

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG072222\  
 Data File : BG054367.D  
 Acq On : 22 Jul 2022 20:03  
 Operator : CG/JU  
 Sample : N3814-10  
 Misc :  
 ALS Vial : 12 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 ENV-DE-1(0-5)

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: BG054367.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         | 5.567       | 363           | 371         | 381          | rBV      | 65361          | 113868        | 10.97%          | 1.097%        |
| 2         | 7.171       | 638           | 644         | 653          | rBV      | 74463          | 133912        | 12.91%          | 1.290%        |
| 3         | 7.994       | 777           | 784         | 791          | rBB      | 72240          | 130425        | 12.57%          | 1.257%        |
| 4         | 9.151       | 975           | 981         | 989          | rBV      | 48952          | 90062         | 8.68%           | 0.868%        |
| 5         | 10.802      | 1254          | 1262        | 1268         | rBV      | 96994          | 183246        | 17.66%          | 1.766%        |
| 6         | 13.247      | 1672          | 1678        | 1687         | rBB      | 158929         | 248114        | 23.91%          | 2.391%        |
| 7         | 14.351      | 1860          | 1866        | 1877         | rBV      | 10587          | 19481         | 1.88%           | 0.188%        |
| 8         | 14.627      | 1905          | 1913        | 1919         | rBV      | 175572         | 292540        | 28.19%          | 2.819%        |
| 9         | 14.686      | 1919          | 1923        | 1931         | rVB      | 12262          | 18848         | 1.82%           | 0.182%        |
| 10        | 15.021      | 1975          | 1980        | 1986         | rBB      | 8086           | 12487         | 1.20%           | 0.120%        |
| 11        | 15.673      | 2084          | 2091        | 2099         | rBB      | 19260          | 31010         | 2.99%           | 0.299%        |
| 12        | 16.114      | 2160          | 2166        | 2176         | rBV3     | 145141         | 223736        | 21.56%          | 2.156%        |
| 13        | 16.678      | 2249          | 2262        | 2267         | rBV4     | 6958           | 17202         | 1.66%           | 0.166%        |
| 14        | 16.901      | 2291          | 2300        | 2304         | rBV6     | 4932           | 11412         | 1.10%           | 0.110%        |
| 15        | 17.025      | 2315          | 2321        | 2327         | rVB      | 9940           | 16335         | 1.57%           | 0.157%        |
| 16        | 17.119      | 2327          | 2337        | 2342         | rBV7     | 9381           | 20813         | 2.01%           | 0.201%        |
| 17        | 17.189      | 2345          | 2349        | 2353         | rVB      | 12171          | 16569         | 1.60%           | 0.160%        |
| 18        | 17.371      | 2373          | 2380        | 2383         | rBV      | 225801         | 341434        | 32.91%          | 3.290%        |
| 19        | 17.412      | 2383          | 2387        | 2394         | rVV      | 285091         | 427661        | 41.22%          | 4.121%        |
| 20        | 17.506      | 2397          | 2403        | 2408         | rVV      | 63324          | 98364         | 9.48%           | 0.948%        |
| 21        | 17.771      | 2438          | 2448        | 2453         | rBV2     | 23184          | 46853         | 4.52%           | 0.451%        |
| 22        | 17.953      | 2474          | 2479        | 2486         | rVB6     | 8981           | 14969         | 1.44%           | 0.144%        |
| 23        | 18.035      | 2486          | 2493        | 2499         | rBV8     | 4623           | 11150         | 1.07%           | 0.107%        |
| 24        | 18.247      | 2524          | 2529        | 2533         | rVV      | 49648          | 81869         | 7.89%           | 0.789%        |
| 25        | 18.294      | 2533          | 2537        | 2545         | rVV      | 63574          | 103560        | 9.98%           | 0.998%        |
| 26        | 18.376      | 2545          | 2551        | 2556         | rVV      | 27279          | 41217         | 3.97%           | 0.397%        |
| 27        | 18.441      | 2556          | 2562        | 2566         | rVV3     | 70711          | 136330        | 13.14%          | 1.314%        |
| 28        | 18.476      | 2566          | 2568        | 2574         | rVB      | 34094          | 45448         | 4.38%           | 0.438%        |
| 29        | 18.552      | 2574          | 2581        | 2585         | rBV6     | 7703           | 14963         | 1.44%           | 0.144%        |
| 30        | 18.740      | 2608          | 2613        | 2617         | rBV      | 39649          | 53220         | 5.13%           | 0.513%        |
| 31        | 18.787      | 2617          | 2621        | 2627         | rVB3     | 20090          | 31831         | 3.07%           | 0.307%        |
| 32        | 18.987      | 2647          | 2655        | 2659         | rBV2     | 13989          | 26044         | 2.51%           | 0.251%        |
| 33        | 19.040      | 2661          | 2664        | 2667         | rBV2     | 10401          | 10410         | 1.00%           | 0.100%        |
| 34        | 19.075      | 2667          | 2670        | 2674         | rVB3     | 10857          | 14833         | 1.43%           | 0.143%        |
| 35        | 19.163      | 2678          | 2685        | 2689         | rBV      | 35779          | 68303         | 6.58%           | 0.658%        |
| 36        | 19.222      | 2691          | 2695        | 2698         | rBV4     | 10510          | 17737         | 1.71%           | 0.171%        |

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 ALS Vial : 12 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 ENV-DE-1(0-5)

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 | 19.304 | 2704 | 2709 | 2716 | rBV2 | 28332  | 53239   | 5.13%   | 0.513% |
| 38 | 19.428 | 2723 | 2730 | 2736 | rBV  | 603429 | 890420  | 85.81%  | 8.580% |
| 39 | 19.481 | 2736 | 2739 | 2742 | rBV  | 10277  | 14199   | 1.37%   | 0.137% |
| 40 | 19.516 | 2742 | 2745 | 2751 | rVB2 | 16374  | 24506   | 2.36%   | 0.236% |
| 41 | 19.575 | 2751 | 2755 | 2758 | rBV  | 10508  | 15145   | 1.46%   | 0.146% |
| 42 | 19.669 | 2768 | 2771 | 2774 | rVB  | 15854  | 15653   | 1.51%   | 0.151% |
| 43 | 19.757 | 2780 | 2786 | 2787 | rBV  | 106552 | 151799  | 14.63%  | 1.463% |
| 44 | 19.786 | 2787 | 2791 | 2798 | rVV  | 493546 | 768805  | 74.09%  | 7.408% |
| 45 | 19.857 | 2799 | 2803 | 2811 | rVB3 | 31143  | 53474   | 5.15%   | 0.515% |
| 46 | 19.980 | 2816 | 2824 | 2828 | rBV2 | 313894 | 452647  | 43.62%  | 4.362% |
| 47 | 20.027 | 2828 | 2832 | 2837 | rVB  | 31429  | 48337   | 4.66%   | 0.466% |
| 48 | 20.103 | 2840 | 2845 | 2846 | rBV2 | 45979  | 62927   | 6.06%   | 0.606% |
| 49 | 20.156 | 2852 | 2854 | 2858 | rBV  | 8639   | 13375   | 1.29%   | 0.129% |
| 50 | 20.280 | 2864 | 2875 | 2880 | rVB  | 89207  | 190824  | 18.39%  | 1.839% |
| 51 | 20.379 | 2887 | 2892 | 2895 | rBV  | 57582  | 90215   | 8.69%   | 0.869% |
| 52 | 20.432 | 2896 | 2901 | 2905 | rVV2 | 48862  | 91118   | 8.78%   | 0.878% |
| 53 | 20.509 | 2912 | 2914 | 2920 | rVB4 | 15047  | 20119   | 1.94%   | 0.194% |
| 54 | 20.573 | 2922 | 2925 | 2929 | rVV  | 34726  | 43774   | 4.22%   | 0.422% |
| 55 | 20.620 | 2930 | 2933 | 2937 | rVB  | 31370  | 43486   | 4.19%   | 0.419% |
| 56 | 21.038 | 3000 | 3004 | 3009 | rVB  | 61801  | 86108   | 8.30%   | 0.830% |
| 57 | 21.214 | 3029 | 3034 | 3039 | rBV3 | 48018  | 100334  | 9.67%   | 0.967% |
| 58 | 21.261 | 3039 | 3042 | 3046 | rVV  | 53581  | 89815   | 8.66%   | 0.865% |
| 59 | 21.314 | 3047 | 3051 | 3056 | rVV2 | 55259  | 108317  | 10.44%  | 1.044% |
| 60 | 21.372 | 3056 | 3061 | 3067 | rVB2 | 57375  | 111634  | 10.76%  | 1.076% |
| 61 | 21.625 | 3097 | 3104 | 3110 | rBV3 | 413998 | 1037607 | 100.00% | 9.998% |
| 62 | 21.684 | 3110 | 3114 | 3125 | rVB  | 282487 | 491545  | 47.37%  | 4.736% |
| 63 | 21.831 | 3135 | 3139 | 3145 | rBV  | 47137  | 90721   | 8.74%   | 0.874% |
| 64 | 22.360 | 3226 | 3229 | 3234 | rVB  | 50530  | 79361   | 7.65%   | 0.765% |
| 65 | 22.430 | 3238 | 3241 | 3247 | rVB  | 29245  | 48930   | 4.72%   | 0.471% |
| 66 | 23.828 | 3471 | 3479 | 3486 | rBV  | 209452 | 677682  | 65.31%  | 6.530% |
| 67 | 24.539 | 3594 | 3600 | 3612 | rVB  | 120364 | 345710  | 33.32%  | 3.331% |
| 68 | 24.686 | 3618 | 3625 | 3635 | rVB  | 168938 | 480592  | 46.32%  | 4.631% |
| 69 | 24.839 | 3643 | 3651 | 3659 | rBV  | 146596 | 419285  | 40.41%  | 4.040% |

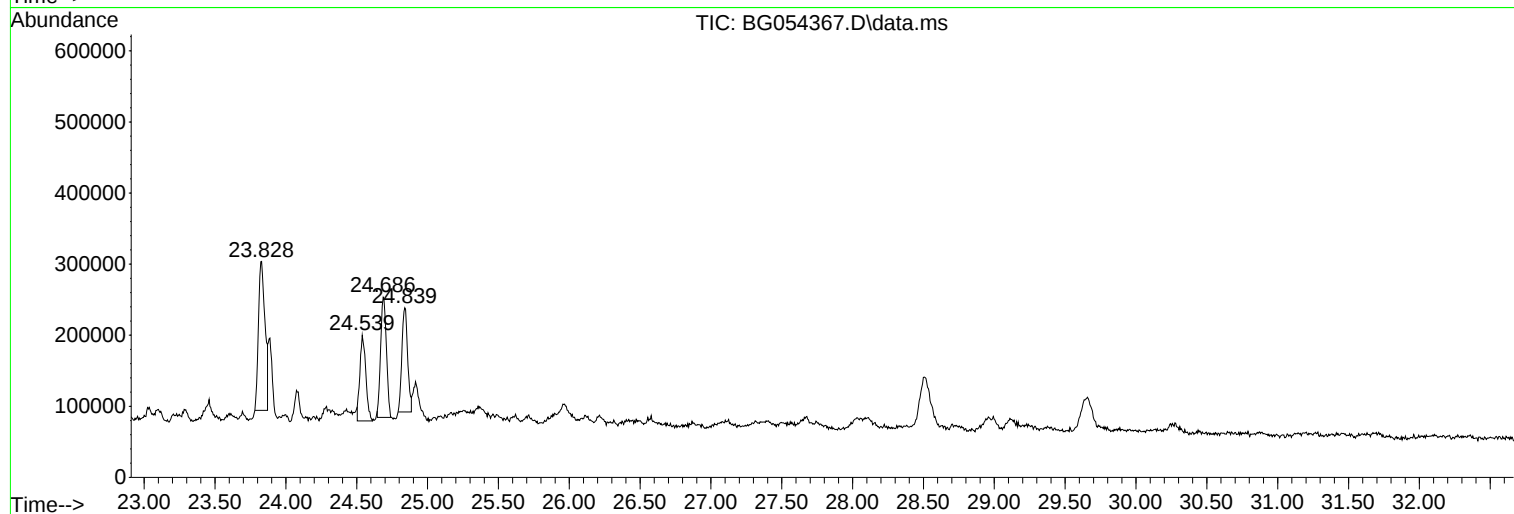
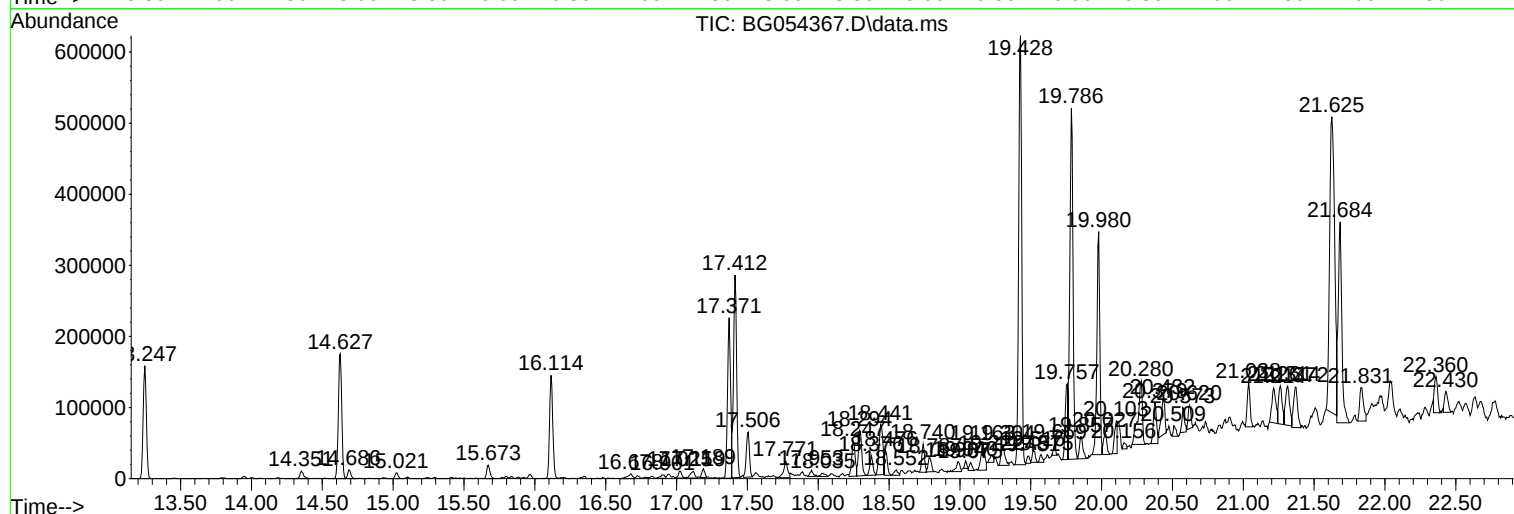
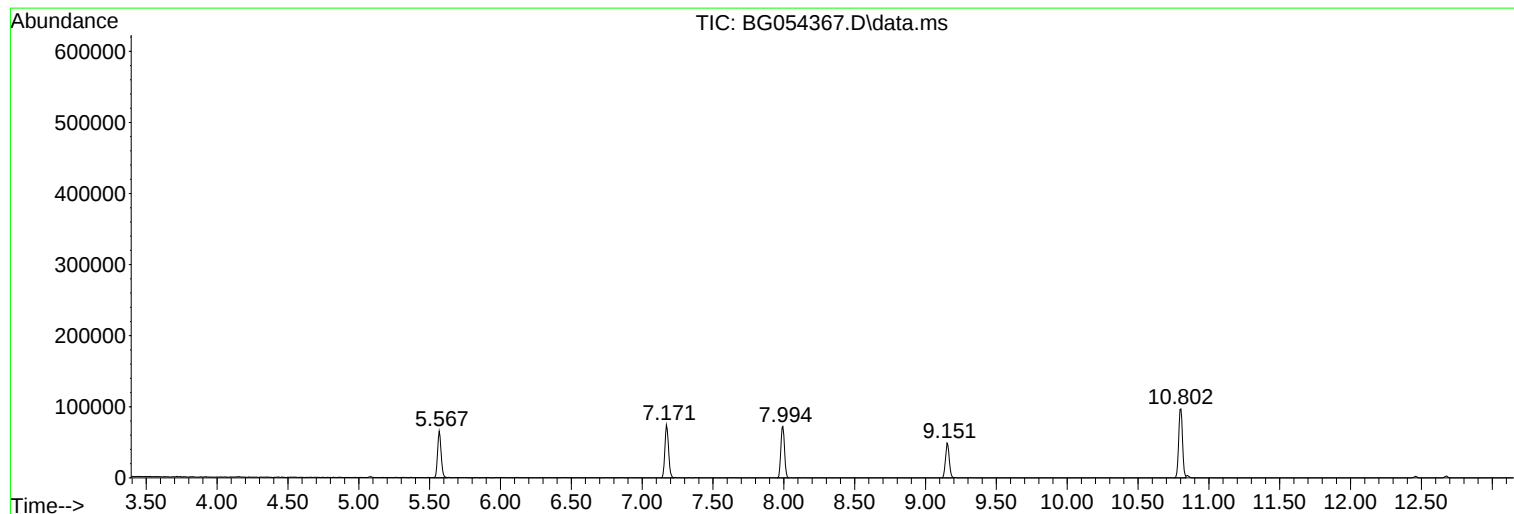
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Sample : N3814-10  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
ENV-DE-1(0-5)

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\BG072222\  
Data File : BG054367.D  
Acq On : 22 Jul 2022 20:03  
Operator : CG/JU  
Sample : N3814-10  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
ENV-DE-1(0-5)

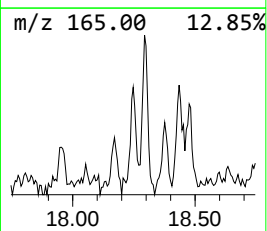
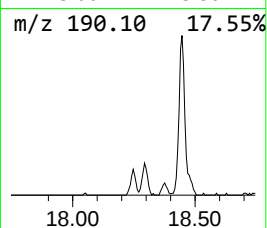
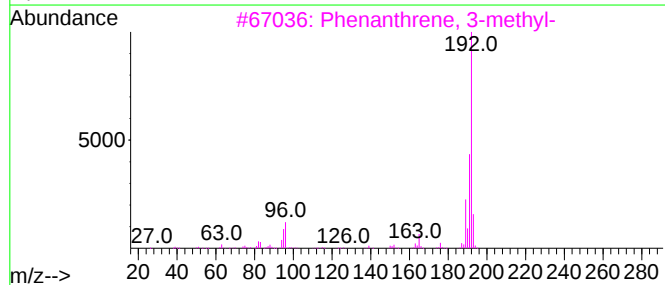
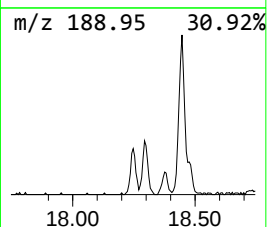
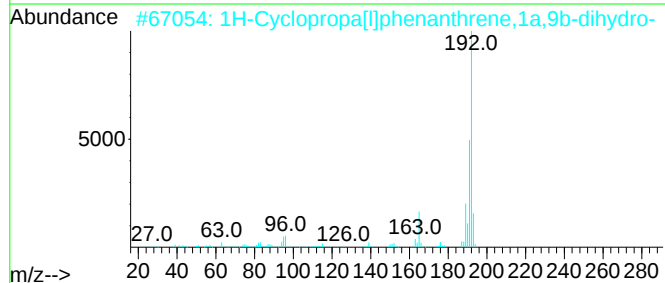
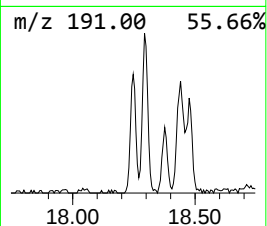
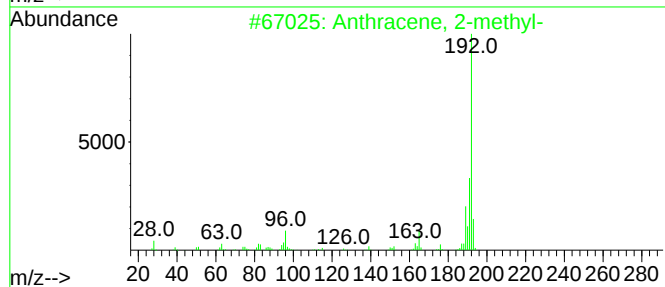
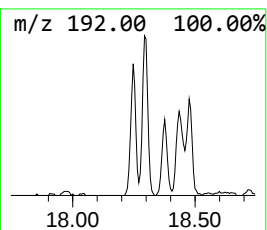
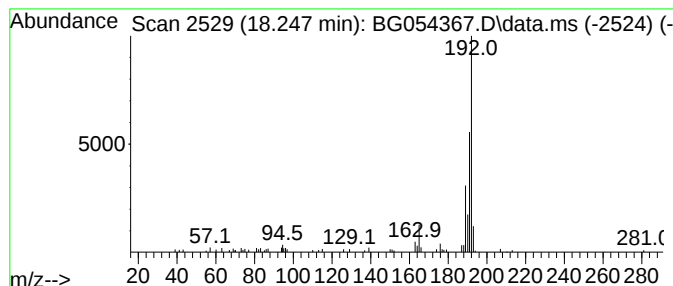
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG071522.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 2 Anthracene, 2-methyl- Concentration Rank 4

| R.T.   | EstConc | Area  | Relative to ISTD | R.T.   |
|--------|---------|-------|------------------|--------|
| 18.247 | 4.80 ng | 81869 | Phenanthrene-d10 | 17.371 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Anthracene, 2-methyl-               | 192 | C15H12  | 000613-12-7 | 94   |
| 2    |    |   | 1H-Cyclopropa[1]phenanthrene,1a,... | 192 | C15H12  | 000949-41-7 | 91   |
| 3    |    |   | Phenanthrene, 3-methyl-             | 192 | C15H12  | 000832-71-3 | 91   |
| 4    |    |   | Phenanthrene, 1-methyl-             | 192 | C15H12  | 000832-69-9 | 91   |
| 5    |    |   | Phenanthrene, 2-methyl-             | 192 | C15H12  | 002531-84-2 | 89   |



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Sample : N3814-10  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
ENV-DE-1(0-5)

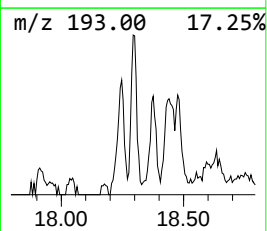
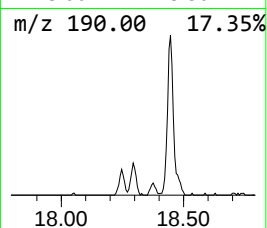
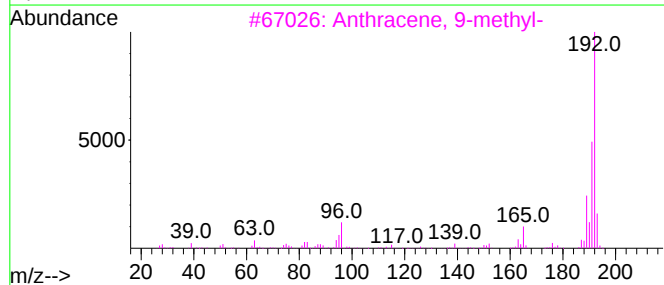
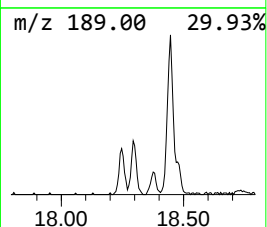
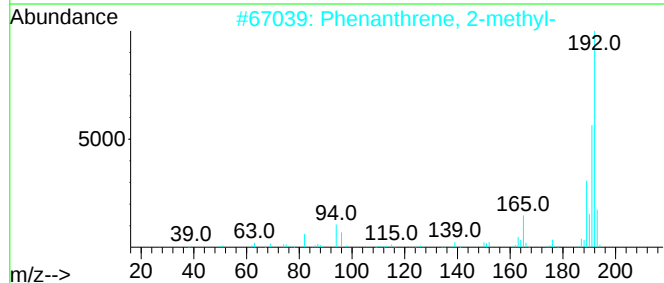
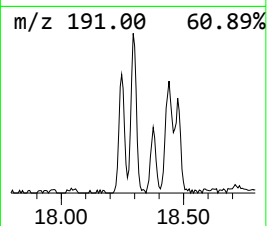
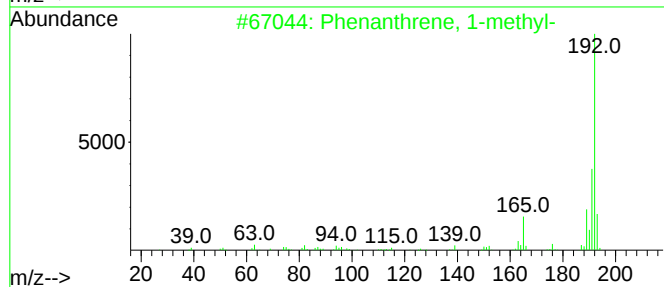
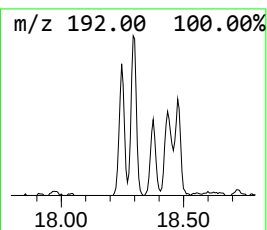
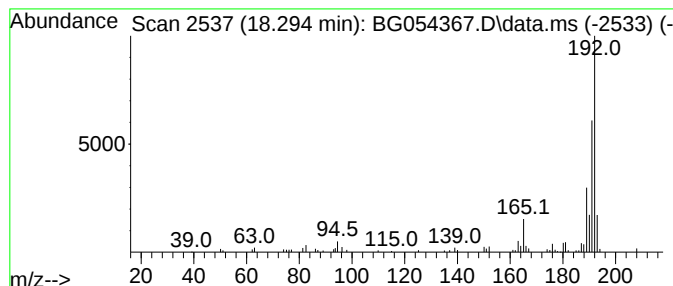
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG071522.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 3 Phenanthrene, 1-methyl- Concentration Rank 3

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 18.294 | 6.07 ng | 103560 | Phenanthrene-d10 | 17.371 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Phenanthrene, 1-methyl- | 192 | C15H12  | 000832-69-9 | 96   |
| 2    |    |   | Phenanthrene, 2-methyl- | 192 | C15H12  | 002531-84-2 | 95   |
| 3    |    |   | Anthracene, 9-methyl-   | 192 | C15H12  | 000779-02-2 | 94   |
| 4    |    |   | Anthracene, 1-methyl-   | 192 | C15H12  | 000610-48-0 | 93   |
| 5    |    |   | 1H-Indene, 2-phenyl-    | 192 | C15H12  | 004505-48-0 | 93   |



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Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
ENV-DE-1(0-5)

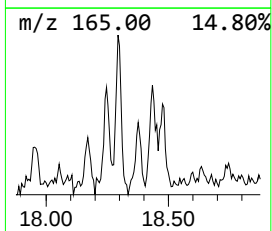
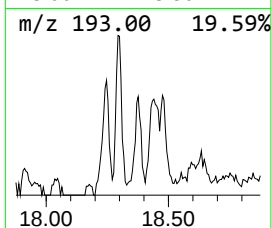
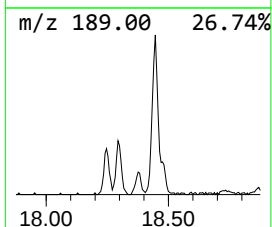
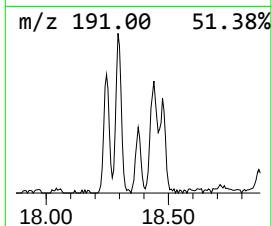
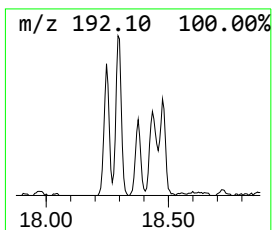
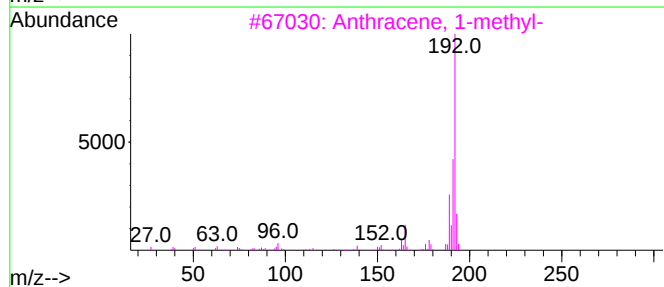
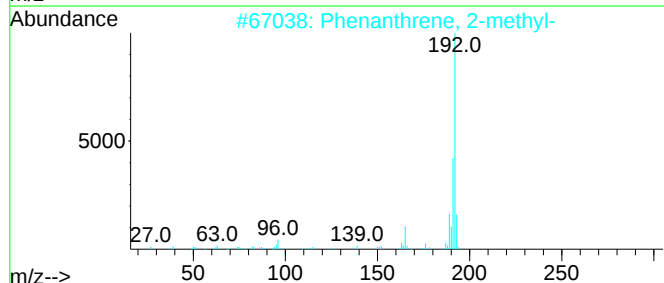
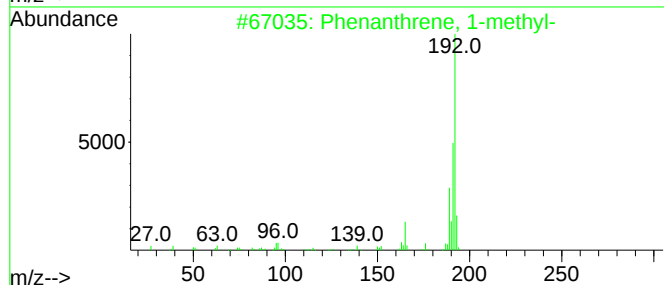
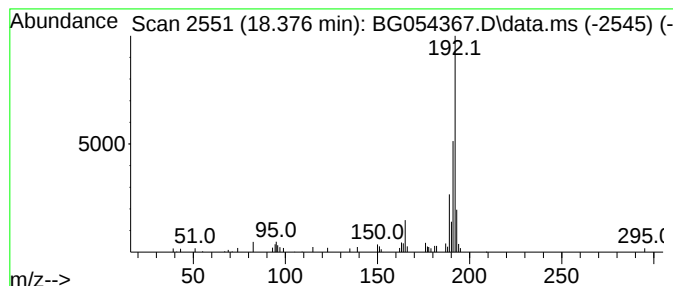
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG071522.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 4 Phenanthrene, 2-methyl- Concentration Rank 10

| R.T.   | EstConc | Area  | Relative to ISTD | R.T.   |
|--------|---------|-------|------------------|--------|
| 18.376 | 2.41 ng | 41217 | Phenanthrene-d10 | 17.371 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Phenanthrene, 1-methyl-             | 192 | C15H12  | 000832-69-9 | 94   |
| 2    |    |   | Phenanthrene, 2-methyl-             | 192 | C15H12  | 002531-84-2 | 93   |
| 3    |    |   | Anthracene, 1-methyl-               | 192 | C15H12  | 000610-48-0 | 91   |
| 4    |    |   | 1H-Cyclopropa[1]phenanthrene,1a,... | 192 | C15H12  | 000949-41-7 | 91   |
| 5    |    |   | Phenanthrene, 4-methyl-             | 192 | C15H12  | 000832-64-4 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG072222\  
Data File : BG054367.D  
Acq On : 22 Jul 2022 20:03  
Operator : CG/JU  
Sample : N3814-10  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
ENV-DE-1(0-5)

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

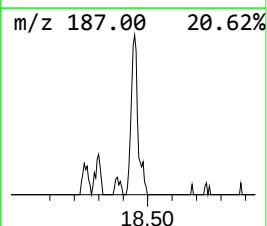
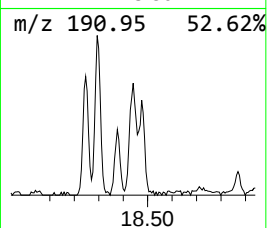
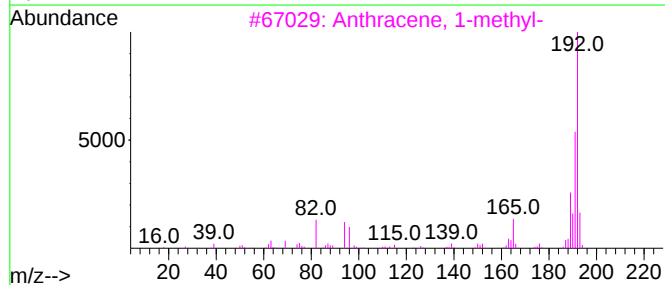
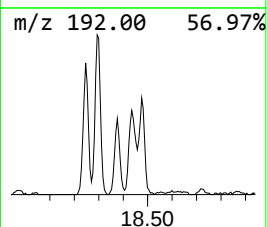
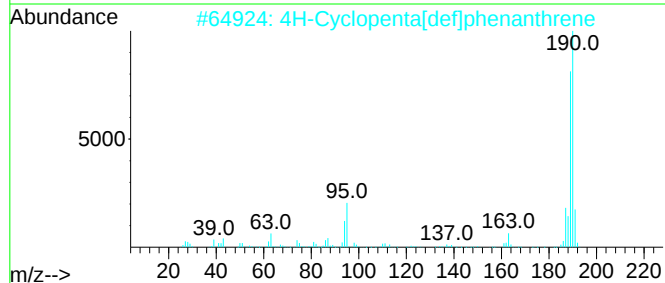
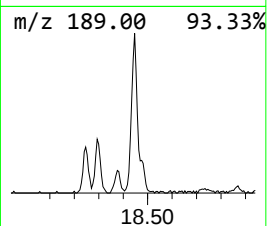
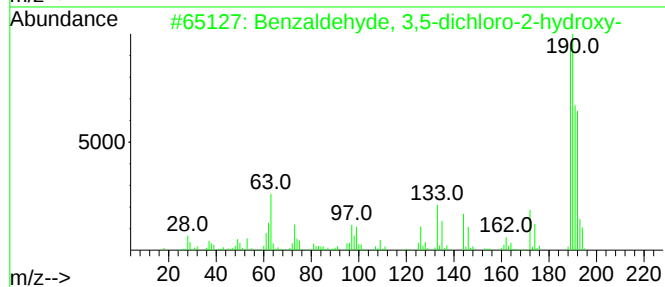
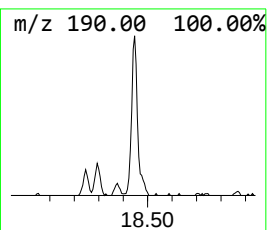
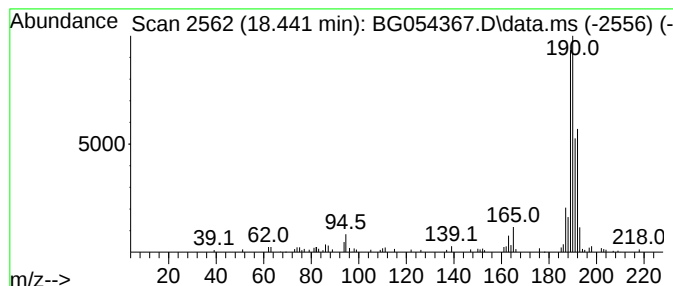
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TIC Integration Parameters: LSCINT.P

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Peak Number 5 Benzaldehyde, 3,5-dichloro-... Concentration Rank 2

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 18.441 | 7.99 ng | 136330 | Phenanthrene-d10 | 17.371 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|-----------|-------------|------|
| 1    |    |   | Benzaldehyde, 3,5-dichloro-2-hyd... | 190 | C7H4Cl2O2 | 000090-60-8 | 72   |
| 2    |    |   | 4H-Cyclopenta[def]phenanthrene      | 190 | C15H10    | 000203-64-5 | 64   |
| 3    |    |   | Anthracene, 1-methyl-               | 192 | C15H12    | 000610-48-0 | 35   |
| 4    |    |   | 5H-Dibenzo[a,d]cycloheptene         | 192 | C15H12    | 000256-81-5 | 30   |
| 5    |    |   | 8,9-Dihydrocyclopenta[def]phenan... | 192 | C15H12    | 027410-55-5 | 30   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG072222\  
Data File : BG054367.D  
Acq On : 22 Jul 2022 20:03  
Operator : CG/JU  
Sample : N3814-10  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
ENV-DE-1(0-5)

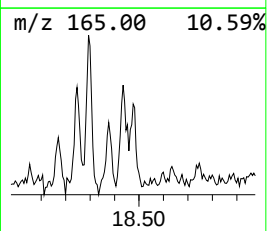
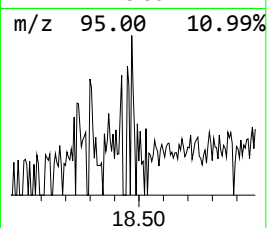
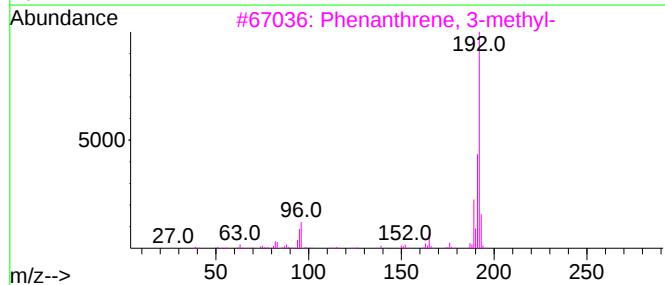
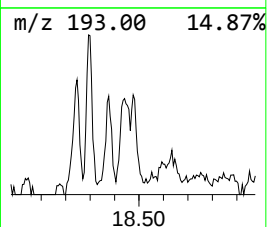
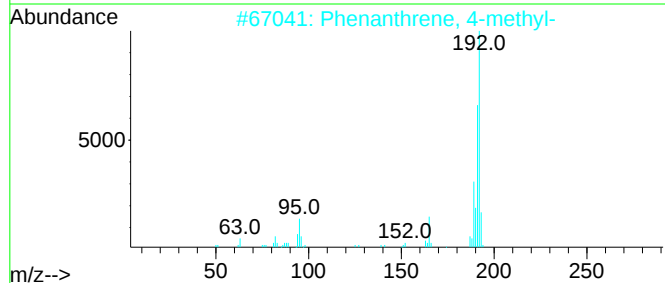
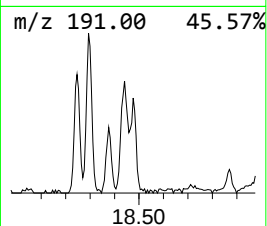
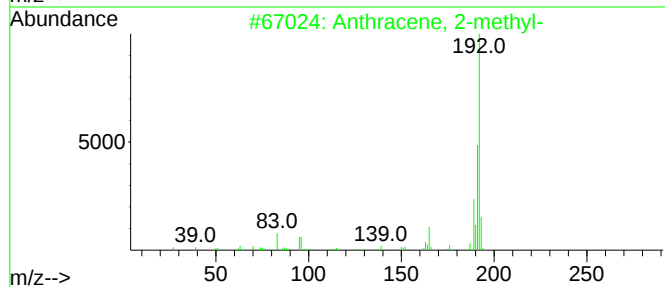
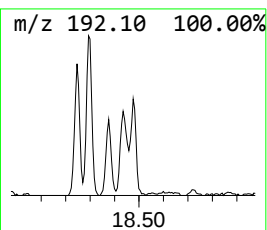
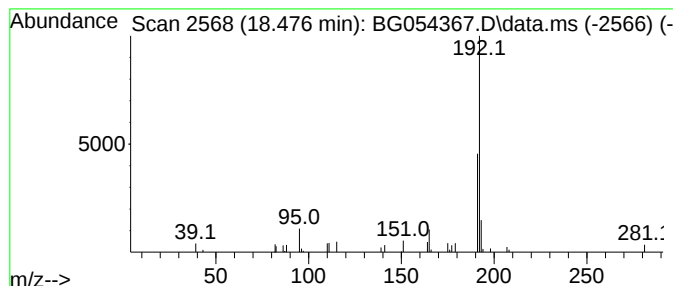
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 6 Phenanthrene, 4-methyl- Concentration Rank 9

| R.T.   | EstConc | Area  | Relative to ISTD | R.T.   |
|--------|---------|-------|------------------|--------|
| 18.476 | 2.66 ng | 45448 | Phenanthrene-d10 | 17.371 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Anthracene, 2-methyl-   | 192 | C15H12  | 000613-12-7 | 87   |
| 2    |    |   | Phenanthrene, 4-methyl- | 192 | C15H12  | 000832-64-4 | 86   |
| 3    |    |   | Phenanthrene, 3-methyl- | 192 | C15H12  | 000832-71-3 | 86   |
| 4    |    |   | Anthracene, 1-methyl-   | 192 | C15H12  | 000610-48-0 | 86   |
| 5    |    |   | Phenanthrene, 9-methyl- | 192 | C15H12  | 000883-20-5 | 86   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG072222\  
Data File : BG054367.D  
Acq On : 22 Jul 2022 20:03  
Operator : CG/JU  
Sample : N3814-10  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
ENV-DE-1(0-5)

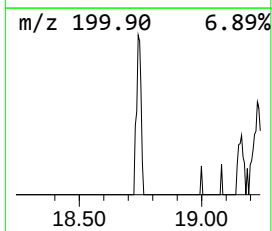
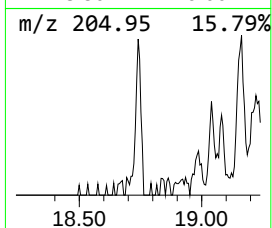
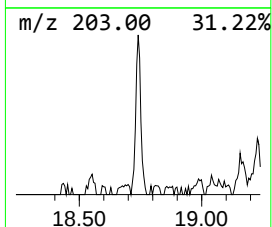
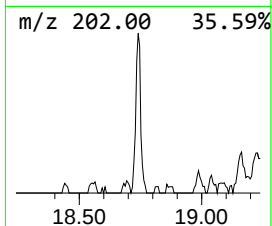
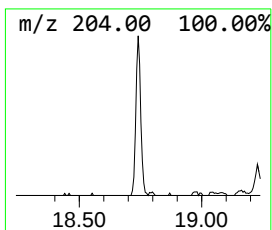
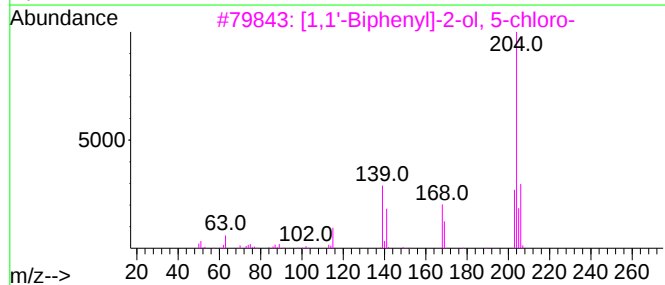
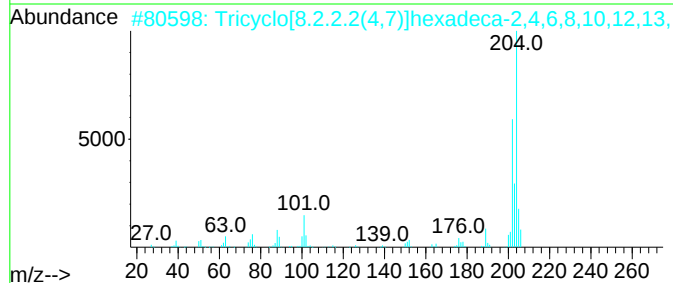
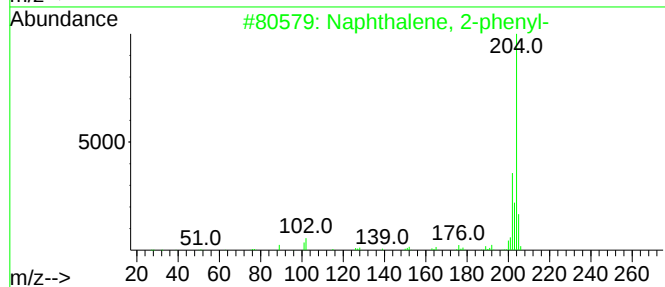
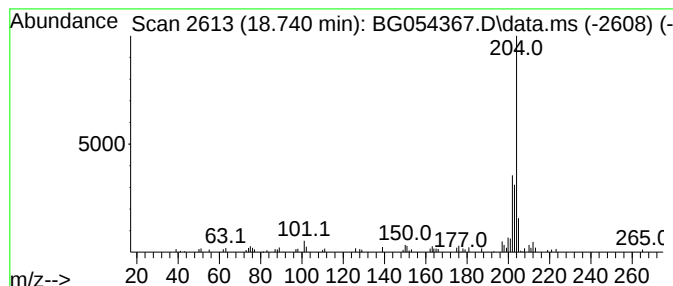
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 7 Naphthalene, 2-phenyl- Concentration Rank 8

| R.T.   | EstConc | Area  | Relative to ISTD | R.T.   |
|--------|---------|-------|------------------|--------|
| 18.740 | 3.12 ng | 53220 | Phenanthrene-d10 | 17.371 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Naphthalene, 2-phenyl-              | 204 | C16H12   | 000612-94-2 | 68   |
| 2    |    |   | Tricyclo[8.2.2.2(4,7)]hexadeca-2... | 204 | C16H12   | 006572-60-7 | 59   |
| 3    |    |   | [1,1'-Biphenyl]-2-ol, 5-chloro-     | 204 | C12H9ClO | 000607-12-5 | 53   |
| 4    |    |   | [1,1'-Biphenyl]-2,2'-dicarbonitrile | 204 | C14H8N2  | 004341-02-0 | 50   |
| 5    |    |   | 6-Cyanophenanthridine               | 204 | C14H8N2  | 002176-28-5 | 49   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG072222\  
Data File : BG054367.D  
Acq On : 22 Jul 2022 20:03  
Operator : CG/JU  
Sample : N3814-10  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
ENV-DE-1(0-5)

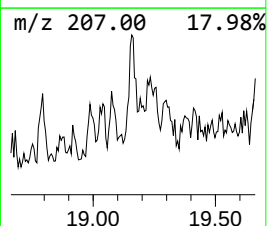
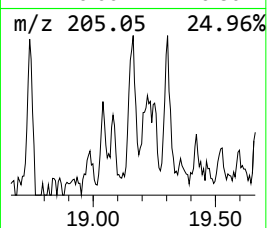
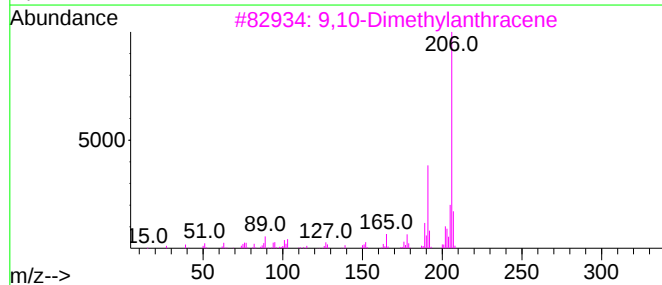
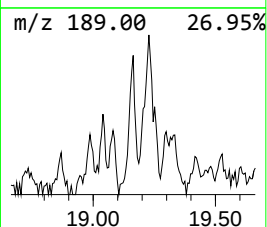
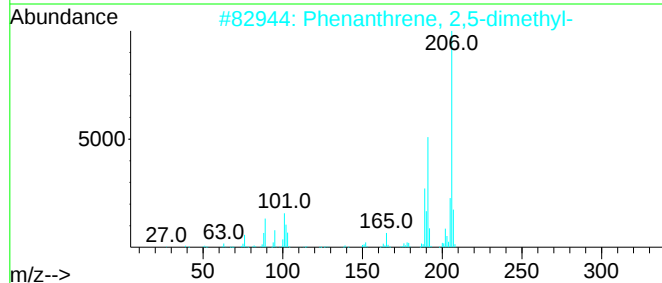
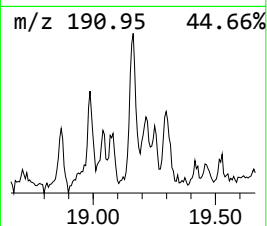
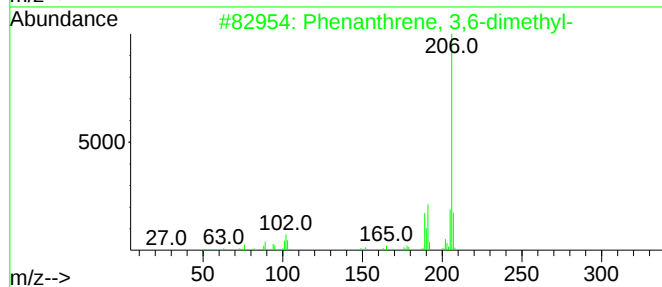
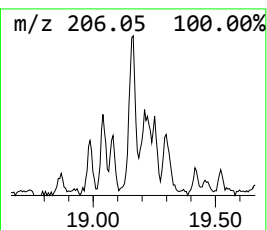
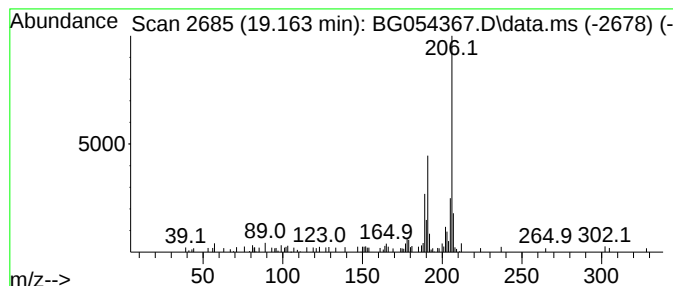
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 8 Phenanthrene, 3,6-dimethyl- Concentration Rank 5

| R.T.   | EstConc | Area  | Relative to ISTD | R.T.   |
|--------|---------|-------|------------------|--------|
| 19.163 | 4.00 ng | 68303 | Phenanthrene-d10 | 17.371 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Phenanthrene, 3,6-dimethyl- | 206 | C16H14  | 001576-67-6 | 91   |
| 2    |    |   | Phenanthrene, 2,5-dimethyl- | 206 | C16H14  | 003674-66-6 | 90   |
| 3    |    |   | 9,10-Dimethylanthracene     | 206 | C16H14  | 000781-43-1 | 87   |
| 4    |    |   | Phenanthrene, 2,7-dimethyl- | 206 | C16H14  | 001576-69-8 | 87   |
| 5    |    |   | Anthracene, 1,4-dimethyl-   | 206 | C16H14  | 000781-92-0 | 87   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG072222\  
Data File : BG054367.D  
Acq On : 22 Jul 2022 20:03  
Operator : CG/JU  
Sample : N3814-10  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

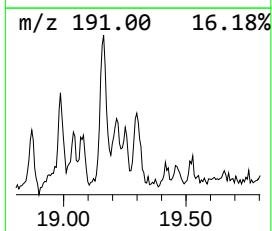
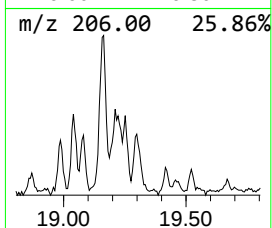
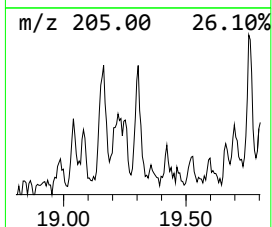
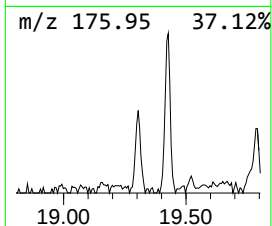
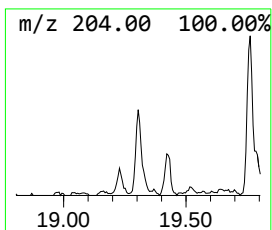
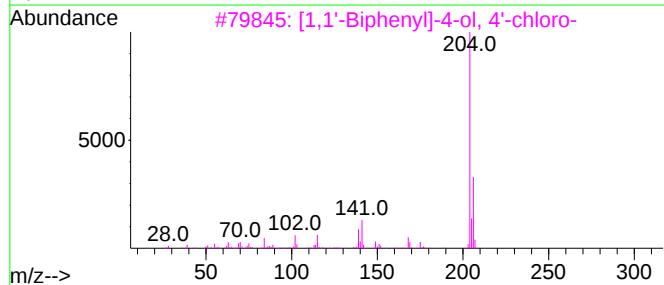
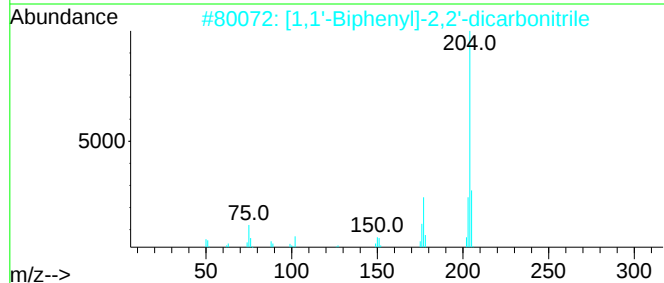
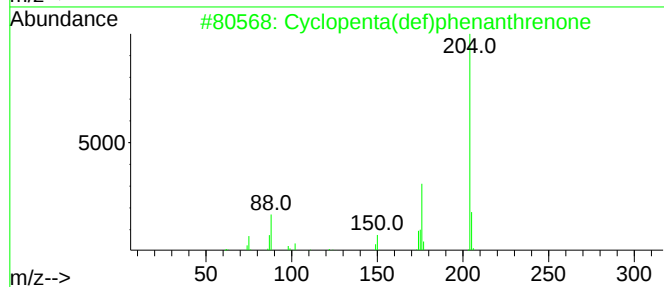
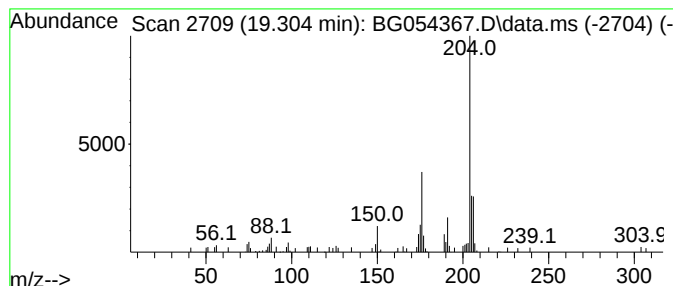
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ClientSampleId :  
ENV-DE-1(0-5)

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 9 Cyclopenta(def)phenanthrenone Concentration Rank 7

| R.T.      | EstConc                             | Area  | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|-------|------------------|--------------|------|
| 19.304    | 3.12 ng                             | 53239 | Phenanthrene-d10 | 17.371       |      |
| Hit# of 5 | Tentative ID                        | MW    | MolForm          | CAS#         | Qual |
| 1         | Cyclopenta(def)phenanthrenone       | 204   | C15H8O           | 005737-13-3  | 68   |
| 2         | [1,1'-Biphenyl]-2,2'-dicarbonitrile | 204   | C14H8N2          | 004341-02-0  | 53   |
| 3         | [1,1'-Biphenyl]-4-ol, 4'-chloro-    | 204   | C12H9ClO         | 028034-99-3  | 47   |
| 4         | 1-Aminocyclopentanecarboxylic ac... | 319   | C15H26ClNO4      | 1000392-62-4 | 47   |
| 5         | L-Rhamnose, (R,R,S,S)-, 4TMS der... | 452   | C18H44O5Si4      | 108392-01-4  | 47   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG072222\  
Data File : BG054367.D  
Acq On : 22 Jul 2022 20:03  
Operator : CG/JU  
Sample : N3814-10  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
ENV-DE-1(0-5)

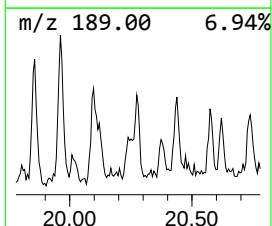
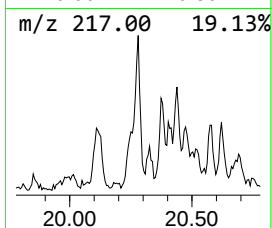
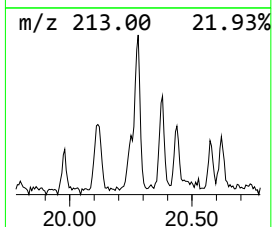
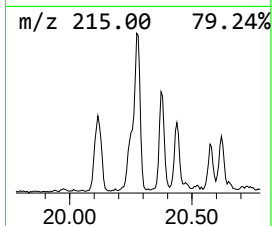
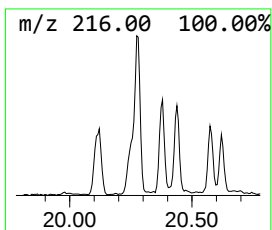
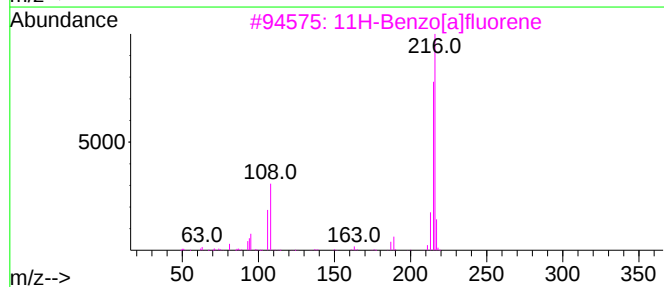
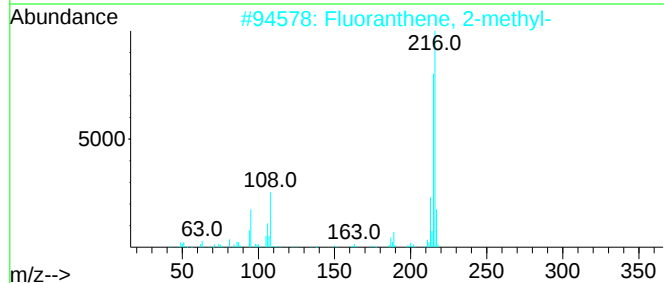
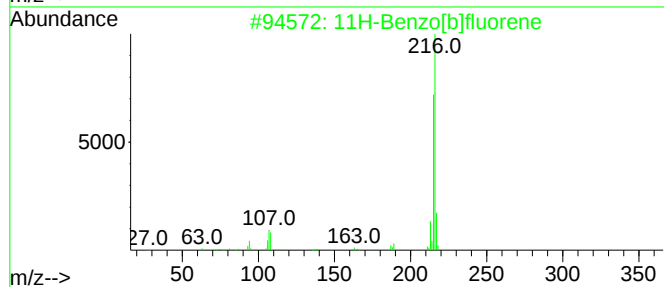
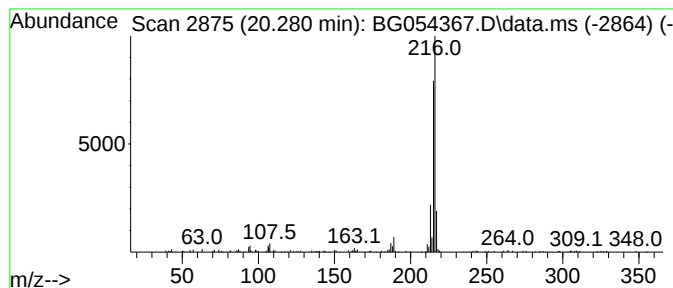
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 10 11H-Benzo[b]fluorene Concentration Rank 6

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 20.280 | 3.68 ng | 190824 | Chrysene-d12     | 21.637 |

| Hit# | of 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------|-----|---------|-------------|------|
| 1    |      | 11H-Benzo[b]fluorene    | 216 | C17H12  | 000243-17-4 | 94   |
| 2    |      | Fluoranthene, 2-methyl- | 216 | C17H12  | 033543-31-6 | 93   |
| 3    |      | 11H-Benzo[a]fluorene    | 216 | C17H12  | 000238-84-6 | 91   |
| 4    |      | Pyrene, 1-methyl-       | 216 | C17H12  | 002381-21-7 | 91   |
| 5    |      | 7H-Benzo[c]fluorene     | 216 | C17H12  | 000205-12-9 | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG072222\  
Data File : BG054367.D  
Acq On : 22 Jul 2022 20:03  
Operator : CG/JU  
Sample : N3814-10  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
ENV-DE-1(0-5)

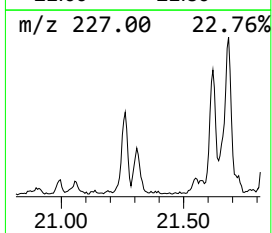
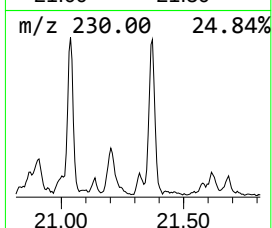
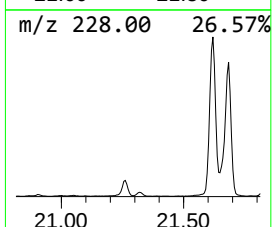
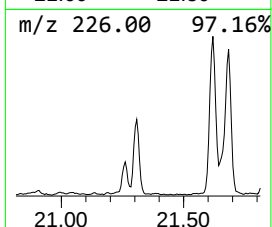
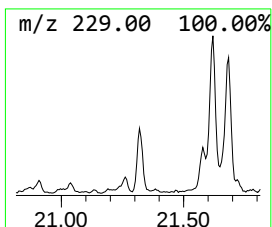
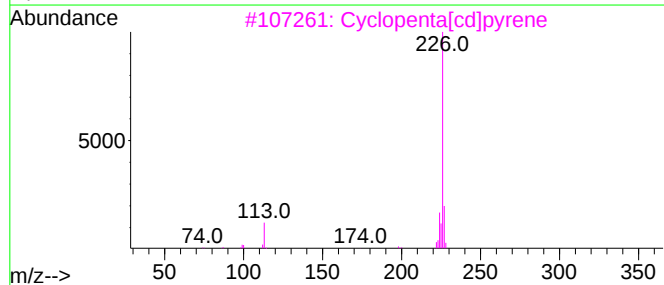
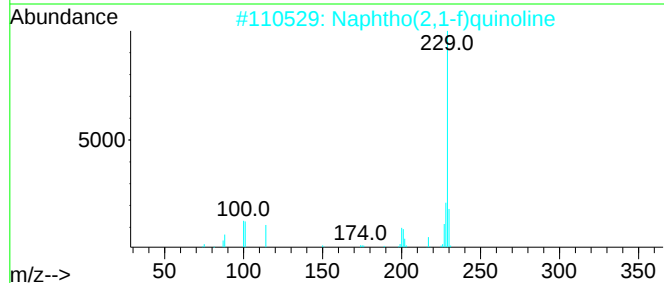
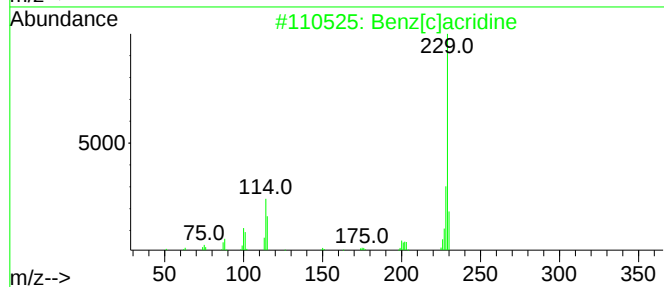
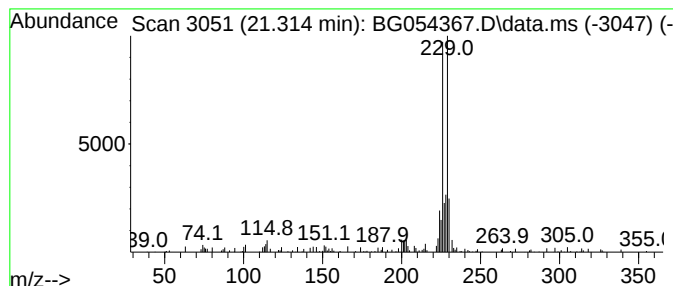
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 11 Benz[c]acridine Concentration Rank 13

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 21.314 | 2.09 ng | 108317 | Chrysene-d12     | 21.637 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Benz[c]acridine                     | 229 | C17H11N   | 000225-51-4  | 50   |
| 2    |    |   | Naphtho(2,1-f)quinoline             | 229 | C17H11N   | 000218-08-6  | 43   |
| 3    |    |   | Cyclopenta[cd]pyrene                | 226 | C18H10    | 027208-37-3  | 38   |
| 4    |    |   | Benzimidazole-5-amine, 1-methyl-... | 229 | C12H11N3S | 021444-73-5  | 35   |
| 5    |    |   | 5-Bromo-6-methoxy-7-azaindole       | 226 | C8H7BrN2O | 1190321-63-1 | 30   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG072222\  
Data File : BG054367.D  
Acq On : 22 Jul 2022 20:03  
Operator : CG/JU  
Sample : N3814-10  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
ENV-DE-1(0-5)

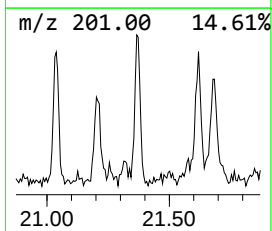
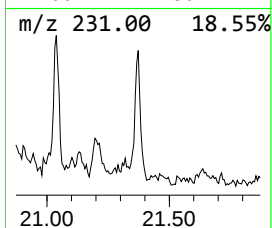
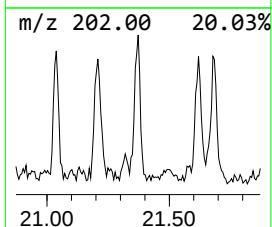
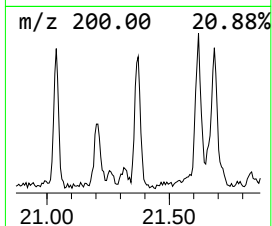
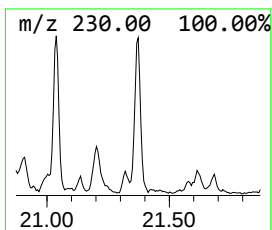
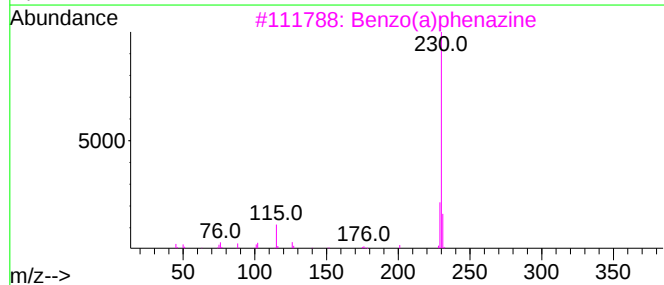
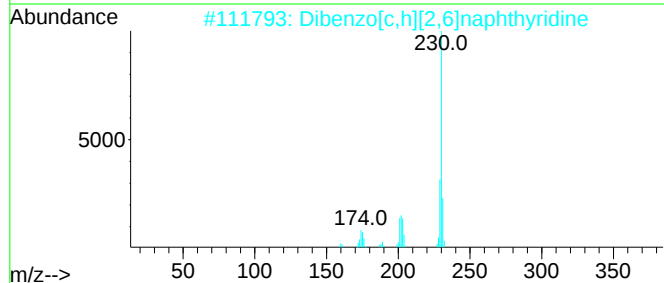
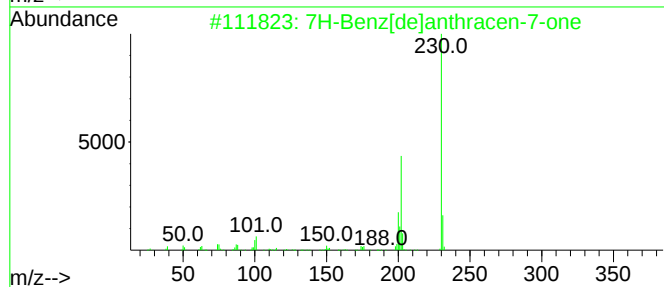
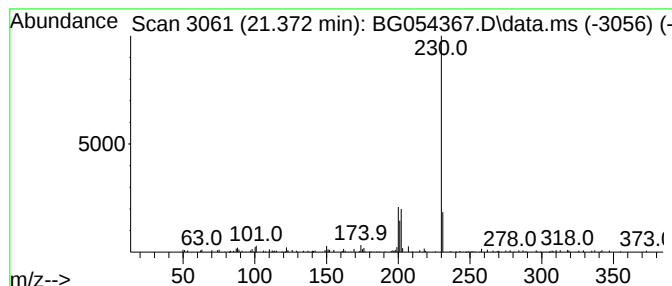
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG071522.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 12 7H-Benz[de]anthracen-7-one Concentration Rank 11

| R.T.      | EstConc                        | Area   | Relative to ISTD |          | R.T.        |      |
|-----------|--------------------------------|--------|------------------|----------|-------------|------|
| 21.372    | 2.15 ng                        | 111634 | Chrysene-d12     |          | 21.637      |      |
| Hit# of 5 | Tentative ID                   |        | MW               | MolForm  | CAS#        | Qual |
| 1         | 7H-Benz[de]anthracen-7-one     |        | 230              | C17H10O  | 000082-05-3 | 64   |
| 2         | Dibenzo[c,h][2,6]naphthyridine |        | 230              | C16H10N2 | 000218-30-4 | 53   |
| 3         | Benzo(a)phenazine              |        | 230              | C16H10N2 | 000225-61-6 | 47   |
| 4         | p-Terphenyl                    |        | 230              | C18H14   | 000092-94-4 | 47   |
| 5         | 11H-Benzo[a]fluoren-11-one     |        | 230              | C17H10O  | 000479-79-8 | 43   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG072222\  
Data File : BG054367.D  
Acq On : 22 Jul 2022 20:03  
Operator : CG/JU  
Sample : N3814-10  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
ENV-DE-1(0-5)

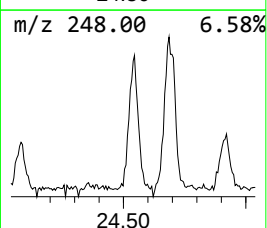
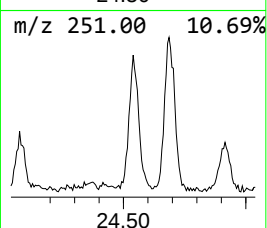
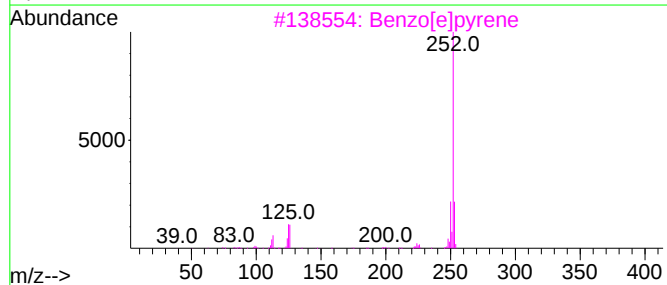
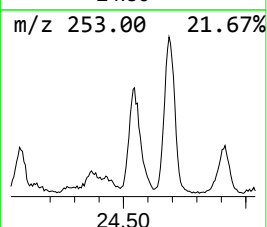
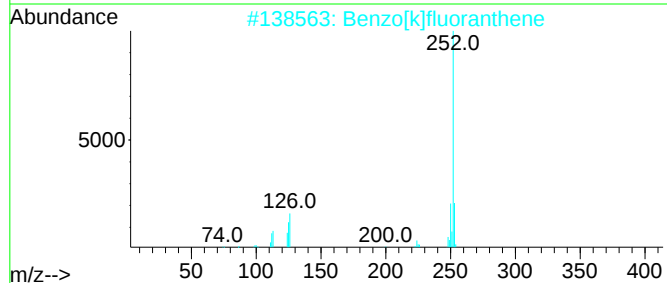
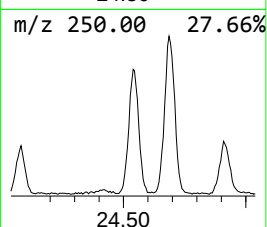
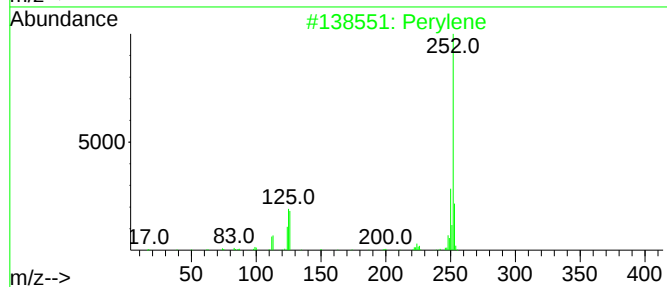
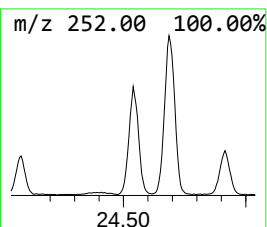
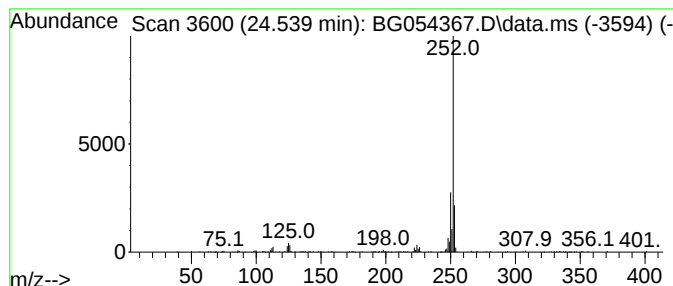
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG071522.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 13 Perylene Concentration Rank 1

| R.T.   | EstConc  | Area   | Relative to ISTD | R.T.   |
|--------|----------|--------|------------------|--------|
| 24.539 | 16.49 ng | 345710 | Perylene-d12     | 24.839 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm  | CAS#        | Qual |
|------|----|---|----------------------------|-----|----------|-------------|------|
| 1    |    |   | Perylene                   | 252 | C20H12   | 000198-55-0 | 93   |
| 2    |    |   | Benzo[k]fluoranthene       | 252 | C20H12   | 000207-08-9 | 89   |
| 3    |    |   | Benzo[e]pyrene             | 252 | C20H12   | 000192-97-2 | 87   |
| 4    |    |   | Benzo[b]fluoranthene       | 252 | C20H12   | 000205-99-2 | 81   |
| 5    |    |   | (1-Cyclopentenyl)ferrocene | 252 | C15H16Fe | 012260-67-2 | 78   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG072222\  
Data File : BG054367.D  
Acq On : 22 Jul 2022 20:03  
Operator : CG/JU  
Sample : N3814-10  
Misc :  
ALS Vial : 12 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
ENV-DE-1(0-5)

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 |
| Anthracene, 2-m... | 18.247 | 4.8     | ng    | 81869    | 4                     | 17.371 | 341434  | 20.0 |
| Phenanthrene, 1... | 18.294 | 6.1     | ng    | 103560   | 4                     | 17.371 | 341434  | 20.0 |
| Phenanthrene, 2... | 18.376 | 2.4     | ng    | 41217    | 4                     | 17.371 | 341434  | 20.0 |
| Benzaldehyde, 3... | 18.441 | 8.0     | ng    | 136330   | 4                     | 17.371 | 341434  | 20.0 |
| Phenanthrene, 4... | 18.476 | 2.7     | ng    | 45448    | 4                     | 17.371 | 341434  | 20.0 |
| Naphthalene, 2-... | 18.740 | 3.1     | ng    | 53220    | 4                     | 17.371 | 341434  | 20.0 |
| Phenanthrene, 3... | 19.163 | 4.0     | ng    | 68303    | 4                     | 17.371 | 341434  | 20.0 |
| Cyclopenta(def)... | 19.304 | 3.1     | ng    | 53239    | 4                     | 17.371 | 341434  | 20.0 |
| 11H-Benzo[b]flu... | 20.280 | 3.7     | ng    | 190824   | 5                     | 21.637 | 1037610 | 20.0 |
| Benz[c]acridine    | 21.314 | 2.1     | ng    | 108317   | 5                     | 21.637 | 1037610 | 20.0 |
| 7H-Benz[de]anth... | 21.372 | 2.1     | ng    | 111634   | 5                     | 21.637 | 1037610 | 20.0 |
| Perylene           | 24.539 | 16.5    | ng    | 345710   | 6                     | 24.839 | 419285  | 20.0 |