 | 83   |
| 3    |    |   | Phenanthrene, 2,5-dimethyl- | 206 | C16H14  | 003674-66-6 | 64   |
| 4    |    |   | Phenanthrene, 3,6-dimethyl- | 206 | C16H14  | 001576-67-6 | 64   |
| 5    |    |   | Phenanthrene, 2,3-dimethyl- | 206 | C16H14  | 003674-65-5 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP081125\  
Data File : BP025355.D  
Acq On : 11 Aug 2025 19:34  
Operator : CG/JU  
Sample : Q2795-03  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
COMP-3

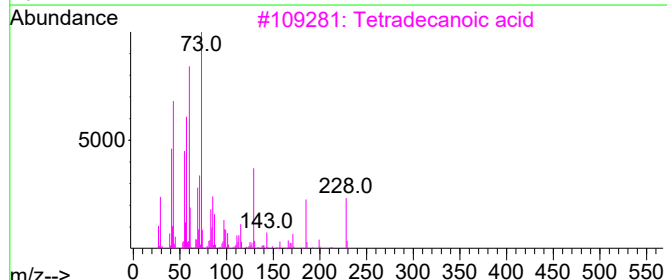
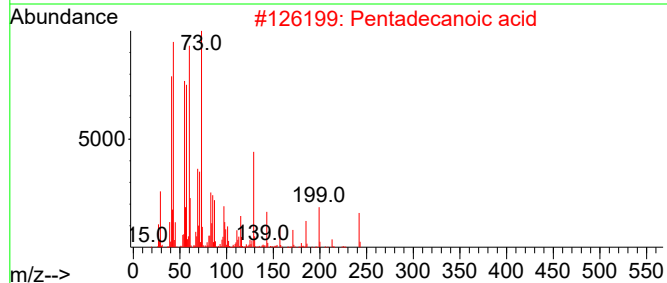
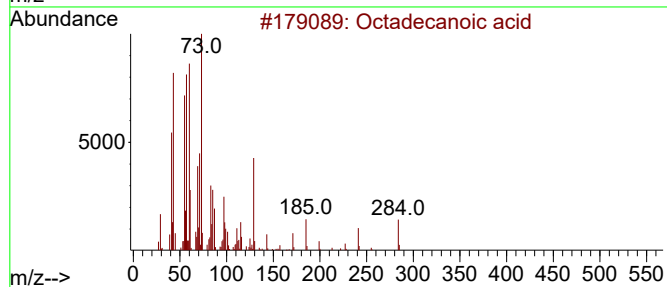
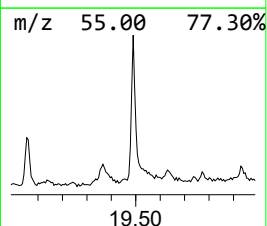
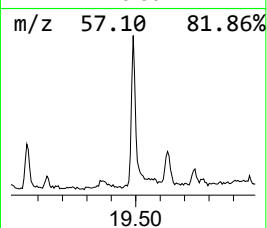
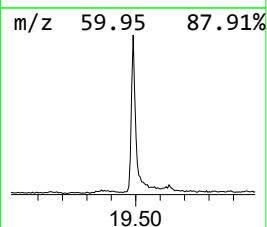
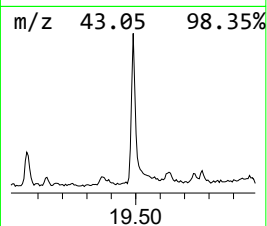
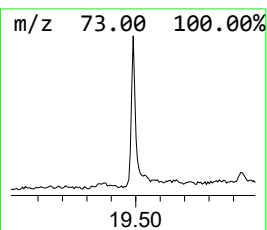
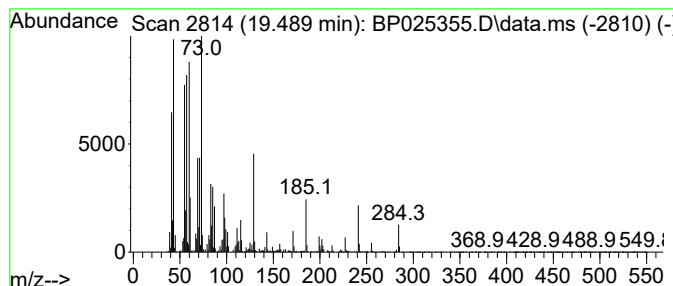
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 8 Octadecanoic acid Concentration Rank 8

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 19.489 | 3.36 ng | 735806 | Chrysene-d12     | 21.660 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Octadecanoic acid                   | 284 | C18H36O2 | 000057-11-4 | 99   |
| 2    |    |   | Pentadecanoic acid                  | 242 | C15H30O2 | 001002-84-2 | 91   |
| 3    |    |   | Tetradecanoic acid                  | 228 | C14H28O2 | 000544-63-8 | 87   |
| 4    |    |   | Octadecanoic acid, 2-(2-hydroxye... | 372 | C22H44O4 | 000106-11-6 | 86   |
| 5    |    |   | n-Hexadecanoic acid                 | 256 | C16H32O2 | 000057-10-3 | 81   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP081125\  
Data File : BP025355.D  
Acq On : 11 Aug 2025 19:34  
Operator : CG/JU  
Sample : Q2795-03  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
COMP-3

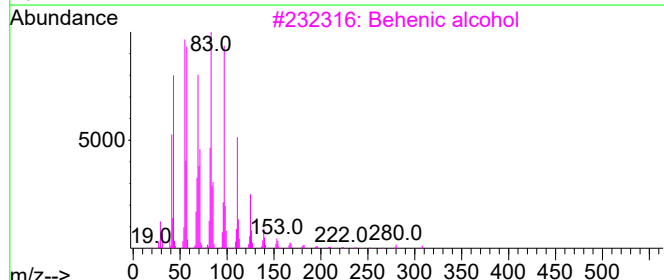
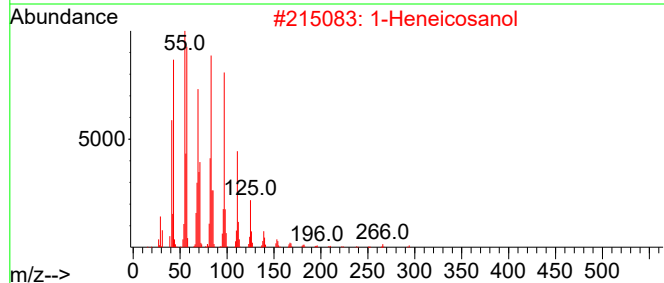
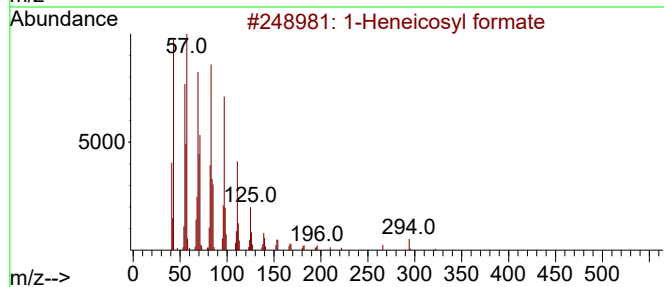
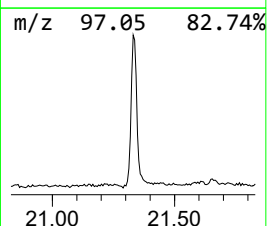
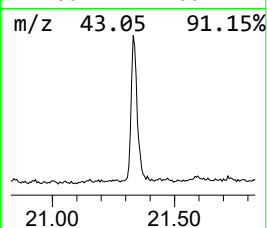
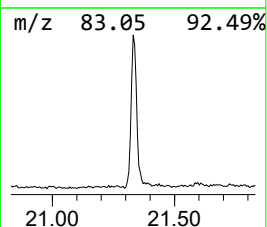
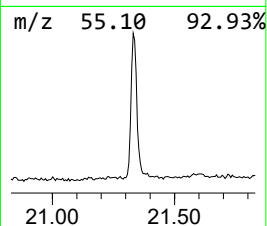
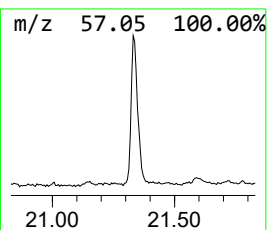
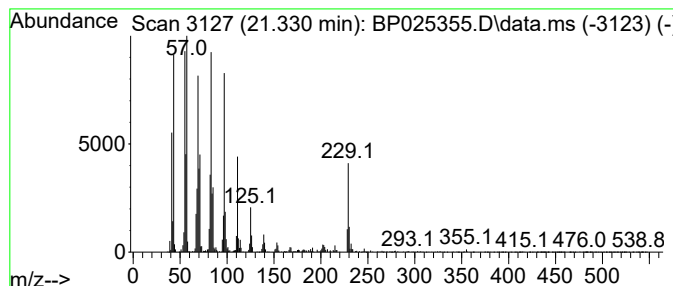
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 9 1-Heneicosyl formate Concentration Rank 4

| R.T.   | EstConc | Area    | Relative to ISTD | R.T.   |
|--------|---------|---------|------------------|--------|
| 21.330 | 7.14 ng | 1563970 | Chrysene-d12     | 21.660 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|------------|-------------|------|
| 1    |    |   | 1-Heneicosyl formate                | 340 | C22H44O2   | 077899-03-7 | 95   |
| 2    |    |   | 1-Heneicosanol                      | 312 | C21H44O    | 015594-90-8 | 91   |
| 3    |    |   | Behenic alcohol                     | 326 | C22H46O    | 000661-19-8 | 91   |
| 4    |    |   | n-Tetracosanol-1                    | 354 | C24H50O    | 000506-51-4 | 91   |
| 5    |    |   | Heptafluorobutyric acid, n-tetra... | 410 | C18H29F7O2 | 007365-36-8 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP081125\  
Data File : BP025355.D  
Acq On : 11 Aug 2025 19:34  
Operator : CG/JU  
Sample : Q2795-03  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
COMP-3

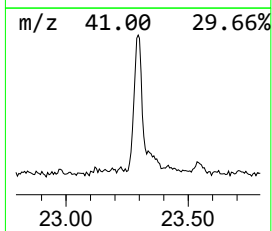
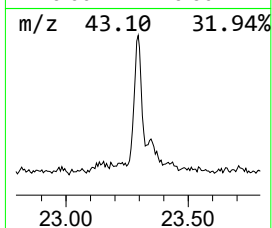
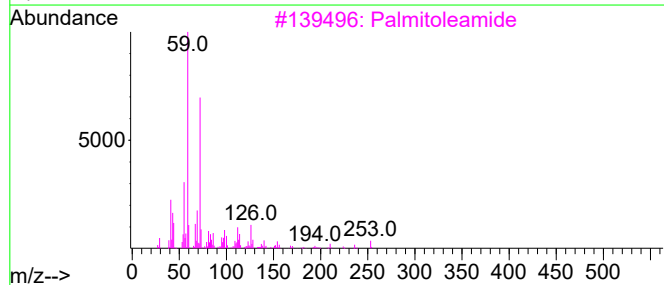
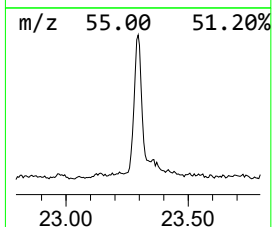
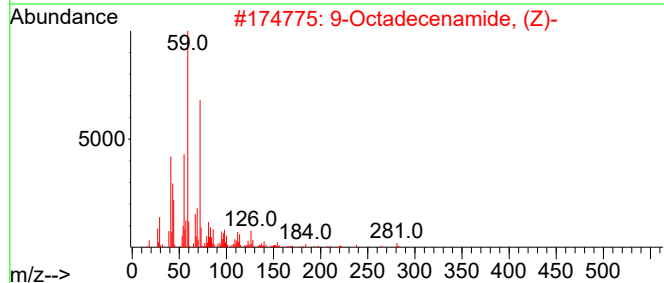
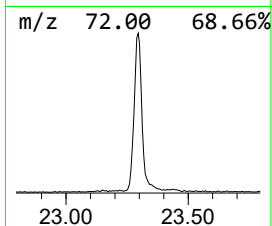
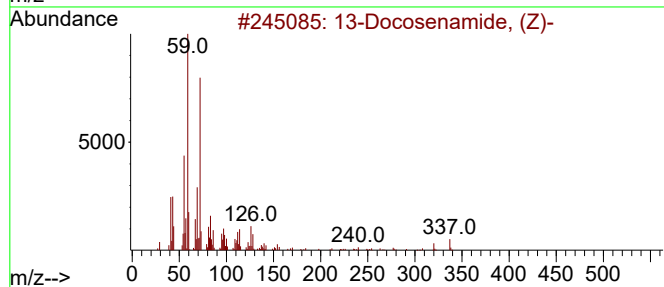
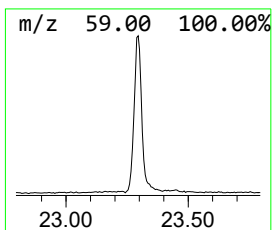
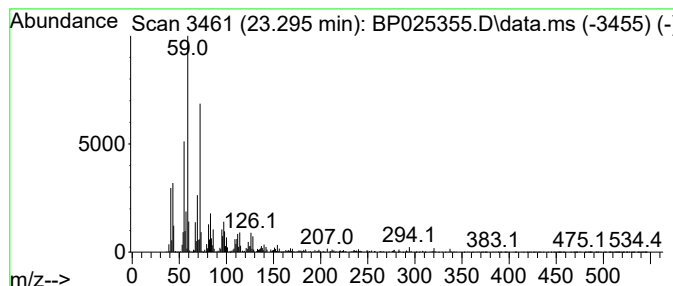
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP080525.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 10 13-Docosenamide, (Z)- Concentration Rank 3

| R.T.   | EstConc | Area    | Relative to ISTD | R.T.   |
|--------|---------|---------|------------------|--------|
| 23.295 | 7.41 ng | 1625120 | Chrysene-d12     | 21.660 |

| Hit# | of 5 | Tentative ID           | MW  | MolForm  | CAS#        | Qual |
|------|------|------------------------|-----|----------|-------------|------|
| 1    |      | 13-Docosenamide, (Z)-  | 337 | C22H43NO | 000112-84-5 | 96   |
| 2    |      | 9-Octadecenamide, (Z)- | 281 | C18H35NO | 000301-02-0 | 91   |
| 3    |      | Palmitoleamide         | 253 | C16H31NO | 106010-22-4 | 64   |
| 4    |      | Tetradecanamide        | 227 | C14H29NO | 000638-58-4 | 59   |
| 5    |      | Dodecanamide           | 199 | C12H25NO | 001120-16-7 | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_P\Data\BP081125\  
Data File : BP025355.D  
Acq On : 11 Aug 2025 19:34  
Operator : CG/JU  
Sample : Q2795-03  
Misc :  
ALS Vial : 14 Sample Multiplier: 1

Instrument :  
BNA\_P  
ClientSampleId :  
COMP-3

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_P\Methods\8270E-BP080525.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 |
| Cyclohexane         | 2.967  | 4.2     | ng    | 362934   | 1                     | 7.802  | 1743930 | 20.0 |
| Butane, 2-metho...  | 3.049  | 32.6    | ng    | 2844450  | 1                     | 7.802  | 1743930 | 20.0 |
| 2-Pentanone, 4-...  | 4.949  | 7.1     | ng    | 621788   | 1                     | 7.802  | 1743930 | 20.0 |
| Benzophenone        | 15.795 | 6.8     | ng    | 872446   | 3                     | 14.419 | 2558440 | 20.0 |
| n-Hexadecanoic ...  | 18.166 | 15.4    | ng    | 1983630  | 4                     | 17.219 | 2573740 | 20.0 |
| 4H-Cyclopenta[d]... | 18.325 | 2.9     | ng    | 371156   | 4                     | 17.219 | 2573740 | 20.0 |
| Phenanthrene, 2...  | 19.060 | 2.9     | ng    | 367911   | 4                     | 17.219 | 2573740 | 20.0 |
| Octadecanoic acid   | 19.489 | 3.4     | ng    | 735806   | 5                     | 21.660 | 4383500 | 20.0 |
| 1-Heneicosyl fo...  | 21.330 | 7.1     | ng    | 1563970  | 5                     | 21.660 | 4383500 | 20.0 |
| 13-Docosenamide...  | 23.295 | 7.4     | ng    | 1625120  | 5                     | 21.660 | 4383500 | 20.0 |