

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF111318\  
 Data File : BF110727.D  
 Acq On : 13 Nov 2018 18:59  
 Operator : JU/SJ  
 Sample : J5925-11  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 TP-5S

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:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF110818.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         | 2.287       | 30            | 34          | 51           | rVB      | 139669         | 232884        | 8.90%           | 0.937%        |
| 2         | 5.198       | 525           | 529         | 539          | rBV      | 74423          | 107697        | 4.12%           | 0.434%        |
| 3         | 5.569       | 588           | 592         | 612          | rBV      | 2282522        | 2430271       | 92.93%          | 9.783%        |
| 4         | 6.575       | 757           | 763         | 772          | rBV      | 2349377        | 2615285       | 100.00%         | 10.528%       |
| 5         | 6.716       | 782           | 787         | 790          | rBV      | 2792759        | 2552553       | 97.60%          | 10.275%       |
| 6         | 6.945       | 822           | 826         | 832          | rBV      | 629641         | 580346        | 22.19%          | 2.336%        |
| 7         | 7.098       | 848           | 852         | 866          | rBV      | 2070453        | 1941757       | 74.25%          | 7.817%        |
| 8         | 7.510       | 917           | 922         | 925          | rBV      | 1554528        | 1580110       | 60.42%          | 6.361%        |
| 9         | 8.228       | 1040          | 1044        | 1058         | rBV      | 700887         | 744546        | 28.47%          | 2.997%        |
| 10        | 9.298       | 1221          | 1226        | 1229         | rBV      | 2909041        | 2418582       | 92.48%          | 9.736%        |
| 11        | 9.692       | 1289          | 1293        | 1300         | rVB      | 298564         | 282880        | 10.82%          | 1.139%        |
| 12        | 9.986       | 1339          | 1343        | 1347         | rBV      | 750326         | 723709        | 27.67%          | 2.913%        |
| 13        | 10.780      | 1473          | 1478        | 1490         | rBV      | 1306023        | 1354896       | 51.81%          | 5.454%        |
| 14        | 11.480      | 1593          | 1597        | 1599         | rBV      | 577972         | 578631        | 22.12%          | 2.329%        |
| 15        | 11.504      | 1599          | 1601        | 1605         | rVB      | 108120         | 93603         | 3.58%           | 0.377%        |
| 16        | 11.980      | 1679          | 1682        | 1685         | rBV      | 144358         | 131934        | 5.04%           | 0.531%        |
| 17        | 12.010      | 1685          | 1687        | 1692         | rVB2     | 60930          | 54061         | 2.07%           | 0.218%        |
| 18        | 12.092      | 1698          | 1701        | 1703         | rBV2     | 69196          | 73914         | 2.83%           | 0.298%        |
| 19        | 12.116      | 1703          | 1705        | 1709         | rVB      | 57750          | 46723         | 1.79%           | 0.188%        |
| 20        | 12.274      | 1728          | 1732        | 1735         | rBV      | 40118          | 37555         | 1.44%           | 0.151%        |
| 21        | 12.421      | 1752          | 1757        | 1760         | rBV3     | 25985          | 30858         | 1.18%           | 0.124%        |
| 22        | 12.527      | 1772          | 1775        | 1778         | rBV      | 72773          | 76304         | 2.92%           | 0.307%        |
| 23        | 12.627      | 1787          | 1792        | 1798         | rVB4     | 30416          | 60155         | 2.30%           | 0.242%        |
| 24        | 12.698      | 1800          | 1804        | 1807         | rBV      | 187561         | 187203        | 7.16%           | 0.754%        |
| 25        | 12.763      | 1812          | 1815        | 1819         | rBV4     | 37904          | 43158         | 1.65%           | 0.174%        |
| 26        | 12.910      | 1837          | 1840        | 1841         | rBV      | 61367          | 56153         | 2.15%           | 0.226%        |
| 27        | 12.927      | 1841          | 1843        | 1848         | rVV      | 280893         | 266771        | 10.20%          | 1.074%        |
| 28        | 12.974      | 1848          | 1851        | 1855         | rVB2     | 35771          | 39727         | 1.52%           | 0.160%        |
| 29        | 13.057      | 1861          | 1865        | 1869         | rBV      | 1570395        | 1495381       | 57.18%          | 6.020%        |
| 30        | 13.145      | 1876          | 1880        | 1885         | rBV2     | 46117          | 80524         | 3.08%           | 0.324%        |
| 31        | 13.257      | 1896          | 1899        | 1904         | rVB      | 68602          | 83616         | 3.20%           | 0.337%        |
| 32        | 13.321      | 1904          | 1910        | 1912         | rBV2     | 46704          | 54106         | 2.07%           | 0.218%        |
| 33        | 13.363      | 1912          | 1917        | 1919         | rBV3     | 60901          | 74001         | 2.83%           | 0.298%        |
| 34        | 13.451      | 1930          | 1932        | 1935         | rBV      | 56032          | 53084         | 2.03%           | 0.214%        |

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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:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF110818.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |        |        |        |        |
|----|--------|------|------|------|------|--------|--------|--------|--------|
| 35 | 13.486 | 1935 | 1938 | 1940 | rVV  | 62852  | 56847  | 2.17%  | 0.229% |
| 36 | 13.527 | 1940 | 1945 | 1948 | rVV  | 174992 | 162010 | 6.19%  | 0.652% |
| 37 | 13.739 | 1978 | 1981 | 1983 | rBV4 | 23923  | 28103  | 1.07%  | 0.113% |
| 38 | 13.768 | 1983 | 1986 | 1989 | rVB2 | 42747  | 48178  | 1.84%  | 0.194% |
| 39 | 13.880 | 2003 | 2005 | 2007 | rBV  | 33989  | 29718  | 1.14%  | 0.120% |
| 40 | 13.904 | 2007 | 2009 | 2011 | rVV  | 48260  | 36484  | 1.40%  | 0.147% |
| 41 | 13.933 | 2011 | 2014 | 2020 | rVV4 | 108983 | 133691 | 5.11%  | 0.538% |
| 42 | 13.980 | 2020 | 2022 | 2028 | rVB  | 40861  | 47165  | 1.80%  | 0.190% |
| 43 | 14.127 | 2042 | 2047 | 2050 | rBV2 | 608868 | 756304 | 28.92% | 3.045% |
| 44 | 14.157 | 2050 | 2052 | 2058 | rVB  | 214923 | 193123 | 7.38%  | 0.777% |
| 45 | 14.239 | 2063 | 2066 | 2067 | rBV  | 31788  | 31806  | 1.22%  | 0.128% |
| 46 | 14.286 | 2072 | 2074 | 2082 | rVB7 | 28538  | 57719  | 2.21%  | 0.232% |
| 47 | 14.480 | 2104 | 2107 | 2109 | rBV  | 30982  | 35316  | 1.35%  | 0.142% |
| 48 | 14.515 | 2110 | 2113 | 2116 | rVV  | 96410  | 90979  | 3.48%  | 0.366% |
| 49 | 14.551 | 2116 | 2119 | 2123 | rVV2 | 77767  | 95385  | 3.65%  | 0.384% |
| 50 | 14.615 | 2128 | 2130 | 2133 | rVB2 | 34632  | 29733  | 1.14%  | 0.120% |
| 51 | 14.645 | 2133 | 2135 | 2137 | rBV  | 56779  | 53936  | 2.06%  | 0.217% |
| 52 | 14.874 | 2171 | 2174 | 2179 | rVB  | 105155 | 112571 | 4.30%  | 0.453% |
| 53 | 15.227 | 2226 | 2234 | 2238 | rBV2 | 166954 | 333388 | 12.75% | 1.342% |
| 54 | 15.315 | 2247 | 2249 | 2254 | rVB  | 36981  | 46054  | 1.76%  | 0.185% |
| 55 | 15.527 | 2282 | 2285 | 2289 | rVB  | 63692  | 71421  | 2.73%  | 0.288% |
| 56 | 15.592 | 2292 | 2296 | 2298 | rVV  | 105403 | 143635 | 5.49%  | 0.578% |
| 57 | 15.621 | 2298 | 2301 | 2303 | rVV  | 150470 | 188920 | 7.22%  | 0.761% |
| 58 | 15.656 | 2303 | 2307 | 2311 | rVB  | 313516 | 391635 | 14.97% | 1.577% |
| 59 | 16.074 | 2373 | 2378 | 2383 | rVB  | 124306 | 184418 | 7.05%  | 0.742% |
| 60 | 16.356 | 2422 | 2426 | 2430 | rBV4 | 38385  | 65050  | 2.49%  | 0.262% |
| 61 | 16.598 | 2463 | 2467 | 2476 | rVB  | 72808  | 134494 | 5.14%  | 0.541% |
| 62 | 17.227 | 2570 | 2574 | 2581 | rVB4 | 51272  | 100462 | 3.84%  | 0.404% |

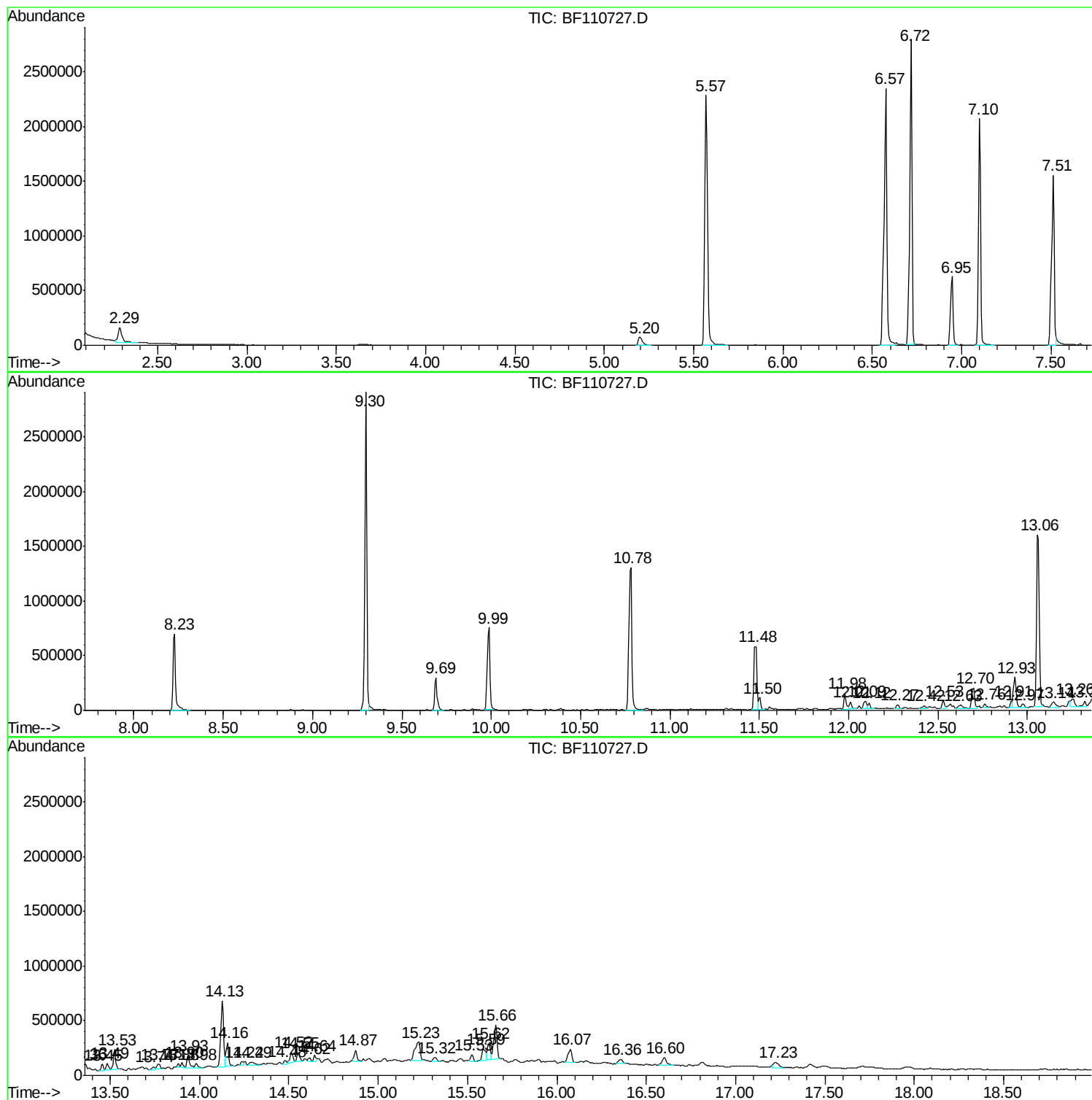
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Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampled :  
TP-5S

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.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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Data File : BF110727.D  
Acq On : 13 Nov 2018 18:59  
Operator : JU/SJ  
Sample : J5925-11  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-5S

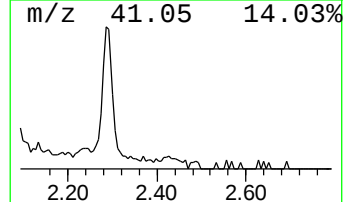
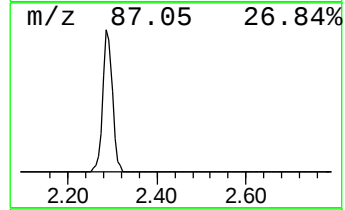
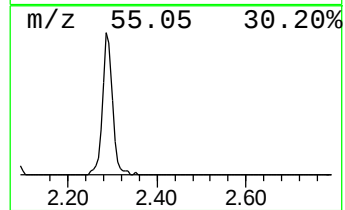
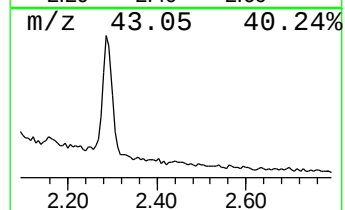
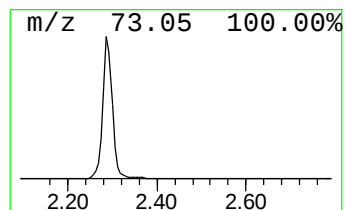
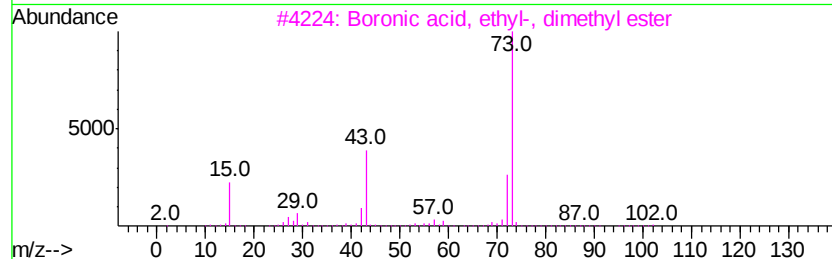
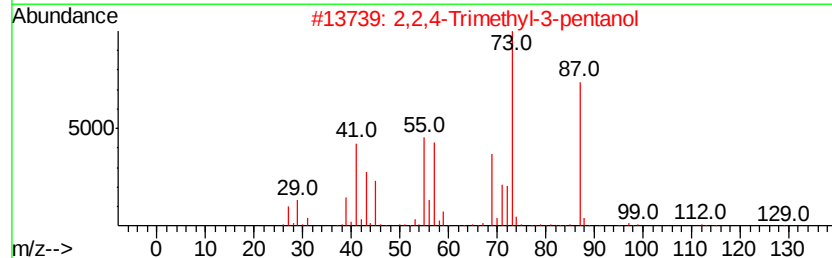
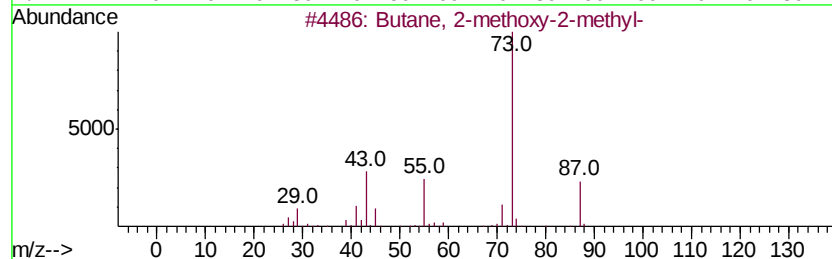
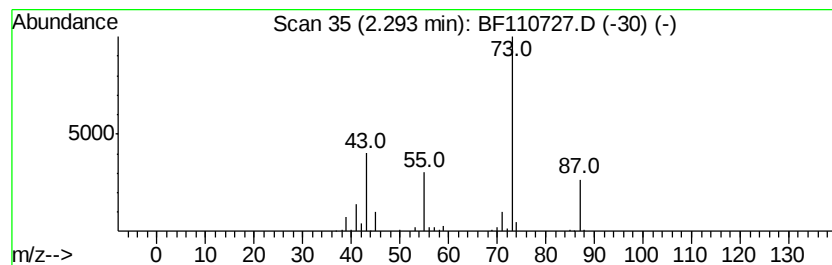
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 4

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 2.29 | 8.03 ng | 232884 | 1,4-Dichlorobenzene-d4 | 6.95 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Butane, 2-methoxy-2-methyl-         | 102 | C6H14O   | 000994-05-8 | 83   |
| 2    |    | 2,2,4-Trimethyl-3-pentanol          | 130 | C8H18O   | 005162-48-1 | 39   |
| 3    |    | Boronic acid, ethyl-, dimethyl e... | 102 | C4H11BO2 | 007318-82-3 | 17   |
| 4    |    | Pentane, 3-methoxy-                 | 102 | C6H14O   | 036839-67-5 | 12   |
| 5    |    | Silane, tetramethyl-                | 88  | C4H12Si  | 000075-76-3 | 12   |



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Sample : J5925-11  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-5S

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

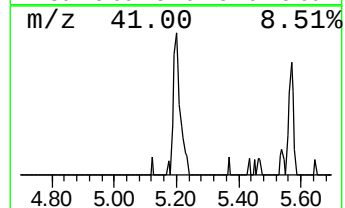
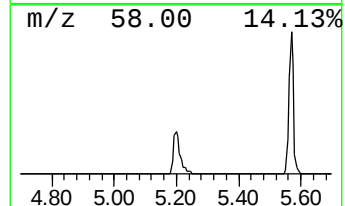
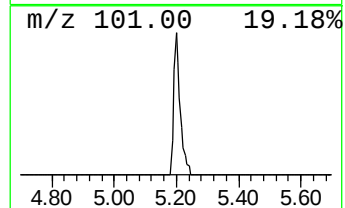
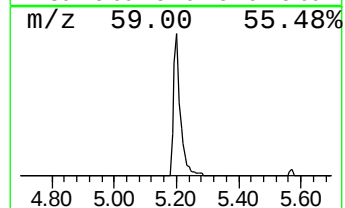
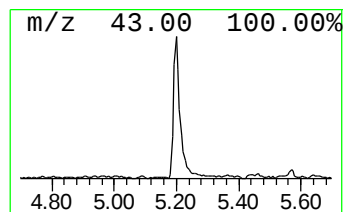
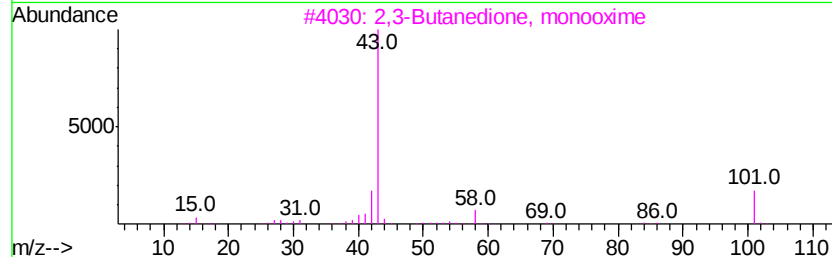
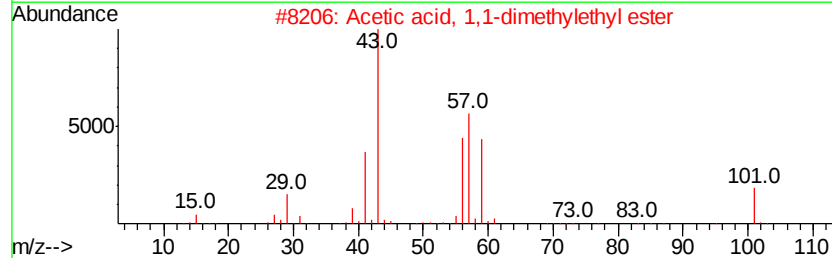
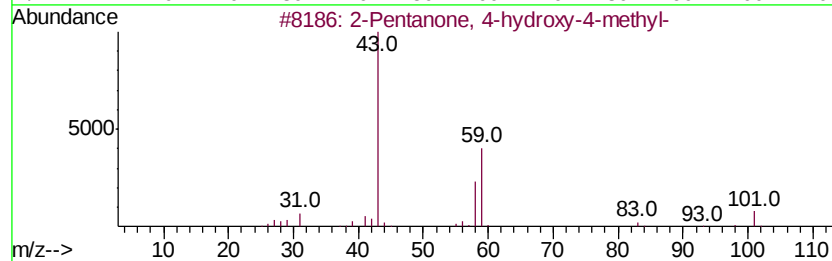
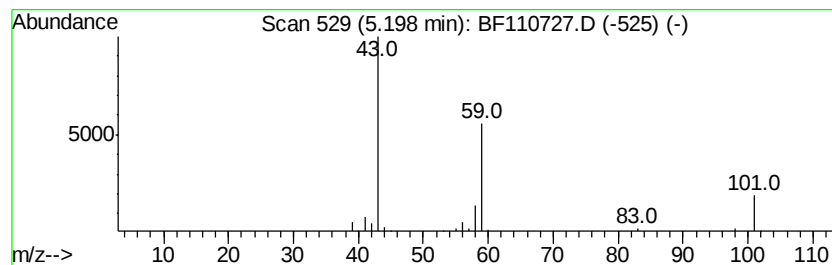
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 8

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 5.20 | 3.71 ng | 107697 | 1,4-Dichlorobenzene-d4 | 6.95 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 2-Pentanone, 4-hydroxy-4-methyl-    | 116 | C6H12O2  | 000123-42-2 | 42   |
| 2    |    | Acetic acid, 1,1-dimethylethyl e... | 116 | C6H12O2  | 000540-88-5 | 39   |
| 3    |    | 2,3-Butanedione, monooxime          | 101 | C4H7NO2  | 000057-71-6 | 25   |
| 4    |    | Morpholine, 4-methyl-               | 101 | C5H11NO  | 000109-02-4 | 9    |
| 5    |    | Butanamide, 2-hydroxy-2,N-dimethyl- | 117 | C5H11NO2 | 039961-72-3 | 9    |



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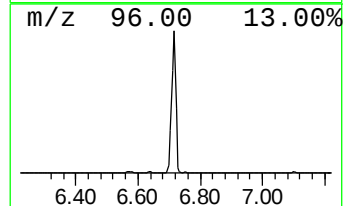
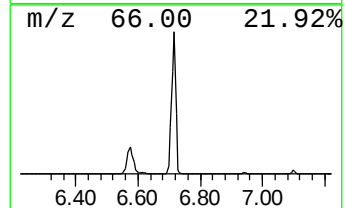
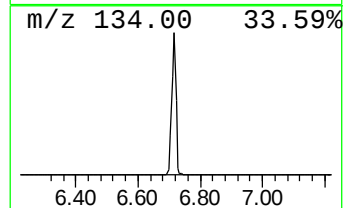
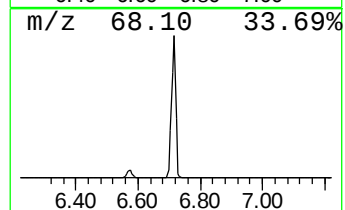
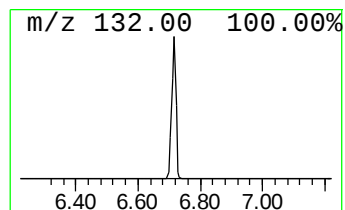
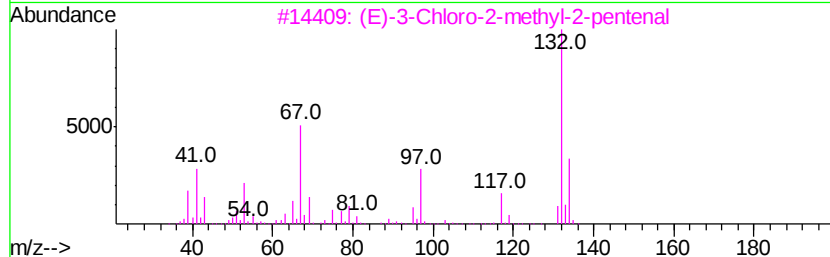
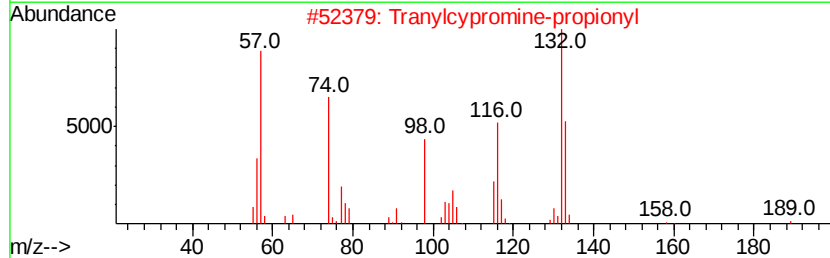
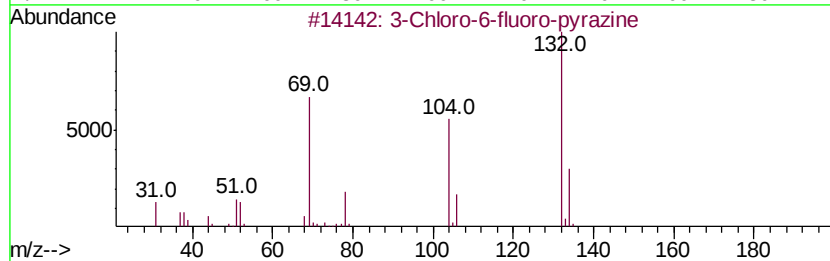
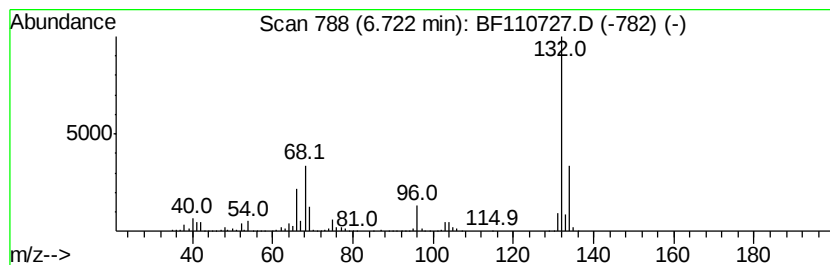
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 3 unknown6.72 Concentration Rank 1

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.72 | 87.97 ng | 2552550 | 1,4-Dichlorobenzene-d4 | 6.95 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 3-Chloro-6-fluoro-pyrazine          | 132 | C4H2ClFN2 | 1000146-10-7 | 27   |
| 2    |    | Tranylcypromine-propionyl           | 189 | C12H15NO  | 1000123-86-3 | 12   |
| 3    |    | (E)-3-Chloro-2-methyl-2-pentenal    | 132 | C6H9ClO   | 031357-76-3  | 12   |
| 4    |    | Bicyclo[4.2.0]octa-1,3,5-triene,... | 132 | C10H12    | 028749-81-7  | 9    |
| 5    |    | 4-Chloro-2-fluoro-pyrimidine        | 132 | C4H2ClFN2 | 051422-00-5  | 9    |



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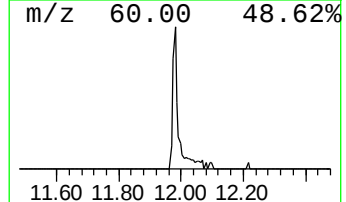
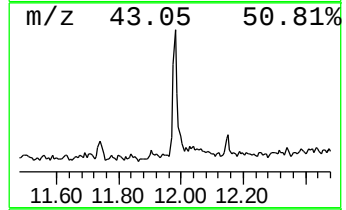
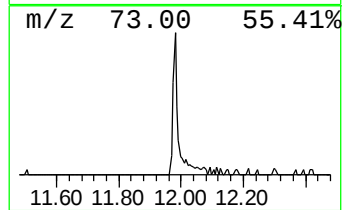
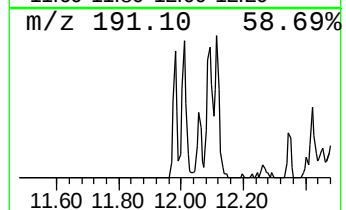
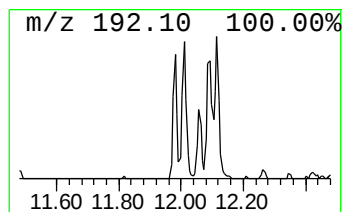
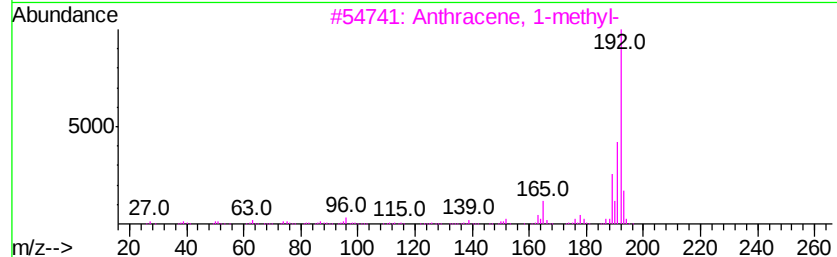
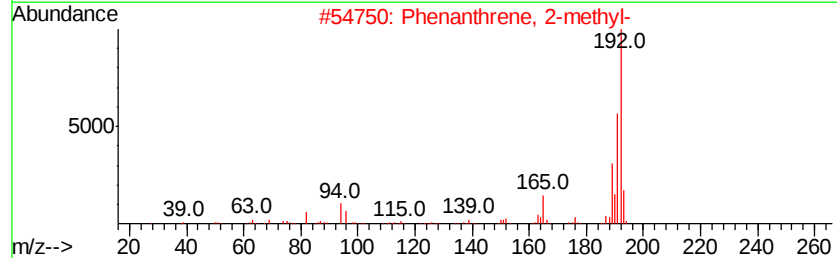
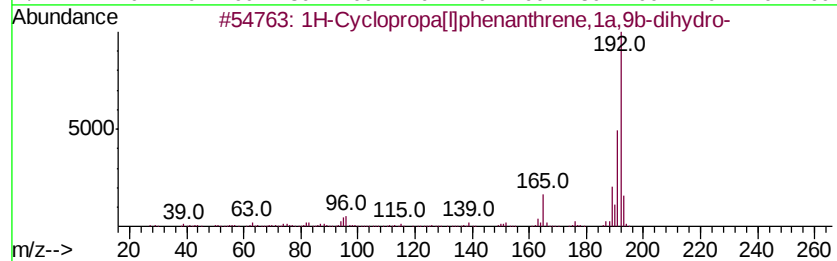
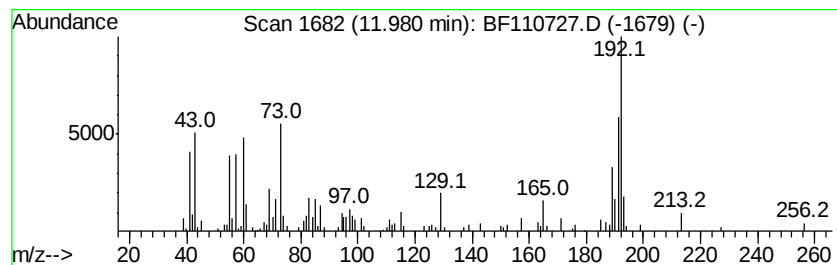
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ClientSampleId :  
TP-5S

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 5 1H-Cyclopropa[l]phenanthren... Concentration Rank 6

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|-------------|------|
| 11.98     | 4.56 ng                             | 131934 | Phenanthrene-d10 | 11.48       |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual |
| 1         | 1H-Cyclopropa[1]phenanthrene,1a,... | 192    | C15H12           | 000949-41-7 | 95   |
| 2         | Phenanthrene, 2-methyl-             | 192    | C15H12           | 002531-84-2 | 95   |
| 3         | Anthracene, 1-methyl-               | 192    | C15H12           | 000610-48-0 | 95   |
| 4         | Phenanthrene, 1-methyl-             | 192    | C15H12           | 000832-69-9 | 95   |
| 5         | Anthracene, 2-methyl-               | 192    | C15H12           | 000613-12-7 | 95   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111318\  
Data File : BF110727.D  
Acq On : 13 Nov 2018 18:59  
Operator : JU/SJ  
Sample : J5925-11  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-5S

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

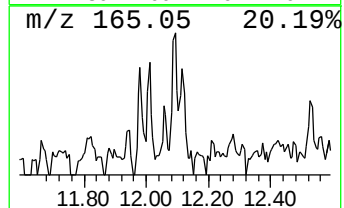
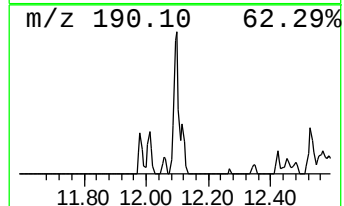
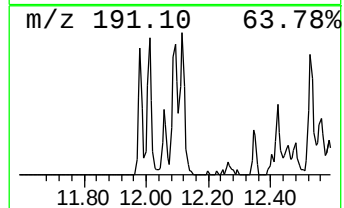
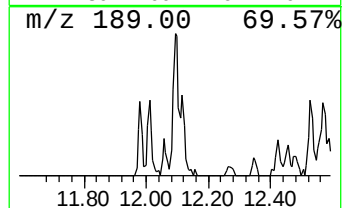
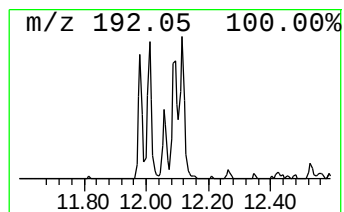
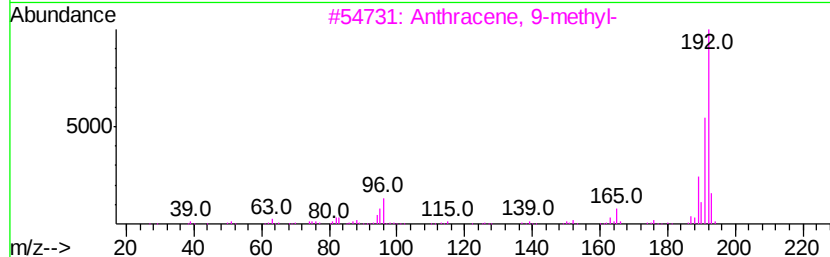
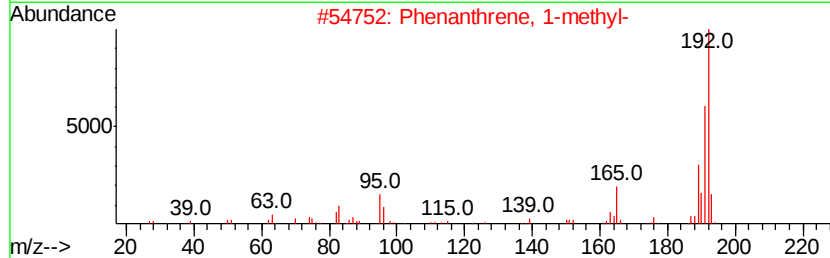
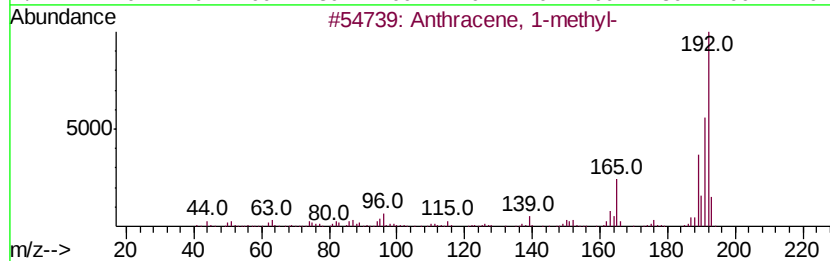
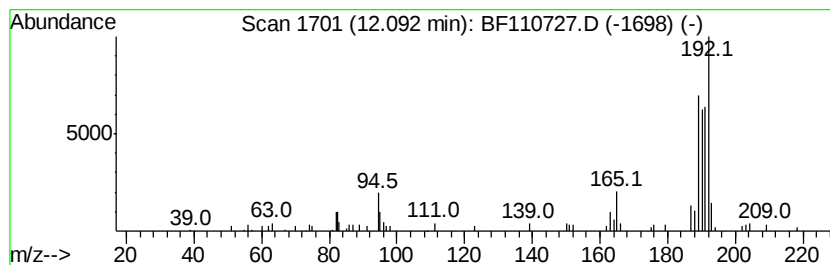
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 6 Anthracene, 1-methyl- Concentration Rank 14

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 12.09 | 2.55 ng | 73914 | Phenanthrene-d10 | 11.48 |

| Hit# | of | Tentative ID                         | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Anthracene, 1-methyl-                | 192 | C15H12  | 000610-48-0 | 90   |
| 2    |    | Phenanthrene, 1-methyl-              | 192 | C15H12  | 000832-69-9 | 78   |
| 3    |    | Anthracene, 9-methyl-                | 192 | C15H12  | 000779-02-2 | 64   |
| 4    |    | Naphtho[2,3-b]norbornadiene          | 192 | C15H12  | 107426-38-0 | 60   |
| 5    |    | 1H-Cyclopropa[1]phenanthrene, 1a,... | 192 | C15H12  | 000949-41-7 | 60   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111318\  
Data File : BF110727.D  
Acq On : 13 Nov 2018 18:59  
Operator : JU/SJ  
Sample : J5925-11  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-5S

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

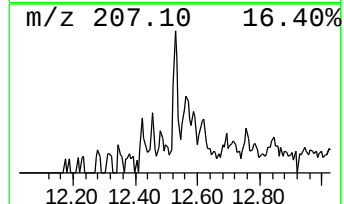
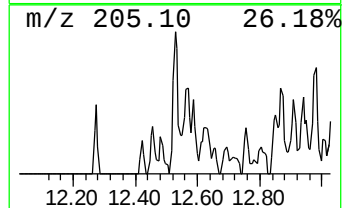
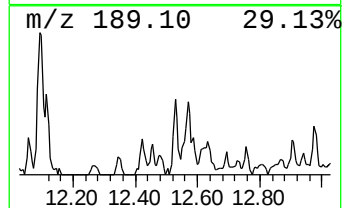
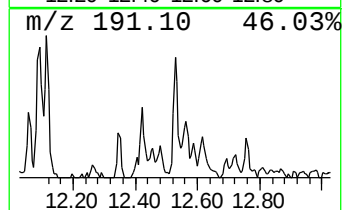
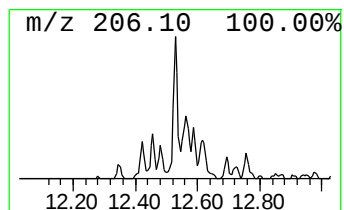
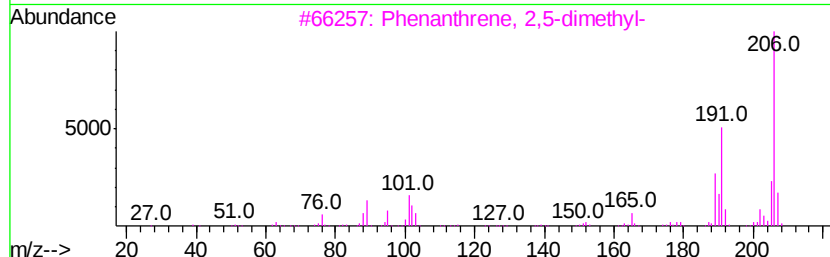
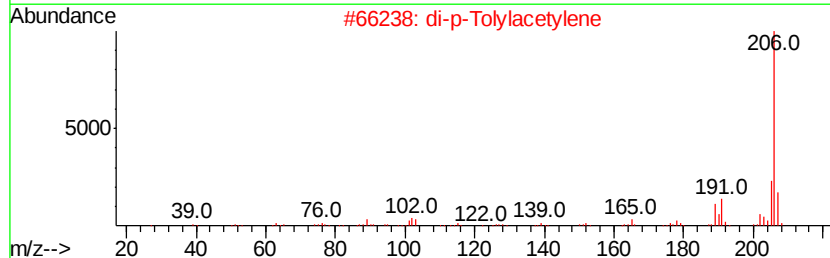
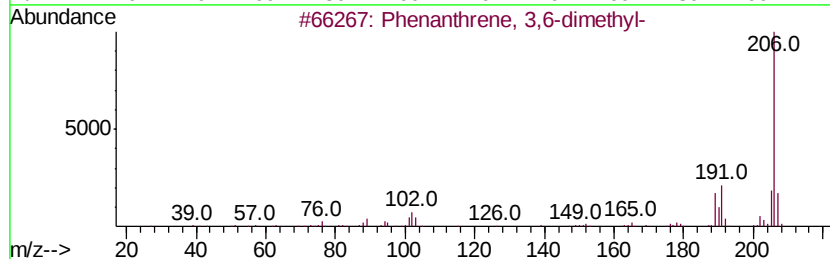
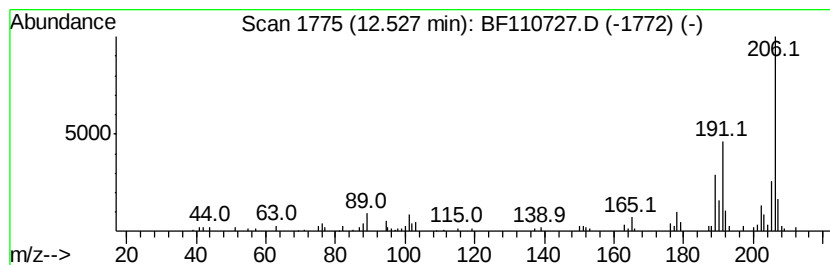
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 Phenanthrene, 3,6-dimethyl- Concentration Rank 13

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 12.53 | 2.64 ng | 76304 | Phenanthrene-d10 | 11.48 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | Phenanthrene, 3,6-dimethyl- | 206 | C16H14  | 001576-67-6 | 94   |
| 2    |    | di-p-Tolylacetylene         | 206 | C16H14  | 002789-88-0 | 94   |
| 3    |    | Phenanthrene, 2,5-dimethyl- | 206 | C16H14  | 003674-66-6 | 93   |
| 4    |    | Phenanthrene, 2,7-dimethyl- | 206 | C16H14  | 001576-69-8 | 93   |
| 5    |    | 9,10-Dimethylantracene      | 206 | C16H14  | 000781-43-1 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111318\  
Data File : BF110727.D  
Acq On : 13 Nov 2018 18:59  
Operator : JU/SJ  
Sample : J5925-11  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-5S

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

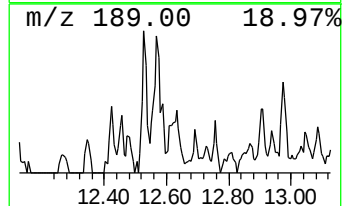
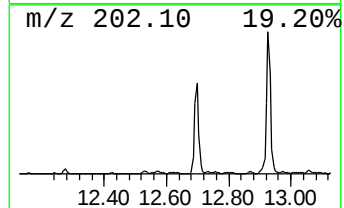
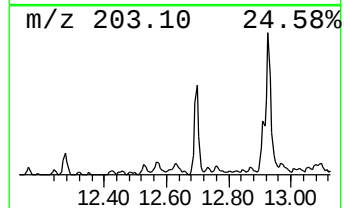
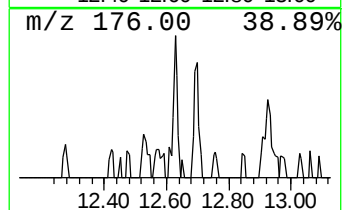
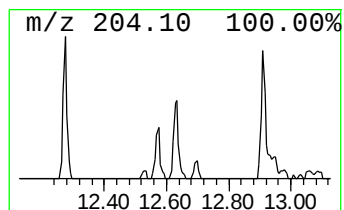
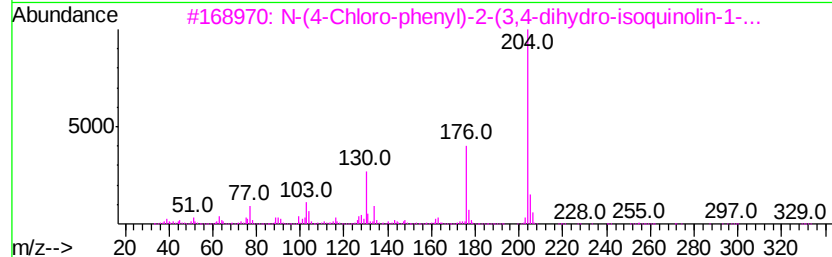
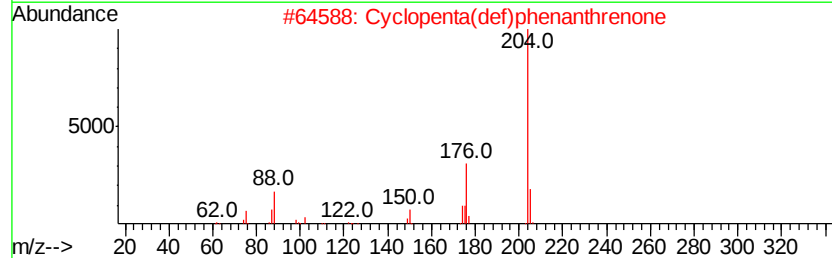
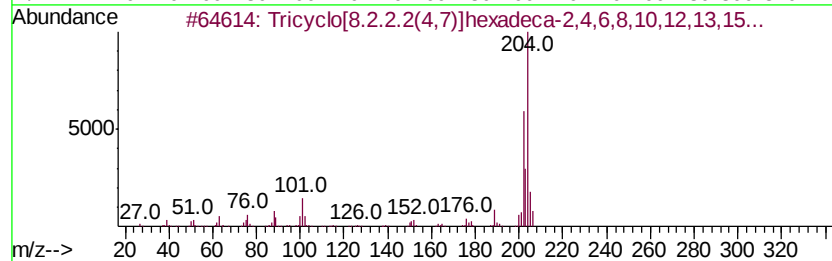
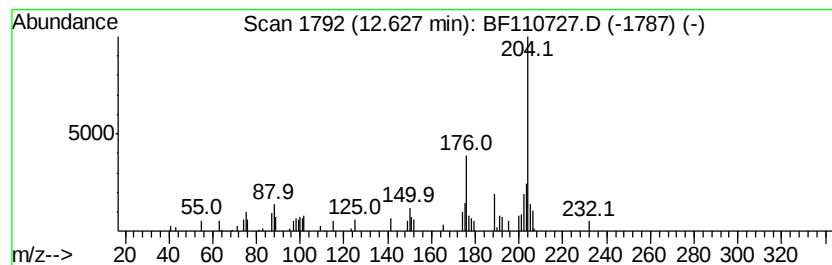
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 Tricyclo[8.2.2.2(4,7)]hexad... Concentration Rank 20

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 12.63 | 2.08 ng | 60155 | Phenanthrene-d10 | 11.48 |

| Hit# | of | Tentative ID                        | MW  | MolForm      | CAS#         | Qual |
|------|----|-------------------------------------|-----|--------------|--------------|------|
| 1    | 5  | Tricyclo[8.2.2.2(4,7)]hexadeca-2... | 204 | C16H12       | 006572-60-7  | 78   |
| 2    |    | Cyclopenta(def)phenanthrenone       | 204 | C15H8O       | 005737-13-3  | 64   |
| 3    |    | N-(4-Chloro-phenyl)-2-(3,4-dihyd... | 330 | C17H15ClN2OS | 1000296-34-3 | 53   |
| 4    |    | 3,4-Biphenyldicarbonitrile          | 204 | C14H8N2      | 004128-63-6  | 53   |
| 5    |    | [1,1'-Biphenyl]-2,2'-dicarbonitrile | 204 | C14H8N2      | 004341-02-0  | 49   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111318\  
Data File : BF110727.D  
Acq On : 13 Nov 2018 18:59  
Operator : JU/SJ  
Sample : J5925-11  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-5S

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

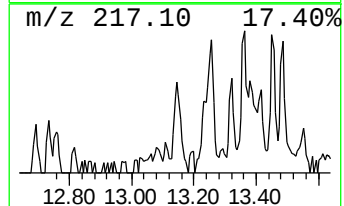
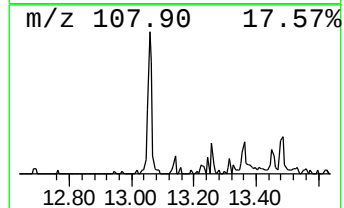
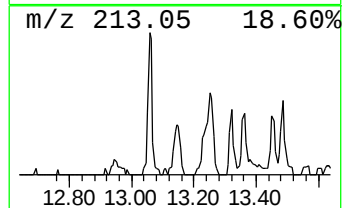
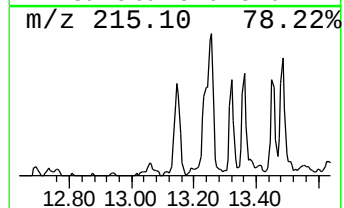
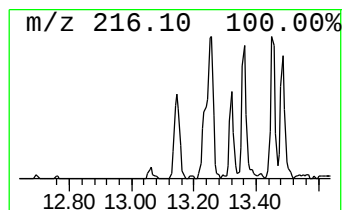
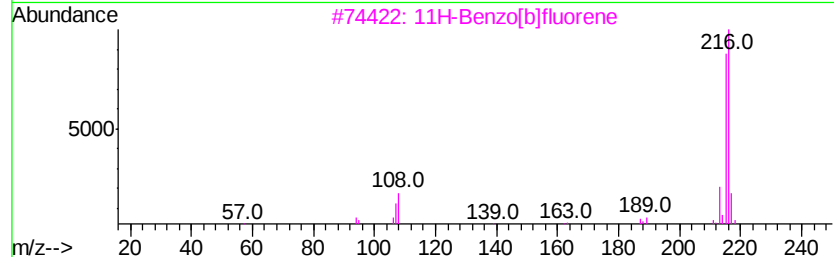
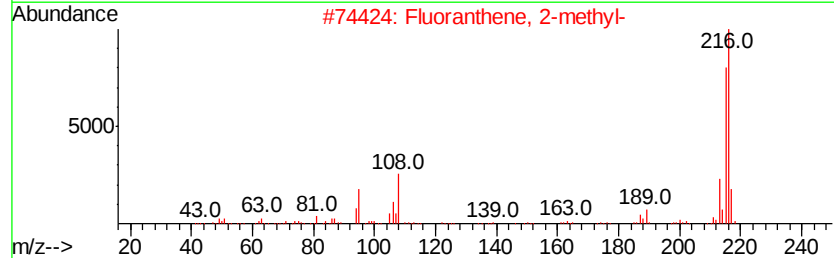
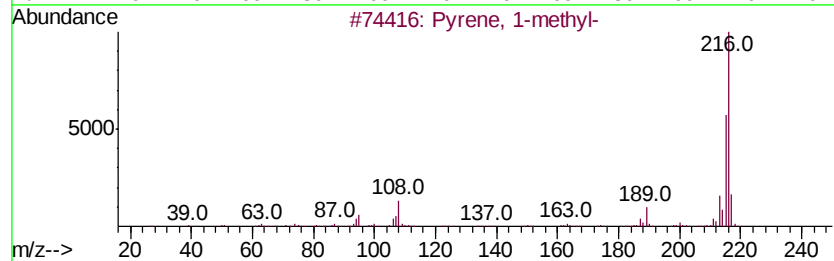
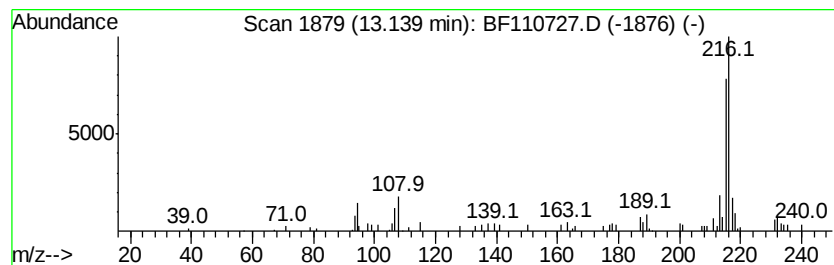
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 Pyrene, 1-methyl- Concentration Rank 19

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 13.14 | 2.13 ng | 80524 | Chrysene-d12     | 14.13 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | Pyrene, 1-methyl-       | 216 | C17H12  | 002381-21-7 | 93   |
| 2    |    | Fluoranthene, 2-methyl- | 216 | C17H12  | 033543-31-6 | 93   |
| 3    |    | 11H-Benzo[b]fluorene    | 216 | C17H12  | 000243-17-4 | 91   |
| 4    |    | 7H-Benzo[c]fluorene     | 216 | C17H12  | 000205-12-9 | 90   |
| 5    |    | Pyrene, 2-methyl-       | 216 | C17H12  | 003442-78-2 | 90   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111318\  
Data File : BF110727.D  
Acq On : 13 Nov 2018 18:59  
Operator : JU/SJ  
Sample : J5925-11  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-5S

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

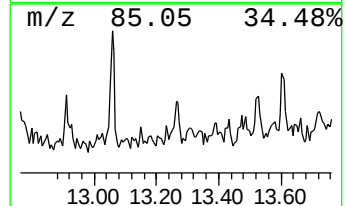
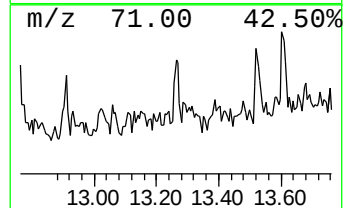
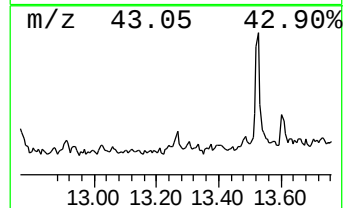
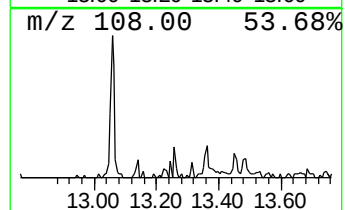
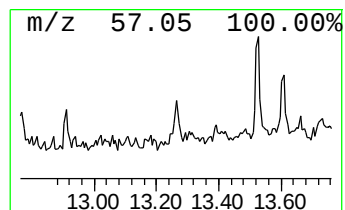
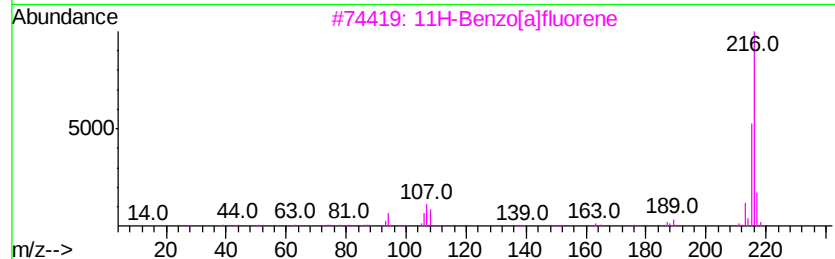
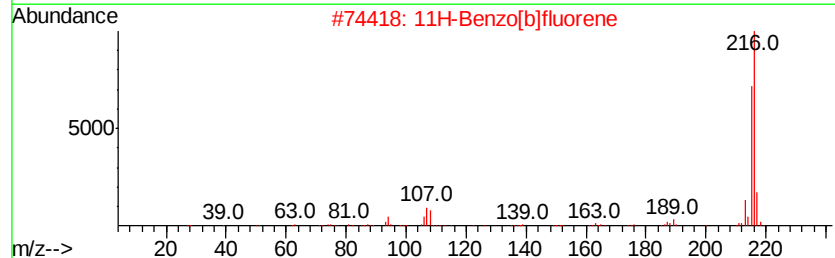
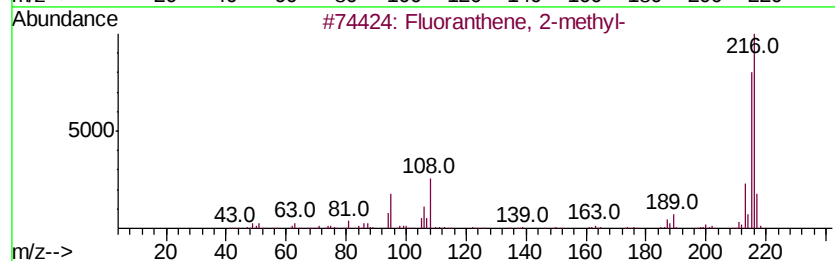
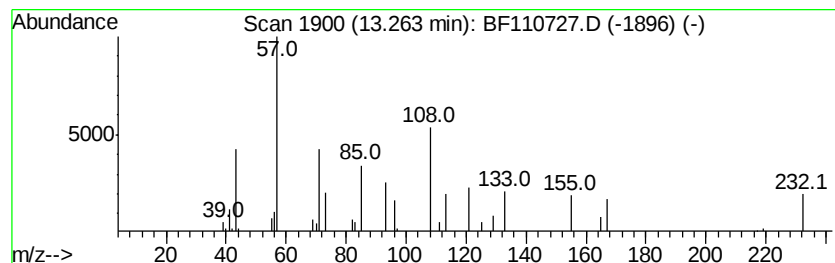
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 Fluoranthene, 2-methyl- Concentration Rank 18

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 13.26 | 2.21 ng | 83616 | Chrysene-d12     | 14.13 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------|-----|------------|--------------|------|
| 1    | 5  | Fluoranthene, 2-methyl-             | 216 | C17H12     | 033543-31-6  | 86   |
| 2    |    | 11H-Benzo[b]fluorene                | 216 | C17H12     | 000243-17-4  | 83   |
| 3    |    | 11H-Benzo[a]fluorene                | 216 | C17H12     | 000238-84-6  | 83   |
| 4    |    | Pvrene, 1-methyl-                   | 216 | C17H12     | 002381-21-7  | 72   |
| 5    |    | 4-Hydroxy-6-methyl-1-pyridin-3-y... | 216 | C12H12N2O2 | 1000318-25-3 | 72   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111318\  
Data File : BF110727.D  
Acq On : 13 Nov 2018 18:59  
Operator : JU/SJ  
Sample : J5925-11  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-5S

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

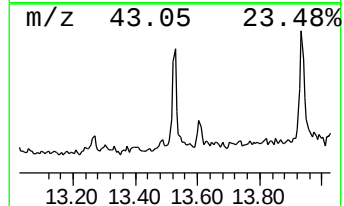
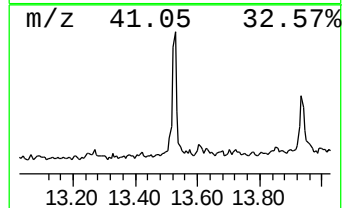
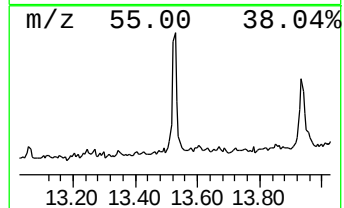
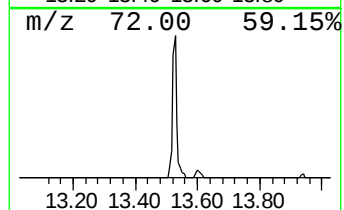
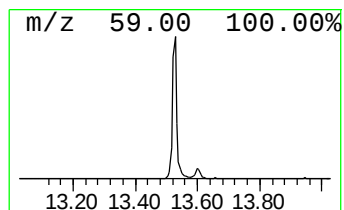
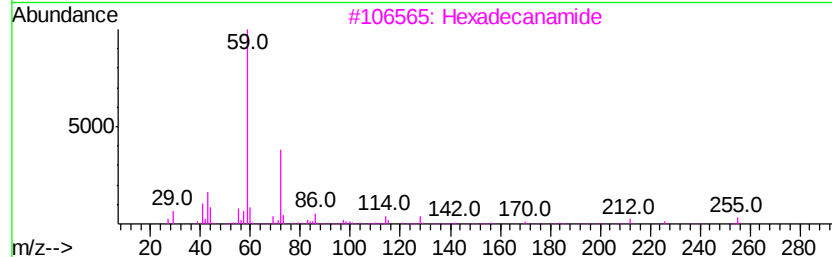
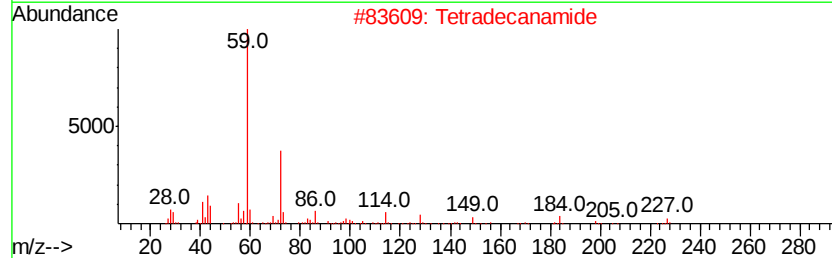
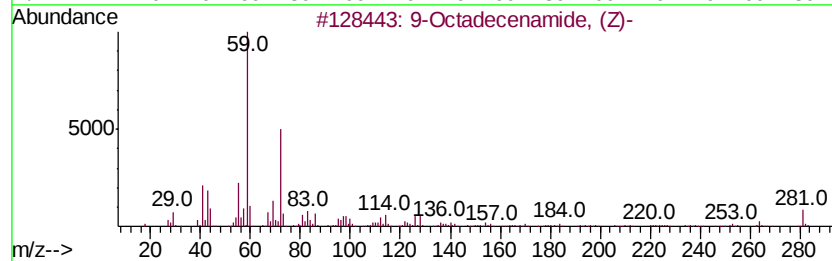
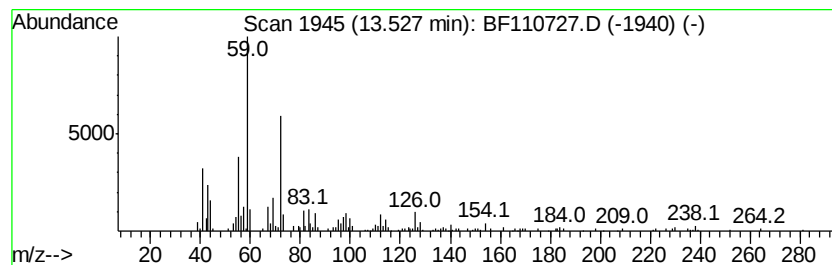
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TIC Integration Parameters: LSCINT.P

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Peak Number 11 9-Octadecenamide, (Z)- Concentration Rank 7

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.53 | 4.28 ng | 162010 | Chrysene-d12     | 14.13 |

| Hit# | of | 5 | Tentative ID           | MW  | MolForm  | CAS#        | Qual |
|------|----|---|------------------------|-----|----------|-------------|------|
| 1    |    |   | 9-Octadecenamide, (Z)- | 281 | C18H35NO | 000301-02-0 | 91   |
| 2    |    |   | Tetradecanamide        | 227 | C14H29NO | 000638-58-4 | 53   |
| 3    |    |   | Hexadecanamide         | 255 | C16H33NO | 000629-54-9 | 53   |
| 4    |    |   | Dodecanamide           | 199 | C12H25NO | 001120-16-7 | 53   |
| 5    |    |   | Octanamide             | 143 | C8H17NO  | 000629-01-6 | 53   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111318\  
Data File : BF110727.D  
Acq On : 13 Nov 2018 18:59  
Operator : JU/SJ  
Sample : J5925-11  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-5S

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

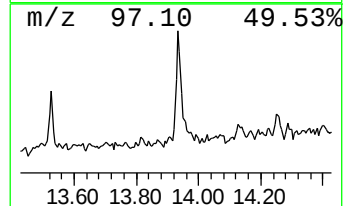
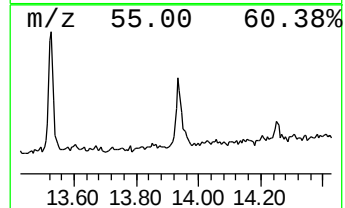
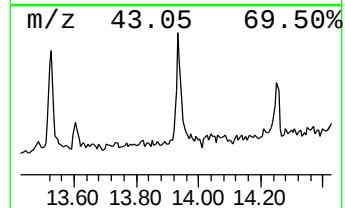
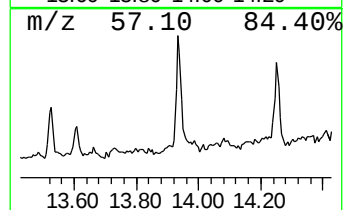
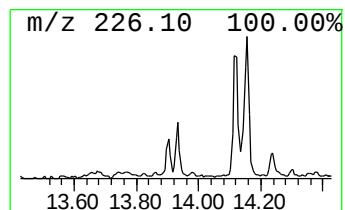
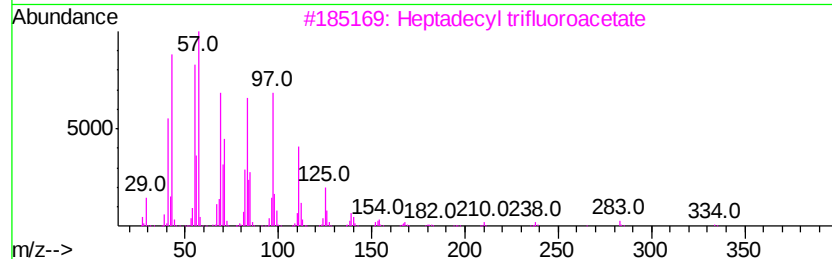
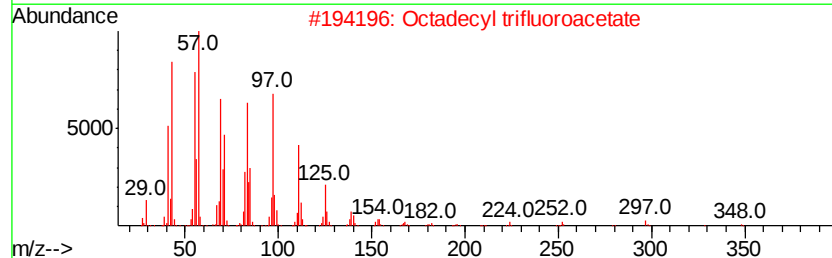
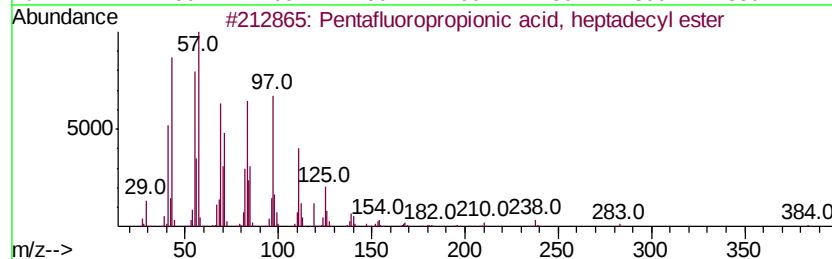
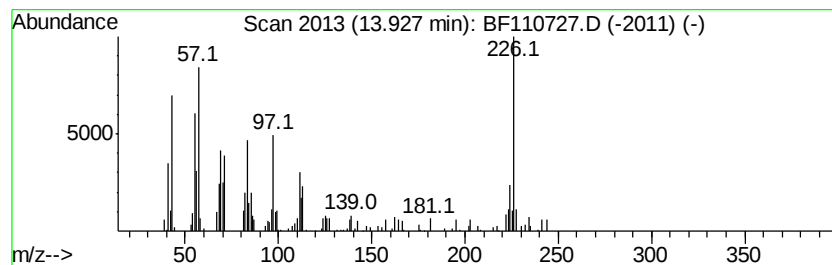
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 Pentafluoropropionic acid, ... Concentration Rank 10

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.93 | 3.54 ng | 133691 | Chrysene-d12     | 14.13 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------|-----|------------|--------------|------|
| 1    | 5  | Pentafluoropropionic acid, hepta... | 402 | C20H35F5O2 | 959218-78-1  | 90   |
| 2    |    | Octadecyl trifluoroacetate          | 366 | C20H37F3O2 | 079392-43-1  | 90   |
| 3    |    | Heptadecyl trifluoroacetate         | 352 | C19H35F3O2 | 1000351-87-0 | 90   |
| 4    |    | Cyclohexadecane                     | 224 | C16H32     | 000295-65-8  | 89   |
| 5    |    | Heptadecyl heptafluorobutyrate      | 452 | C21H35F7O2 | 959085-66-6  | 78   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111318\  
Data File : BF110727.D  
Acq On : 13 Nov 2018 18:59  
Operator : JU/SJ  
Sample : J5925-11  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-5S

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

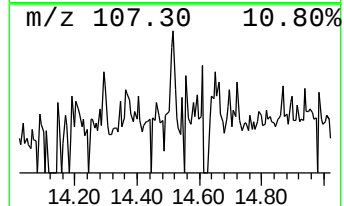
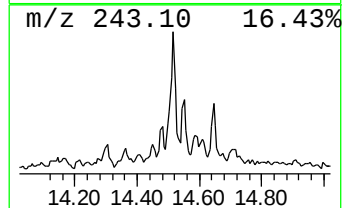
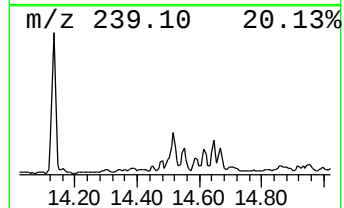
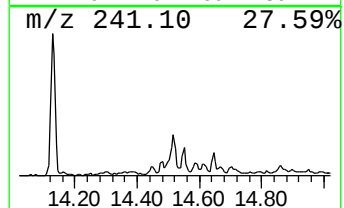
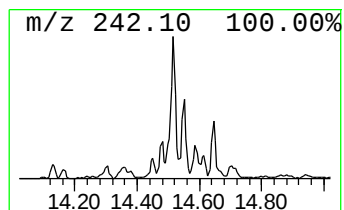
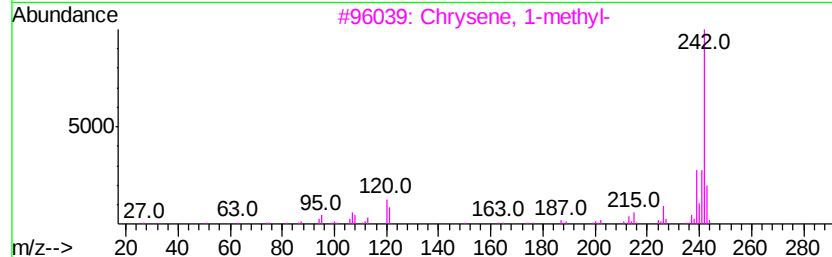
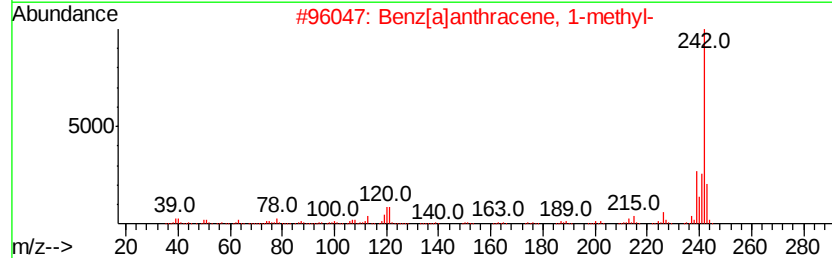
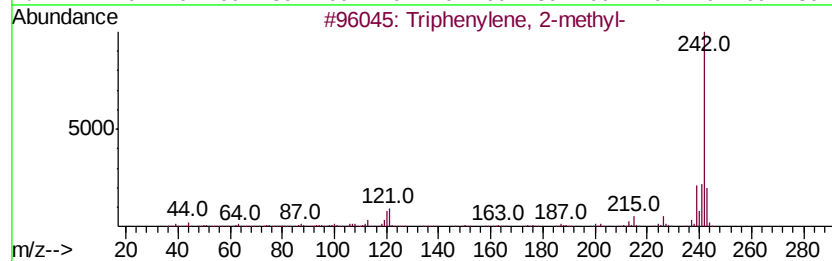
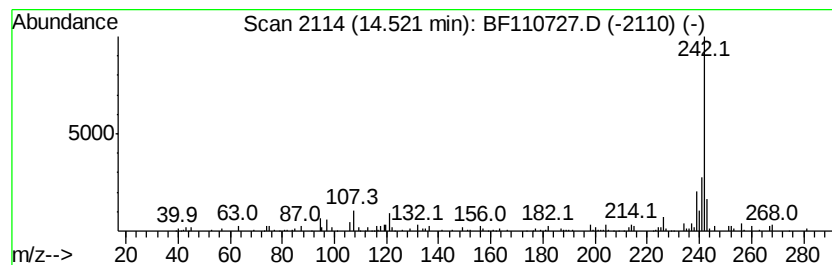
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 Triphenylene, 2-methyl- Concentration Rank 16

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 14.52 | 2.41 ng | 90979 | Chrysene-d12     | 14.13 |

| Hit# | of | Tentative ID                          | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Triphenylene, 2-methyl-               | 242 | C19H14  | 001705-84-6 | 98   |
| 2    |    | Benz[ <i>a</i> ]anthracene, 1-methyl- | 242 | C19H14  | 002498-77-3 | 96   |
| 3    |    | Chrysene, 1-methyl-                   | 242 | C19H14  | 003351-28-8 | 94   |
| 4    |    | 2-Methylchrysene                      | 242 | C19H14  | 003351-32-4 | 94   |
| 5    |    | Benz[ <i>a</i> ]anthracene, 7-methyl- | 242 | C19H14  | 002541-69-7 | 94   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111318\  
Data File : BF110727.D  
Acq On : 13 Nov 2018 18:59  
Operator : JU/SJ  
Sample : J5925-11  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

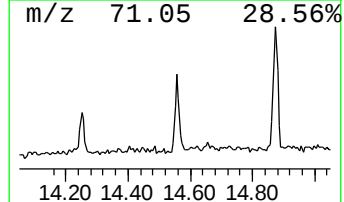
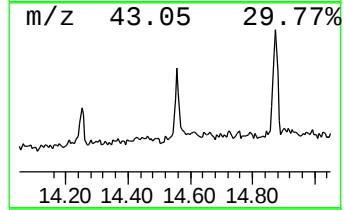
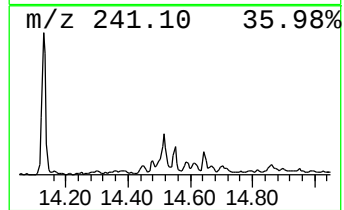
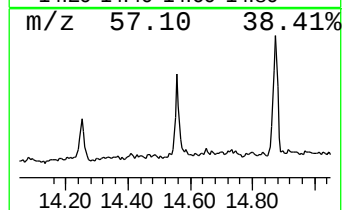
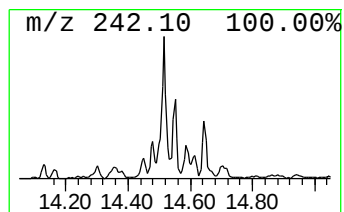
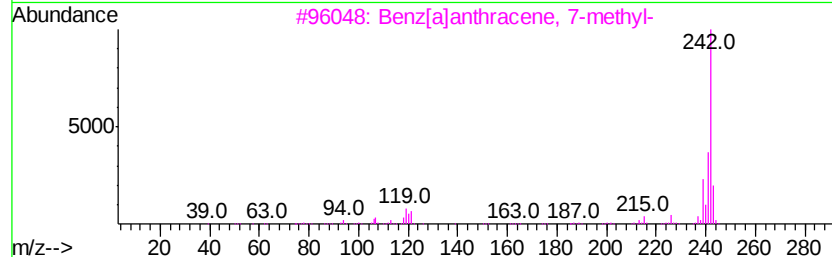
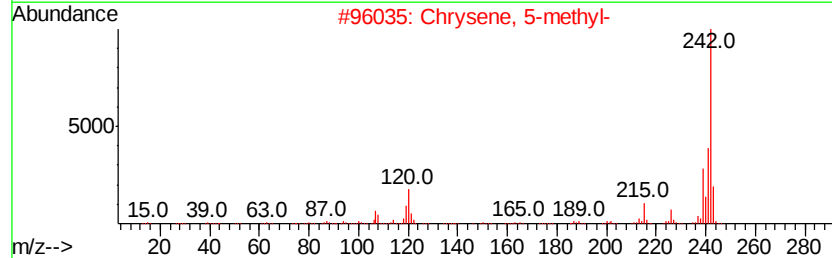
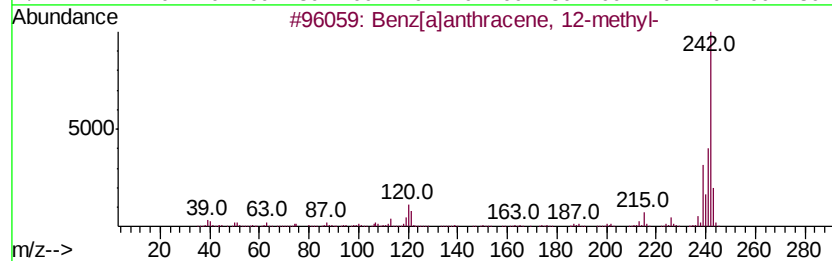
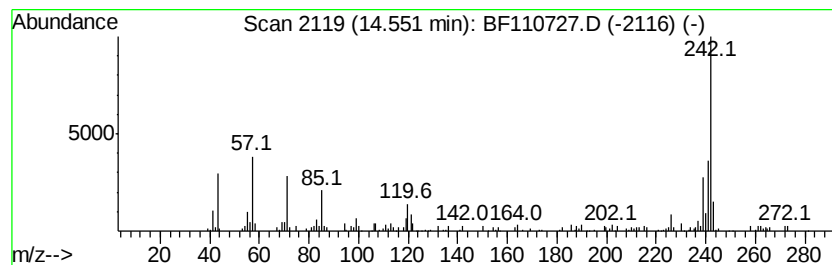
Instrument :  
BNA\_F  
ClientSampleId :  
TP-5S

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 14 Benz[a]anthracene, 12-methyl- Concentration Rank 15

| R.T.    | EstConc | Area                          | Relative to ISTD | R.T.           |
|---------|---------|-------------------------------|------------------|----------------|
| 14.55   | 2.52 ng | 95385                         | Chrysene-d12     | 14.13          |
| Hit# of | 5       | Tentative ID                  | MW MolForm       | CAS# Qual      |
| 1       |         | Benz[a]anthracene, 12-methyl- | 242 C19H14       | 002422-79-9 70 |
| 2       |         | Chrysene, 5-methyl-           | 242 C19H14       | 003697-24-3 70 |
| 3       |         | Benz[a]anthracene, 7-methyl-  | 242 C19H14       | 002541-69-7 60 |
| 4       |         | Chrysene, 1-methyl-           | 242 C19H14       | 003351-28-8 58 |
| 5       |         | Triphenylene, 2-methyl-       | 242 C19H14       | 001705-84-6 55 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111318\  
Data File : BF110727.D  
Acq On : 13 Nov 2018 18:59  
Operator : JU/SJ  
Sample : J5925-11  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-5S

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

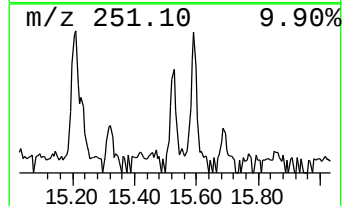
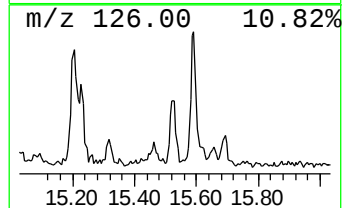
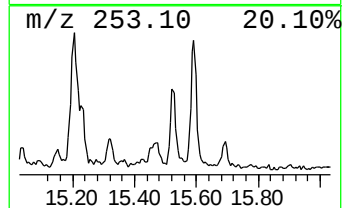
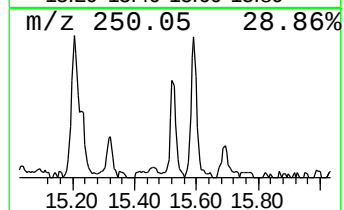
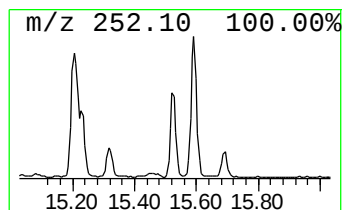
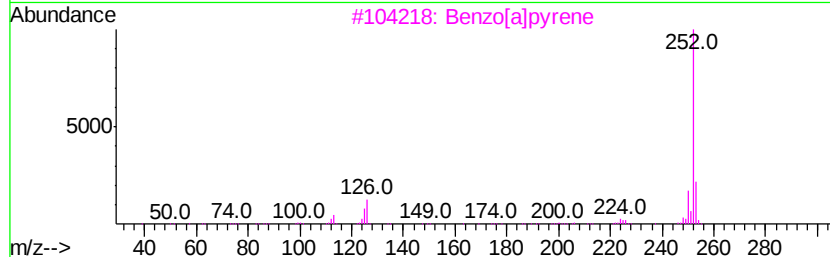
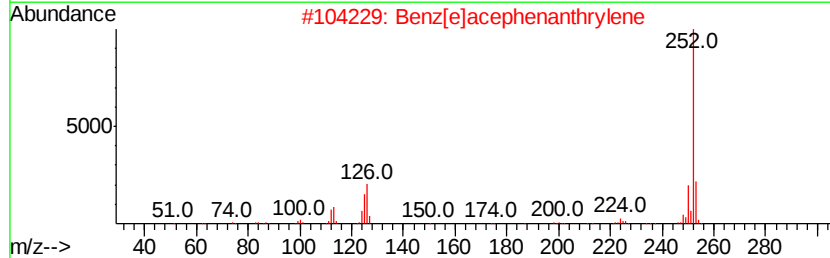
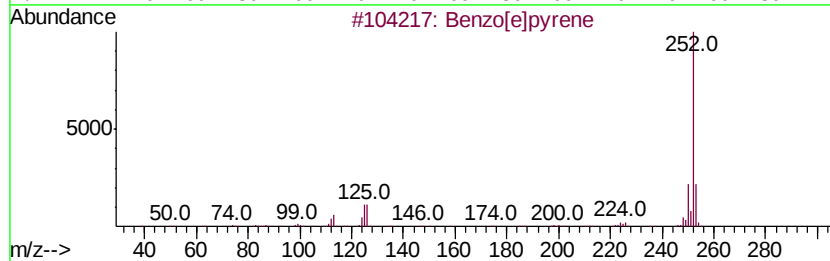
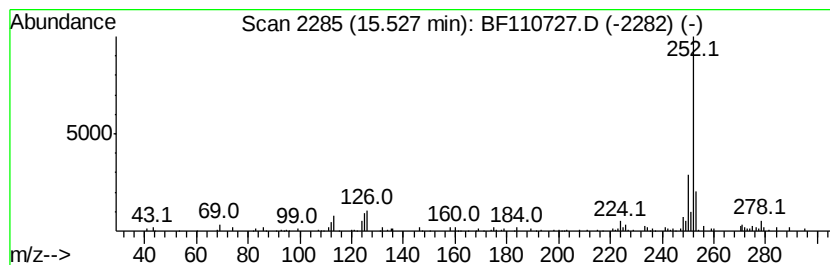
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 17 Benzo[e]pyrene Concentration Rank 9

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 15.53 | 3.65 ng | 71421 | Perylene-d12     | 15.66 |

| Hit# | of | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzo[e]bpvrene           | 252 | C20H12  | 000192-97-2 | 93   |
| 2    |    | Benzo[e]acephenanthrylene | 252 | C20H12  | 000205-99-2 | 91   |
| 3    |    | Benzo[a]bpvrene           | 252 | C20H12  | 000050-32-8 | 90   |
| 4    |    | Benzo[i]lfluoranthene     | 252 | C20H12  | 000205-82-3 | 90   |
| 5    |    | Perylene                  | 252 | C20H12  | 000198-55-0 | 81   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111318\  
Data File : BF110727.D  
Acq On : 13 Nov 2018 18:59  
Operator : JU/SJ  
Sample : J5925-11  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-5S

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

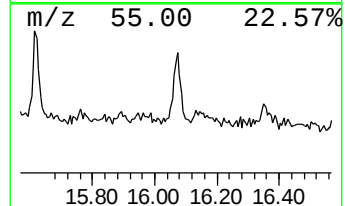
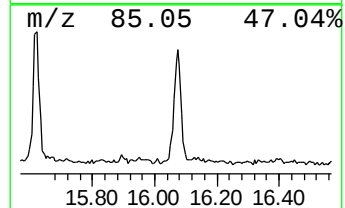
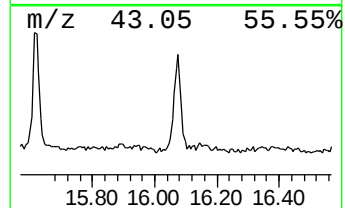
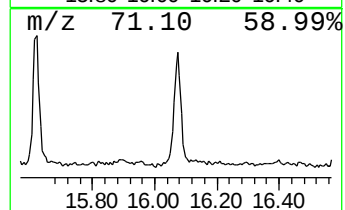
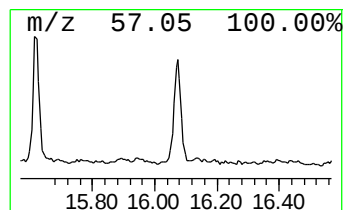
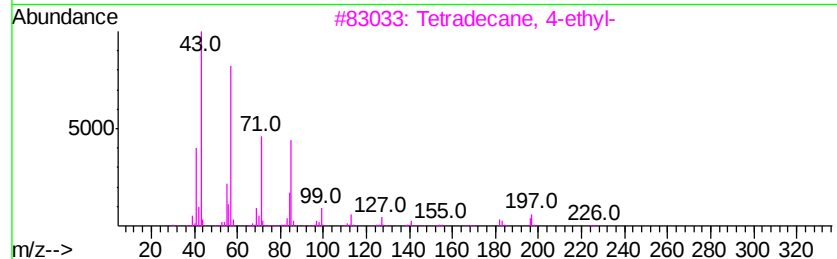
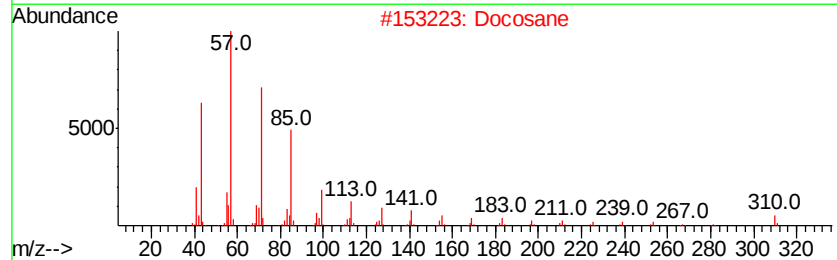
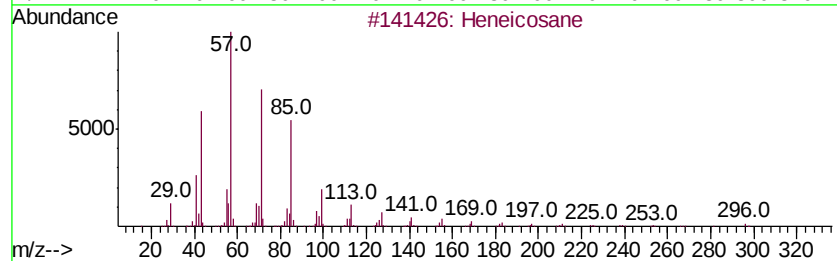
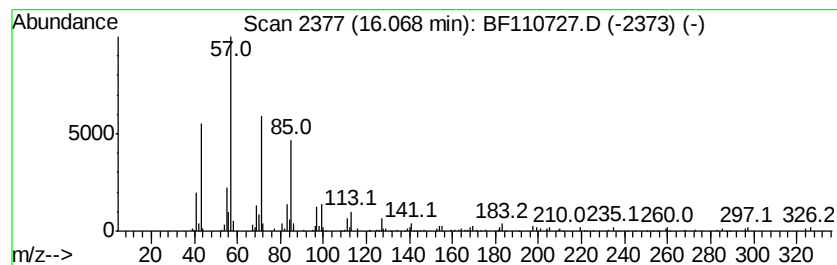
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 Heneicosane Concentration Rank 3

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 16.07 | 9.42 ng | 184418 | Perylene-d12     | 15.66 |

| Hit# of | 5                      | Tentative ID | MW  | MolForm | CAS#        | Qual |
|---------|------------------------|--------------|-----|---------|-------------|------|
| 1       | Heneicosane            |              | 296 | C21H44  | 000629-94-7 | 97   |
| 2       | Docosane               |              | 310 | C22H46  | 000629-97-0 | 96   |
| 3       | Tetradecane, 4-ethyl-  |              | 226 | C16H34  | 055045-14-2 | 90   |
| 4       | Tetratetracontane      |              | 619 | C44H90  | 007098-22-8 | 87   |
| 5       | Heneicosane, 11-decyl- |              | 437 | C31H64  | 055320-06-4 | 87   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111318\  
Data File : BF110727.D  
Acq On : 13 Nov 2018 18:59  
Operator : JU/SJ  
Sample : J5925-11  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-5S

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

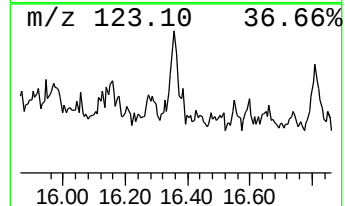
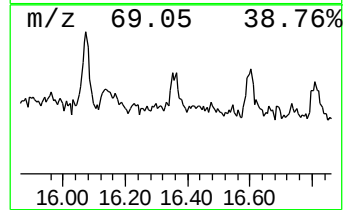
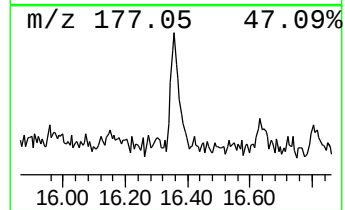
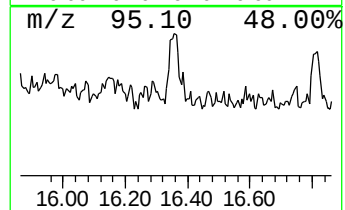
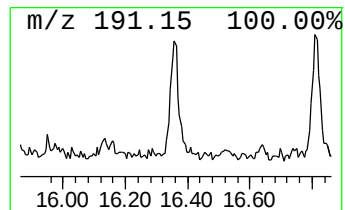
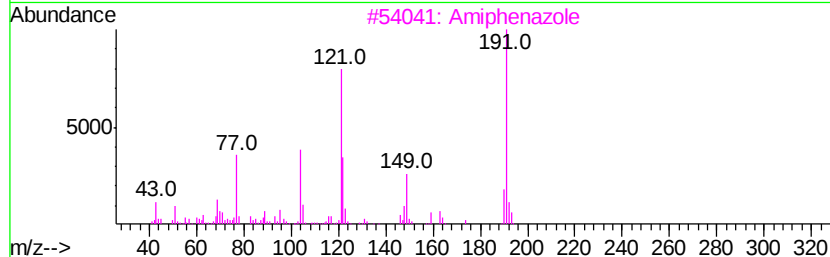
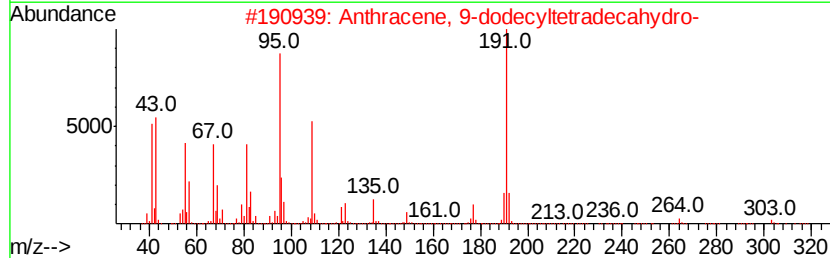
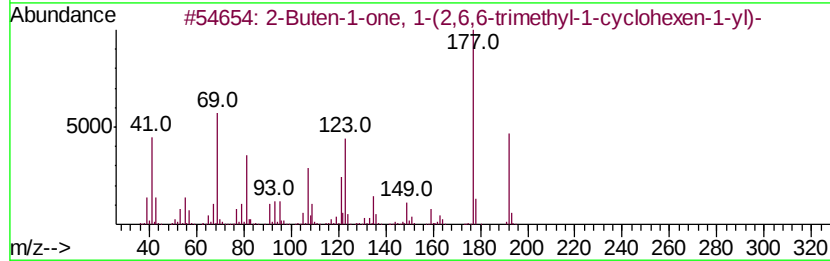
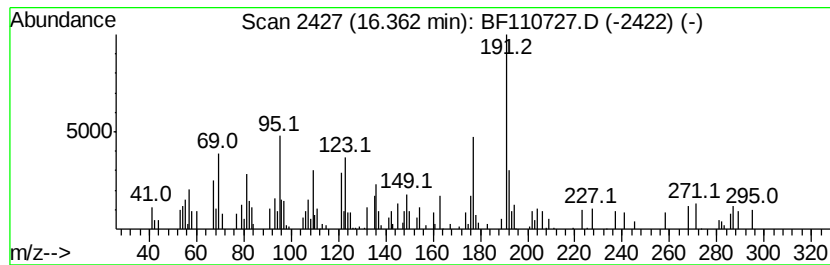
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 unknown16.36 Concentration Rank 11

| R.T.  | EstConc | Area  | Relative to ISTD | R.T.  |
|-------|---------|-------|------------------|-------|
| 16.36 | 3.32 ng | 65050 | Perylene-d12     | 15.66 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|-------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | 2-Buten-1-one, 1-(2,6,6-trimethy... | 192 | C13H20O   | 035044-68-9 | 41   |
| 2    |    | Anthracene, 9-dodecyltetradecahy... | 360 | C26H48    | 055401-75-7 | 38   |
| 3    |    | Amiphenazole                        | 191 | C9H9N3S   | 000490-55-1 | 25   |
| 4    |    | 1-Penten-3-one, 1-(2,6,6-trimeth... | 206 | C14H22O   | 000127-43-5 | 22   |
| 5    |    | 2-Chloro-4,6-dimethylquinoline      | 191 | C11H10ClN | 003913-18-6 | 22   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF111318\  
Data File : BF110727.D  
Acq On : 13 Nov 2018 18:59  
Operator : JU/SJ  
Sample : J5925-11  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-5S

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

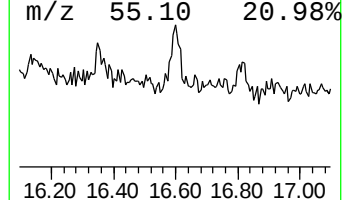
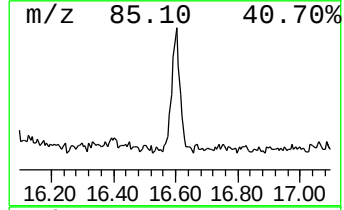
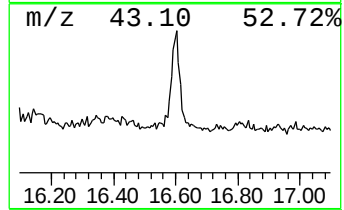
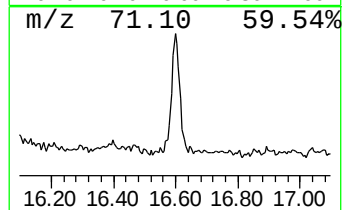
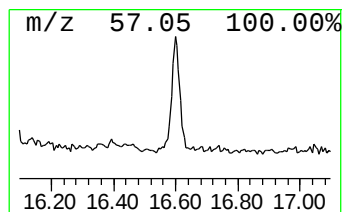
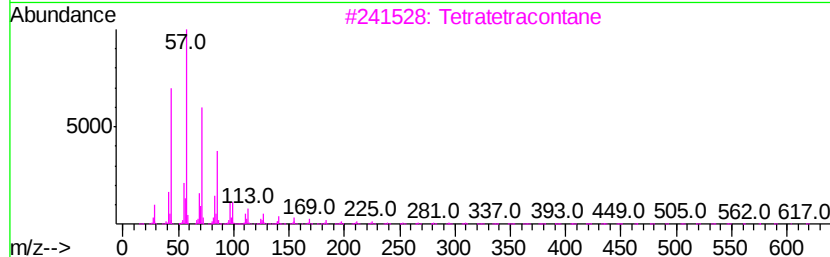
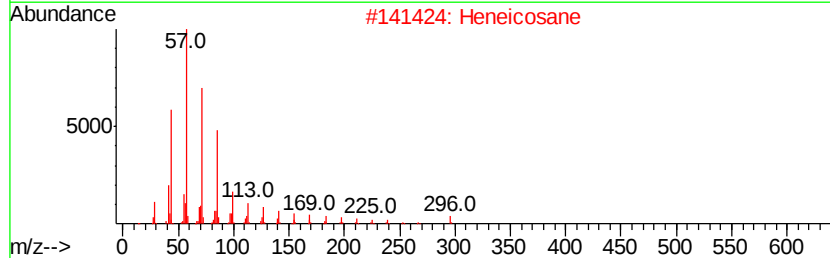
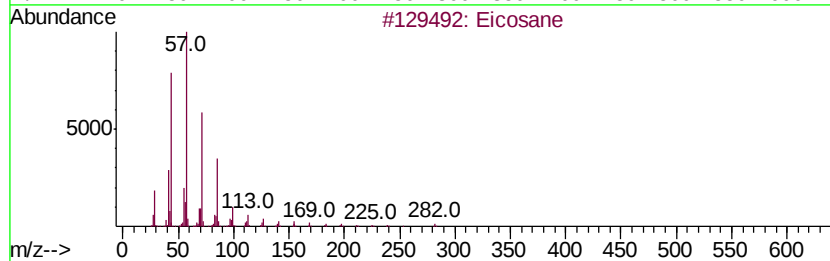
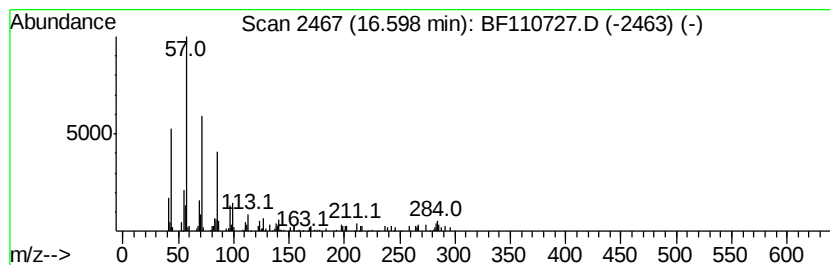
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 Eicosane Concentration Rank 5

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 16.60 | 6.87 ng | 134494 | Perylene-d12     | 15.66 |

| Hit# of | 5                  | Tentative ID | MW  | MolForm | CAS#         | Qual |
|---------|--------------------|--------------|-----|---------|--------------|------|
| 1       | Eicosane           |              | 282 | C20H42  | 000112-95-8  | 90   |
| 2       | Heneicosane        |              | 296 | C21H44  | 000629-94-7  | 89   |
| 3       | Tetratetracontane  |              | 619 | C44H90  | 007098-22-8  | 83   |
| 4       | 2-methyloctacosane |              | 408 | C29H60  | 1000376-72-8 | 80   |
| 5       | Hexatriacontane    |              | 507 | C36H74  | 000630-06-8  | 80   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF111318\  
 Data File : BF110727.D  
 Acq On : 13 Nov 2018 18:59  
 Operator : JU/SJ  
 Sample : J5925-11  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 TP-5S

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF110818.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 |
| Butane, 2-methoxy...  | 2.29  | 8.0     | ng    | 232884   | 1                     | 6.95  | 580346 | 20.0 |
| 2-Pentanone, 4-hy...  | 5.20  | 3.7     | ng    | 107697   | 1                     | 6.95  | 580346 | 20.0 |
| unknown6.72           | 6.72  | 88.0    | ng    | 2552550  | 1                     | 6.95  | 580346 | 20.0 |
| 1H-Cyclopropa[1]p...  | 11.98 | 4.6     | ng    | 131934   | 4                     | 11.48 | 578631 | 20.0 |
| Anthracene, 1-met...  | 12.09 | 2.5     | ng    | 73914    | 4                     | 11.48 | 578631 | 20.0 |
| Phenanthrene, 3,6...  | 12.53 | 2.6     | ng    | 76304    | 4                     | 11.48 | 578631 | 20.0 |
| Tricyclo[8.2.2.2(...  | 12.63 | 2.1     | ng    | 60155    | 4                     | 11.48 | 578631 | 20.0 |
| Pyrene, 1-methyl-     | 13.14 | 2.1     | ng    | 80524    | 5                     | 14.13 | 756304 | 20.0 |
| Fluoranthene, 2-m...  | 13.26 | 2.2     | ng    | 83616    | 5                     | 14.13 | 756304 | 20.0 |
| 9-Octadecenamide, ... | 13.53 | 4.3     | ng    | 162010   | 5                     | 14.13 | 756304 | 20.0 |
| Pentafluoropropio...  | 13.93 | 3.5     | ng    | 133691   | 5                     | 14.13 | 756304 | 20.0 |
| Triphenylene, 2-m...  | 14.52 | 2.4     | ng    | 90979    | 5                     | 14.13 | 756304 | 20.0 |
| Benz[a]anthracene...  | 14.55 | 2.5     | ng    | 95385    | 5                     | 14.13 | 756304 | 20.0 |
| Benzo[e]pyrene        | 15.53 | 3.6     | ng    | 71421    | 6                     | 15.66 | 391635 | 20.0 |
| Heneicosane           | 16.07 | 9.4     | ng    | 184418   | 6                     | 15.66 | 391635 | 20.0 |
| unknown16.36          | 16.36 | 3.3     | ng    | 65050    | 6                     | 15.66 | 391635 | 20.0 |
| Eicosane              | 16.60 | 6.9     | ng    | 134494   | 6                     | 15.66 | 391635 | 20.0 |