

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG110121\  
 Data File : BG050799.D  
 Acq On : 1 Nov 2021 15:12  
 Operator : CG/JU  
 Sample : M4422-01 2X  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 OR-03-102921

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-BG100621.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG050799.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         | 5.774       | 383           | 389         | 398          | rBV      | 306131         | 524283        | 23.86%          | 1.765%        |
| 2         | 7.384       | 654           | 663         | 672          | rBV      | 316162         | 576398        | 26.23%          | 1.940%        |
| 3         | 8.236       | 802           | 808         | 815          | rVB      | 155390         | 276189        | 12.57%          | 0.930%        |
| 4         | 9.411       | 1000          | 1008        | 1020         | rVB      | 197642         | 367253        | 16.71%          | 1.236%        |
| 5         | 10.809      | 1240          | 1246        | 1255         | rBV2     | 22487          | 45099         | 2.05%           | 0.152%        |
| 6         | 11.003      | 1271          | 1279        | 1283         | rBV2     | 14765          | 26327         | 1.20%           | 0.089%        |
| 7         | 11.068      | 1283          | 1290        | 1295         | rVV      | 216222         | 417502        | 19.00%          | 1.405%        |
| 8         | 11.115      | 1295          | 1298        | 1305         | rVV      | 38618          | 63735         | 2.90%           | 0.215%        |
| 9         | 12.431      | 1517          | 1522        | 1529         | rBV      | 26034          | 40123         | 1.83%           | 0.135%        |
| 10        | 12.701      | 1564          | 1568        | 1574         | rVB2     | 28040          | 49431         | 2.25%           | 0.166%        |
| 11        | 12.919      | 1598          | 1605        | 1615         | rBV2     | 27495          | 57563         | 2.62%           | 0.194%        |
| 12        | 13.371      | 1676          | 1682        | 1688         | rBV3     | 17495          | 28338         | 1.29%           | 0.095%        |
| 13        | 13.483      | 1692          | 1701        | 1712         | rVB      | 452548         | 705107        | 32.08%          | 2.374%        |
| 14        | 13.659      | 1724          | 1731        | 1735         | rBV      | 42912          | 73506         | 3.34%           | 0.247%        |
| 15        | 14.182      | 1815          | 1820        | 1825         | rBV2     | 45539          | 77490         | 3.53%           | 0.261%        |
| 16        | 14.229      | 1825          | 1828        | 1833         | rVB2     | 25720          | 40475         | 1.84%           | 0.136%        |
| 17        | 14.311      | 1838          | 1842        | 1846         | rVB3     | 31258          | 46407         | 2.11%           | 0.156%        |
| 18        | 14.423      | 1856          | 1861        | 1870         | rVB3     | 25692          | 53898         | 2.45%           | 0.181%        |
| 19        | 14.587      | 1883          | 1889        | 1897         | rBV2     | 62819          | 126827        | 5.77%           | 0.427%        |
| 20        | 14.722      | 1908          | 1912        | 1916         | rVB2     | 48274          | 64772         | 2.95%           | 0.218%        |
| 21        | 14.858      | 1921          | 1935        | 1941         | rBV      | 331646         | 584324        | 26.59%          | 1.967%        |
| 22        | 14.922      | 1941          | 1946        | 1953         | rVV      | 94514          | 153974        | 7.01%           | 0.518%        |
| 23        | 14.993      | 1953          | 1958        | 1964         | rVB7     | 11309          | 23591         | 1.07%           | 0.079%        |
| 24        | 15.063      | 1964          | 1970        | 1978         | rVB7     | 18794          | 42785         | 1.95%           | 0.144%        |
| 25        | 15.163      | 1980          | 1987        | 1992         | rBV9     | 14133          | 37913         | 1.73%           | 0.128%        |
| 26        | 15.251      | 1995          | 2002        | 2010         | rVB2     | 73974          | 148258        | 6.75%           | 0.499%        |
| 27        | 15.322      | 2010          | 2014        | 2023         | rBV5     | 37125          | 69257         | 3.15%           | 0.233%        |
| 28        | 15.469      | 2032          | 2039        | 2043         | rBV3     | 34918          | 73886         | 3.36%           | 0.249%        |
| 29        | 15.522      | 2044          | 2048        | 2052         | rVB      | 21650          | 33816         | 1.54%           | 0.114%        |
| 30        | 15.639      | 2061          | 2068        | 2070         | rBV5     | 22059          | 49122         | 2.24%           | 0.165%        |
| 31        | 15.668      | 2070          | 2073        | 2078         | rVB2     | 76993          | 98808         | 4.50%           | 0.333%        |
| 32        | 15.903      | 2107          | 2113        | 2123         | rVB2     | 127840         | 225604        | 10.27%          | 0.759%        |
| 33        | 16.062      | 2136          | 2140        | 2145         | rVV5     | 41395          | 73834         | 3.36%           | 0.249%        |
| 34        | 16.191      | 2154          | 2162        | 2170         | rVB4     | 33251          | 83556         | 3.80%           | 0.281%        |
| 35        | 16.344      | 2182          | 2188        | 2195         | rVB      | 368108         | 570386        | 25.95%          | 1.920%        |
| 36        | 16.532      | 2215          | 2220        | 2222         | rBV      | 77198          | 114194        | 5.20%           | 0.384%        |

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 Sample : M4422-01 2X  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 OR-03-102921

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-BG100621.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 37 | 16.697 | 2245 | 2248 | 2253 | rBV4 | 17977   | 30860   | 1.40%  | 0.104% |
| 38 | 16.902 | 2275 | 2283 | 2287 | rBV2 | 48687   | 103455  | 4.71%  | 0.348% |
| 39 | 16.955 | 2289 | 2292 | 2296 | rVB2 | 28142   | 39109   | 1.78%  | 0.132% |
| 40 | 17.049 | 2305 | 2308 | 2312 | rVB3 | 29347   | 47350   | 2.15%  | 0.159% |
| 41 | 17.126 | 2314 | 2321 | 2325 | rBV5 | 25651   | 73898   | 3.36%  | 0.249% |
| 42 | 17.261 | 2340 | 2344 | 2348 | rVB4 | 21985   | 37362   | 1.70%  | 0.126% |
| 43 | 17.319 | 2350 | 2354 | 2359 | rVV2 | 88255   | 131378  | 5.98%  | 0.442% |
| 44 | 17.372 | 2360 | 2363 | 2367 | rVV  | 40175   | 57404   | 2.61%  | 0.193% |
| 45 | 17.425 | 2367 | 2372 | 2377 | rVB  | 71655   | 116503  | 5.30%  | 0.392% |
| 46 | 17.607 | 2397 | 2403 | 2406 | rBV  | 380959  | 634056  | 28.85% | 2.134% |
| 47 | 17.648 | 2406 | 2410 | 2416 | rVV  | 905288  | 1362886 | 62.02% | 4.588% |
| 48 | 17.737 | 2420 | 2425 | 2430 | rVV  | 229327  | 367388  | 16.72% | 1.237% |
| 49 | 17.795 | 2432 | 2435 | 2439 | rVB2 | 41447   | 58442   | 2.66%  | 0.197% |
| 50 | 17.866 | 2444 | 2447 | 2452 | rBV6 | 14231   | 27668   | 1.26%  | 0.093% |
| 51 | 18.007 | 2466 | 2471 | 2476 | rBV2 | 43090   | 81620   | 3.71%  | 0.275% |
| 52 | 18.060 | 2476 | 2480 | 2486 | rVB  | 74233   | 105245  | 4.79%  | 0.354% |
| 53 | 18.113 | 2486 | 2489 | 2494 | rBV  | 42935   | 58709   | 2.67%  | 0.198% |
| 54 | 18.183 | 2497 | 2501 | 2506 | rVV  | 68198   | 103335  | 4.70%  | 0.348% |
| 55 | 18.236 | 2506 | 2510 | 2517 | rVB  | 230559  | 295884  | 13.46% | 0.996% |
| 56 | 18.336 | 2521 | 2527 | 2534 | rVB  | 42831   | 88195   | 4.01%  | 0.297% |
| 57 | 18.477 | 2545 | 2551 | 2555 | rBV  | 201097  | 343456  | 15.63% | 1.156% |
| 58 | 18.524 | 2555 | 2559 | 2564 | rVB  | 255269  | 391900  | 17.83% | 1.319% |
| 59 | 18.606 | 2568 | 2573 | 2577 | rBV  | 92742   | 138366  | 6.30%  | 0.466% |
| 60 | 18.671 | 2577 | 2584 | 2588 | rBV2 | 294773  | 622423  | 28.32% | 2.095% |
| 61 | 18.706 | 2588 | 2590 | 2594 | rVB  | 153065  | 175721  | 8.00%  | 0.592% |
| 62 | 18.747 | 2594 | 2597 | 2600 | rBV  | 51919   | 56820   | 2.59%  | 0.191% |
| 63 | 18.871 | 2614 | 2618 | 2621 | rVB  | 38455   | 52960   | 2.41%  | 0.178% |
| 64 | 18.965 | 2629 | 2634 | 2638 | rBV  | 156018  | 225350  | 10.25% | 0.759% |
| 65 | 19.017 | 2639 | 2643 | 2652 | rVB5 | 73587   | 150874  | 6.87%  | 0.508% |
| 66 | 19.094 | 2652 | 2656 | 2658 | rBV  | 20122   | 30989   | 1.41%  | 0.104% |
| 67 | 19.205 | 2668 | 2675 | 2680 | rBV5 | 73835   | 149981  | 6.82%  | 0.505% |
| 68 | 19.258 | 2680 | 2684 | 2688 | rBV  | 81476   | 107511  | 4.89%  | 0.362% |
| 69 | 19.299 | 2688 | 2691 | 2695 | rVB2 | 63729   | 79964   | 3.64%  | 0.269% |
| 70 | 19.370 | 2698 | 2703 | 2705 | rBV  | 934821  | 1240859 | 56.46% | 4.177% |
| 71 | 19.399 | 2705 | 2708 | 2712 | rVB2 | 544944  | 681112  | 30.99% | 2.293% |
| 72 | 19.529 | 2724 | 2730 | 2743 | rBV4 | 202663  | 415204  | 18.89% | 1.398% |
| 73 | 19.646 | 2744 | 2750 | 2756 | rBV  | 1307183 | 2041675 | 92.90% | 6.873% |
| 74 | 19.699 | 2756 | 2759 | 2762 | rVV2 | 57252   | 76599   | 3.49%  | 0.258% |
| 75 | 19.740 | 2762 | 2766 | 2771 | rVB2 | 59795   | 82556   | 3.76%  | 0.278% |
| 76 | 19.793 | 2773 | 2775 | 2779 | rVB  | 33313   | 38222   | 1.74%  | 0.129% |

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 Sample : M4422-01 2X  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 OR-03-102921

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-BG100621.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |         |        |
|----|--------|------|------|------|------|---------|---------|---------|--------|
| 77 | 19.887 | 2786 | 2791 | 2798 | rBV3 | 77751   | 170070  | 7.74%   | 0.573% |
| 78 | 19.975 | 2802 | 2806 | 2807 | rVV  | 203309  | 254491  | 11.58%  | 0.857% |
| 79 | 20.010 | 2807 | 2812 | 2819 | rVV  | 1419812 | 2197647 | 100.00% | 7.398% |
| 80 | 20.075 | 2819 | 2823 | 2828 | rVB  | 121045  | 176380  | 8.03%   | 0.594% |
| 81 | 20.193 | 2838 | 2843 | 2847 | rVB  | 679551  | 894273  | 40.69%  | 3.010% |
| 82 | 20.234 | 2847 | 2850 | 2858 | rVB3 | 53899   | 110776  | 5.04%   | 0.373% |
| 83 | 20.334 | 2860 | 2867 | 2873 | rBV4 | 150965  | 353647  | 16.09%  | 1.190% |
| 84 | 20.492 | 2885 | 2894 | 2901 | rVB  | 291033  | 696646  | 31.70%  | 2.345% |
| 85 | 20.592 | 2907 | 2911 | 2915 | rBV  | 204894  | 275322  | 12.53%  | 0.927% |
| 86 | 20.651 | 2917 | 2921 | 2925 | rVB  | 194890  | 266120  | 12.11%  | 0.896% |
| 87 | 20.792 | 2941 | 2945 | 2949 | rBV  | 193726  | 275685  | 12.54%  | 0.928% |
| 88 | 20.839 | 2949 | 2953 | 2961 | rVB2 | 180749  | 320578  | 14.59%  | 1.079% |
| 89 | 21.274 | 3023 | 3027 | 3033 | rVB2 | 105484  | 189810  | 8.64%   | 0.639% |
| 90 | 21.468 | 3057 | 3060 | 3063 | rBV  | 91015   | 112793  | 5.13%   | 0.380% |
| 91 | 21.509 | 3064 | 3067 | 3072 | rVB2 | 136576  | 202403  | 9.21%   | 0.681% |
| 92 | 21.562 | 3072 | 3076 | 3081 | rBV3 | 101810  | 170598  | 7.76%   | 0.574% |
| 93 | 21.891 | 3125 | 3132 | 3139 | rBV2 | 671350  | 1799930 | 81.90%  | 6.059% |
| 94 | 21.955 | 3139 | 3143 | 3155 | rVB  | 540305  | 983960  | 44.77%  | 3.312% |
| 95 | 22.114 | 3166 | 3170 | 3175 | rBV  | 85535   | 137782  | 6.27%   | 0.464% |
| 96 | 24.229 | 3522 | 3530 | 3538 | rBV  | 283593  | 1006012 | 45.78%  | 3.387% |
| 97 | 24.999 | 3654 | 3661 | 3676 | rVB2 | 207026  | 648296  | 29.50%  | 2.182% |
| 98 | 25.157 | 3679 | 3688 | 3700 | rVB  | 321697  | 989157  | 45.01%  | 3.330% |
| 99 | 25.316 | 3707 | 3715 | 3723 | rBV2 | 174661  | 506644  | 23.05%  | 1.706% |

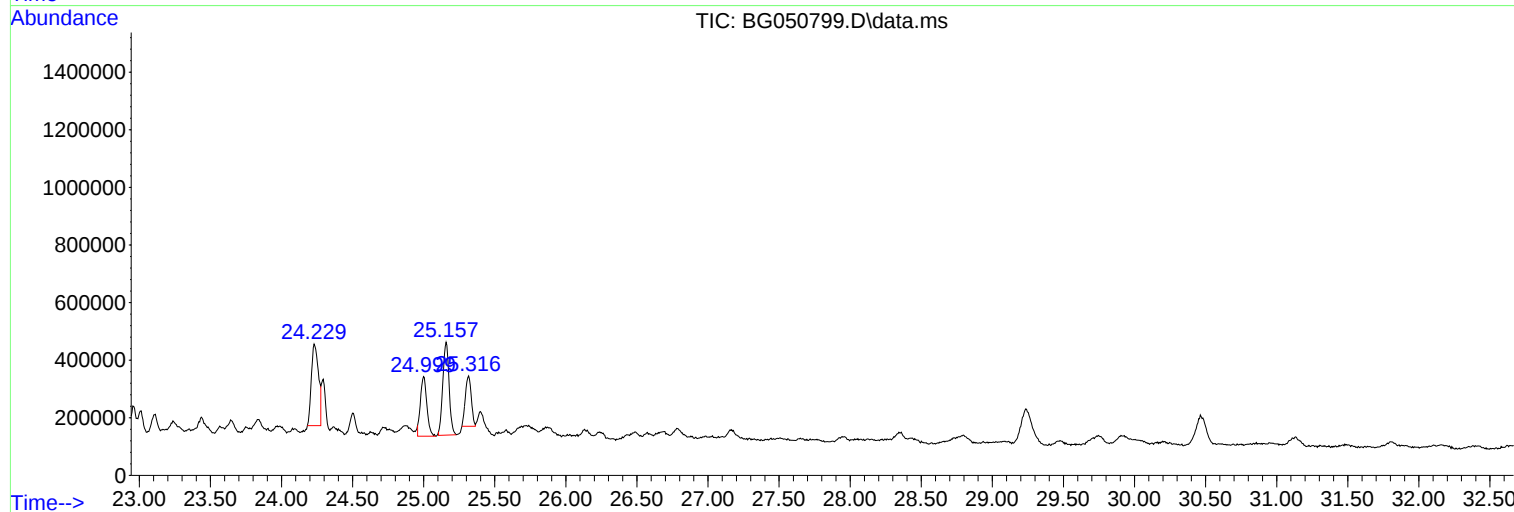
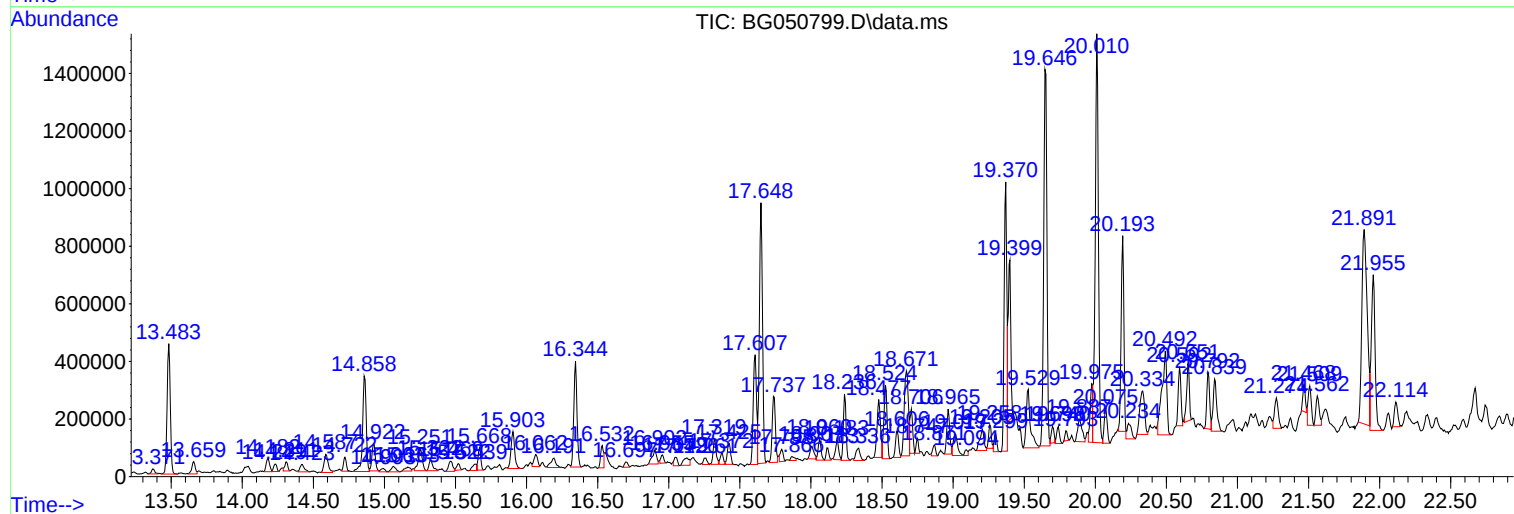
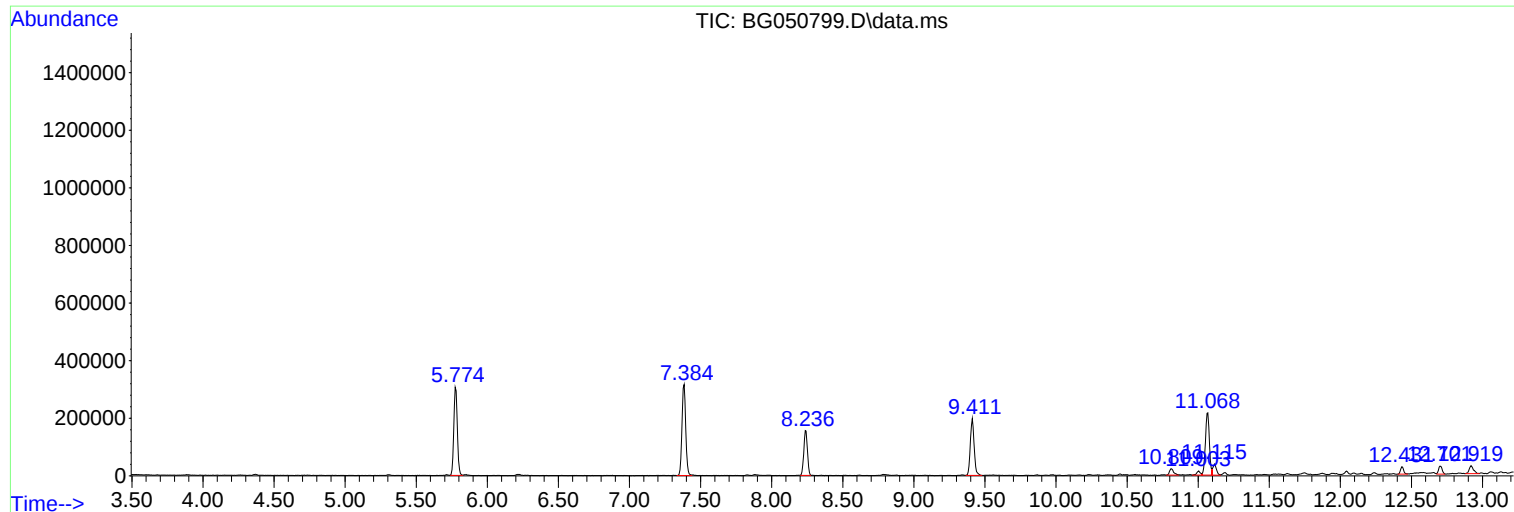
Sum of corrected areas: 29706340

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG110121\  
Data File : BG050799.D  
Acq On : 1 Nov 2021 15:12  
Operator : CG/JU  
Sample : M4422-01 2X  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
OR-03-102921

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG100621.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\BG110121\  
Data File : BG050799.D  
Acq On : 1 Nov 2021 15:12  
Operator : CG/JU  
Sample : M4422-01 2X  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
OR-03-102921

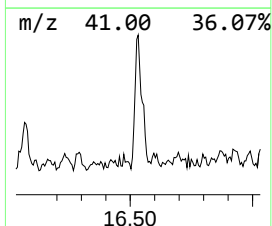
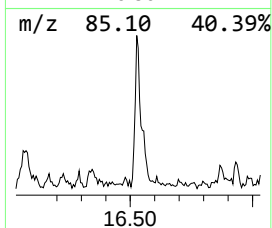
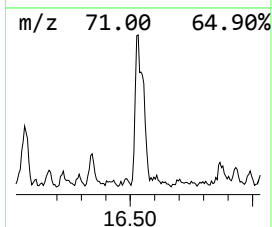
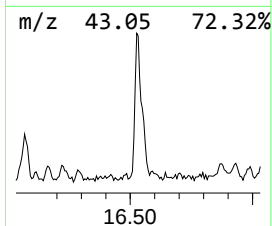
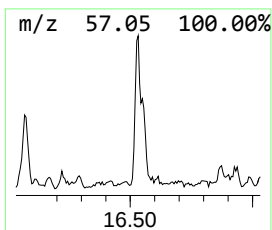
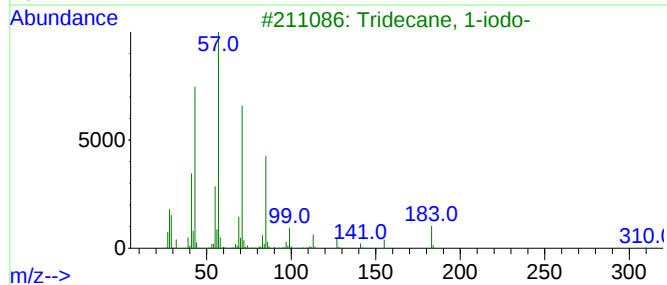
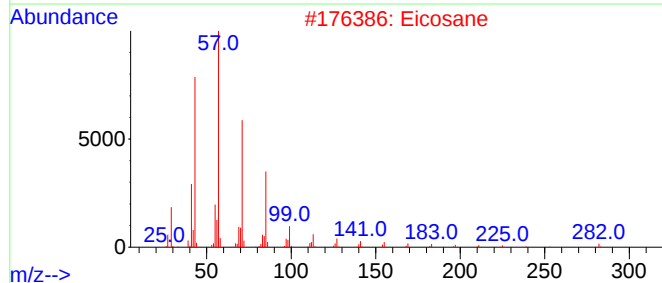
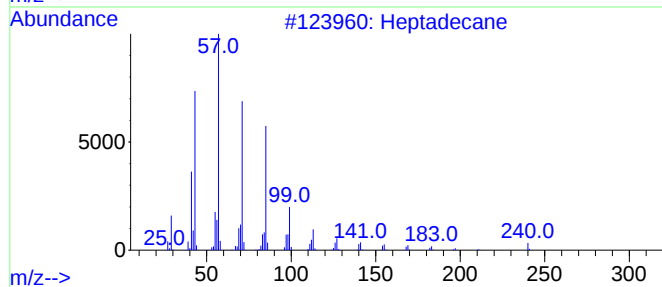
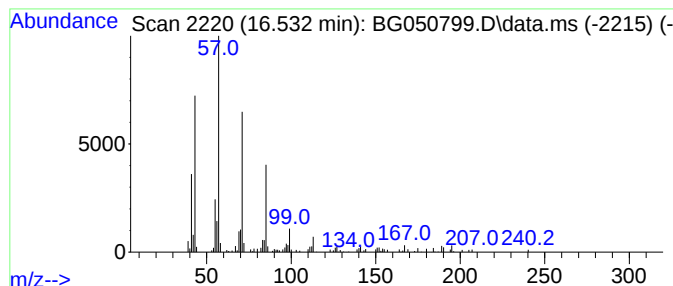
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG100621.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 1 Heptadecane Concentration Rank 18

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 16.532 | 3.60 ng | 114194 | Phenanthrene-d10 | 17.607 |

| Hit# | of | 5 | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------------|-----|---------|-------------|------|
| 1    |    |   | Heptadecane                  | 240 | C17H36  | 000629-78-7 | 95   |
| 2    |    |   | Eicosane                     | 282 | C20H42  | 000112-95-8 | 90   |
| 3    |    |   | Tridecane, 1-iodo-           | 310 | C13H27I | 035599-77-0 | 90   |
| 4    |    |   | Heptacosane                  | 380 | C27H56  | 000593-49-7 | 90   |
| 5    |    |   | Dodecane, 2-methyl-6-propyl- | 226 | C16H34  | 055045-08-4 | 86   |



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ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
OR-03-102921

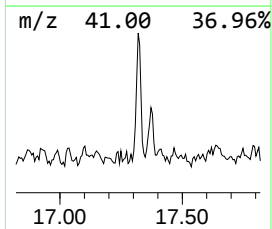
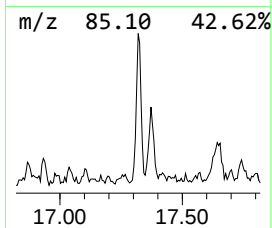
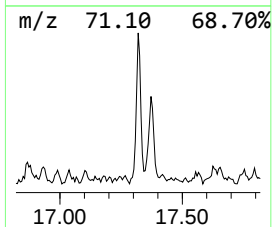
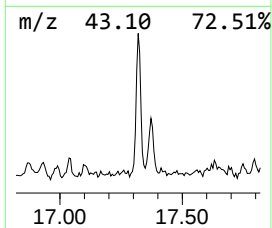
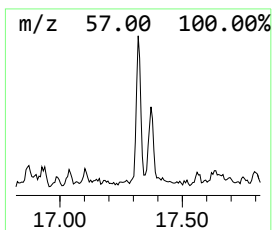
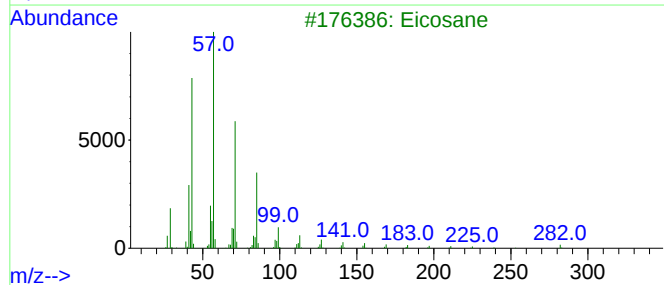
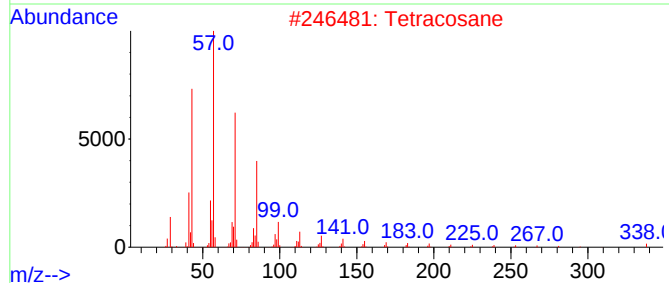
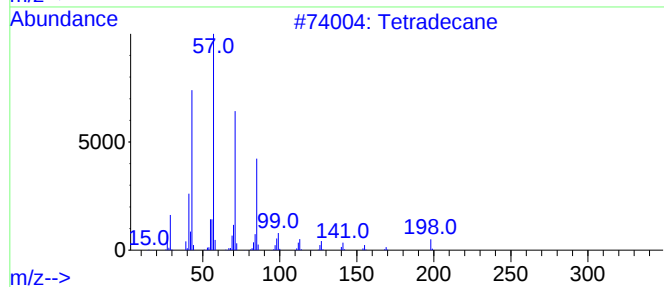
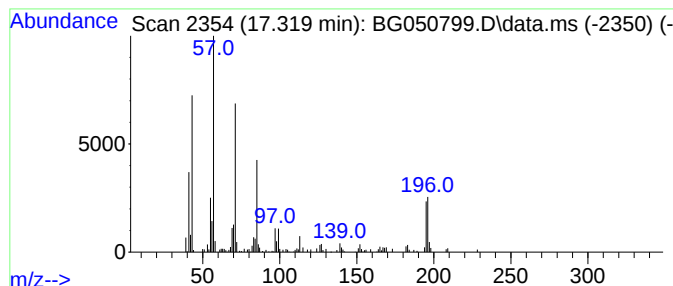
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG100621.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 2 Tetradecane Concentration Rank 15

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 17.320 | 4.14 ng | 131378 | Phenanthrene-d10 | 17.607 |

| Hit# | of | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------|-----|---------|-------------|------|
| 1    |    |   | Tetradecane  | 198 | C14H30  | 000629-59-4 | 64   |
| 2    |    |   | Tetracosane  | 338 | C24H50  | 000646-31-1 | 62   |
| 3    |    |   | Eicosane     | 282 | C20H42  | 000112-95-8 | 58   |
| 4    |    |   | Nonadecane   | 268 | C19H40  | 000629-92-5 | 58   |
| 5    |    |   | Octacosane   | 394 | C28H58  | 000630-02-4 | 58   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG110121\  
Data File : BG050799.D  
Acq On : 1 Nov 2021 15:12  
Operator : CG/JU  
Sample : M4422-01 2X  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
OR-03-102921

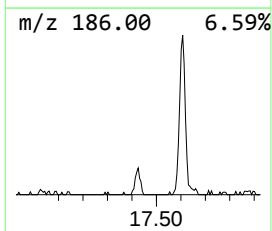
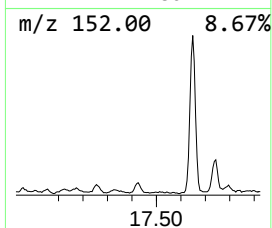
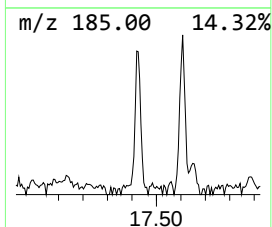
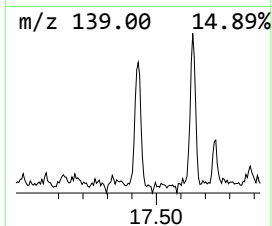
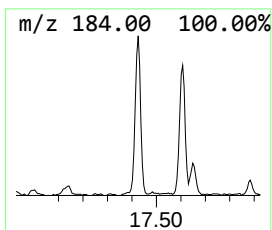
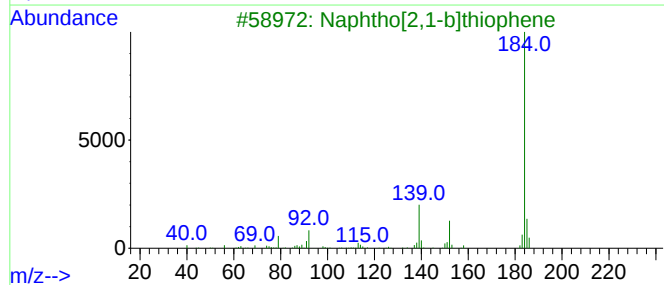
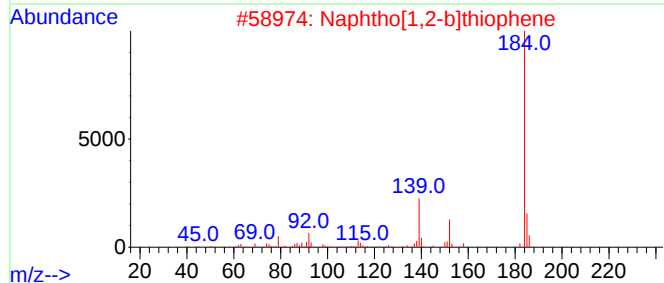
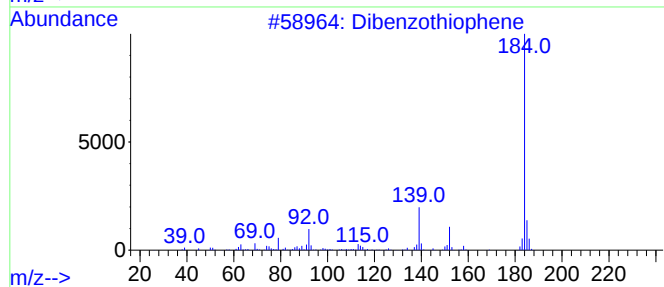
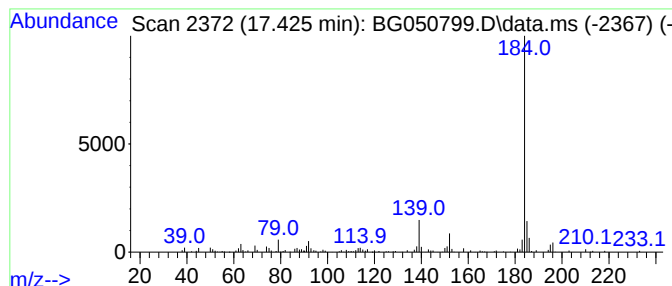
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 3 Dibenzothiophene Concentration Rank 17

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 17.425 | 3.67 ng | 116503 | Phenanthrene-d10 | 17.607 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Dibenzothiophene        | 184 | C12H8S  | 000132-65-0 | 95   |
| 2    |    |   | Naphtho[1,2-b]thiophene | 184 | C12H8S  | 000234-41-3 | 94   |
| 3    |    |   | Naphtho[2,1-b]thiophene | 184 | C12H8S  | 000233-02-3 | 93   |
| 4    |    |   | Azuleno(2,1-b)thiophene | 184 | C12H8S  | 000248-13-5 | 87   |
| 5    |    |   | Naphtho[2,3-b]thiophene | 184 | C12H8S  | 000268-77-9 | 87   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG110121\  
Data File : BG050799.D  
Acq On : 1 Nov 2021 15:12  
Operator : CG/JU  
Sample : M4422-01 2X  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
OR-03-102921

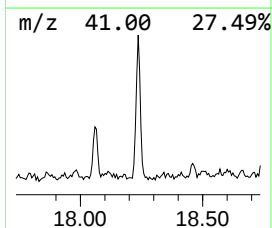
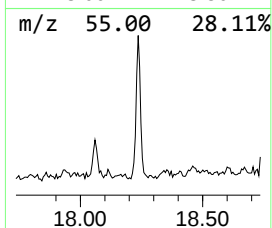
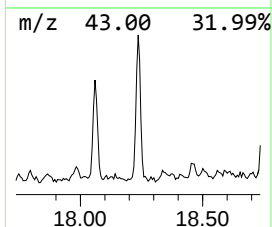
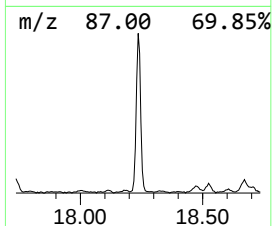
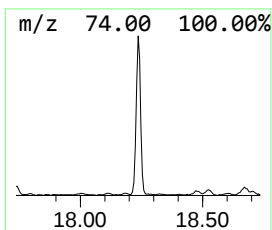
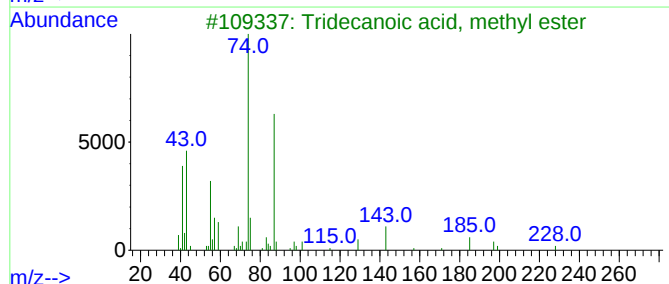
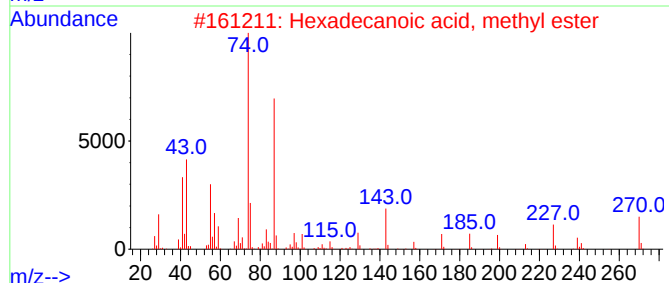
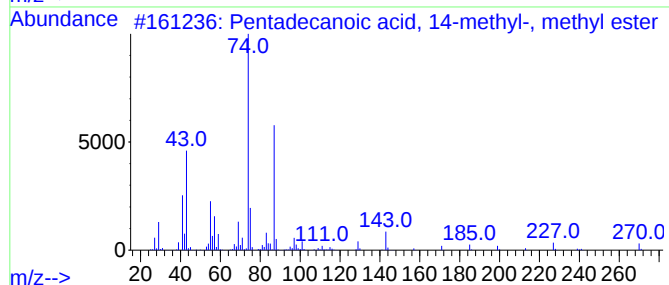
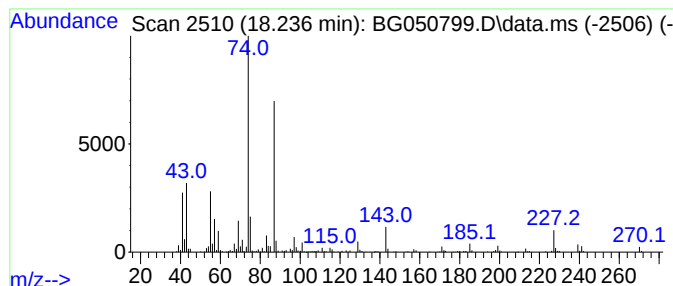
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 4 Pentadecanoic acid, 14-meth... Concentration Rank 8

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 18.236 | 9.33 ng | 295884 | Phenanthrene-d10 | 17.607 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Pentadecanoic acid, 14-methyl-, ... | 270 | C17H34O2 | 005129-60-2 | 96   |
| 2    |    |   | Hexadecanoic acid, methyl ester     | 270 | C17H34O2 | 000112-39-0 | 91   |
| 3    |    |   | Tridecanoic acid, methyl ester      | 228 | C14H28O2 | 001731-88-0 | 90   |
| 4    |    |   | Methyl tetradecanoate               | 242 | C15H30O2 | 000124-10-7 | 86   |
| 5    |    |   | Dodecanoic acid, methyl ester       | 214 | C13H26O2 | 000111-82-0 | 86   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG110121\  
Data File : BG050799.D  
Acq On : 1 Nov 2021 15:12  
Operator : CG/JU  
Sample : M4422-01 2X  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
OR-03-102921

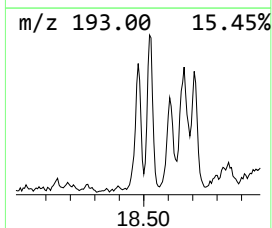
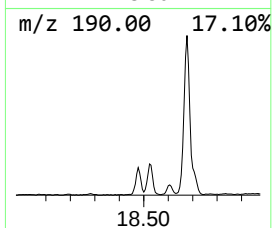
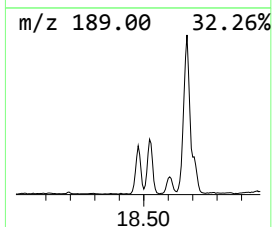
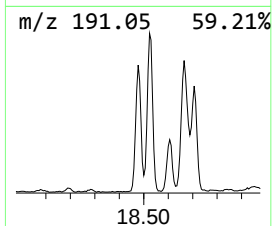
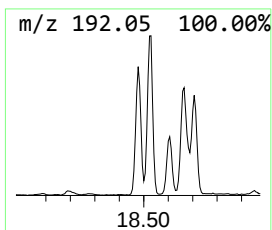
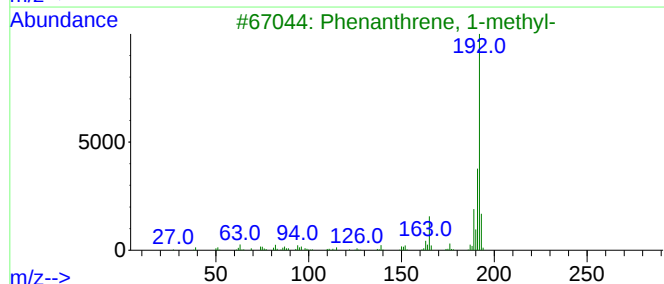
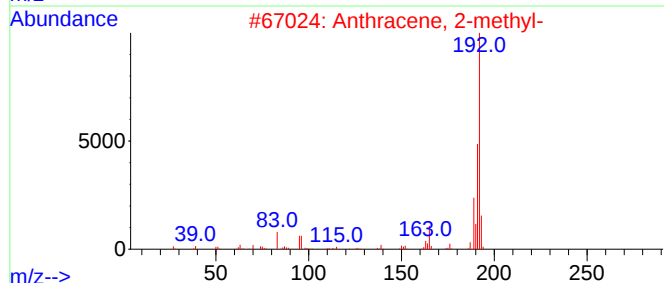
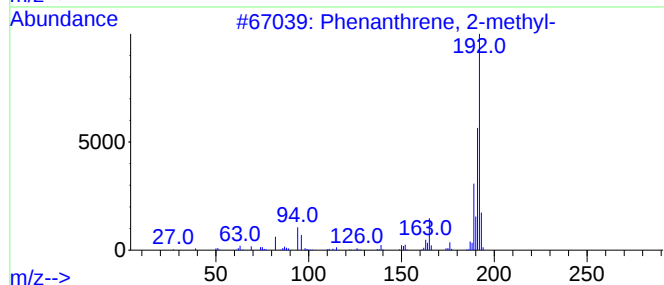
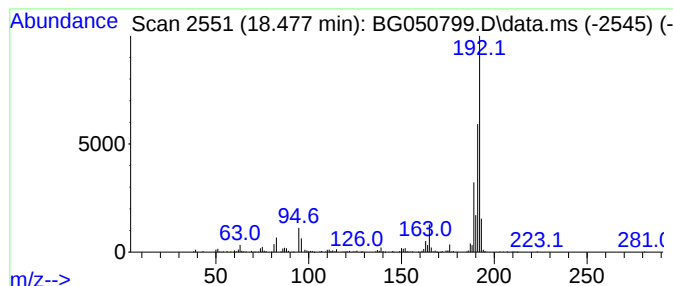
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 5 Phenanthrene, 2-methyl- Concentration Rank 7

| R.T.   | EstConc  | Area   | Relative to ISTD | R.T.   |
|--------|----------|--------|------------------|--------|
| 18.477 | 10.83 ng | 343456 | Phenanthrene-d10 | 17.607 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Phenanthrene, 2-methyl-             | 192 | C15H12  | 002531-84-2 | 97   |
| 2    |    |   | Anthracene, 2-methyl-               | 192 | C15H12  | 000613-12-7 | 96   |
| 3    |    |   | Phenanthrene, 1-methyl-             | 192 | C15H12  | 000832-69-9 | 95   |
| 4    |    |   | 1H-Cyclopropa[1]phenanthrene,1a,... | 192 | C15H12  | 000949-41-7 | 95   |
| 5    |    |   | Anthracene, 9-methyl-               | 192 | C15H12  | 000779-02-2 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG110121\  
Data File : BG050799.D  
Acq On : 1 Nov 2021 15:12  
Operator : CG/JU  
Sample : M4422-01 2X  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
OR-03-102921

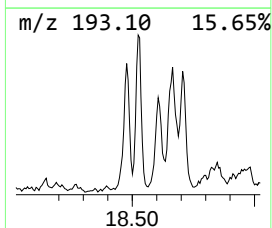
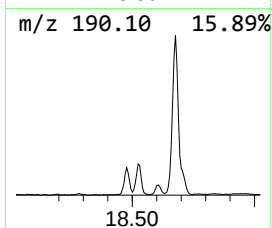
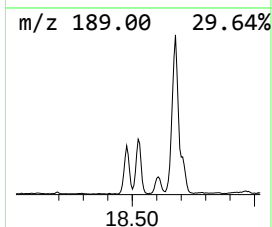
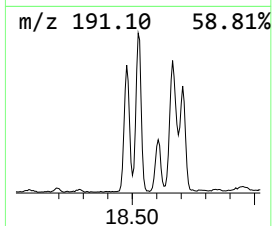
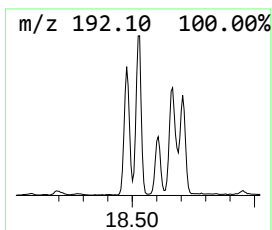
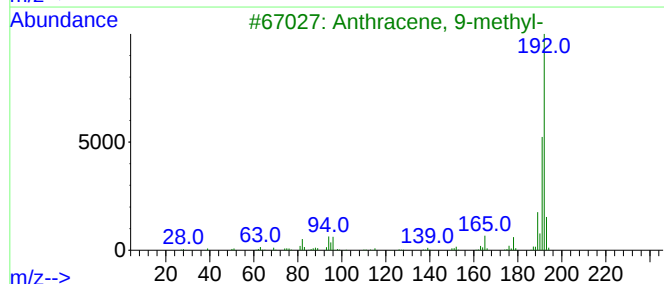
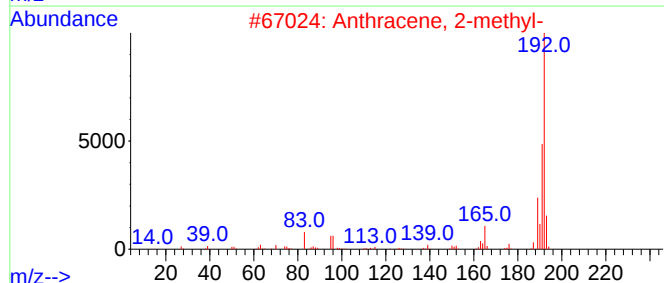
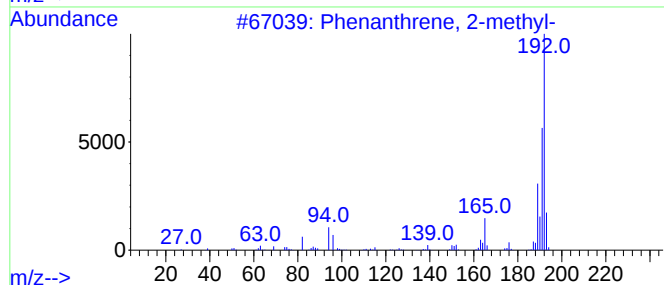
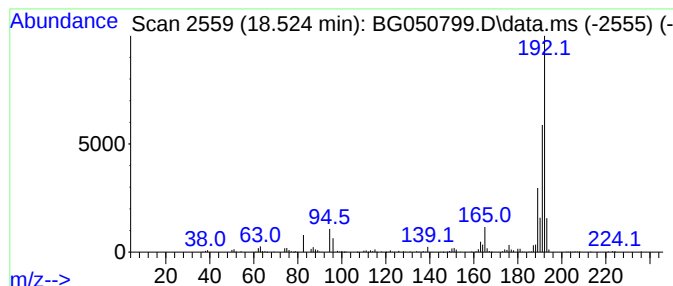
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 6 Anthracene, 2-methyl- Concentration Rank 6

| R.T.   | EstConc  | Area   | Relative to ISTD | R.T.   |
|--------|----------|--------|------------------|--------|
| 18.524 | 12.36 ng | 391900 | Phenanthrene-d10 | 17.607 |

| Hit# | of 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------|-----|---------|-------------|------|
| 1    |      | Phenanthrene, 2-methyl- | 192 | C15H12  | 002531-84-2 | 97   |
| 2    |      | Anthracene, 2-methyl-   | 192 | C15H12  | 000613-12-7 | 96   |
| 3    |      | Anthracene, 9-methyl-   | 192 | C15H12  | 000779-02-2 | 93   |
| 4    |      | Phenanthrene, 1-methyl- | 192 | C15H12  | 000832-69-9 | 93   |
| 5    |      | Phenanthrene, 4-methyl- | 192 | C15H12  | 000832-64-4 | 93   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG110121\  
Data File : BG050799.D  
Acq On : 1 Nov 2021 15:12  
Operator : CG/JU  
Sample : M4422-01 2X  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
OR-03-102921

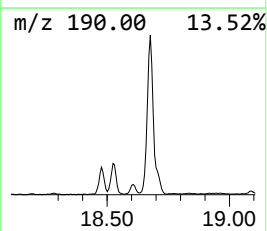
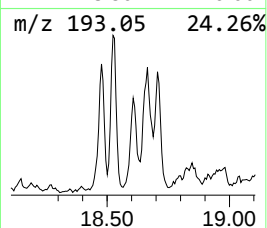
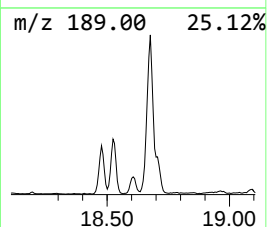
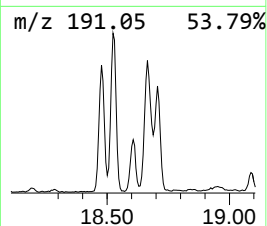
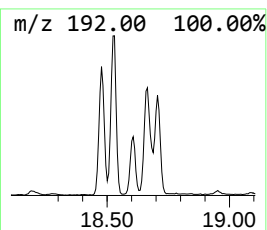
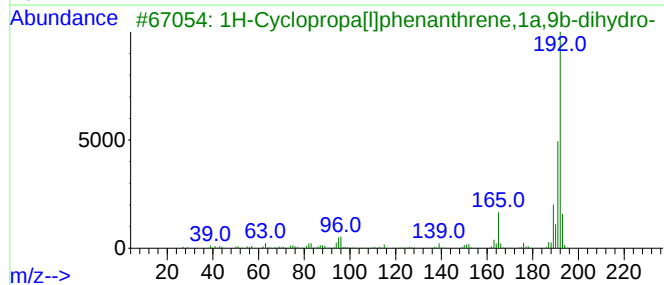
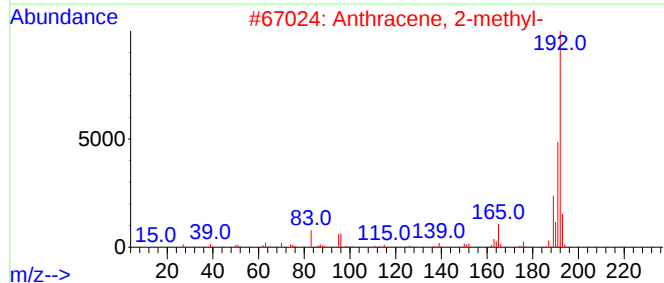
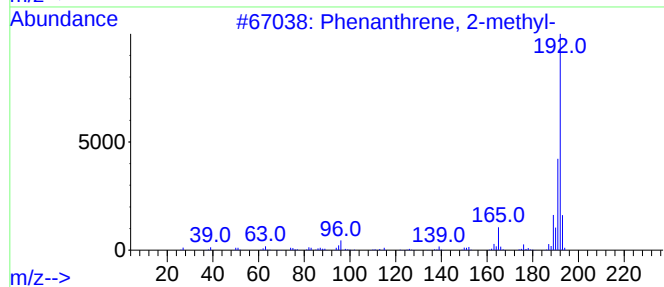
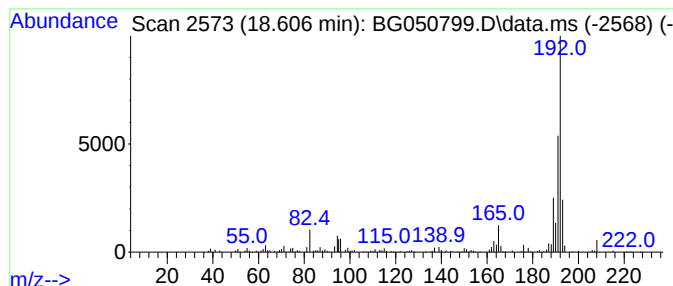
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 7 1H-Cyclopropa[l]phenanthren... Concentration Rank 14

| R.T.      | EstConc                             | Area   | Relative to ISTD |         |             | R.T.   |
|-----------|-------------------------------------|--------|------------------|---------|-------------|--------|
| 18.606    | 4.36 ng                             | 138366 | Phenanthrene-d10 |         |             | 17.607 |
| Hit# of 5 | Tentative ID                        |        | MW               | MolForm | CAS#        | Qual   |
| 1         | Phenanthrene, 2-methyl-             |        | 192              | C15H12  | 002531-84-2 | 96     |
| 2         | Anthracene, 2-methyl-               |        | 192              | C15H12  | 000613-12-7 | 95     |
| 3         | 1H-Cyclopropa[1]phenanthrene,1a,... |        | 192              | C15H12  | 000949-41-7 | 95     |
| 4         | Phenanthrene, 1-methyl-             |        | 192              | C15H12  | 000832-69-9 | 95     |
| 5         | Anthracene, 9-methyl-               |        | 192              | C15H12  | 000779-02-2 | 90     |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG110121\  
Data File : BG050799.D  
Acq On : 1 Nov 2021 15:12  
Operator : CG/JU  
Sample : M4422-01 2X  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
OR-03-102921

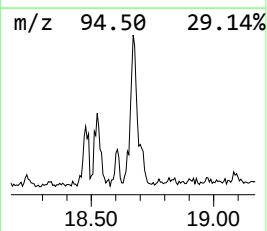
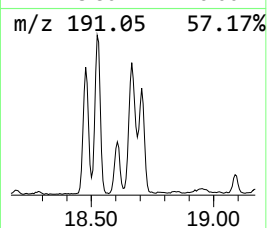
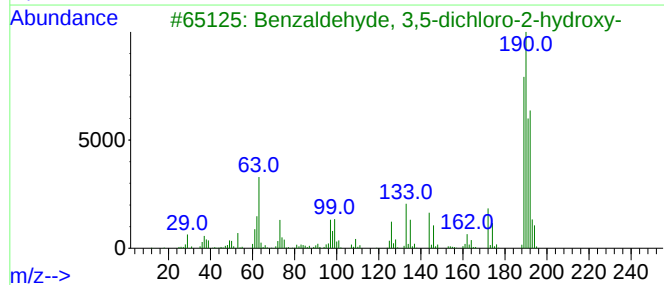
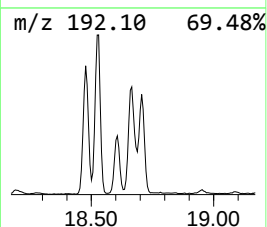
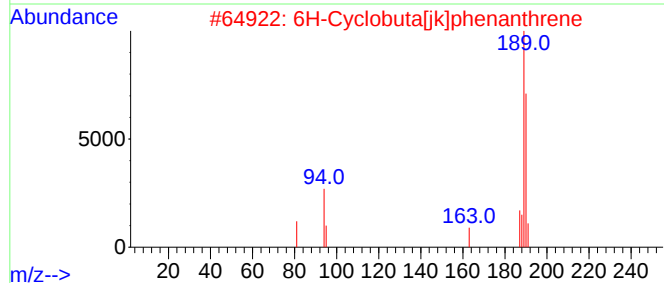
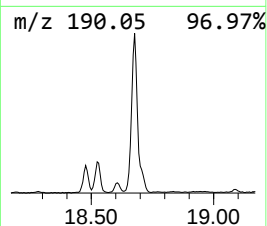
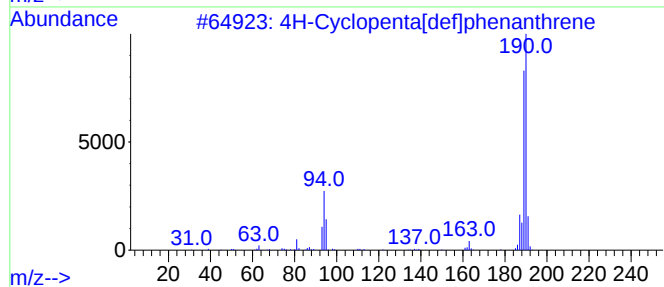
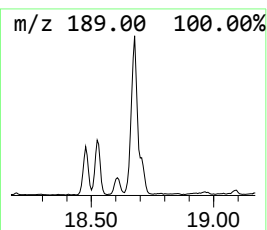
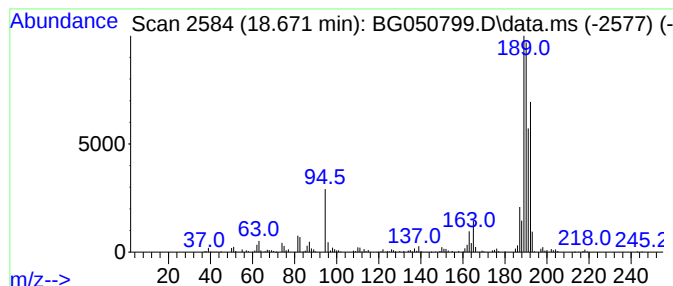
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 8 4H-Cyclopenta[def]phenanthrene Concentration Rank 4

| R.T.      | EstConc                             | Area   | Relative to ISTD |           |             | R.T.   |
|-----------|-------------------------------------|--------|------------------|-----------|-------------|--------|
| 18.671    | 19.63 ng                            | 622423 | Phenanthrene-d10 |           |             | 17.607 |
| Hit# of 5 | Tentative ID                        |        | MW               | MolForm   | CAS#        | Qual   |
| 1         | 4H-Cyclopenta[def]phenanthrene      |        | 190              | C15H10    | 000203-64-5 | 70     |
| 2         | 6H-Cyclobuta[jk]phenanthrene        |        | 190              | C15H10    | 083469-43-6 | 58     |
| 3         | Benzaldehyde, 3,5-dichloro-2-hyd... |        | 190              | C7H4Cl2O2 | 000090-60-8 | 47     |
| 4         | 1a,9b-Dihydro-1H-cyclopropa[a]an... |        | 192              | C15H12    | 006109-24-6 | 30     |
| 5         | 3-Bromo-2-furoic acid               |        | 190              | C5H3BrO3  | 014903-90-3 | 27     |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG110121\  
Data File : BG050799.D  
Acq On : 1 Nov 2021 15:12  
Operator : CG/JU  
Sample : M4422-01 2X  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
OR-03-102921

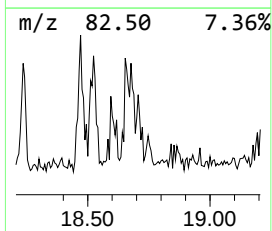
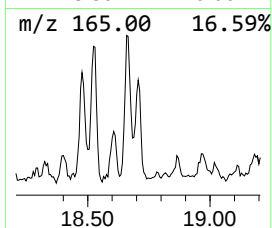
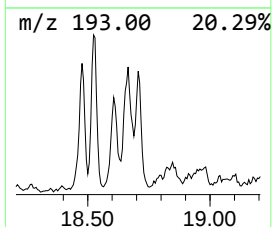
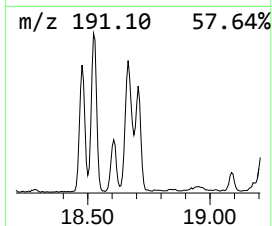
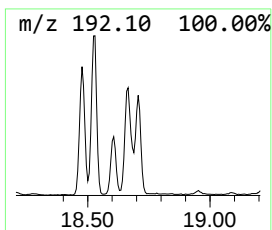
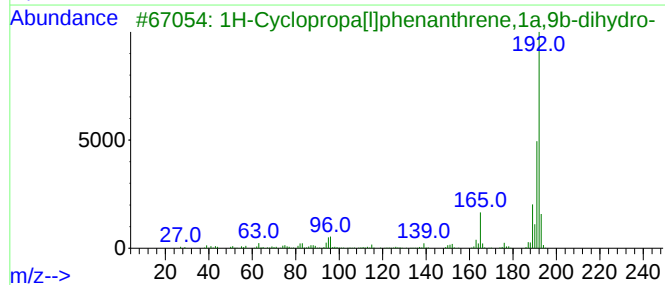
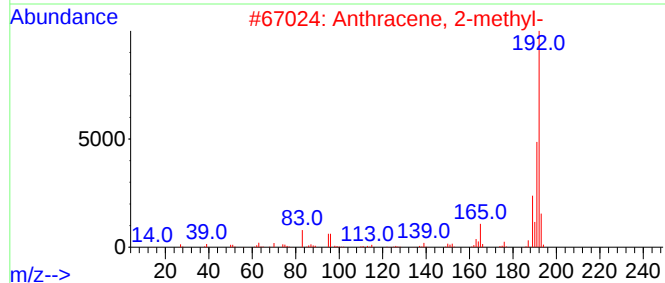
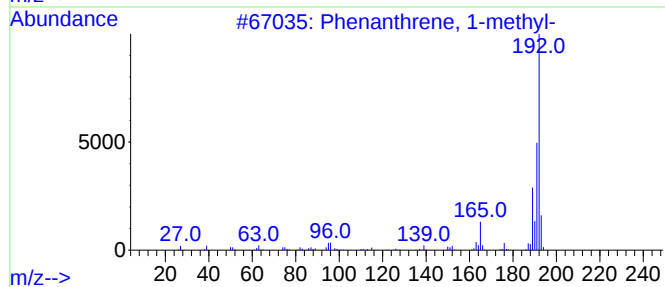
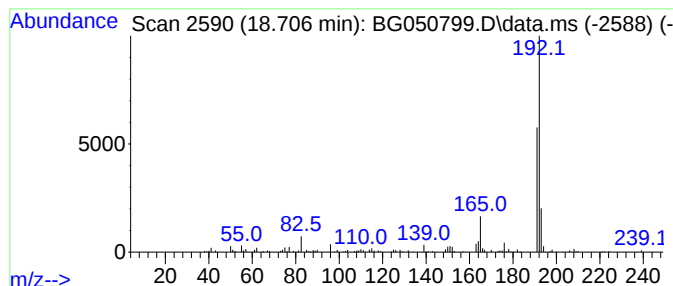
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 9 Phenanthrene, 1-methyl- Concentration Rank 11

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|-------------|------|
| 18.706    | 5.54 ng                             | 175721 | Phenanthrene-d10 | 17.607      |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual |
| 1         | Phenanthrene, 1-methyl-             | 192    | C15H12           | 000832-69-9 | 95   |
| 2         | Anthracene, 2-methyl-               | 192    | C15H12           | 000613-12-7 | 91   |
| 3         | 1H-Cyclopropa[1]phenanthrene,1a,... | 192    | C15H12           | 000949-41-7 | 91   |
| 4         | Phenanthrene, 2-methyl-             | 192    | C15H12           | 002531-84-2 | 91   |
| 5         | Anthracene, 1-methyl-               | 192    | C15H12           | 000610-48-0 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG110121\  
Data File : BG050799.D  
Acq On : 1 Nov 2021 15:12  
Operator : CG/JU  
Sample : M4422-01 2X  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
OR-03-102921

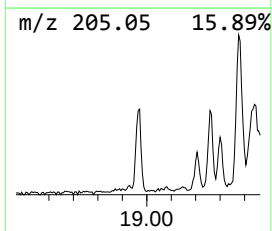
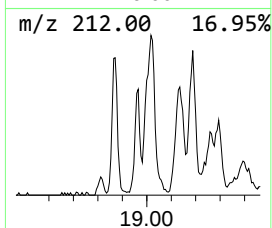
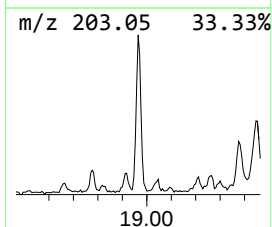
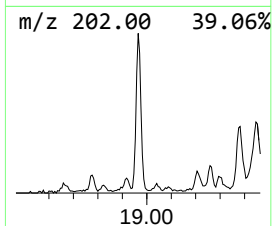
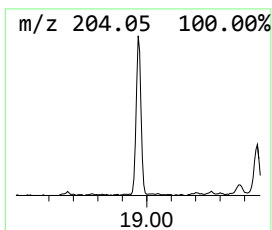
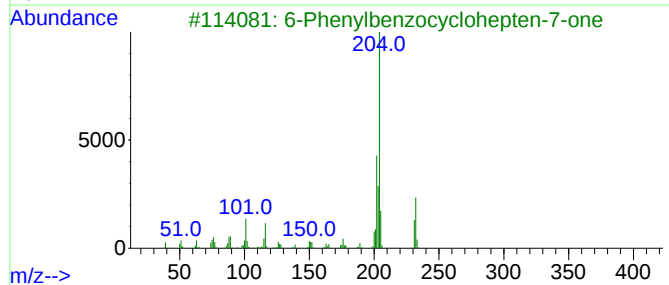
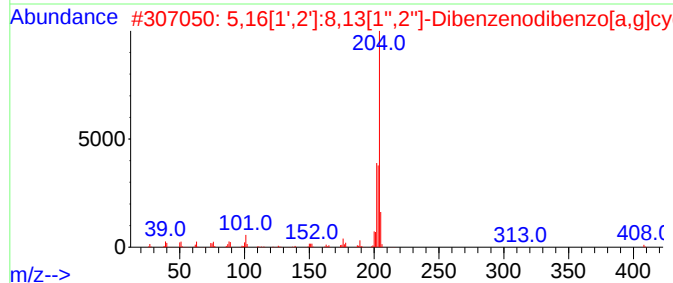
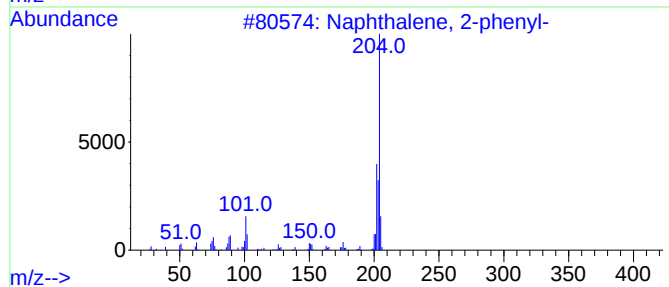
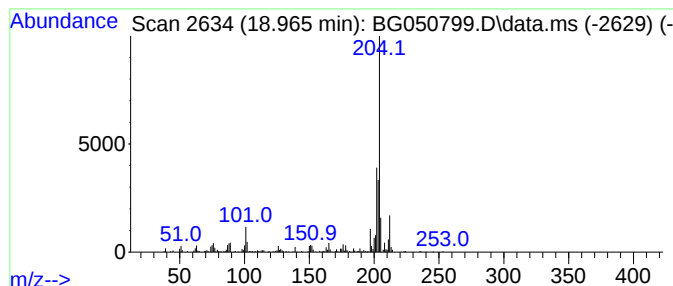
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 18.965 | 7.11 ng | 225350 | Phenanthrene-d10 | 17.607 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|-----------|-------------|------|
| 1    |    |   | Naphthalene, 2-phenyl-              | 204 | C16H12    | 000612-94-2 | 90   |
| 2    |    |   | 5,16[1',2']:8,13[1'',2'']-Dibenz... | 408 | C32H24    | 005672-97-9 | 74   |
| 3    |    |   | 6-Phenylbenzocyclohepten-7-one      | 232 | C17H12O   | 093327-56-1 | 72   |
| 4    |    |   | Tricyclo[8.2.2.2(4,7)]hexadeca-2... | 204 | C16H12    | 006572-60-7 | 60   |
| 5    |    |   | 9,10-Bis(bromomethyl)anthracene     | 362 | C16H12Br2 | 034373-96-1 | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG110121\  
Data File : BG050799.D  
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Operator : CG/JU  
Sample : M4422-01 2X  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

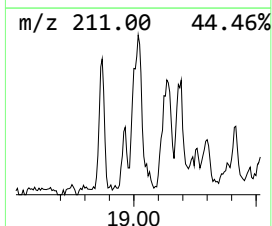
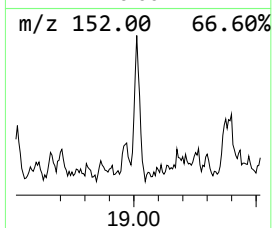
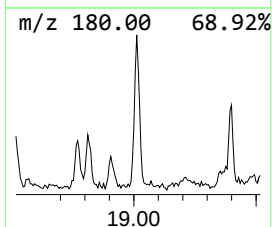
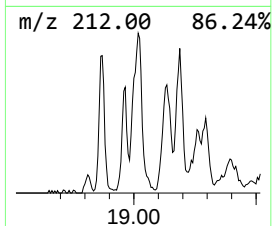
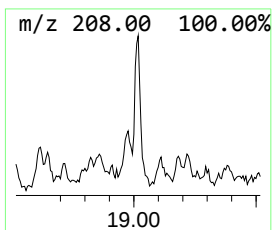
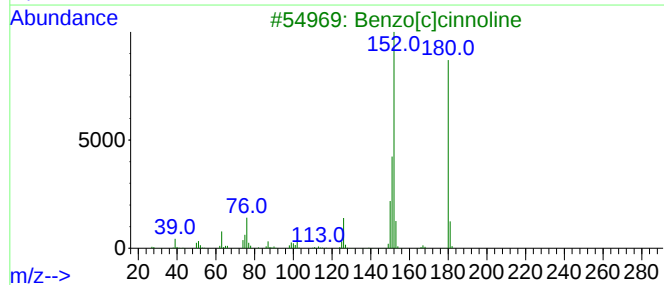
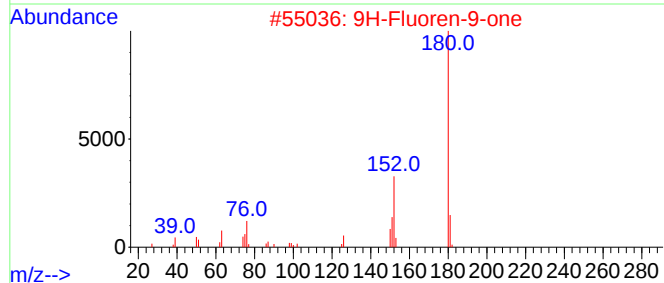
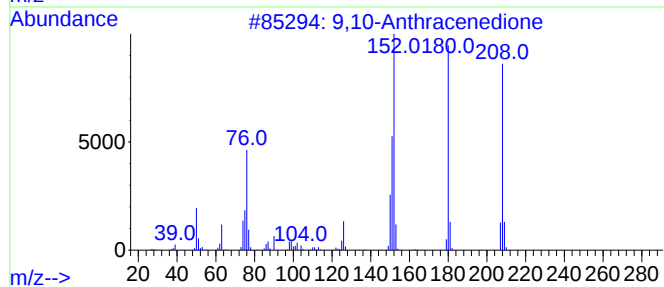
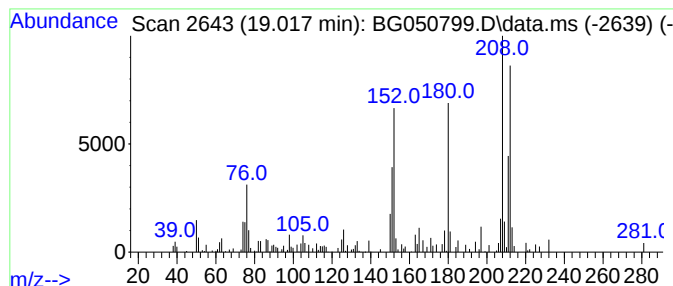
Instrument :  
BNA\_G  
ClientSampleId :  
OR-03-102921

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG100621.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 11 9,10-Anthracenedione Concentration Rank 12

| R.T.   | EstConc                            | Area         | Relative to ISTD |             |      | R.T.   |
|--------|------------------------------------|--------------|------------------|-------------|------|--------|
| 19.017 | 4.76 ng                            | 150874       | Phenanthrene-d10 |             |      | 17.607 |
| Hit#   | of 5                               | Tentative ID | MW               | MolForm     | CAS# | Qual   |
| 1      | 9,10-Anthracenedione               | 208          | C14H8O2          | 000084-65-1 | 90   |        |
| 2      | 9H-Fluoren-9-one                   | 180          | C13H8O           | 000486-25-9 | 47   |        |
| 3      | Benzo[c]cinnoline                  | 180          | C12H8N2          | 000230-17-1 | 38   |        |
| 4      | 5,6,7,8-Tetrahydro-1-phenylapht... | 208          | C16H16           | 067064-63-5 | 30   |        |
| 5      | 1H-Phenalen-1-one                  | 180          | C13H8O           | 000548-39-0 | 25   |        |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG110121\  
Data File : BG050799.D  
Acq On : 1 Nov 2021 15:12  
Operator : CG/JU  
Sample : M4422-01 2X  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
OR-03-102921

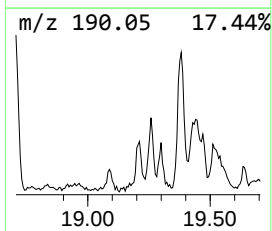
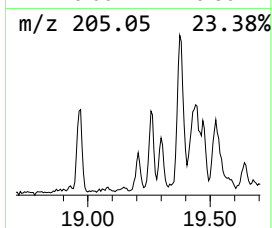
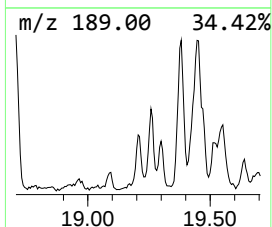
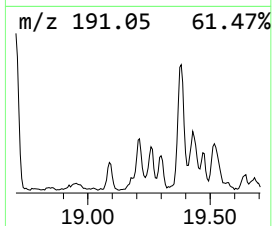
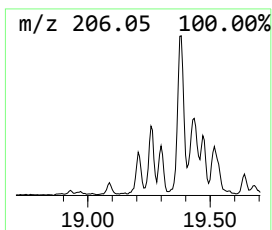
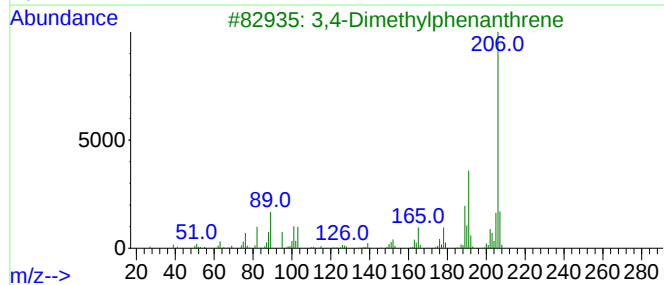
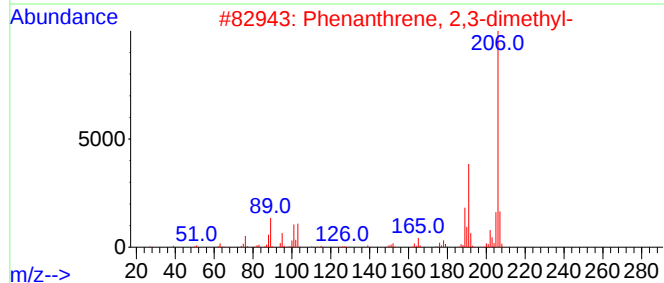
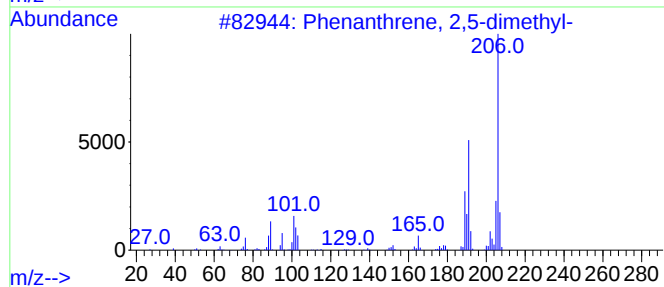
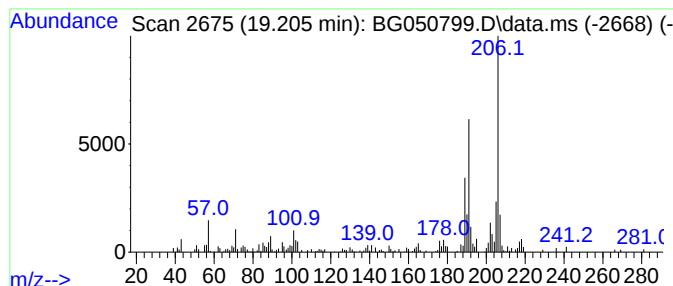
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 12 Phenanthrene, 2,5-dimethyl- Concentration Rank 13

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 19.206 | 4.73 ng | 149981 | Phenanthrene-d10 | 17.607 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#         | Qual |
|------|----|---|-----------------------------|-----|---------|--------------|------|
| 1    |    |   | Phenanthrene, 2,5-dimethyl- | 206 | C16H14  | 003674-66-6  | 94   |
| 2    |    |   | Phenanthrene, 2,3-dimethyl- | 206 | C16H14  | 003674-65-5  | 94   |
| 3    |    |   | 3,4-Dimethylphenanthrene    | 206 | C16H14  | 1000452-24-5 | 94   |
| 4    |    |   | di-p-Tolylacetylene         | 206 | C16H14  | 002789-88-0  | 93   |
| 5    |    |   | Anthracene, 2-ethyl-        | 206 | C16H14  | 052251-71-5  | 93   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG110121\  
Data File : BG050799.D  
Acq On : 1 Nov 2021 15:12  
Operator : CG/JU  
Sample : M4422-01 2X  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
OR-03-102921

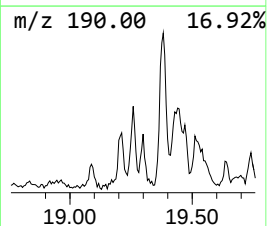
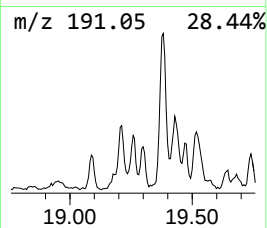
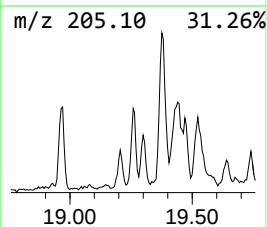
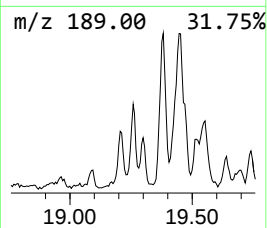
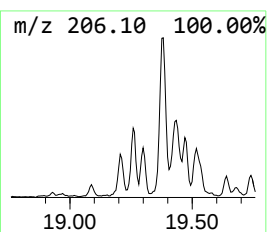
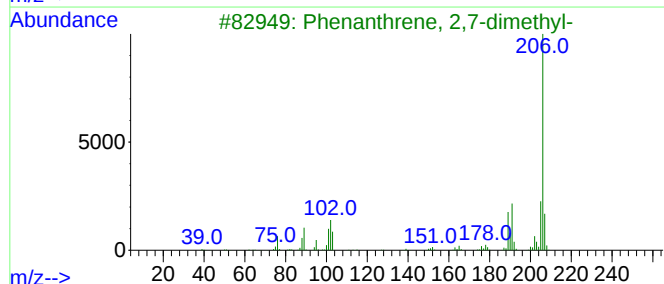
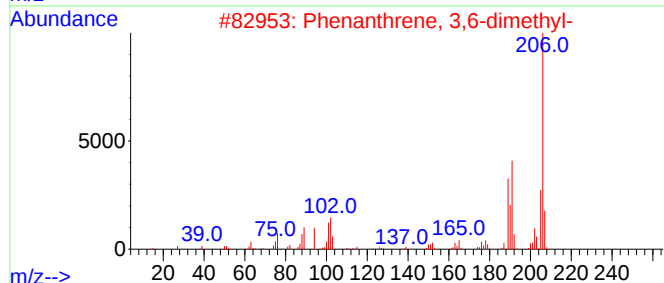
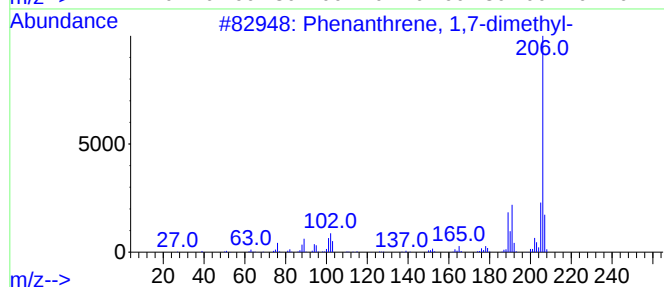
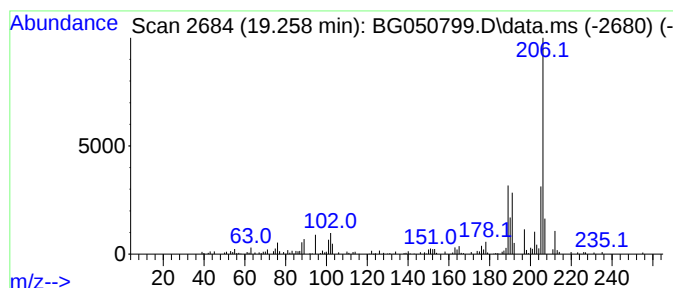
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 13 Phenanthrene, 1,7-dimethyl- Concentration Rank 20

| R.T.      | EstConc                            | Area   | Relative to ISTD | R.T.        |      |
|-----------|------------------------------------|--------|------------------|-------------|------|
| 19.258    | 3.39 ng                            | 107511 | Phenanthrene-d10 | 17.607      |      |
| Hit# of 5 | Tentative ID                       | MW     | MolForm          | CAS#        | Qual |
| 1         | Phenanthrene, 1,7-dimethyl-        | 206    | C16H14           | 000483-87-4 | 95   |
| 2         | Phenanthrene, 3,6-dimethyl-        | 206    | C16H14           | 001576-67-6 | 94   |
| 3         | Phenanthrene, 2,7-dimethyl-        | 206    | C16H14           | 001576-69-8 | 83   |
| 4         | di-p-Tolylacetylene                | 206    | C16H14           | 002789-88-0 | 81   |
| 5         | Naphthalene, 1,2-dihydro-4-phenyl- | 206    | C16H14           | 007469-40-1 | 76   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG110121\  
Data File : BG050799.D  
Acq On : 1 Nov 2021 15:12  
Operator : CG/JU  
Sample : M4422-01 2X  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
OR-03-102921

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG100621.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

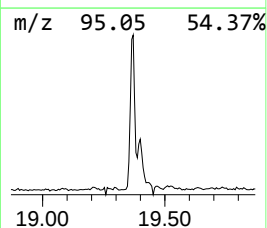
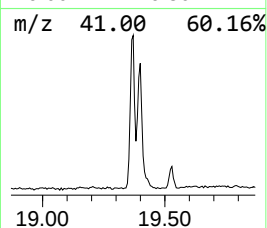
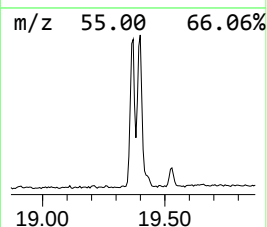
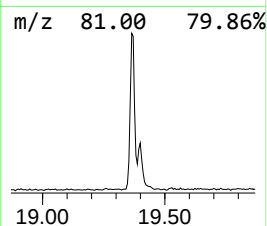
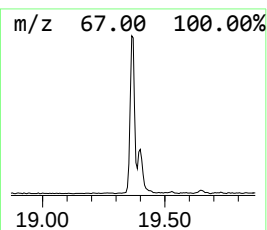
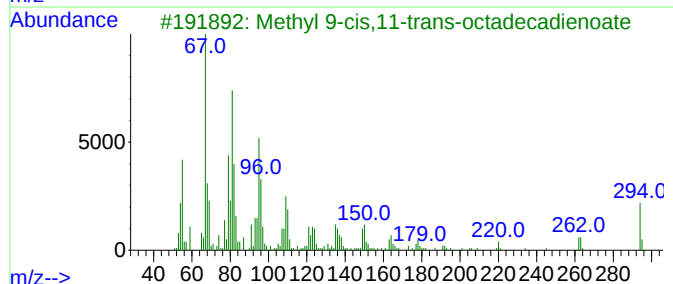
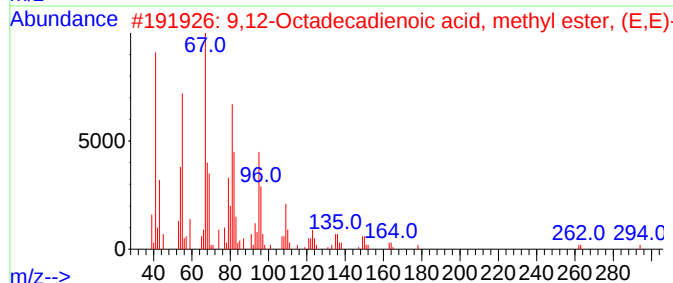
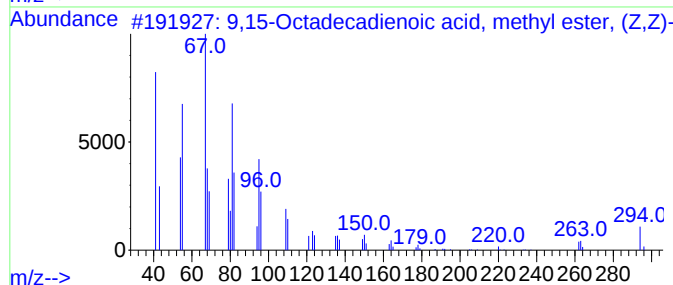
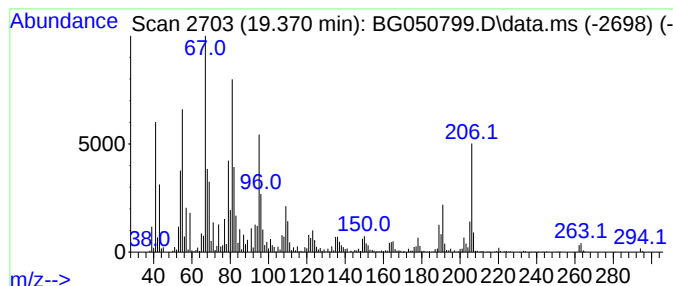
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TIC Integration Parameters: LSCINT.P

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Peak Number 14 9,15-Octadecadienoic acid, ... Concentration Rank 1

| R.T.   | EstConc  | Area    | Relative to ISTD | R.T.   |
|--------|----------|---------|------------------|--------|
| 19.370 | 39.14 ng | 1240860 | Phenanthrene-d10 | 17.607 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 9,15-Octadecadienoic acid, methy... | 294 | C19H34O2 | 017309-05-6 | 99   |
| 2    |    |   | 9,12-Octadecadienoic acid, methy... | 294 | C19H34O2 | 002566-97-4 | 97   |
| 3    |    |   | Methyl 9-cis,11-trans-octadecadi... | 294 | C19H34O2 | 013058-52-1 | 95   |
| 4    |    |   | 9,11-Octadecadienoic acid, methy... | 294 | C19H34O2 | 013038-47-6 | 95   |
| 5    |    |   | 11,14-Octadecadienoic acid, meth... | 294 | C19H34O2 | 056554-61-1 | 95   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG110121\  
Data File : BG050799.D  
Acq On : 1 Nov 2021 15:12  
Operator : CG/JU  
Sample : M4422-01 2X  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
OR-03-102921

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG100621.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

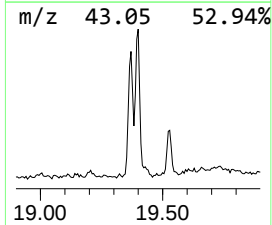
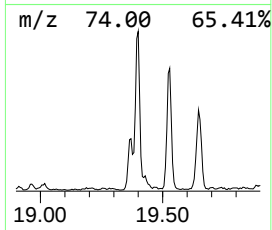
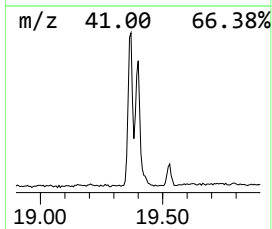
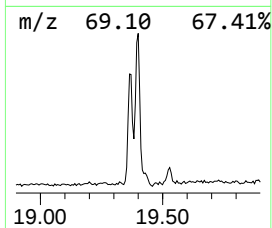
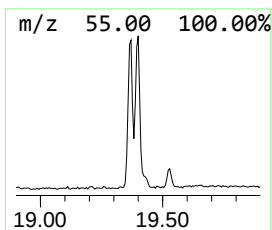
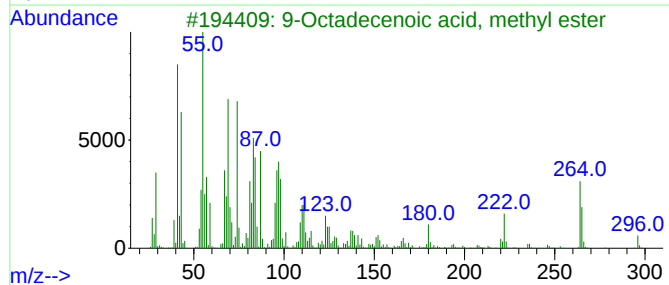
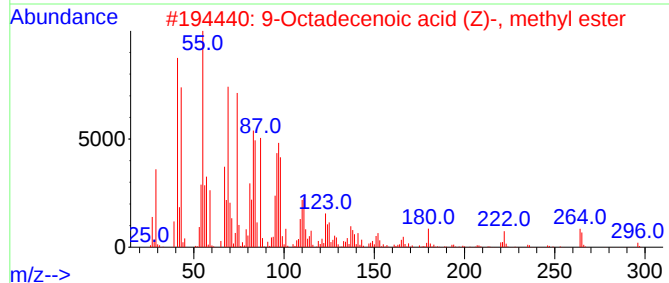
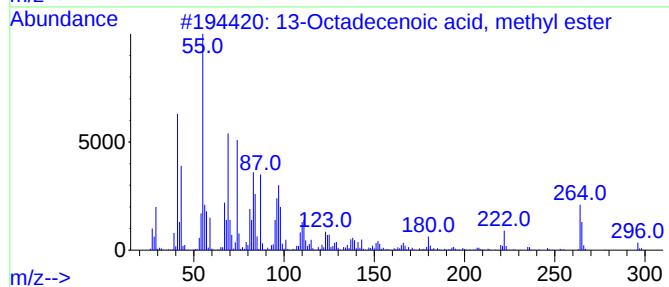
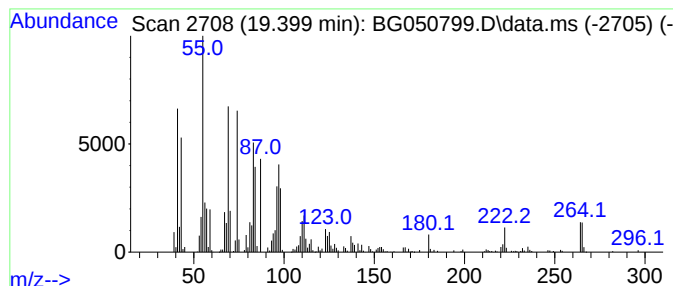
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Peak Number 15 13-Octadecenoic acid, methyl... Concentration Rank 3

| R.T.   | EstConc  | Area   | Relative to ISTD | R.T.   |
|--------|----------|--------|------------------|--------|
| 19.399 | 21.48 ng | 681112 | Phenanthrene-d10 | 17.607 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 13-Octadecenoic acid, methyl ester  | 296 | C19H36O2 | 056554-47-3 | 94   |
| 2    |    |   | 9-Octadecenoic acid (Z)-, methyl... | 296 | C19H36O2 | 000112-62-9 | 91   |
| 3    |    |   | 9-Octadecenoic acid, methyl ester   | 296 | C19H36O2 | 002462-84-2 | 91   |
| 4    |    |   | 8-Octadecenoic acid, methyl ester   | 296 | C19H36O2 | 002345-29-1 | 91   |
| 5    |    |   | 11-Octadecenoic acid, methyl est... | 296 | C19H36O2 | 001937-63-9 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG110121\  
Data File : BG050799.D  
Acq On : 1 Nov 2021 15:12  
Operator : CG/JU  
Sample : M4422-01 2X  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
OR-03-102921

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG100621.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

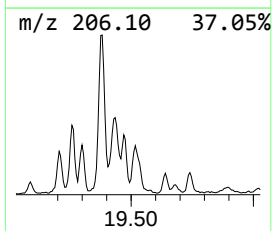
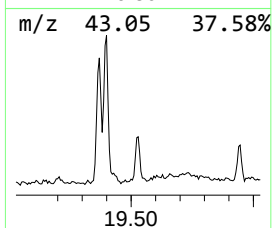
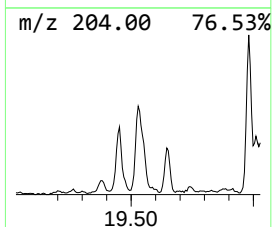
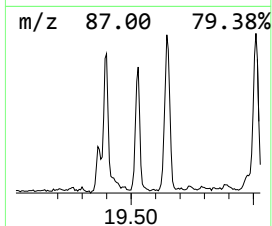
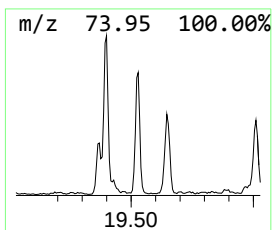
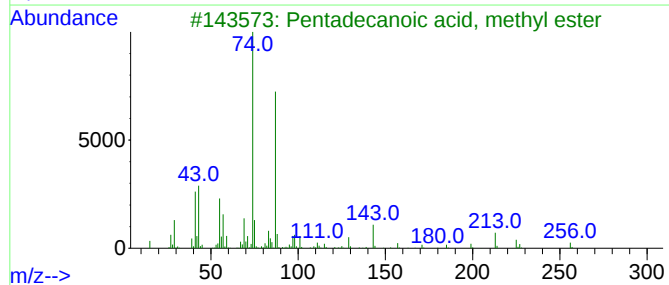
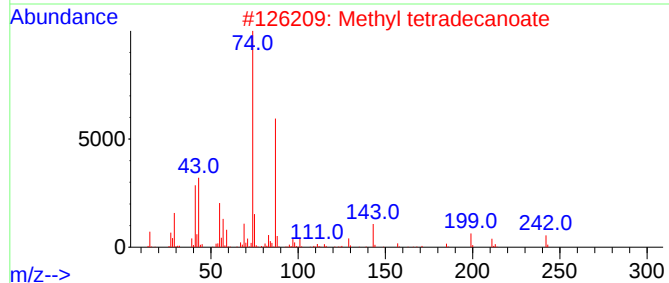
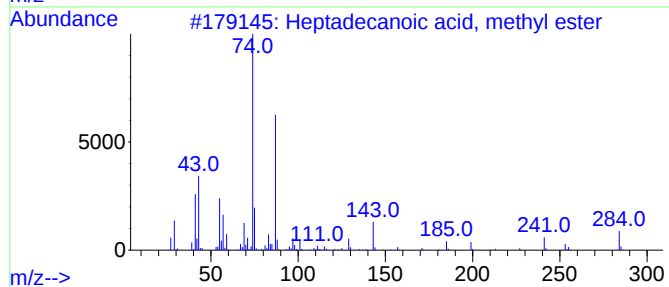
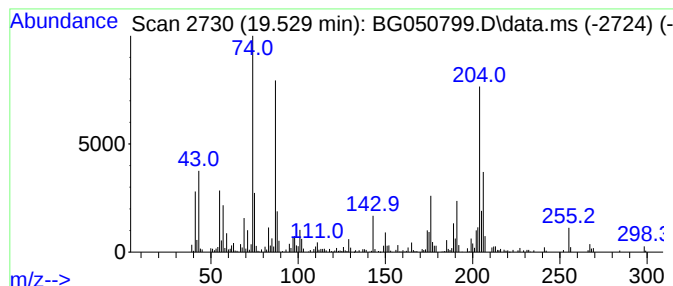
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 16 unknown19.529 Concentration Rank 5

| R.T.   | EstConc  | Area   | Relative to ISTD | R.T.   |
|--------|----------|--------|------------------|--------|
| 19.529 | 13.10 ng | 415204 | Phenanthrene-d10 | 17.607 |

| Hit# | of 5 | Tentative ID                     | MW  | MolForm  | CAS#         | Qual |
|------|------|----------------------------------|-----|----------|--------------|------|
| 1    |      | Heptadecanoic acid, methyl ester | 284 | C18H36O2 | 001731-92-6  | 46   |
| 2    |      | Methyl tetradecanoate            | 242 | C15H30O2 | 000124-10-7  | 46   |
| 3    |      | Pentadecanoic acid, methyl ester | 256 | C16H32O2 | 007132-64-1  | 46   |
| 4    |      | Methyl stearate                  | 298 | C19H38O2 | 000112-61-8  | 41   |
| 5    |      | Methyl 8-methyl-nonanoate        | 186 | C11H22O2 | 1000452-00-9 | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG110121\  
Data File : BG050799.D  
Acq On : 1 Nov 2021 15:12  
Operator : CG/JU  
Sample : M4422-01 2X  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
OR-03-102921

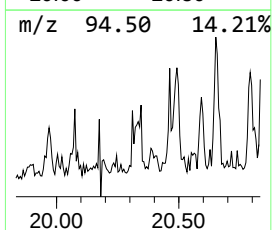
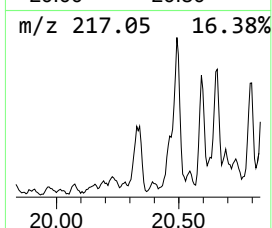
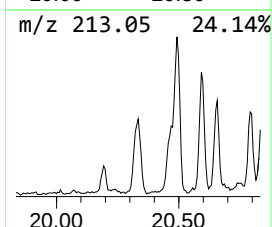
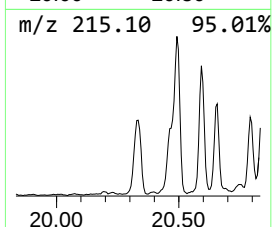
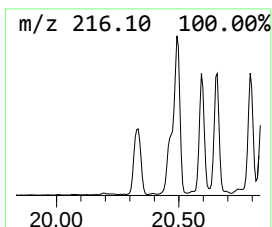
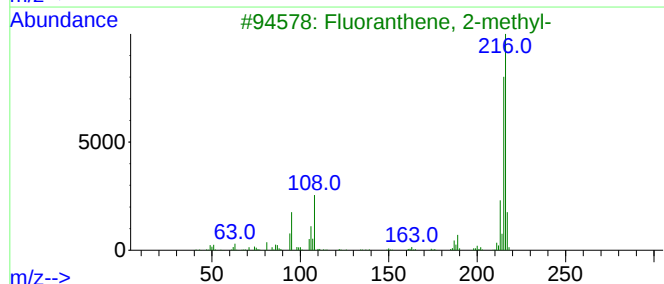
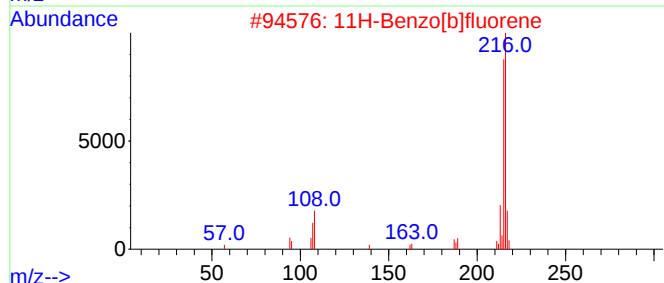
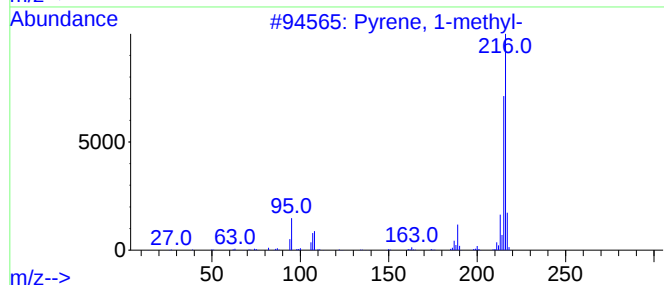
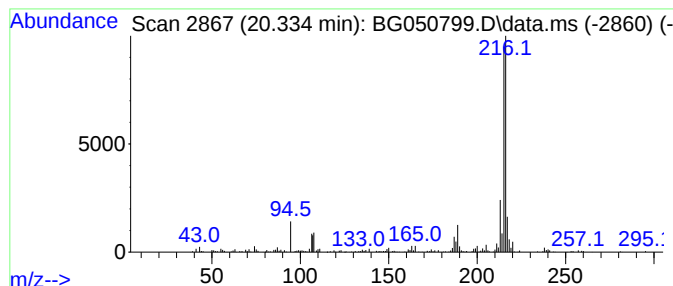
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 17 11H-Benzo[b]fluorene Concentration Rank 16

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 20.334 | 3.93 ng | 353647 | Chrysene-d12     | 21.908 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Pyrene, 1-methyl-       | 216 | C17H12  | 002381-21-7 | 93   |
| 2    |    |   | 11H-Benzo[b]fluorene    | 216 | C17H12  | 000243-17-4 | 93   |
| 3    |    |   | Fluoranthene, 2-methyl- | 216 | C17H12  | 033543-31-6 | 91   |
| 4    |    |   | 7H-Benzo[c]fluorene     | 216 | C17H12  | 000205-12-9 | 83   |
| 5    |    |   | Pyrene, 2-methyl-       | 216 | C17H12  | 003442-78-2 | 76   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG110121\  
Data File : BG050799.D  
Acq On : 1 Nov 2021 15:12  
Operator : CG/JU  
Sample : M4422-01 2X  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
OR-03-102921

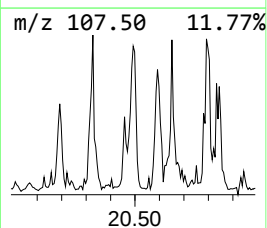
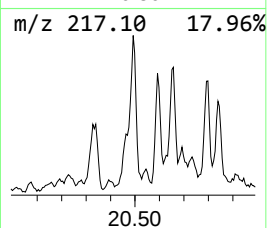
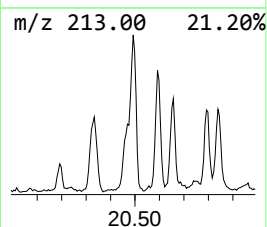
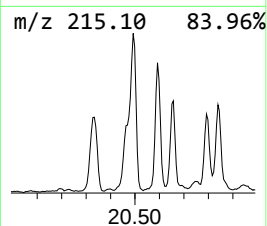
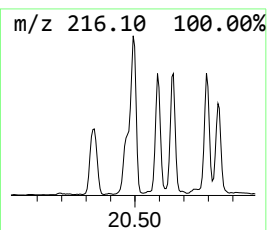
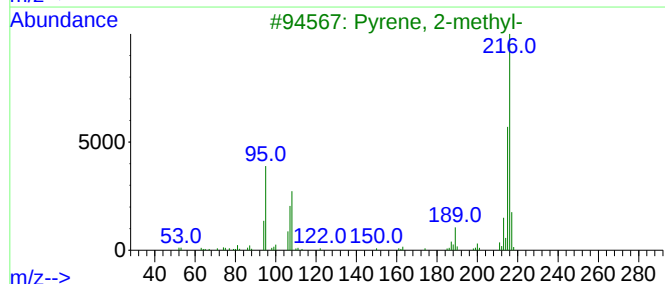
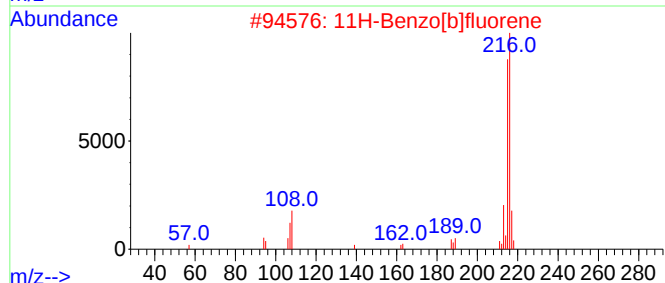
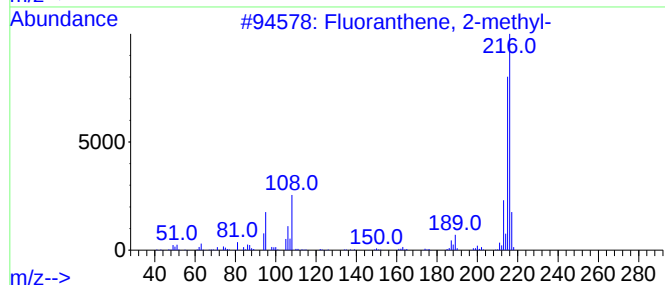
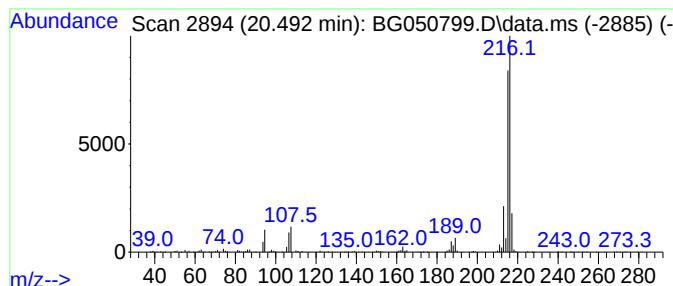
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 18 Fluoranthene, 2-methyl- Concentration Rank 9

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 20.492 | 7.74 ng | 696646 | Chrysene-d12     | 21.908 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Fluoranthene, 2-methyl- | 216 | C17H12  | 033543-31-6 | 95   |
| 2    |    |   | 11H-Benzo[b]fluorene    | 216 | C17H12  | 000243-17-4 | 94   |
| 3    |    |   | Pyrene, 2-methyl-       | 216 | C17H12  | 003442-78-2 | 90   |
| 4    |    |   | Pyrene, 1-methyl-       | 216 | C17H12  | 002381-21-7 | 90   |
| 5    |    |   | 11H-Benzo[a]fluorene    | 216 | C17H12  | 000238-84-6 | 87   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG110121\  
Data File : BG050799.D  
Acq On : 1 Nov 2021 15:12  
Operator : CG/JU  
Sample : M4422-01 2X  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
OR-03-102921

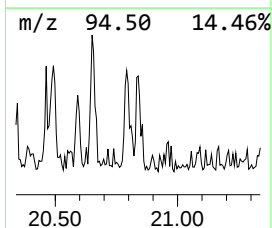
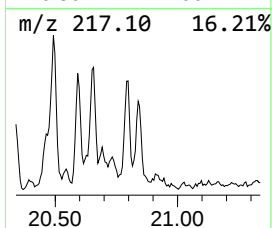
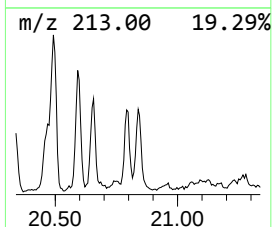
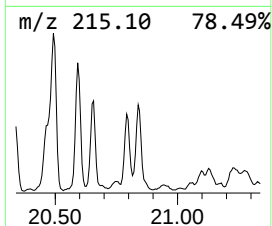
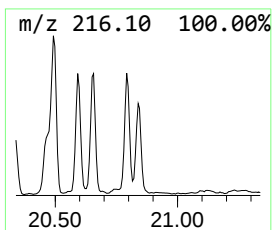
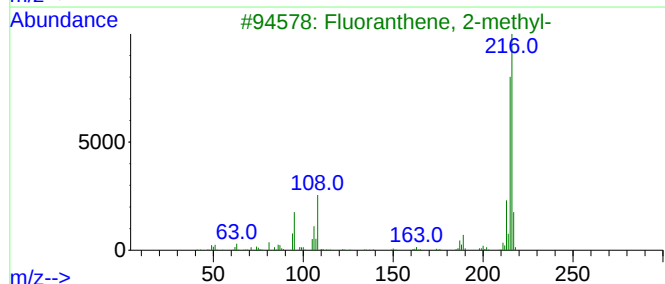
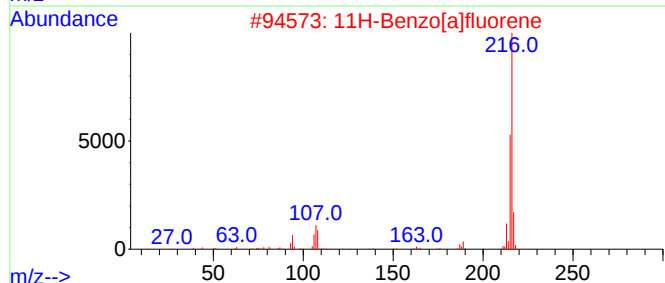
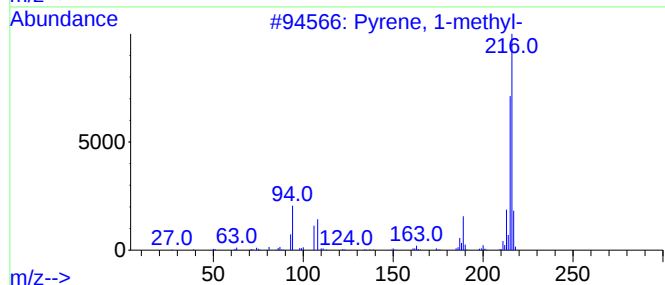
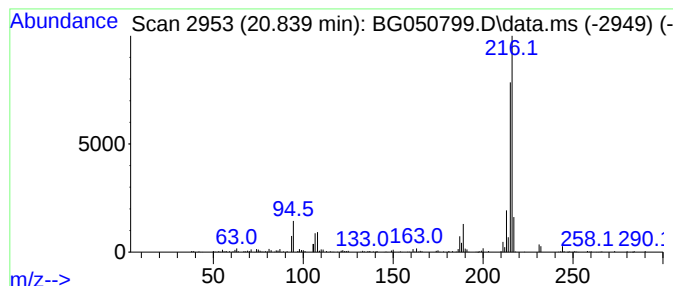
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 19 Pyrene, 1-methyl- Concentration Rank 19

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 20.839 | 3.56 ng | 320578 | Chrysene-d12     | 21.908 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | Pyrene, 1-methyl-       | 216 | C17H12  | 002381-21-7 | 97   |
| 2    |    |   | 11H-Benzo[a]fluorene    | 216 | C17H12  | 000238-84-6 | 94   |
| 3    |    |   | Fluoranthene, 2-methyl- | 216 | C17H12  | 033543-31-6 | 91   |
| 4    |    |   | 11H-Benzo[b]fluorene    | 216 | C17H12  | 000243-17-4 | 87   |
| 5    |    |   | 7H-Benzo[c]fluorene     | 216 | C17H12  | 000205-12-9 | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG110121\  
Data File : BG050799.D  
Acq On : 1 Nov 2021 15:12  
Operator : CG/JU  
Sample : M4422-01 2X  
Misc :  
ALS Vial : 7 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
OR-03-102921

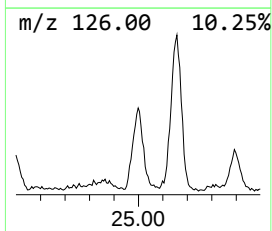
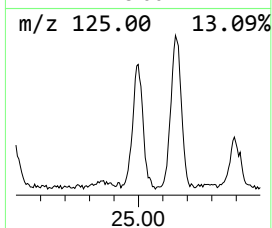
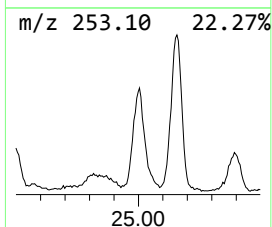
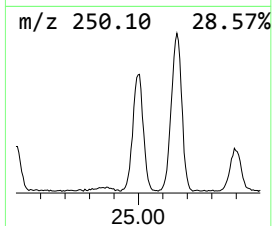
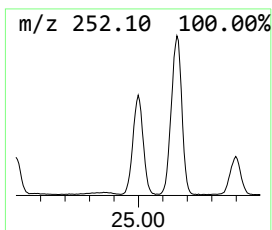
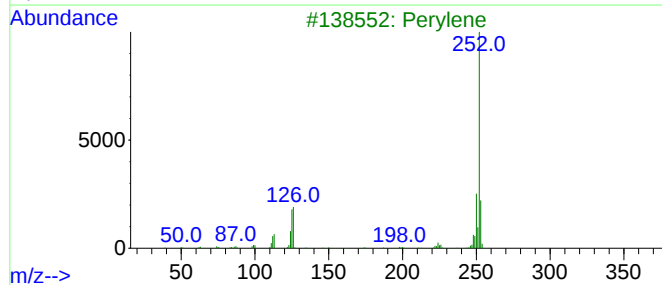
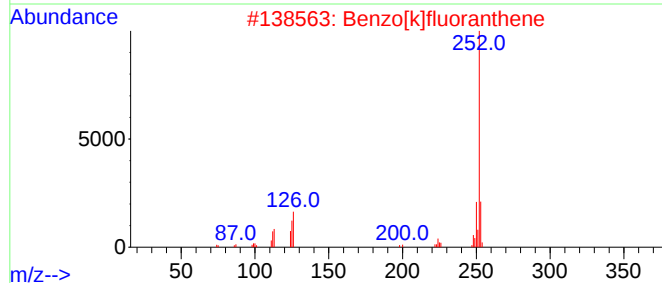
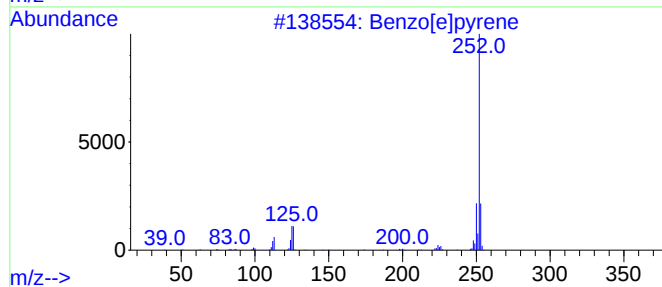
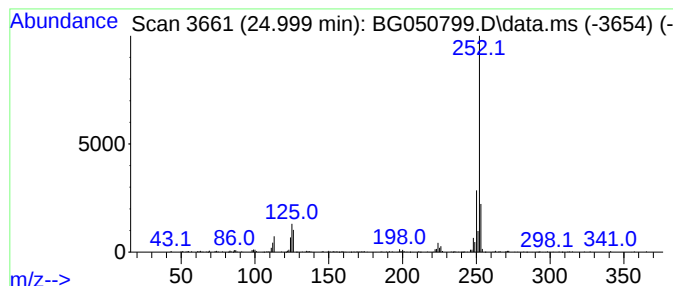
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 20 Benzo[e]pyrene Concentration Rank 2

| R.T.   | EstConc  | Area   | Relative to ISTD | R.T.   |
|--------|----------|--------|------------------|--------|
| 24.999 | 25.59 ng | 648296 | Perylene-d12     | 25.316 |

| 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    |    |   | Perylene             | 252 | C20H12  | 000198-55-0 | 95   |
| 4    |    |   | Benzo[j]fluoranthene | 252 | C20H12  | 000205-82-3 | 94   |
| 5    |    |   | Benzo[b]fluoranthene | 252 | C20H12  | 000205-99-2 | 90   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG110121\  
 Data File : BG050799.D  
 Acq On : 1 Nov 2021 15:12  
 Operator : CG/JU  
 Sample : M4422-01 2X  
 Misc :  
 ALS Vial : 7 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 OR-03-102921

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\8270-BG100621.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 |
| Heptadecane        | 16.532 | 3.6     | ng    | 114194   | 4                     | 17.607 | 634056  | 20.0 |
| Tetradecane        | 17.320 | 4.1     | ng    | 131378   | 4                     | 17.607 | 634056  | 20.0 |
| Dibenzothiophene   | 17.425 | 3.7     | ng    | 116503   | 4                     | 17.607 | 634056  | 20.0 |
| Pentadecanoic a... | 18.236 | 9.3     | ng    | 295884   | 4                     | 17.607 | 634056  | 20.0 |
| Phenanthrene, 2... | 18.477 | 10.8    | ng    | 343456   | 4                     | 17.607 | 634056  | 20.0 |
| Anthracene, 2-m... | 18.524 | 12.4    | ng    | 391900   | 4                     | 17.607 | 634056  | 20.0 |
| 1H-Cyclopropa[l... | 18.606 | 4.4     | ng    | 138366   | 4                     | 17.607 | 634056  | 20.0 |
| 4H-Cyclopenta[d... | 18.671 | 19.6    | ng    | 622423   | 4                     | 17.607 | 634056  | 20.0 |
| Phenanthrene, 1... | 18.706 | 5.5     | ng    | 175721   | 4                     | 17.607 | 634056  | 20.0 |
| Naphthalene, 2-... | 18.965 | 7.1     | ng    | 225350   | 4                     | 17.607 | 634056  | 20.0 |
| 9,10-Anthracene... | 19.017 | 4.8     | ng    | 150874   | 4                     | 17.607 | 634056  | 20.0 |
| Phenanthrene, 2... | 19.206 | 4.7     | ng    | 149981   | 4                     | 17.607 | 634056  | 20.0 |
| Phenanthrene, 1... | 19.258 | 3.4     | ng    | 107511   | 4                     | 17.607 | 634056  | 20.0 |
| 9,15-Octadecadi... | 19.370 | 39.1    | ng    | 1240860  | 4                     | 17.607 | 634056  | 20.0 |
| 13-Octadecenoic... | 19.399 | 21.5    | ng    | 681112   | 4                     | 17.607 | 634056  | 20.0 |
| unknown19.529      | 19.529 | 13.1    | ng    | 415204   | 4                     | 17.607 | 634056  | 20.0 |
| 11H-Benzo[b]flu... | 20.334 | 3.9     | ng    | 353647   | 5                     | 21.908 | 1799930 | 20.0 |
| Fluoranthene, 2... | 20.492 | 7.7     | ng    | 696646   | 5                     | 21.908 | 1799930 | 20.0 |
| Pyrene, 1-methyl-  | 20.839 | 3.6     | ng    | 320578   | 5                     | 21.908 | 1799930 | 20.0 |
| Benzo[e]pyrene     | 24.999 | 25.6    | ng    | 648296   | 6                     | 25.316 | 506644  | 20.0 |