

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF050317\  
 Data File : BF094848.D  
 Acq On : 3 May 2017 22:42  
 Operator : SJ/MA  
 Sample : I2914-01  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SB-03-2.5-3.5

Integration Parameters: rteint.p  
 Integrator: RTE  
 Smoothing : ON  
 Sampling : 1  
 Start Thrs: 0.2  
 Stop Thrs : 0  
 Filtering: 5  
 Min Area: 3 % of largest Peak  
 Max Peaks: 100  
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >  
 Peak separation: 5

Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF050217.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

| 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         | 5.043       | 524           | 528         | 545          | rBV      | 286996         | 599137        | 19.29%          | 1.831%        |
| 2         | 5.678       | 631           | 636         | 652          | rBV      | 2012133        | 2858254       | 92.03%          | 8.733%        |
| 3         | 6.248       | 729           | 733         | 741          | rVB      | 25685          | 37891         | 1.22%           | 0.116%        |
| 4         | 7.272       | 901           | 907         | 921          | rBV      | 2023637        | 3069064       | 98.82%          | 9.377%        |
| 5         | 7.389       | 921           | 927         | 939          | rVB      | 2378591        | 3105749       | 100.00%         | 9.489%        |
| 6         | 7.725       | 980           | 984         | 991          | rBV      | 506993         | 556922        | 17.93%          | 1.702%        |
| 7         | 7.966       | 1019          | 1025        | 1035         | rBV      | 1960318        | 2527347       | 81.38%          | 7.722%        |
| 8         | 8.625       | 1131          | 1137        | 1152         | rBV      | 1432876        | 1926220       | 62.02%          | 5.885%        |
| 9         | 9.742       | 1322          | 1327        | 1340         | rBV      | 667301         | 800251        | 25.77%          | 2.445%        |
| 10        | 11.519      | 1623          | 1629        | 1633         | rBV      | 2498707        | 3063322       | 98.63%          | 9.360%        |
| 11        | 12.201      | 1741          | 1745        | 1752         | rBV      | 528962         | 645889        | 20.80%          | 1.973%        |
| 12        | 12.342      | 1766          | 1769        | 1776         | rVB      | 35619          | 47659         | 1.53%           | 0.146%        |
| 13        | 12.571      | 1803          | 1808        | 1823         | rBV2     | 706699         | 897216        | 28.89%          | 2.741%        |
| 14        | 12.819      | 1845          | 1850        | 1857         | rBV3     | 29892          | 43504         | 1.40%           | 0.133%        |
| 15        | 13.419      | 1949          | 1952        | 1956         | rVB      | 43246          | 47395         | 1.53%           | 0.145%        |
| 16        | 13.783      | 2005          | 2014        | 2018         | rBV5     | 21942          | 44890         | 1.45%           | 0.137%        |
| 17        | 13.866      | 2023          | 2028        | 2041         | rBV      | 1239420        | 1713206       | 55.16%          | 5.235%        |
| 18        | 14.107      | 2063          | 2069        | 2073         | rBV      | 43471          | 57640         | 1.86%           | 0.176%        |
| 19        | 14.189      | 2080          | 2083        | 2086         | rBV      | 40653          | 57118         | 1.84%           | 0.175%        |
| 20        | 14.618      | 2150          | 2156        | 2160         | rBV7     | 23709          | 48909         | 1.57%           | 0.149%        |
| 21        | 14.924      | 2203          | 2208        | 2211         | rVB2     | 32789          | 38687         | 1.25%           | 0.118%        |
| 22        | 14.971      | 2211          | 2216        | 2220         | rBV      | 652714         | 801988        | 25.82%          | 2.450%        |
| 23        | 15.007      | 2220          | 2222        | 2228         | rVB      | 88885          | 107369        | 3.46%           | 0.328%        |
| 24        | 15.095      | 2234          | 2237        | 2243         | rBV      | 39515          | 45779         | 1.47%           | 0.140%        |
| 25        | 15.195      | 2248          | 2254        | 2262         | rBV7     | 73358          | 185075        | 5.96%           | 0.565%        |
| 26        | 15.683      | 2331          | 2337        | 2346         | rBV8     | 61576          | 198490        | 6.39%           | 0.606%        |
| 27        | 15.754      | 2346          | 2349        | 2353         | rVB      | 43299          | 46947         | 1.51%           | 0.143%        |
| 28        | 15.795      | 2353          | 2356        | 2358         | rBV      | 47778          | 56268         | 1.81%           | 0.172%        |
| 29        | 15.907      | 2371          | 2375        | 2378         | rVV2     | 84170          | 114945        | 3.70%           | 0.351%        |
| 30        | 15.942      | 2378          | 2381        | 2390         | rVB2     | 190912         | 239723        | 7.72%           | 0.732%        |
| 31        | 16.195      | 2418          | 2424        | 2426         | rBV4     | 36107          | 58497         | 1.88%           | 0.179%        |
| 32        | 16.412      | 2457          | 2461        | 2465         | rBV4     | 21179          | 32044         | 1.03%           | 0.098%        |
| 33        | 16.548      | 2479          | 2484        | 2488         | rBV      | 100008         | 132071        | 4.25%           | 0.404%        |
| 34        | 16.589      | 2488          | 2491        | 2495         | rVB2     | 51579          | 68450         | 2.20%           | 0.209%        |

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 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SB-03-2.5-3.5

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF050217.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 35 | 16.665 | 2500 | 2504 | 2510 | rVB2 | 47564   | 63528   | 2.05%  | 0.194% |
| 36 | 16.748 | 2514 | 2518 | 2524 | rVV  | 402035  | 437956  | 14.10% | 1.338% |
| 37 | 16.848 | 2532 | 2535 | 2538 | rBV4 | 37255   | 46637   | 1.50%  | 0.142% |
| 38 | 16.883 | 2538 | 2541 | 2544 | rVB  | 45887   | 47928   | 1.54%  | 0.146% |
| 39 | 17.048 | 2562 | 2569 | 2575 | rVV  | 458112  | 564786  | 18.19% | 1.726% |
| 40 | 17.177 | 2587 | 2591 | 2594 | rBV2 | 30973   | 37179   | 1.20%  | 0.114% |
| 41 | 17.301 | 2606 | 2612 | 2621 | rBV  | 2085993 | 2472132 | 79.60% | 7.553% |
| 42 | 17.377 | 2621 | 2625 | 2628 | rVB2 | 52666   | 73029   | 2.35%  | 0.223% |
| 43 | 17.512 | 2645 | 2648 | 2655 | rVB2 | 141715  | 159158  | 5.12%  | 0.486% |
| 44 | 17.601 | 2660 | 2663 | 2667 | rBV  | 120721  | 116036  | 3.74%  | 0.355% |
| 45 | 17.648 | 2667 | 2671 | 2676 | rBV2 | 74391   | 107234  | 3.45%  | 0.328% |
| 46 | 17.759 | 2688 | 2690 | 2694 | rVB  | 61551   | 61207   | 1.97%  | 0.187% |
| 47 | 17.801 | 2695 | 2697 | 2700 | rVB  | 37121   | 36482   | 1.17%  | 0.111% |
| 48 | 18.095 | 2745 | 2747 | 2750 | rVB3 | 43088   | 39415   | 1.27%  | 0.120% |
| 49 | 18.306 | 2779 | 2783 | 2787 | rBV3 | 88237   | 151931  | 4.89%  | 0.464% |
| 50 | 18.377 | 2793 | 2795 | 2798 | rVB  | 37773   | 34525   | 1.11%  | 0.105% |
| 51 | 18.530 | 2818 | 2821 | 2823 | rBV3 | 91060   | 101339  | 3.26%  | 0.310% |
| 52 | 18.559 | 2823 | 2826 | 2828 | rVV  | 78355   | 120666  | 3.89%  | 0.369% |
| 53 | 18.589 | 2828 | 2831 | 2833 | rVV  | 169130  | 187626  | 6.04%  | 0.573% |
| 54 | 18.636 | 2833 | 2839 | 2842 | rVV2 | 615121  | 935660  | 30.13% | 2.859% |
| 55 | 18.665 | 2842 | 2844 | 2848 | rVV  | 242649  | 283811  | 9.14%  | 0.867% |
| 56 | 18.701 | 2848 | 2850 | 2857 | rVB  | 88063   | 105205  | 3.39%  | 0.321% |
| 57 | 19.142 | 2922 | 2925 | 2929 | rVB2 | 52878   | 63180   | 2.03%  | 0.193% |
| 58 | 19.253 | 2941 | 2944 | 2947 | rVB2 | 47556   | 48983   | 1.58%  | 0.150% |
| 59 | 19.900 | 3049 | 3054 | 3056 | rBV  | 232721  | 333056  | 10.72% | 1.018% |
| 60 | 20.012 | 3070 | 3073 | 3077 | rVB  | 53279   | 65807   | 2.12%  | 0.201% |
| 61 | 20.065 | 3078 | 3082 | 3085 | rBV  | 100517  | 127335  | 4.10%  | 0.389% |
| 62 | 20.189 | 3099 | 3103 | 3109 | rVB  | 185605  | 219657  | 7.07%  | 0.671% |
| 63 | 20.247 | 3109 | 3113 | 3118 | rBV  | 214989  | 279188  | 8.99%  | 0.853% |
| 64 | 20.300 | 3118 | 3122 | 3126 | rBV  | 464854  | 533979  | 17.19% | 1.632% |
| 65 | 20.783 | 3201 | 3204 | 3211 | rVB2 | 68216   | 115468  | 3.72%  | 0.353% |
| 66 | 21.283 | 3285 | 3289 | 3295 | rVB5 | 44626   | 67745   | 2.18%  | 0.207% |
| 67 | 21.594 | 3337 | 3342 | 3348 | rVB2 | 101599  | 192794  | 6.21%  | 0.589% |
| 68 | 21.747 | 3365 | 3368 | 3371 | rBV2 | 18715   | 35180   | 1.13%  | 0.107% |
| 69 | 21.965 | 3400 | 3405 | 3413 | rBV2 | 71362   | 149151  | 4.80%  | 0.456% |
| 70 | 23.883 | 3725 | 3731 | 3735 | rBV6 | 18185   | 39869   | 1.28%  | 0.122% |

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Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-03-2.5-3.5

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:\HPCHEM1\BNA\_F\METHODS\8270-BF050217.M  
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

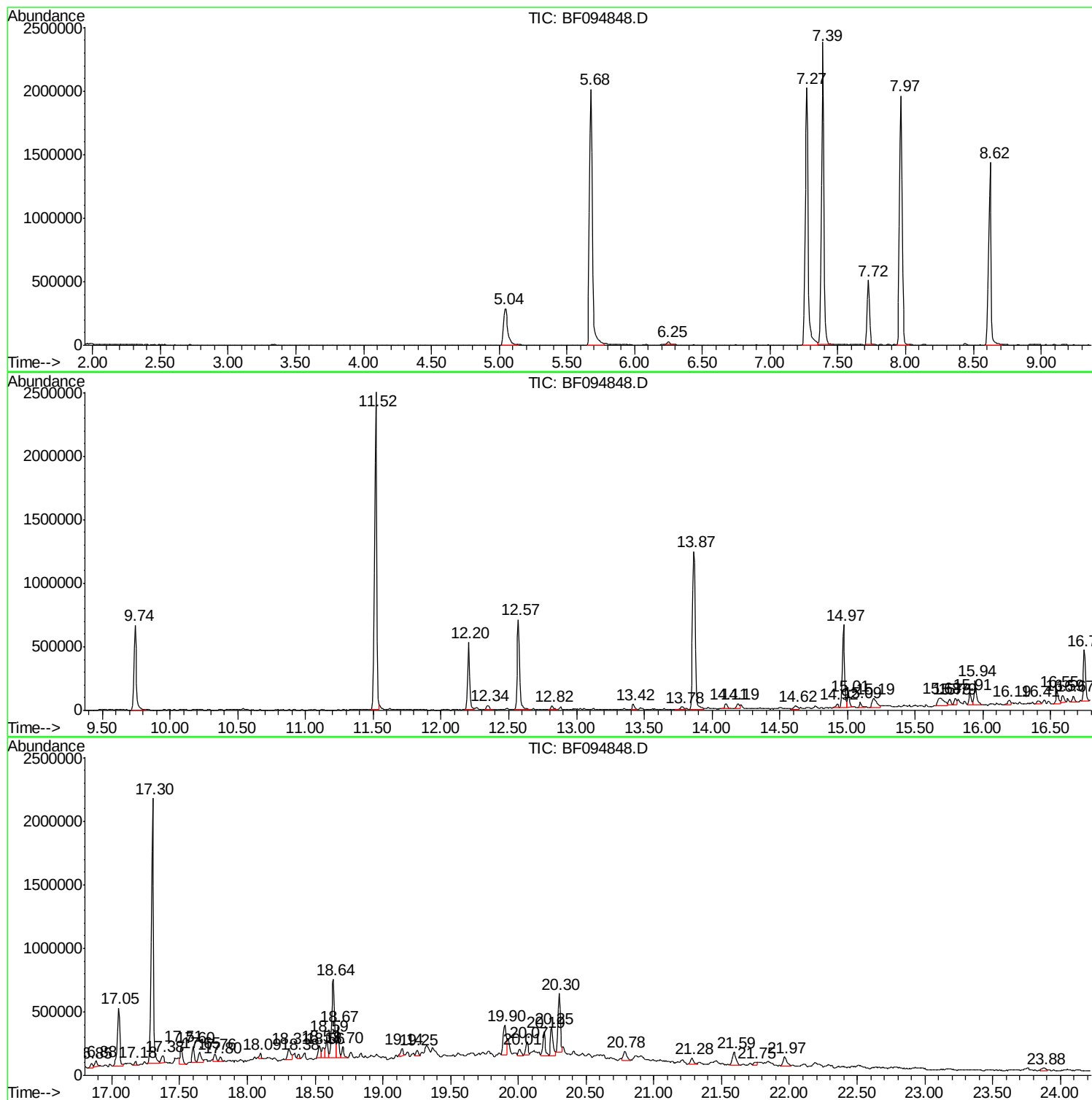
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Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF050317\  
Data File : BF094848.D  
Acq On : 3 May 2017 22:42  
Operator : SJ/MA  
Sample : I2914-01  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-03-2.5-3.5

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF050217.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF050317\  
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Sample : I2914-01  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-03-2.5-3.5

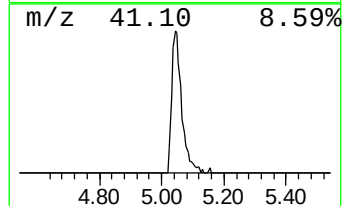
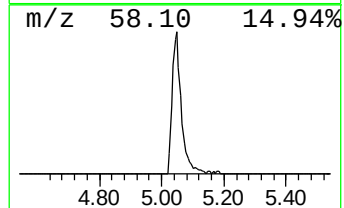
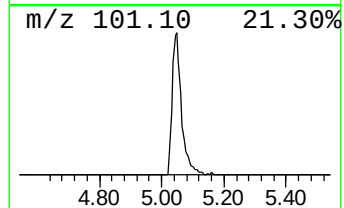
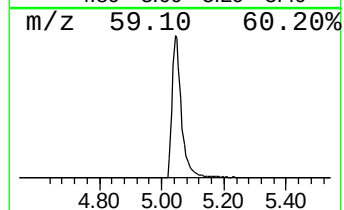
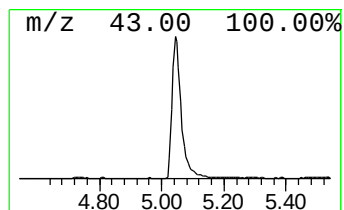
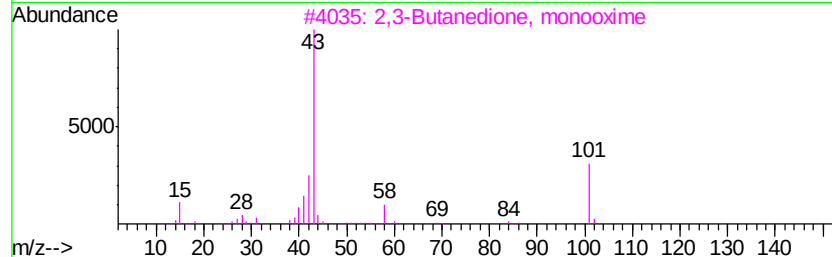
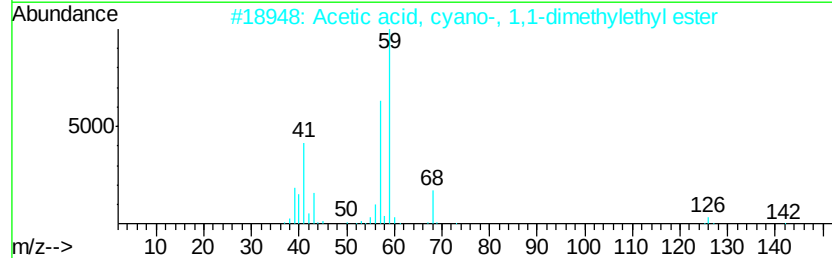
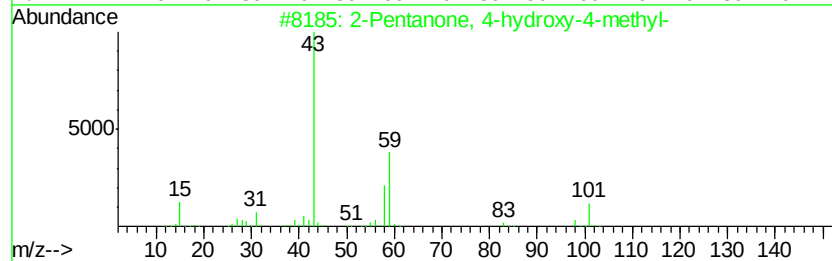
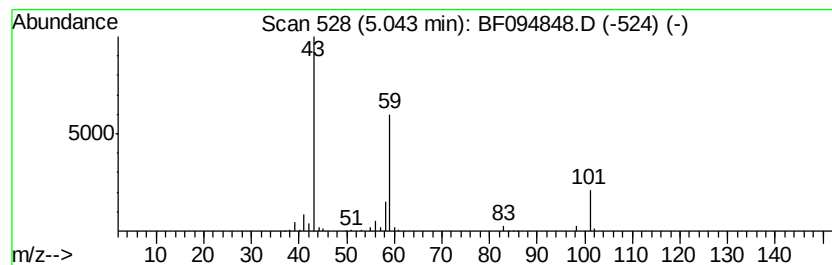
Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF050217.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 5.04 | 21.52 ng | 599137 | 1,4-Dichlorobenzene-d4 | 7.72 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2  | 000123-42-2 | 72   |
| 2    |    | Acetic acid, cyano-, 1,1-dimethy... | 141 | C7H11NO2 | 001116-98-9 | 23   |
| 3    |    | 2,3-Butanedione, monooxime          | 101 | C4H7NO2  | 000057-71-6 | 17   |
| 4    |    | Butane, 1-ethoxy-                   | 102 | C6H14O   | 000628-81-9 | 9    |
| 5    |    | Morpholine, 4-methyl-               | 101 | C5H11NO  | 000109-02-4 | 9    |



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Instrument :  
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ClientSampleId :  
SB-03-2.5-3.5

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF050217.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

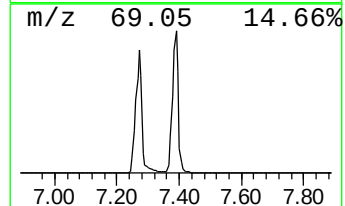
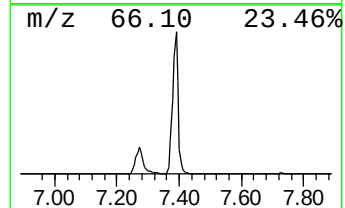
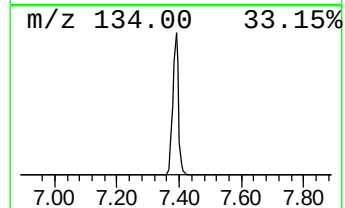
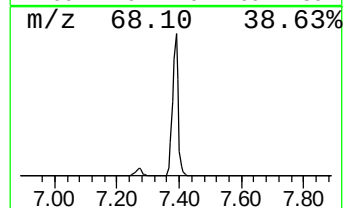
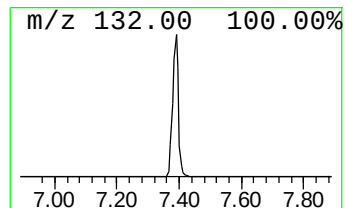
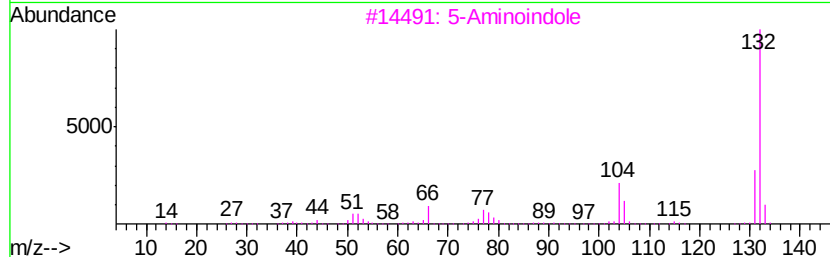
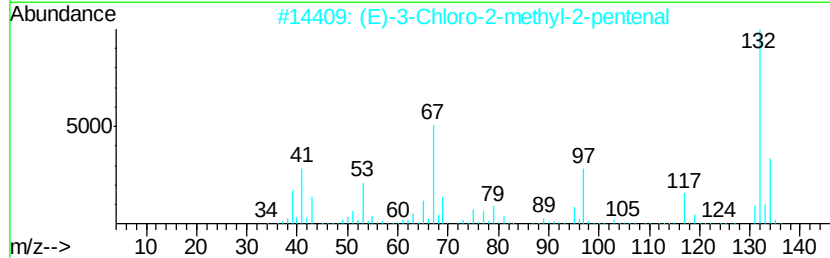
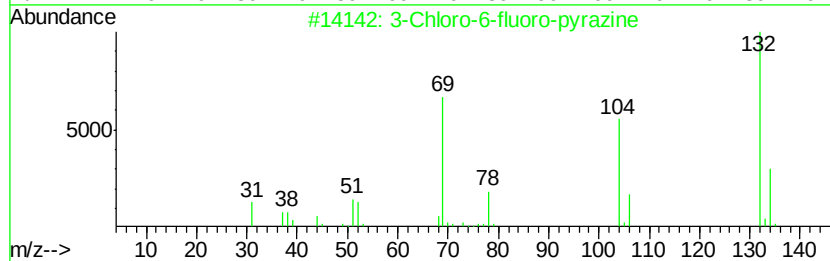
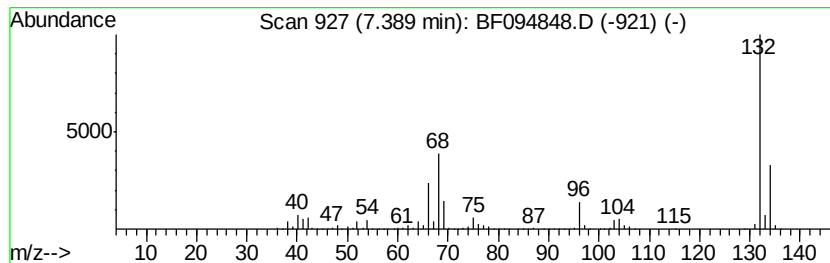
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 unknown7.39 Concentration Rank 1

| R.T. | EstConc   | Area    | Relative to ISTD       | R.T. |
|------|-----------|---------|------------------------|------|
| 7.39 | 111.53 ng | 3105750 | 1,4-Dichlorobenzene-d4 | 7.72 |

| Hit# | of | Tentative ID                     | MW  | MolForm   | CAS#         | Qual |
|------|----|----------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine       | 132 | C4H2ClFN2 | 1000146-10-7 | 27   |
| 2    |    | (E)-3-Chloro-2-methyl-2-pentenal | 132 | C6H9ClO   | 031357-76-3  | 12   |
| 3    |    | 5-Aminoindole                    | 132 | C8H8N2    | 005192-03-0  | 12   |
| 4    |    | Tranylcypromine-propionyl        | 189 | C12H15NO  | 1000123-86-3 | 10   |
| 5    |    | 5-Fluoro-2-chloropyrimidine      | 132 | C4H2ClFN2 | 062802-42-0  | 9    |



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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF050217.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

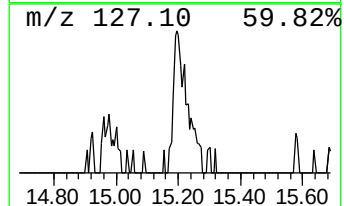
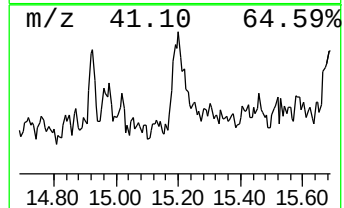
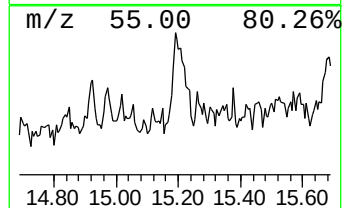
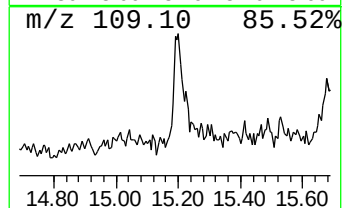
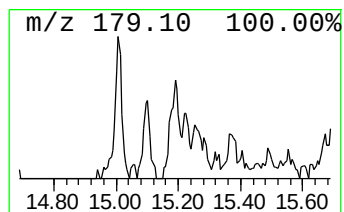
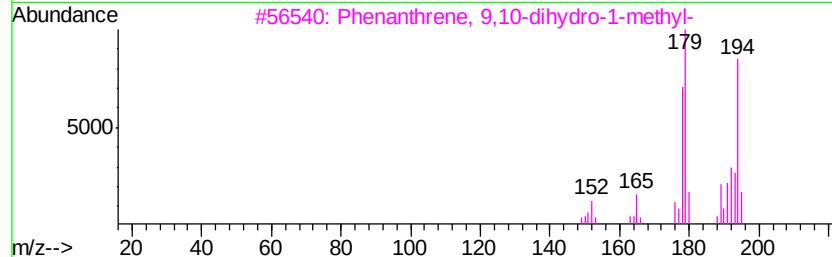
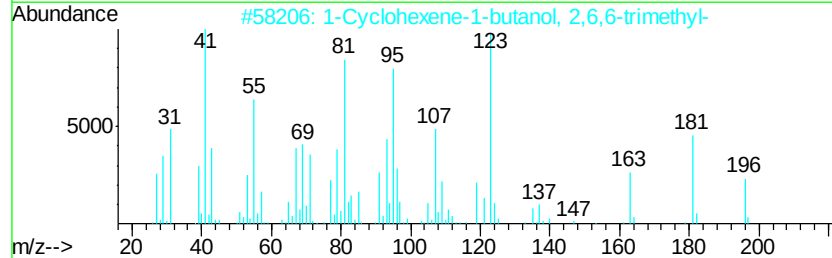
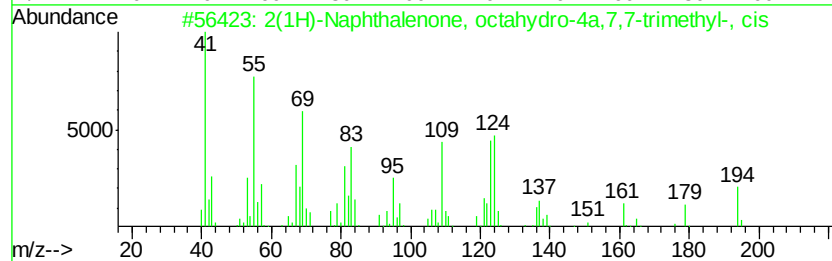
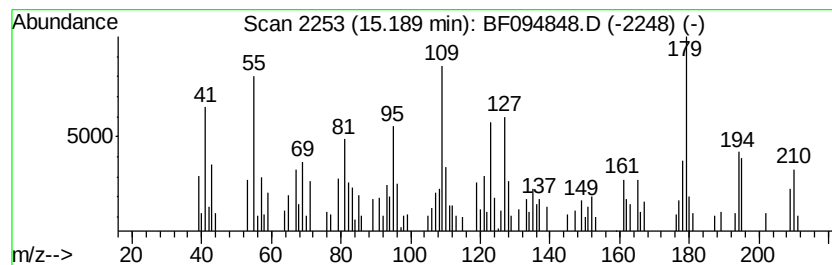
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 4 2(1H)-Naphthalenone, octahy... Concentration Rank 8

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 15.19 | 4.62 ng | 185075 | Phenanthrene-d10 | 14.97 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 2(1H)-Naphthalenone, octahydro-4... | 194 | C13H22O | 007056-56-6 | 50   |
| 2    |    | 1-Cyclohexene-1-butanol, 2,6,6-t... | 196 | C13H24O | 054344-91-1 | 42   |
| 3    |    | Phenanthrene, 9,10-dihydro-1-met... | 194 | C15H14  | 095676-48-5 | 38   |
| 4    |    | 9H-Fluorene, 9,9-dimethyl-          | 194 | C15H14  | 004569-45-3 | 18   |
| 5    |    | 1,4-Methanoazulene-9-methanol, d... | 222 | C15H26O | 001139-17-9 | 15   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF050317\  
Data File : BF094848.D  
Acq On : 3 May 2017 22:42  
Operator : SJ/MA  
Sample : I2914-01  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-03-2.5-3.5

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

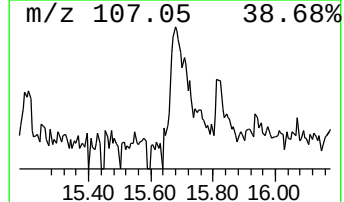
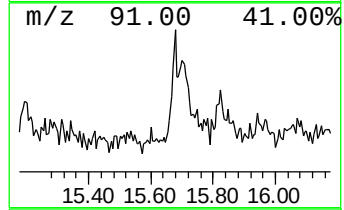
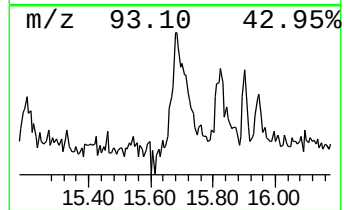
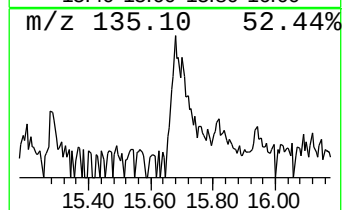
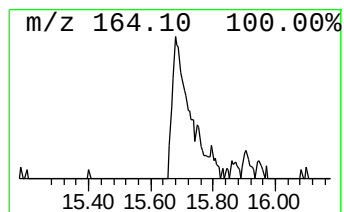
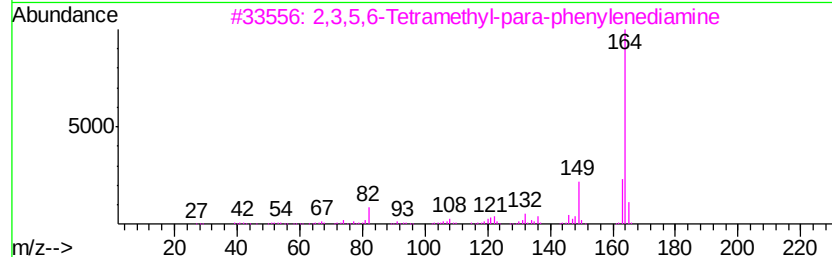
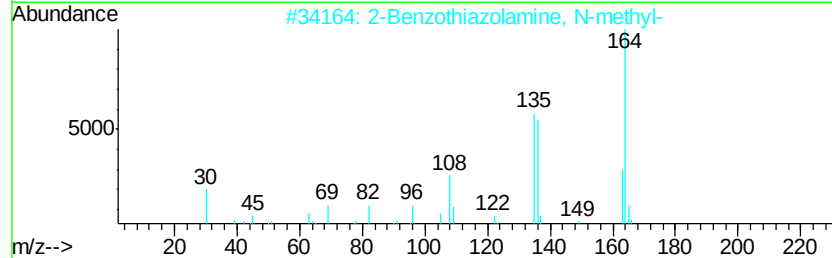
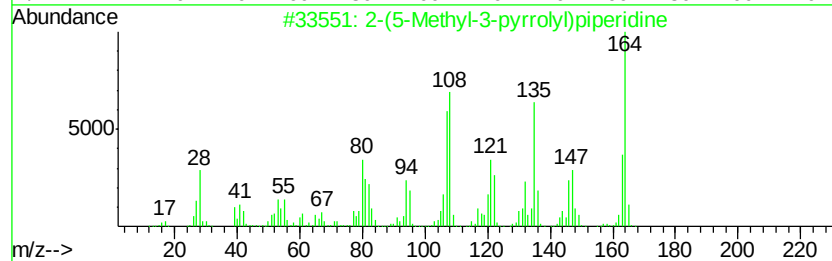
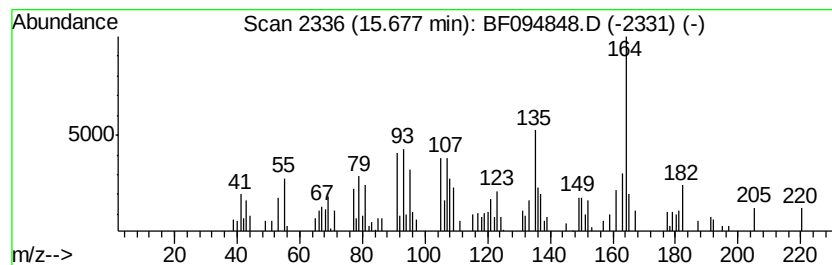
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TIC Integration Parameters: LSCINT.P

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Peak Number 5 2-(5-Methyl-3-pyrrolyl)pipe... Concentration Rank 6

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 15.68 | 4.95 ng | 198490 | Phenanthrene-d10 | 14.97 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | 2-(5-Methyl-3-pyrrolyl)piperidine   | 164 | C10H16N2 | 1000245-18-2 | 55   |
| 2    |    | 2-Benzothiazolamine, N-methyl-      | 164 | C8H8N2S  | 016954-69-1  | 38   |
| 3    |    | 2,3,5,6-Tetramethyl-para-phenyle... | 164 | C10H16N2 | 003102-87-2  | 35   |
| 4    |    | 1H-Indene-1,5(6H)-dione, 2,3,7,7... | 164 | C10H12O2 | 019576-08-0  | 30   |
| 5    |    | s-Triazolo[4,3-a]pyridine, 8-nitro- | 164 | C6H4N4O2 | 031040-09-2  | 30   |





Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF050317\  
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Operator : SJ/MA  
Sample : I2914-01  
Misc :  
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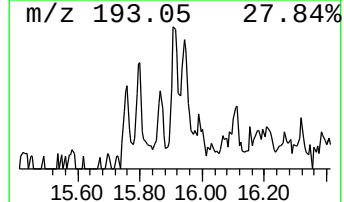
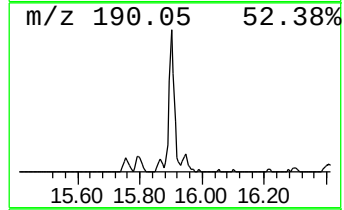
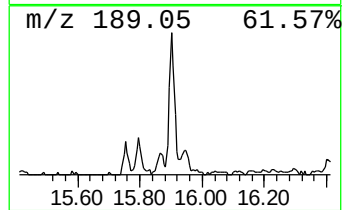
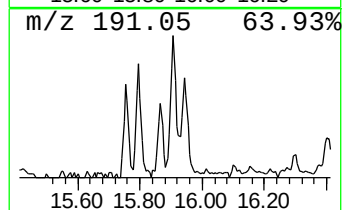
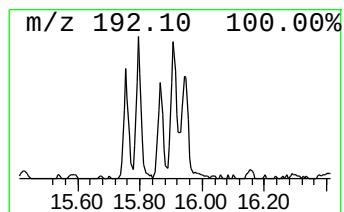
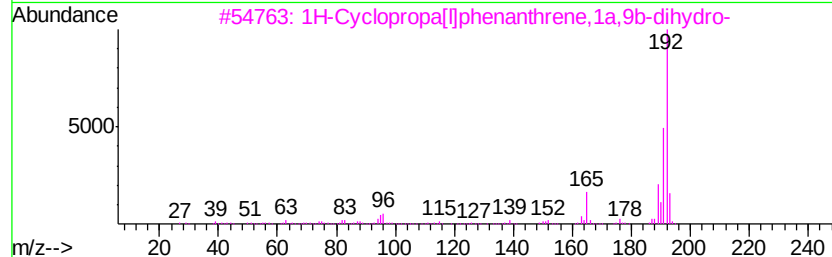
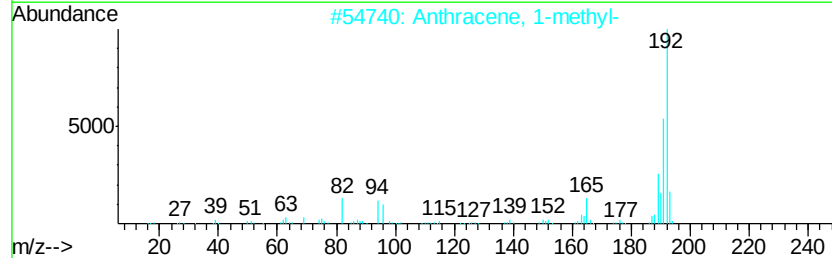
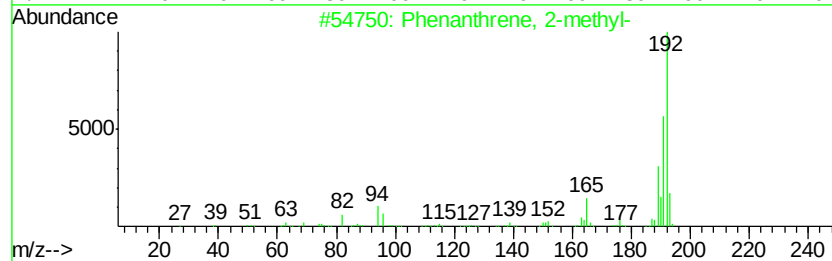
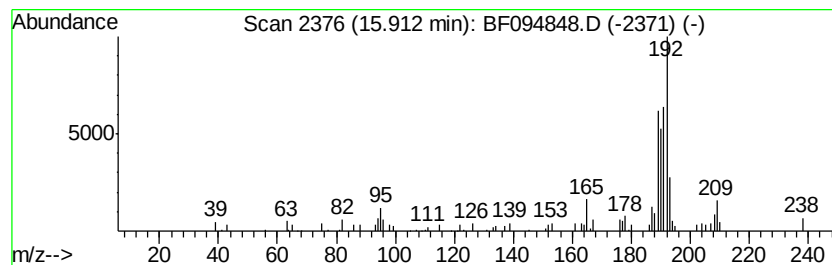
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SB-03-2.5-3.5

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 6 Phenanthrene, 2-methyl- Concentration Rank 13

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|-------------|------|
| 15.91     | 2.87 ng                             | 114945 | Phenanthrene-d10 | 14.97       |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual |
| 1         | Phenanthrene, 2-methyl-             | 192    | C15H12           | 002531-84-2 | 55   |
| 2         | Anthracene, 1-methyl-               | 192    | C15H12           | 000610-48-0 | 50   |
| 3         | 1H-Cyclopropa[l]phenanthrene,1a,... | 192    | C15H12           | 000949-41-7 | 50   |
| 4         | Phenanthrene, 1-methyl-             | 192    | C15H12           | 000832-69-9 | 50   |
| 5         | Phenanthrene, 4-methyl-             | 192    | C15H12           | 000832-64-4 | 50   |



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Operator : SJ/MA  
Sample : I2914-01  
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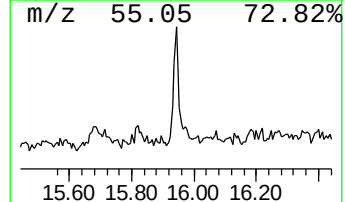
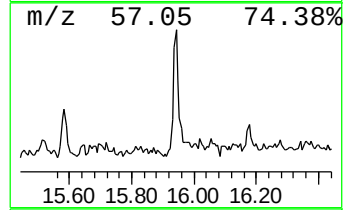
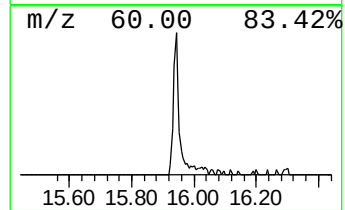
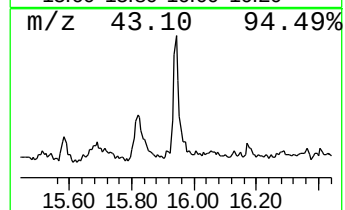
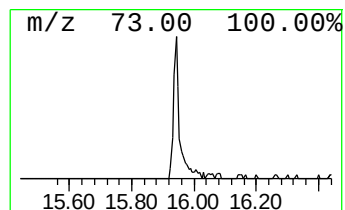
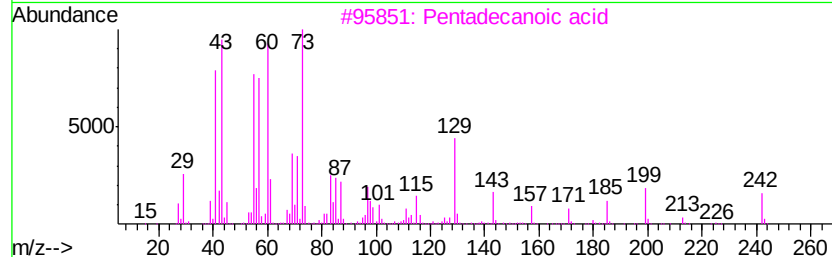
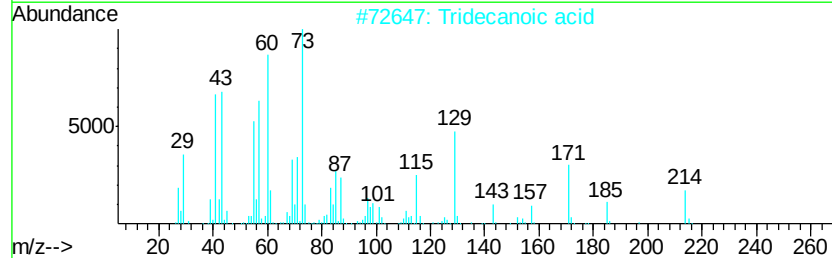
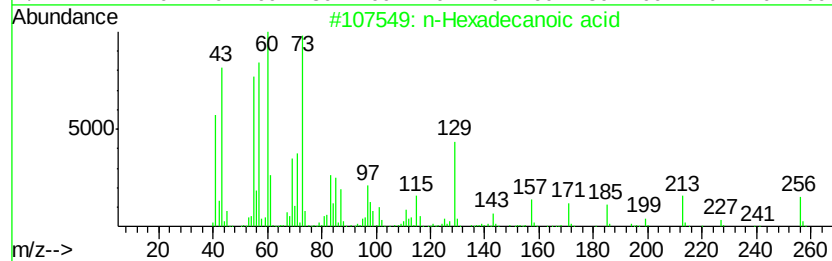
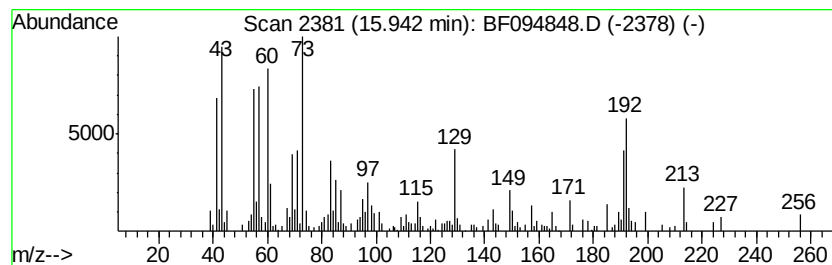
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ClientSampleId :  
SB-03-2.5-3.5

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 n-Hexadecanoic acid Concentration Rank 5

| R.T.      | EstConc             | Area   | Relative to ISTD | R.T.        |      |
|-----------|---------------------|--------|------------------|-------------|------|
| 15.94     | 5.98 ng             | 239723 | Phenanthrene-d10 | 14.97       |      |
| Hit# of 5 | Tentative ID        | MW     | MolForm          | CAS#        | Qual |
| 1         | n-Hexadecanoic acid | 256    | C16H32O2         | 000057-10-3 | 98   |
| 2         | Tridecanoic acid    | 214    | C13H26O2         | 000638-53-9 | 78   |
| 3         | Pentadecanoic acid  | 242    | C15H30O2         | 001002-84-2 | 60   |
| 4         | n-Decanoic acid     | 172    | C10H20O2         | 000334-48-5 | 50   |
| 5         | Tetradecanoic acid  | 228    | C14H28O2         | 000544-63-8 | 49   |



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Operator : SJ/MA  
Sample : I2914-01  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-03-2.5-3.5

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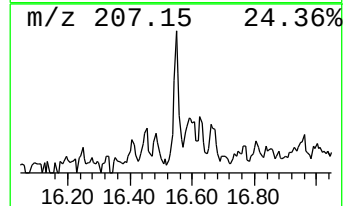
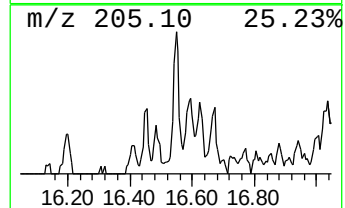
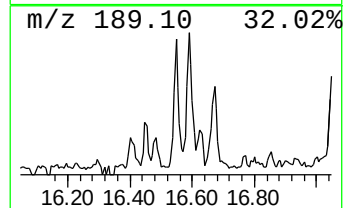
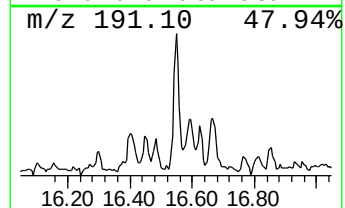
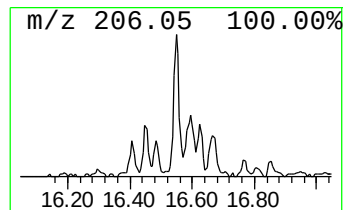
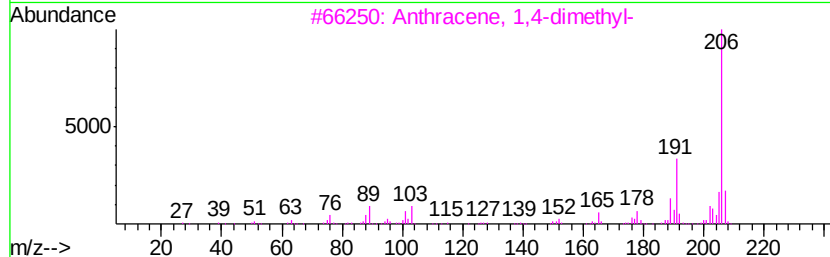
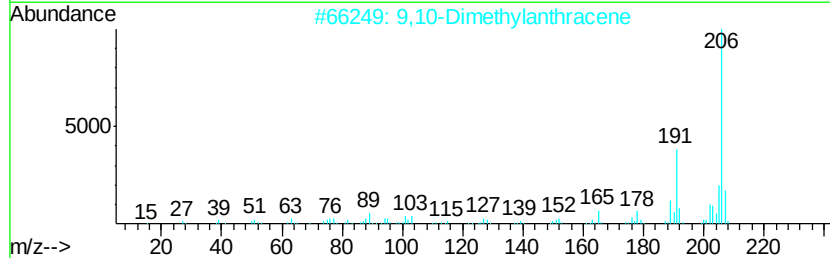
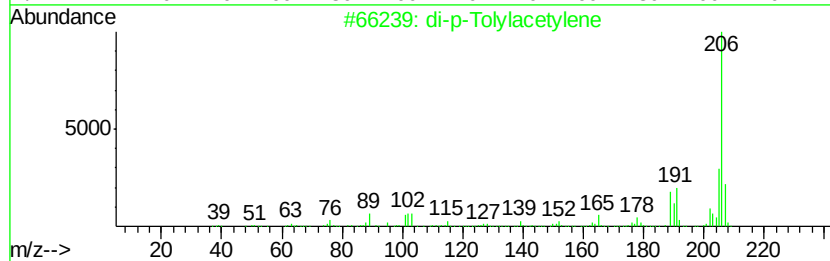
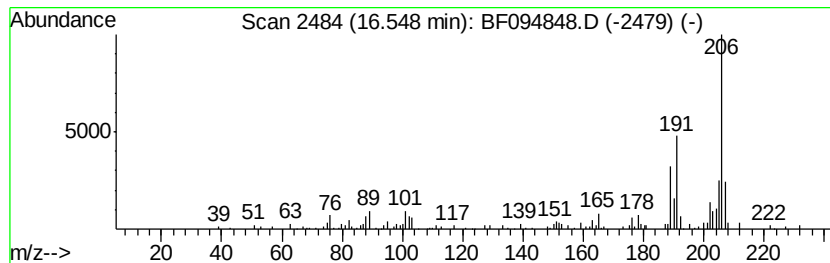
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TIC Integration Parameters: LSCINT.P

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Peak Number 8 di-p-Tolylacetylene Concentration Rank 11

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 16.55 | 3.29 ng | 132071 | Phenanthrene-d10 | 14.97 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | di-p-Tolylacetylene         | 206 | C16H14  | 002789-88-0 | 92   |
| 2    |    | 9,10-Dimethylantracene      | 206 | C16H14  | 000781-43-1 | 91   |
| 3    |    | Anthracene, 1,4-dimethyl-   | 206 | C16H14  | 000781-92-0 | 91   |
| 4    |    | Phenanthrene, 2,5-dimethyl- | 206 | C16H14  | 003674-66-6 | 90   |
| 5    |    | Phenanthrene, 3,6-dimethyl- | 206 | C16H14  | 001576-67-6 | 89   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF050317\  
Data File : BF094848.D  
Acq On : 3 May 2017 22:42  
Operator : SJ/MA  
Sample : I2914-01  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
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ClientSampleId :  
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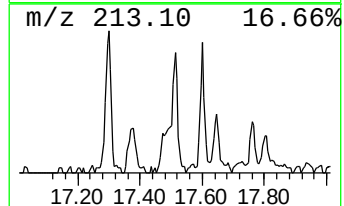
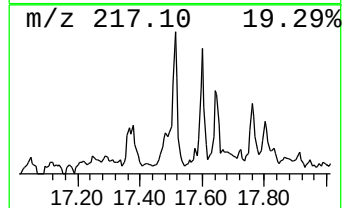
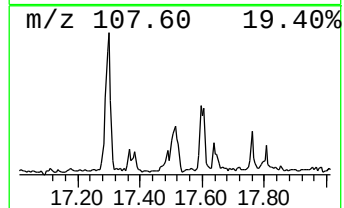
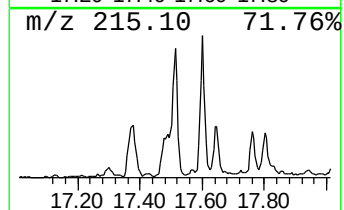
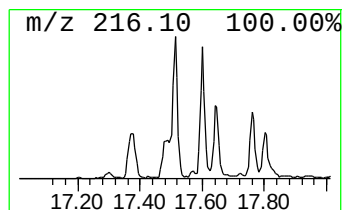
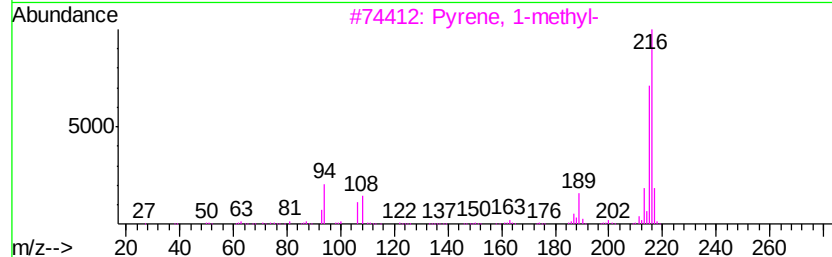
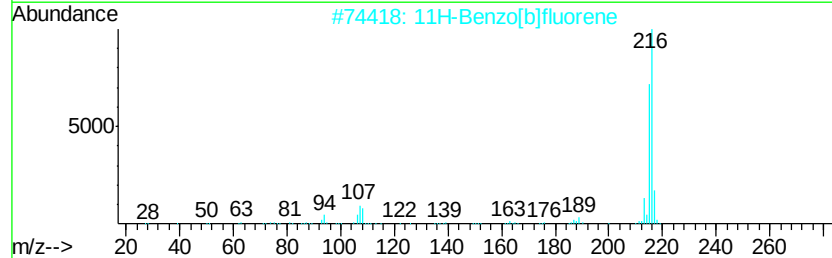
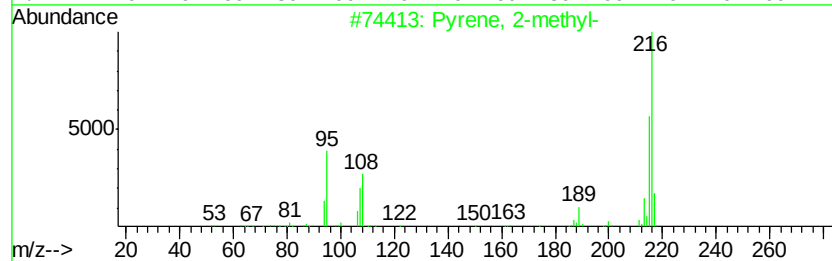
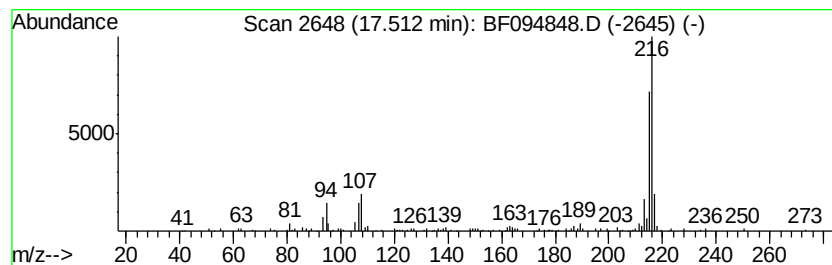
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TIC Integration Parameters: LSCINT.P

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Peak Number 9 Pyrene, 2-methyl- Concentration Rank 10

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 17.51 | 3.40 ng | 159158 | Chrysene-d12     | 18.64 |

| Hit# of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|---------|---|-------------------------|-----|---------|-------------|------|
| 1       |   | Pyrene, 2-methyl-       | 216 | C17H12  | 003442-78-2 | 94   |
| 2       |   | 11H-Benzo[b]fluorene    | 216 | C17H12  | 000243-17-4 | 91   |
| 3       |   | Pyrene, 1-methyl-       | 216 | C17H12  | 002381-21-7 | 90   |
| 4       |   | 11H-Benzo[a]fluorene    | 216 | C17H12  | 000238-84-6 | 90   |
| 5       |   | Fluoranthene, 2-methyl- | 216 | C17H12  | 033543-31-6 | 68   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF050317\  
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Sample : I2914-01  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

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ClientSampleId :  
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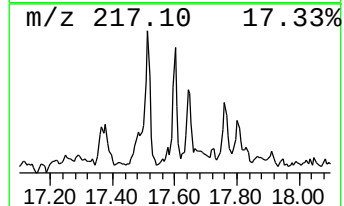
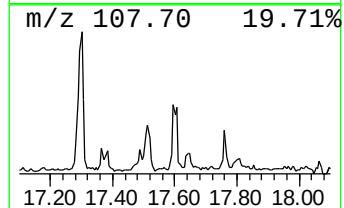
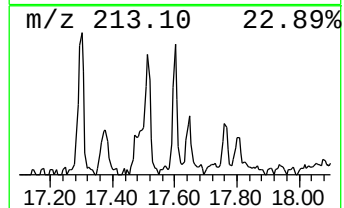
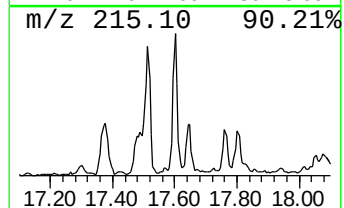
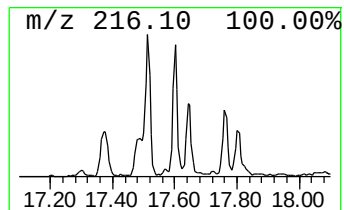
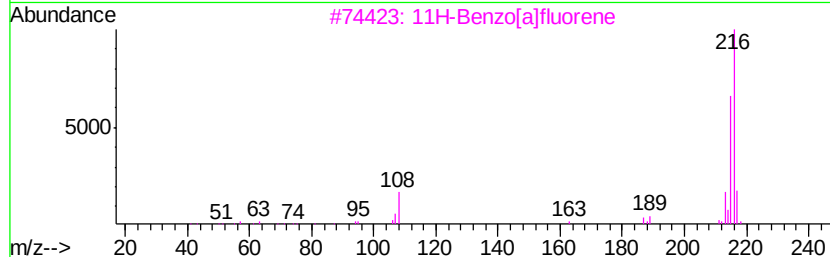
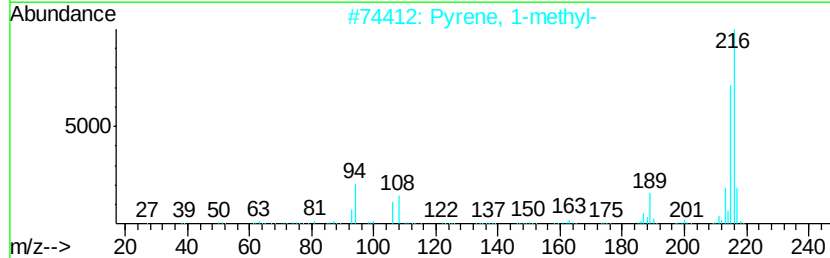
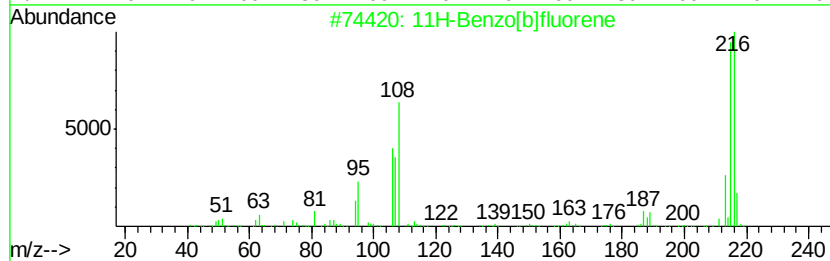
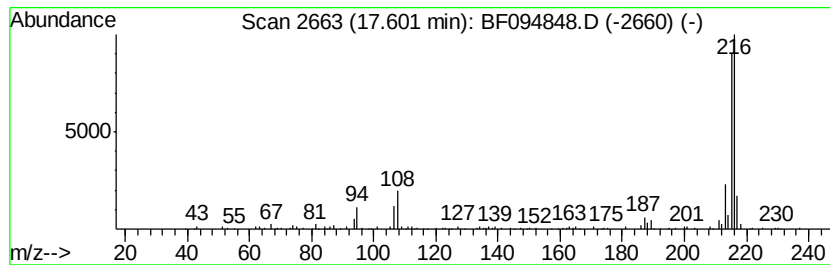
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 11H-Benzo[b]fluorene Concentration Rank 16

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 17.60 | 2.48 ng | 116036 | Chrysene-d12     | 18.64 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | 11H-Benzo[b]fluorene    | 216 | C17H12  | 000243-17-4 | 91   |
| 2    |    |   | Pyrene, 1-methyl-       | 216 | C17H12  | 002381-21-7 | 91   |
| 3    |    |   | 11H-Benzo[a]fluorene    | 216 | C17H12  | 000238-84-6 | 91   |
| 4    |    |   | Fluoranthene, 2-methyl- | 216 | C17H12  | 033543-31-6 | 90   |
| 5    |    |   | 7H-Benzo[c]fluorene     | 216 | C17H12  | 000205-12-9 | 81   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF050317\  
Data File : BF094848.D  
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Sample : I2914-01  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
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ClientSampleId :  
SB-03-2.5-3.5

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

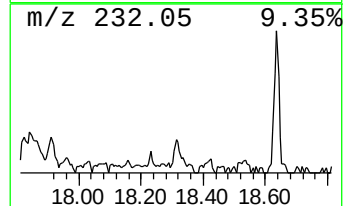
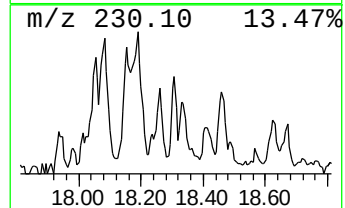
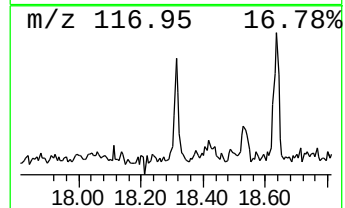
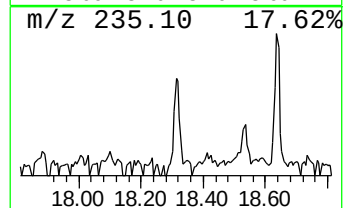
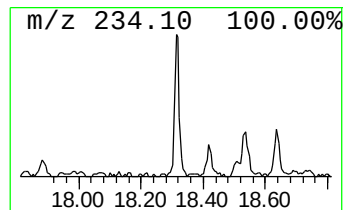
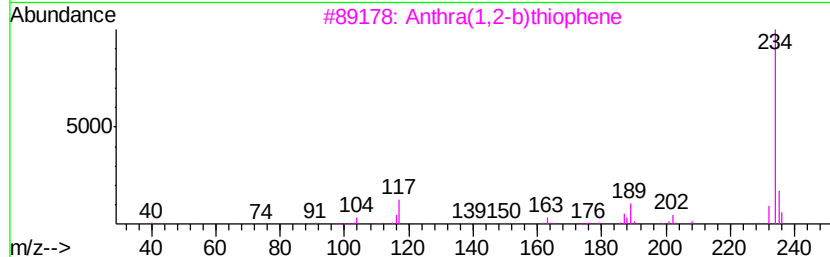
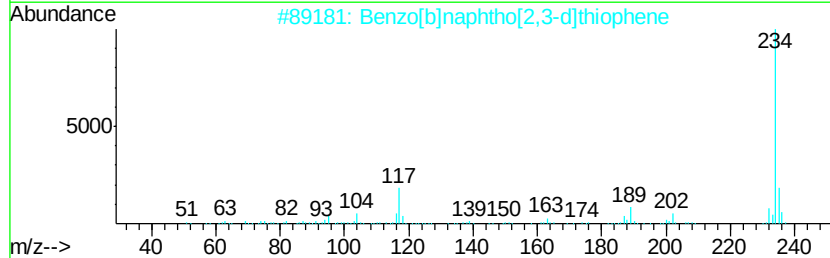
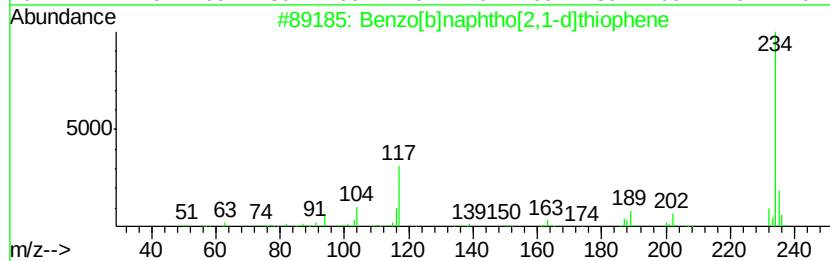
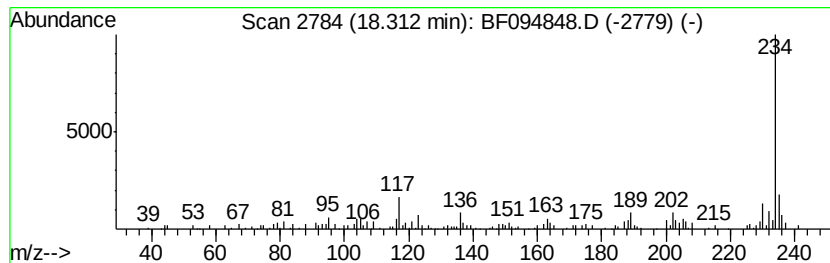
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TIC Integration Parameters: LSCINT.P

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Peak Number 12 unknown18.31 Concentration Rank 12

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 18.31 | 3.25 ng | 151931 | Chrysene-d12     | 18.64 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|------|----|-------------------------------------|-----|------------|-------------|------|
| 1    | 5  | Benzo[b]naphtho[2,1-d]thiophene     | 234 | C16H10S    | 000239-35-0 | 47   |
| 2    |    | Benzo[b]naphtho[2,3-d]thiophene     | 234 | C16H10S    | 000243-46-9 | 38   |
| 3    |    | Anthra(1,2-b)thiophene              | 234 | C16H10S    | 000227-86-1 | 38   |
| 4    |    | 4-Imidazolidinone, 5-(1-methylet... | 234 | C12H14N2OS | 004333-20-4 | 25   |
| 5    |    | Benzo[b]naphtho[1,2-d]thiophene     | 234 | C16H10S    | 000205-43-6 | 25   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF050317\  
Data File : BF094848.D  
Acq On : 3 May 2017 22:42  
Operator : SJ/MA  
Sample : I2914-01  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-03-2.5-3.5

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF050217.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

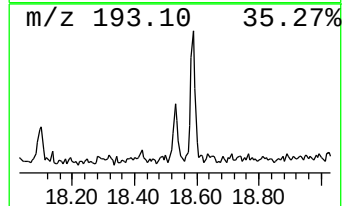
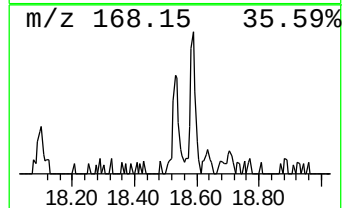
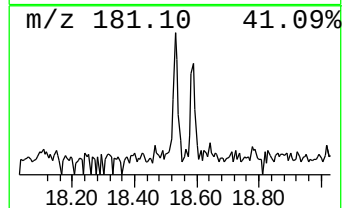
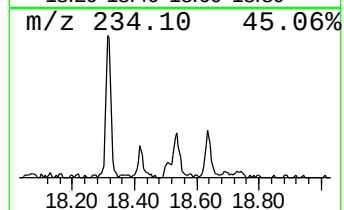
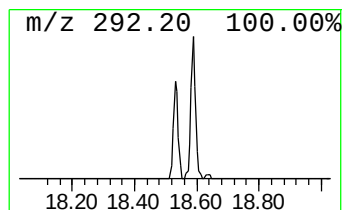
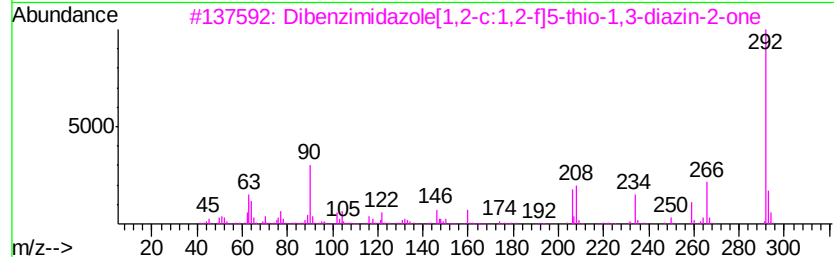
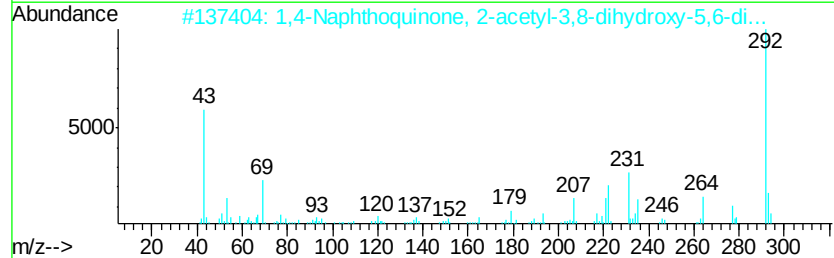
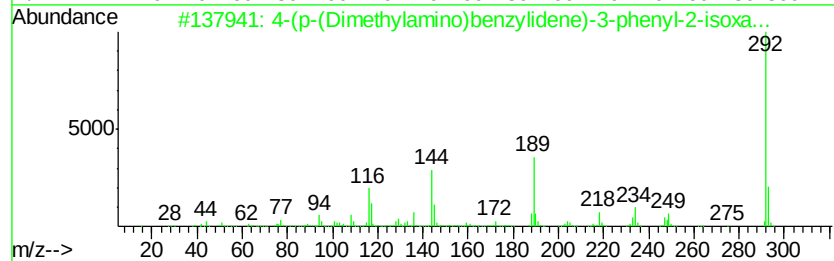
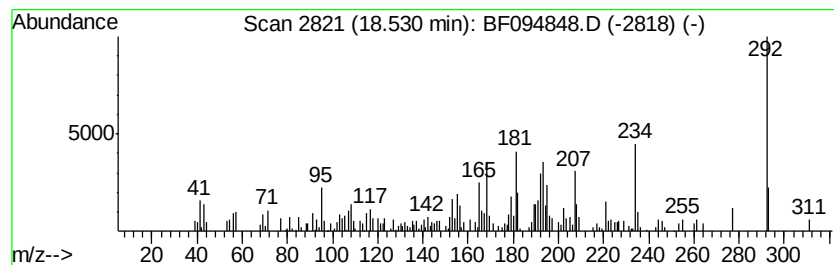
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TIC Integration Parameters: LSCINT.P

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Peak Number 13 unknown18.53 Concentration Rank 19

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 18.53 | 2.17 ng | 101339 | Chrysene-d12     | 18.64 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | 4-(p-(Dimethylamino)benzylidene)... | 292 | C18H16N2O2 | 065156-10-7  | 27   |
| 2    |    |   | 1,4-Naphthoquinone, 2-acetyl-3,8... | 292 | C14H12O7   | 014090-53-0  | 27   |
| 3    |    |   | Dibenzimidazole[1,2-c:1,2-f]5-th... | 292 | C15H8N4OS  | 1000211-46-2 | 27   |
| 4    |    |   | Benzoxazole, 2-[(2,3-dihydro-3,5... | 292 | C18H16N2O2 | 024261-18-5  | 27   |
| 5    |    |   | 1,1'-Biphenyl, 4,4-bis(1-pyrroli... | 292 | C20H24N2   | 1000193-54-6 | 27   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF050317\  
Data File : BF094848.D  
Acq On : 3 May 2017 22:42  
Operator : SJ/MA  
Sample : I2914-01  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-03-2.5-3.5

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF050217.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

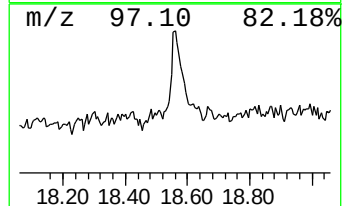
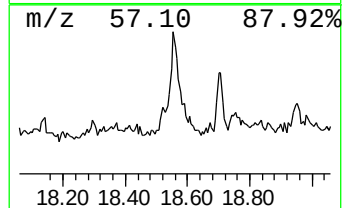
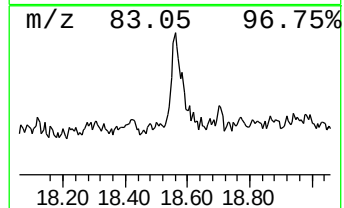
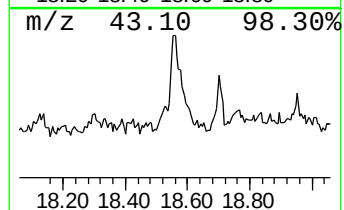
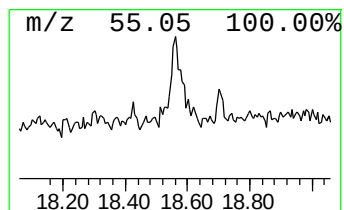
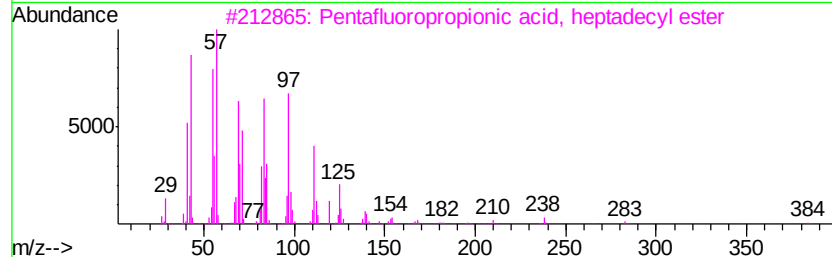
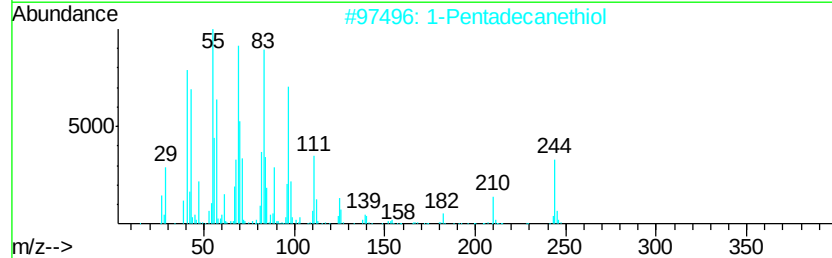
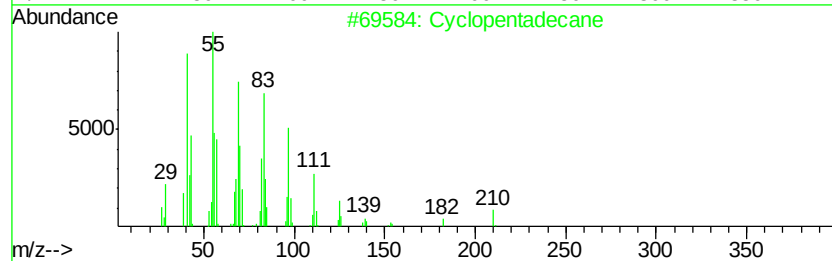
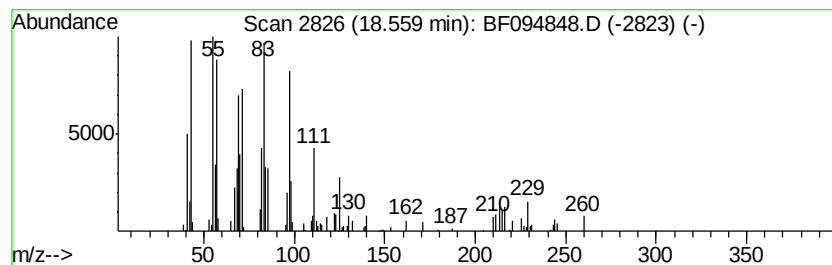
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 14 Cyclopentadecane Concentration Rank 14

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 18.56 | 2.58 ng | 120666 | Chrysene-d12     | 18.64 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|------|----|-------------------------------------|-----|------------|-------------|------|
| 1    | 5  | Cyclopentadecane                    | 210 | C15H30     | 000295-48-7 | 96   |
| 2    |    | 1-Pentadecanethiol                  | 244 | C15H32S    | 025276-70-4 | 93   |
| 3    |    | Pentafluoropropionic acid, hepta... | 402 | C20H35F5O2 | 959218-78-1 | 91   |
| 4    |    | Behenic alcohol                     | 326 | C22H46O    | 000661-19-8 | 87   |
| 5    |    | n-Tetracosanol-1                    | 354 | C24H50O    | 000506-51-4 | 87   |





Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF050317\  
Data File : BF094848.D  
Acq On : 3 May 2017 22:42  
Operator : SJ/MA  
Sample : I2914-01  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-03-2.5-3.5

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF050217.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

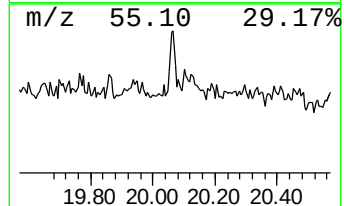
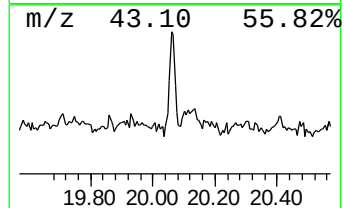
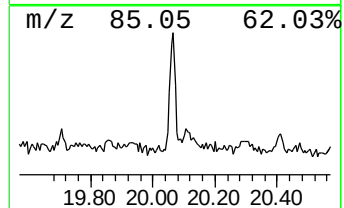
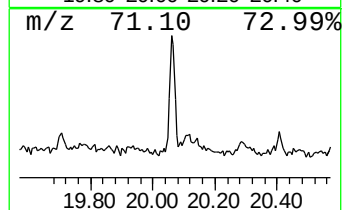
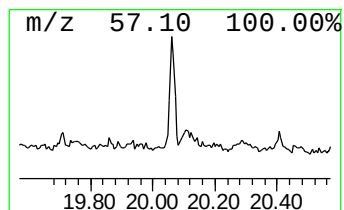
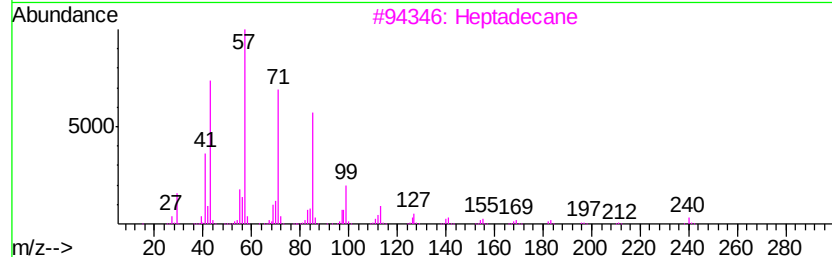
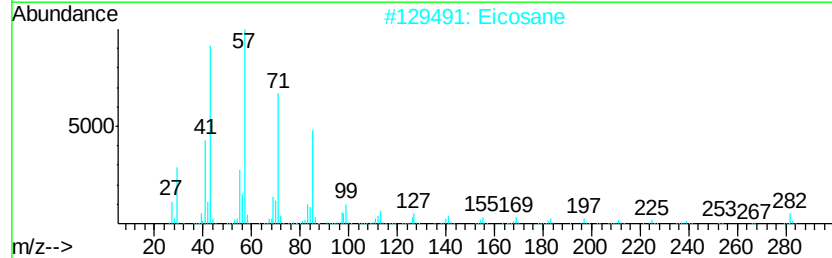
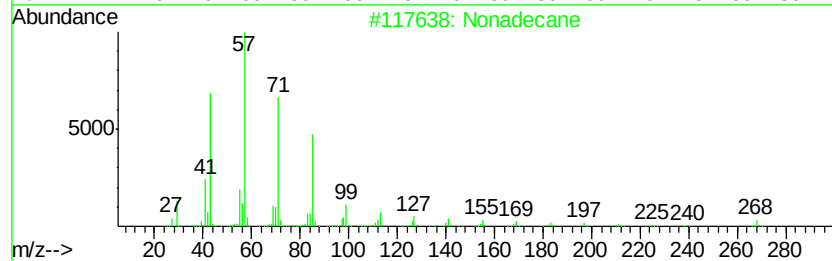
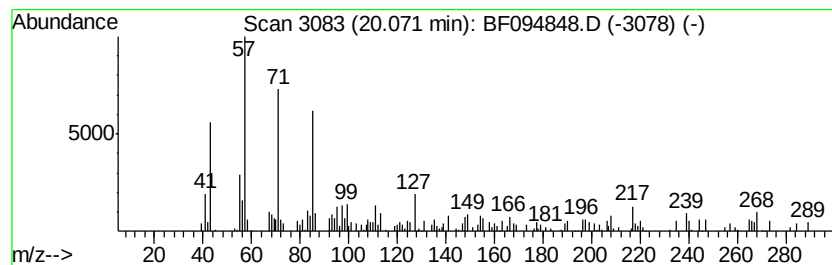
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 16 Nonadecane Concentration Rank 7

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 20.07 | 4.77 ng | 127335 | Perylene-d12     | 20.30 |

| Hit# of | 5           | Tentative ID | MW  | MolForm | CAS#        | Qual |
|---------|-------------|--------------|-----|---------|-------------|------|
| 1       | Nonadecane  |              | 268 | C19H40  | 000629-92-5 | 93   |
| 2       | Eicosane    |              | 282 | C20H42  | 000112-95-8 | 91   |
| 3       | Heptadecane |              | 240 | C17H36  | 000629-78-7 | 91   |
| 4       | Hexadecane  |              | 226 | C16H34  | 000544-76-3 | 86   |
| 5       | Triacontane |              | 422 | C30H62  | 000638-68-6 | 86   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF050317\  
Data File : BF094848.D  
Acq On : 3 May 2017 22:42  
Operator : SJ/MA  
Sample : I2914-01  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-03-2.5-3.5

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF050217.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

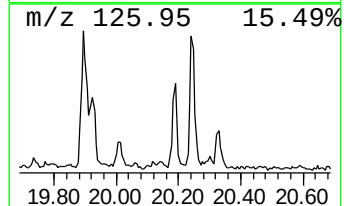
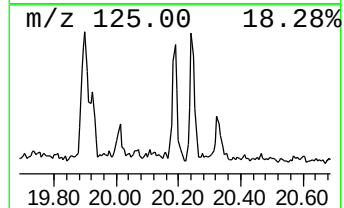
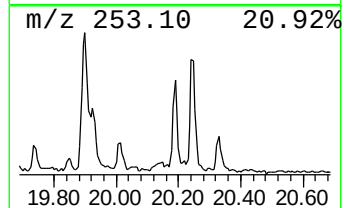
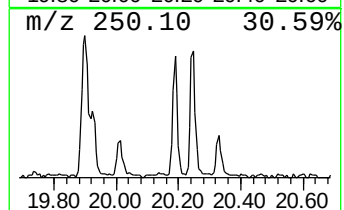
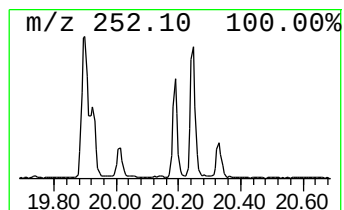
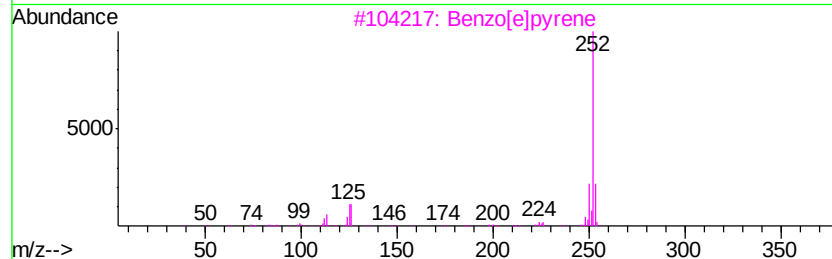
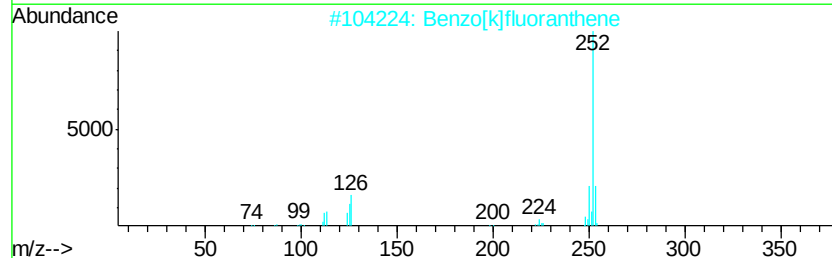
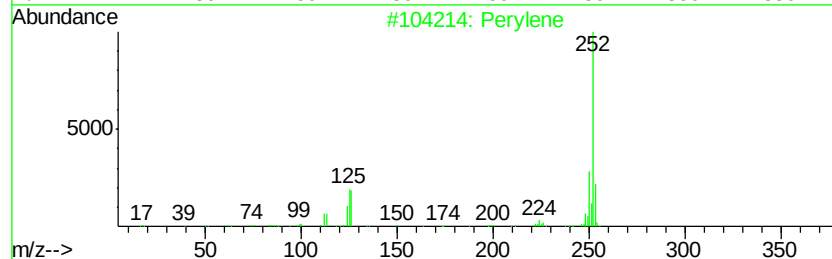
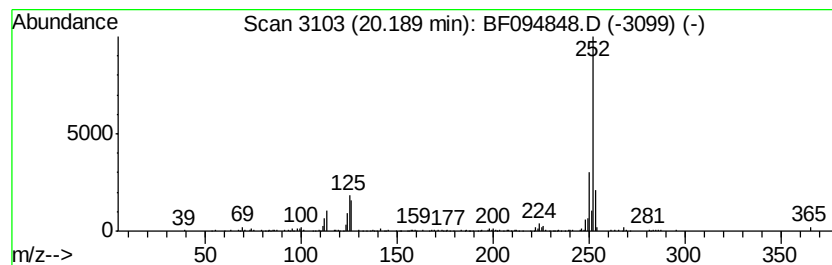
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 17 Perylene Concentration Rank 4

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 20.19 | 8.23 ng | 219657 | Perylene-d12     | 20.30 |

| Hit# | of | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------|-----|---------|-------------|------|
| 1    | 5  | Perylene                 | 252 | C20H12  | 000198-55-0 | 99   |
| 2    |    | Benzo[k]fluoranthene     | 252 | C20H12  | 000207-08-9 | 98   |
| 3    |    | Benzo[e]pyrene           | 252 | C20H12  | 000192-97-2 | 98   |
| 4    |    | Benzo[j]fluoranthene     | 252 | C20H12  | 000205-82-3 | 98   |
| 5    |    | Benz[e]acephenanthrylene | 252 | C20H12  | 000205-99-2 | 96   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF050317\  
Data File : BF094848.D  
Acq On : 3 May 2017 22:42  
Operator : SJ/MA  
Sample : I2914-01  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-03-2.5-3.5

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF050217.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

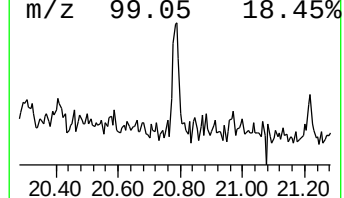
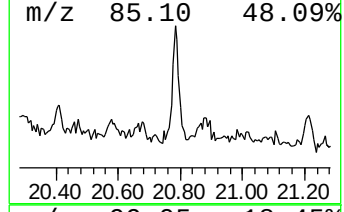
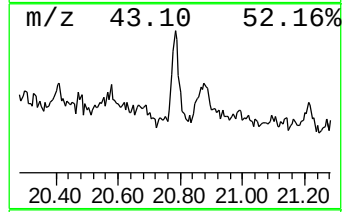
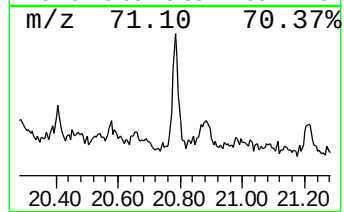
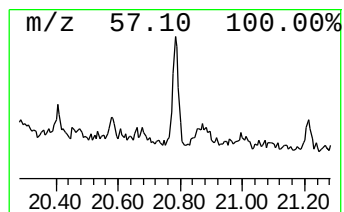
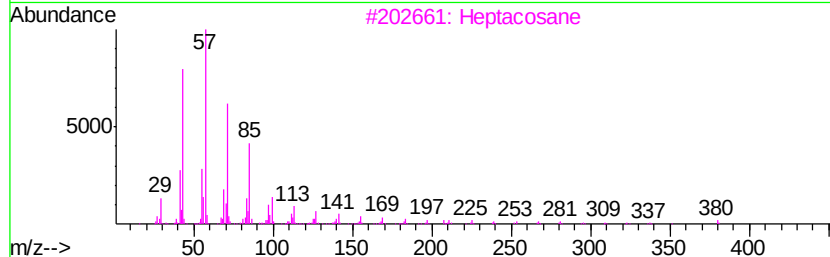
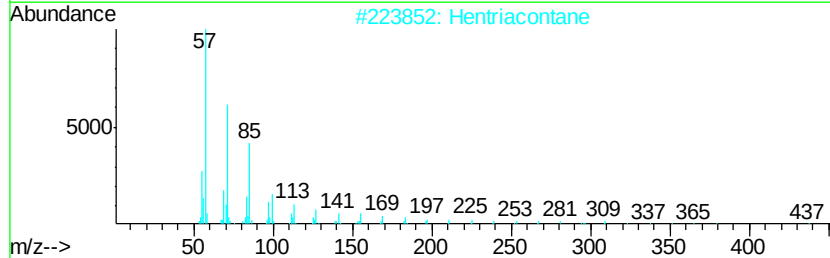
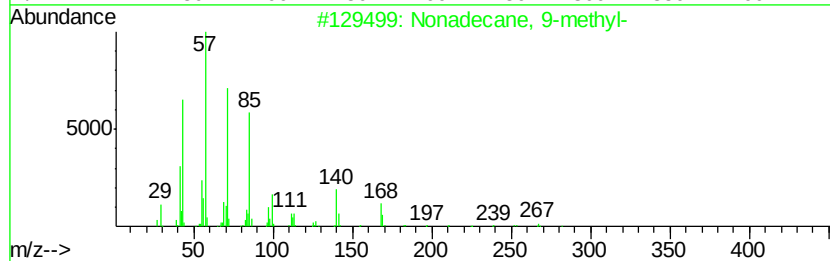
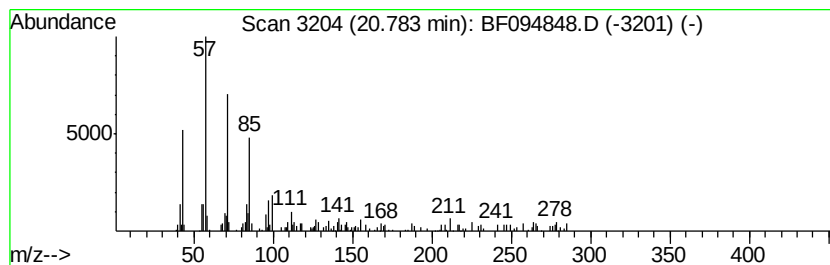
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 Nonadecane, 9-methyl- Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 20.78 | 4.32 ng | 115468 | Perylene-d12     | 20.30 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#         | Qual |
|------|----|----------------------------|-----|---------|--------------|------|
| 1    | 5  | Nonadecane, 9-methyl-      | 282 | C20H42  | 013287-24-6  | 58   |
| 2    |    | Hentriacontane             | 437 | C31H64  | 000630-04-6  | 53   |
| 3    |    | Heptacosane                | 380 | C27H56  | 000593-49-7  | 53   |
| 4    |    | Nonadecane, 2,3-dimethyl-  | 296 | C21H44  | 075163-99-4  | 53   |
| 5    |    | 5-Ethyl-5-methylnonadecane | 310 | C22H46  | 1000360-42-7 | 52   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF050317\  
Data File : BF094848.D  
Acq On : 3 May 2017 22:42  
Operator : SJ/MA  
Sample : I2914-01  
Misc :  
ALS Vial : 11 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
SB-03-2.5-3.5

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF050217.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

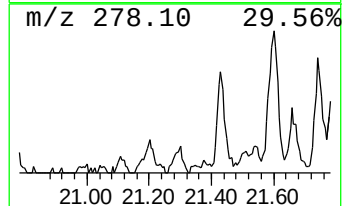
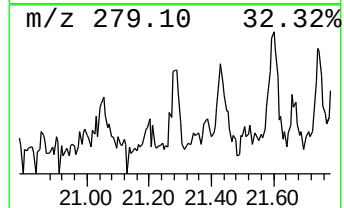
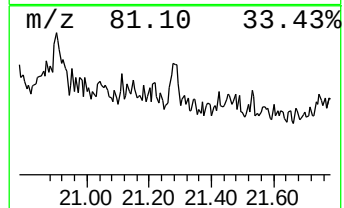
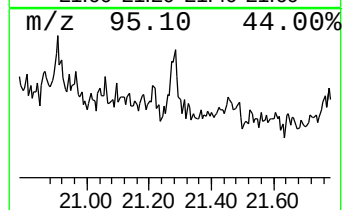
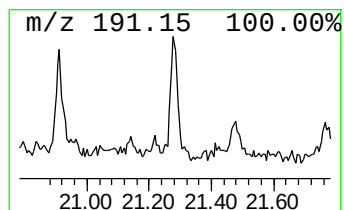
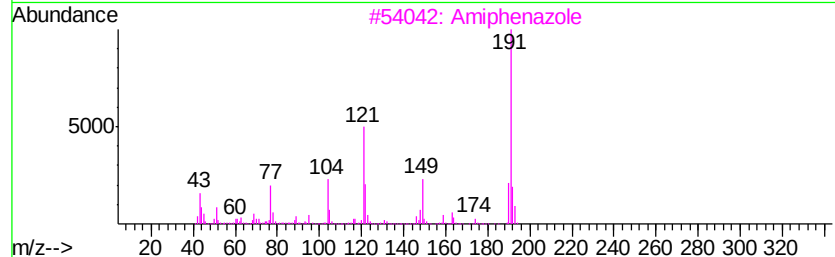
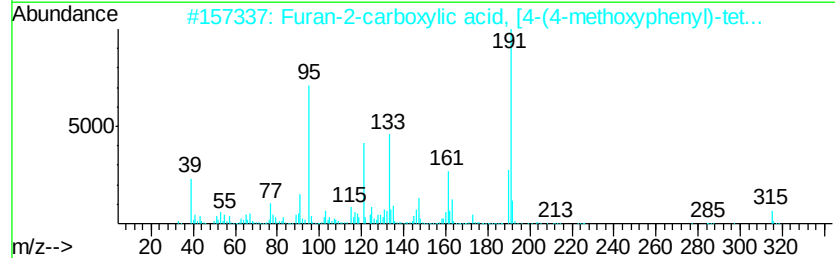
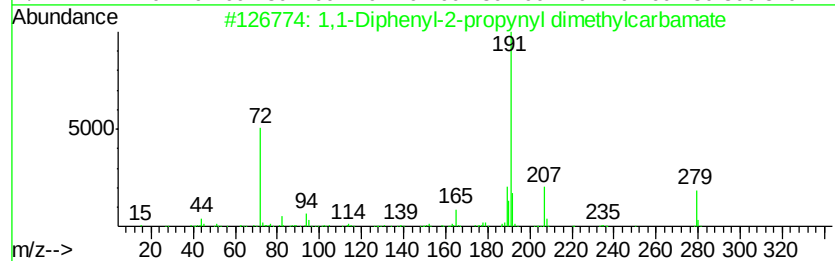
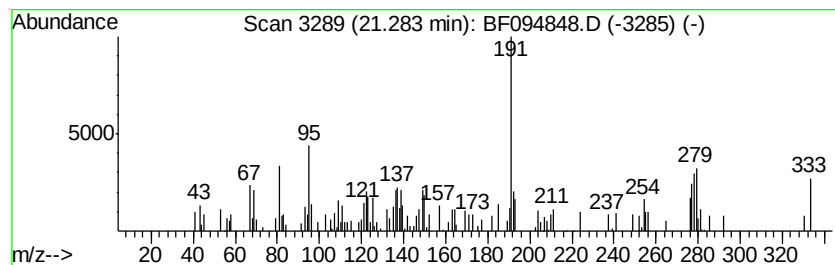
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 unknown21.28 Concentration Rank 15

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 21.28 | 2.54 ng | 67745 | Perylene-d12     | 20.30 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 1,1-Diphenyl-2-propynyl dimethyl... | 279 | C18H17N02 | 010473-88-8  | 35   |
| 2    |    | Furan-2-carboxylic acid, [4-(4-m... | 315 | C18H21N04 | 1000305-72-3 | 32   |
| 3    |    | Amiphenazole                        | 191 | C9H9N3S   | 000490-55-1  | 30   |
| 4    |    | 4H-Benzo[def]carbazole              | 191 | C14H9N    | 000203-65-6  | 27   |
| 5    |    | 9-Anthracenepropanoic acid, 10-e... | 280 | C19H20O2  | 110551-78-5  | 22   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF050317\  
 Data File : BF094848.D  
 Acq On : 3 May 2017 22:42  
 Operator : SJ/MA  
 Sample : I2914-01  
 Misc :  
 ALS Vial : 11 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SB-03-2.5-3.5

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF050217.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
 TIC Integration Parameters: LSCINT.P

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |        |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|--------|------|
|                      |       |         |       |          | #                     | RT    | Resp   | Conc |
| 2-Pentanone, 4-hy... | 5.04  | 21.5    | ng    | 599137   | 1                     | 7.72  | 556922 | 20.0 |
| unknown7.39          | 7.39  | 111.5   | ng    | 3105750  | 1                     | 7.72  | 556922 | 20.0 |
| 2(1H)-Naphthaleno... | 15.19 | 4.6     | ng    | 185075   | 4                     | 14.97 | 801988 | 20.0 |
| 2-(5-Methyl-3-pyr... | 15.68 | 5.0     | ng    | 198490   | 4                     | 14.97 | 801988 | 20.0 |
| Phenanthrene, 2-m... | 15.91 | 2.9     | ng    | 114945   | 4                     | 14.97 | 801988 | 20.0 |
| n-Hexadecanoic acid  | 15.94 | 6.0     | ng    | 239723   | 4                     | 14.97 | 801988 | 20.0 |
| di-p-Tolylacetylene  | 16.55 | 3.3     | ng    | 132071   | 4                     | 14.97 | 801988 | 20.0 |
| Pyrene, 2-methyl-    | 17.51 | 3.4     | ng    | 159158   | 5                     | 18.64 | 935660 | 20.0 |
| 11H-Benzo[b]fluorene | 17.60 | 2.5     | ng    | 116036   | 5                     | 18.64 | 935660 | 20.0 |
| unknown18.31         | 18.31 | 3.3     | ng    | 151931   | 5                     | 18.64 | 935660 | 20.0 |
| unknown18.53         | 18.53 | 2.2     | ng    | 101339   | 5                     | 18.64 | 935660 | 20.0 |
| Cyclopentadecane     | 18.56 | 2.6     | ng    | 120666   | 5                     | 18.64 | 935660 | 20.0 |
| Nonadecane           | 20.07 | 4.8     | ng    | 127335   | 6                     | 20.30 | 533979 | 20.0 |
| Perylene             | 20.19 | 8.2     | ng    | 219657   | 6                     | 20.30 | 533979 | 20.0 |
| Nonadecane, 9-met... | 20.78 | 4.3     | ng    | 115468   | 6                     | 20.30 | 533979 | 20.0 |
| unknown21.28         | 21.28 | 2.5     | ng    | 67745    | 6                     | 20.30 | 533979 | 20.0 |