

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA\_F\DATA\BF090917\  
 Data File : BF098397.D  
 Acq On : 9 Sep 2017 20:30  
 Operator : SJ/JU  
 Sample : I5151-03 50X  
 Misc :  
 ALS Vial : 22 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 NYSEG-PH4-ESMI2-3

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-BF081517.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         | 6.498       | 746           | 750         | 754          | rBV2     | 183782         | 197471        | 2.21%           | 0.199%        |
| 2         | 6.675       | 776           | 780         | 783          | rBV      | 721995         | 592412        | 6.62%           | 0.597%        |
| 3         | 6.851       | 808           | 810         | 814          | rVB      | 303340         | 231814        | 2.59%           | 0.234%        |
| 4         | 6.934       | 820           | 824         | 827          | rBV      | 206985         | 178013        | 1.99%           | 0.179%        |
| 5         | 7.710       | 948           | 956         | 959          | rBV3     | 273597         | 336453        | 3.76%           | 0.339%        |
| 6         | 7.739       | 959           | 961         | 967          | rVB      | 218164         | 259521        | 2.90%           | 0.262%        |
| 7         | 7.998       | 992           | 1005        | 1007         | rBV      | 4963549        | 8951646       | 100.00%         | 9.025%        |
| 8         | 8.034       | 1007          | 1011        | 1014         | rVV      | 524266         | 421323        | 4.71%           | 0.425%        |
| 9         | 8.675       | 1114          | 1120        | 1123         | rBV      | 3343437        | 3502144       | 39.12%          | 3.531%        |
| 10        | 8.769       | 1130          | 1136        | 1139         | rVB      | 2241322        | 2109203       | 23.56%          | 2.127%        |
| 11        | 9.133       | 1191          | 1198        | 1201         | rBV      | 1271024        | 1104848       | 12.34%          | 1.114%        |
| 12        | 9.228       | 1211          | 1214        | 1219         | rVB      | 471256         | 466800        | 5.21%           | 0.471%        |
| 13        | 9.298       | 1219          | 1226        | 1229         | rBV      | 784937         | 1146473       | 12.81%          | 1.156%        |
| 14        | 9.369       | 1233          | 1238        | 1240         | rVV      | 1117803        | 1128602       | 12.61%          | 1.138%        |
| 15        | 9.392       | 1240          | 1242        | 1246         | rVV      | 821688         | 743916        | 8.31%           | 0.750%        |
| 16        | 9.433       | 1246          | 1249        | 1252         | rVV      | 221748         | 182869        | 2.04%           | 0.184%        |
| 17        | 9.486       | 1252          | 1258        | 1266         | rVV      | 541771         | 693179        | 7.74%           | 0.699%        |
| 18        | 9.569       | 1266          | 1272        | 1277         | rVB      | 979421         | 846721        | 9.46%           | 0.854%        |
| 19        | 9.710       | 1287          | 1296        | 1298         | rBV2     | 864724         | 1234951       | 13.80%          | 1.245%        |
| 20        | 9.751       | 1298          | 1303        | 1305         | rVV      | 3963878        | 4455922       | 49.78%          | 4.493%        |
| 21        | 9.916       | 1327          | 1331        | 1334         | rVV      | 2613011        | 2585754       | 28.89%          | 2.607%        |
| 22        | 9.951       | 1334          | 1337        | 1341         | rVV      | 296700         | 256767        | 2.87%           | 0.259%        |
| 23        | 10.022      | 1345          | 1349        | 1352         | rBV2     | 353931         | 434053        | 4.85%           | 0.438%        |
| 24        | 10.051      | 1352          | 1354        | 1360         | rVB2     | 289725         | 278575        | 3.11%           | 0.281%        |
| 25        | 10.116      | 1360          | 1365        | 1366         | rBV2     | 212245         | 272829        | 3.05%           | 0.275%        |
| 26        | 10.157      | 1370          | 1372        | 1375         | rVV      | 362372         | 297762        | 3.33%           | 0.300%        |
| 27        | 10.210      | 1375          | 1381        | 1382         | rVV2     | 142076         | 223567        | 2.50%           | 0.225%        |
| 28        | 10.222      | 1382          | 1383        | 1385         | rVV      | 251835         | 198318        | 2.22%           | 0.200%        |
| 29        | 10.263      | 1385          | 1390        | 1394         | rVB      | 2826312        | 3022171       | 33.76%          | 3.047%        |
| 30        | 10.339      | 1401          | 1403        | 1405         | rVV2     | 193132         | 199462        | 2.23%           | 0.201%        |
| 31        | 10.363      | 1405          | 1407        | 1410         | rVV2     | 393203         | 376179        | 4.20%           | 0.379%        |
| 32        | 10.410      | 1412          | 1415        | 1418         | rVV      | 753714         | 713247        | 7.97%           | 0.719%        |
| 33        | 10.475      | 1422          | 1426        | 1427         | rVV      | 647164         | 544350        | 6.08%           | 0.549%        |
| 34        | 10.492      | 1427          | 1429        | 1433         | rVV      | 864910         | 916197        | 10.23%          | 0.924%        |

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 ALS Vial : 22 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 NYSEG-PH4-ESMI2-3

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-BF081517.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 35 | 10.798 | 1476 | 1481 | 1484 | rBV2 | 509631  | 592624  | 6.62%  | 0.597% |
| 36 | 10.880 | 1492 | 1495 | 1499 | rVV2 | 272325  | 278575  | 3.11%  | 0.281% |
| 37 | 10.927 | 1500 | 1503 | 1505 | rVV  | 279736  | 280119  | 3.13%  | 0.282% |
| 38 | 10.951 | 1505 | 1507 | 1509 | rVV  | 269139  | 253046  | 2.83%  | 0.255% |
| 39 | 10.998 | 1513 | 1515 | 1518 | rVV2 | 261653  | 233011  | 2.60%  | 0.235% |
| 40 | 11.039 | 1518 | 1522 | 1525 | rVV  | 283963  | 314342  | 3.51%  | 0.317% |
| 41 | 11.086 | 1525 | 1530 | 1537 | rVB  | 966237  | 988372  | 11.04% | 0.996% |
| 42 | 11.233 | 1543 | 1555 | 1557 | rBV  | 4360572 | 8082361 | 90.29% | 8.149% |
| 43 | 11.275 | 1557 | 1562 | 1565 | rVV  | 2783369 | 2818026 | 31.48% | 2.841% |
| 44 | 11.304 | 1565 | 1567 | 1570 | rVV  | 339447  | 321063  | 3.59%  | 0.324% |
| 45 | 11.427 | 1580 | 1588 | 1594 | rVV  | 1069823 | 1153380 | 12.88% | 1.163% |
| 46 | 11.486 | 1594 | 1598 | 1602 | rVV3 | 279353  | 351414  | 3.93%  | 0.354% |
| 47 | 11.527 | 1602 | 1605 | 1608 | rVV2 | 136067  | 186108  | 2.08%  | 0.188% |
| 48 | 11.563 | 1608 | 1611 | 1615 | rVV2 | 339287  | 333956  | 3.73%  | 0.337% |
| 49 | 11.692 | 1627 | 1633 | 1635 | rBV  | 875085  | 831161  | 9.29%  | 0.838% |
| 50 | 11.722 | 1635 | 1638 | 1642 | rVB  | 1247445 | 1034592 | 11.56% | 1.043% |
| 51 | 11.769 | 1642 | 1646 | 1649 | rBV  | 718177  | 627835  | 7.01%  | 0.633% |
| 52 | 11.804 | 1649 | 1652 | 1654 | rVV  | 1504348 | 1574006 | 17.58% | 1.587% |
| 53 | 11.827 | 1654 | 1656 | 1660 | rVV  | 511650  | 438430  | 4.90%  | 0.442% |
| 54 | 11.986 | 1680 | 1683 | 1688 | rVB  | 630660  | 562428  | 6.28%  | 0.567% |
| 55 | 12.239 | 1720 | 1726 | 1729 | rBV2 | 373783  | 498806  | 5.57%  | 0.503% |
| 56 | 12.280 | 1729 | 1733 | 1734 | rVV  | 271093  | 339082  | 3.79%  | 0.342% |
| 57 | 12.327 | 1738 | 1741 | 1748 | rVB3 | 169703  | 280888  | 3.14%  | 0.283% |
| 58 | 12.416 | 1749 | 1756 | 1759 | rBV  | 3766963 | 5012786 | 56.00% | 5.054% |
| 59 | 12.498 | 1767 | 1770 | 1773 | rVB  | 928726  | 787033  | 8.79%  | 0.794% |
| 60 | 12.551 | 1773 | 1779 | 1781 | rBV2 | 324183  | 360701  | 4.03%  | 0.364% |
| 61 | 12.645 | 1787 | 1795 | 1798 | rBV2 | 3347001 | 5410488 | 60.44% | 5.455% |
| 62 | 12.680 | 1798 | 1801 | 1808 | rVB4 | 337020  | 425537  | 4.75%  | 0.429% |
| 63 | 12.745 | 1808 | 1812 | 1815 | rBV  | 361951  | 380987  | 4.26%  | 0.384% |
| 64 | 12.786 | 1817 | 1819 | 1823 | rVB  | 269966  | 260768  | 2.91%  | 0.263% |
| 65 | 12.851 | 1823 | 1830 | 1833 | rBV2 | 309509  | 565718  | 6.32%  | 0.570% |
| 66 | 12.933 | 1837 | 1844 | 1845 | rVV4 | 304175  | 364779  | 4.07%  | 0.368% |
| 67 | 12.957 | 1845 | 1848 | 1851 | rVB  | 920879  | 892087  | 9.97%  | 0.899% |
| 68 | 13.021 | 1856 | 1859 | 1862 | rVV  | 889518  | 899995  | 10.05% | 0.907% |
| 69 | 13.057 | 1862 | 1865 | 1869 | rVV2 | 493424  | 606481  | 6.78%  | 0.611% |
| 70 | 13.121 | 1873 | 1876 | 1879 | rVV2 | 403263  | 487820  | 5.45%  | 0.492% |
| 71 | 13.151 | 1879 | 1881 | 1884 | rVV  | 279115  | 284668  | 3.18%  | 0.287% |

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 Misc :  
 ALS Vial : 22 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 NYSEG-PH4-ESMI2-3

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-BF081517.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|     |        |      |      |      |      |         |         |        |        |
|-----|--------|------|------|------|------|---------|---------|--------|--------|
| 72  | 13.180 | 1884 | 1886 | 1893 | rVB2 | 296307  | 398605  | 4.45%  | 0.402% |
| 73  | 13.357 | 1910 | 1916 | 1918 | rVV4 | 120504  | 210647  | 2.35%  | 0.212% |
| 74  | 13.439 | 1926 | 1930 | 1933 | rBV  | 157528  | 223619  | 2.50%  | 0.225% |
| 75  | 13.574 | 1947 | 1953 | 1955 | rBV  | 192256  | 262361  | 2.93%  | 0.265% |
| 76  | 13.598 | 1955 | 1957 | 1959 | rBV  | 344144  | 271041  | 3.03%  | 0.273% |
| 77  | 13.621 | 1959 | 1961 | 1963 | rBV  | 219697  | 192432  | 2.15%  | 0.194% |
| 78  | 13.821 | 1987 | 1995 | 1998 | rBV2 | 2303664 | 3468060 | 38.74% | 3.497% |
| 79  | 13.857 | 1998 | 2001 | 2008 | rVV  | 1840477 | 1793148 | 20.03% | 1.808% |
| 80  | 13.933 | 2011 | 2014 | 2018 | rVV  | 678869  | 588511  | 6.57%  | 0.593% |
| 81  | 14.051 | 2032 | 2034 | 2037 | rVB  | 236588  | 218768  | 2.44%  | 0.221% |
| 82  | 14.168 | 2051 | 2054 | 2056 | rVV2 | 245835  | 269377  | 3.01%  | 0.272% |
| 83  | 14.210 | 2057 | 2061 | 2064 | rVV  | 540351  | 615889  | 6.88%  | 0.621% |
| 84  | 14.245 | 2064 | 2067 | 2069 | rVV2 | 215765  | 242128  | 2.70%  | 0.244% |
| 85  | 14.304 | 2071 | 2077 | 2080 | rVV2 | 306485  | 501960  | 5.61%  | 0.506% |
| 86  | 14.333 | 2080 | 2082 | 2084 | rVV2 | 267861  | 279089  | 3.12%  | 0.281% |
| 87  | 14.357 | 2084 | 2086 | 2092 | rVV2 | 368171  | 528763  | 5.91%  | 0.533% |
| 88  | 14.686 | 2139 | 2142 | 2148 | rBV  | 161853  | 189127  | 2.11%  | 0.191% |
| 89  | 14.845 | 2162 | 2169 | 2177 | rBV  | 1405619 | 3079577 | 34.40% | 3.105% |
| 90  | 14.933 | 2181 | 2184 | 2188 | rVB  | 472611  | 500100  | 5.59%  | 0.504% |
| 91  | 15.115 | 2210 | 2215 | 2221 | rVB  | 815007  | 1176031 | 13.14% | 1.186% |
| 92  | 15.180 | 2221 | 2226 | 2230 | rVV  | 1438267 | 1829017 | 20.43% | 1.844% |
| 93  | 15.233 | 2231 | 2235 | 2237 | rVV  | 467724  | 546966  | 6.11%  | 0.551% |
| 94  | 15.262 | 2237 | 2240 | 2245 | rVB  | 533322  | 652758  | 7.29%  | 0.658% |
| 95  | 16.580 | 2457 | 2464 | 2469 | rVB  | 539468  | 992488  | 11.09% | 1.001% |
| 96  | 16.733 | 2484 | 2490 | 2494 | rBV  | 124525  | 234655  | 2.62%  | 0.237% |
| 97  | 16.986 | 2525 | 2533 | 2543 | rBV  | 398762  | 846660  | 9.46%  | 0.854% |
| 98  | 17.198 | 2563 | 2569 | 2576 | rVB  | 192212  | 381668  | 4.26%  | 0.385% |
| 99  | 19.027 | 2871 | 2880 | 2893 | rVB2 | 79349   | 262907  | 2.94%  | 0.265% |
| 100 | 19.203 | 2897 | 2910 | 2915 | rBV2 | 62189   | 211443  | 2.36%  | 0.213% |

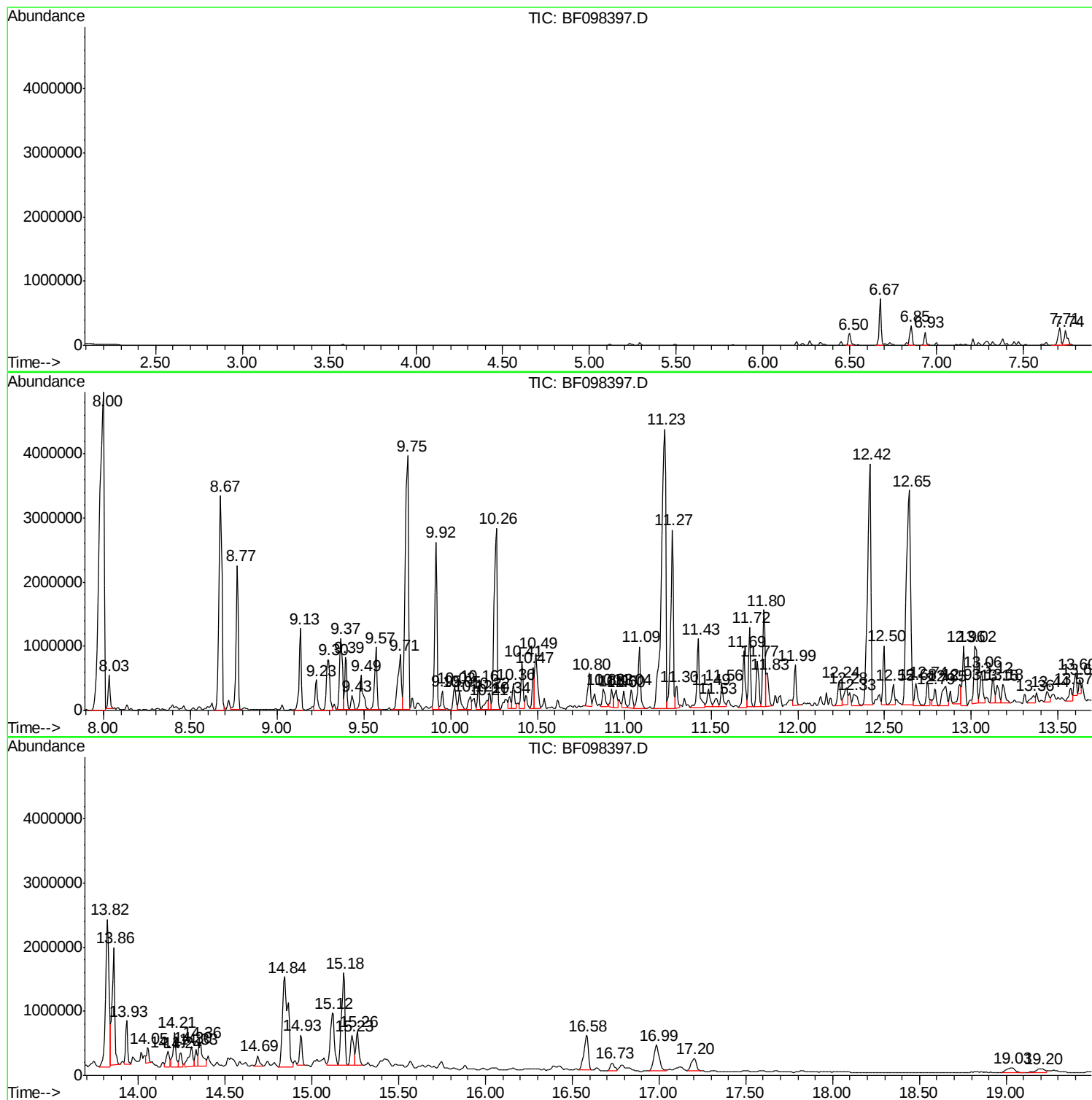
Sum of corrected areas: 99184780

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ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
NYSEG-PH4-ESMI2-3

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF090917\  
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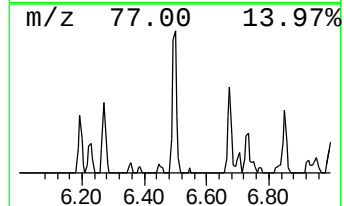
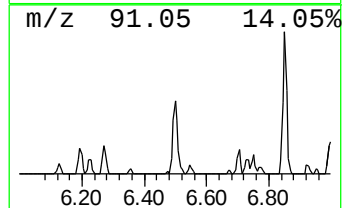
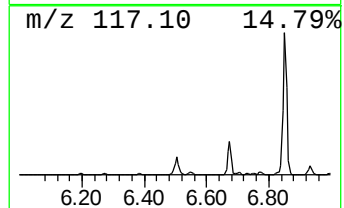
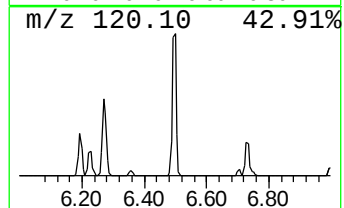
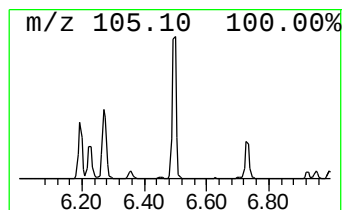
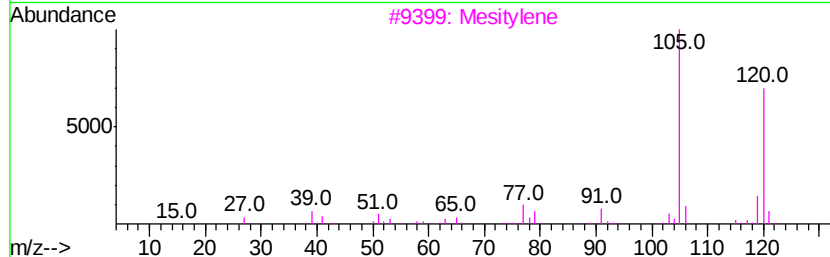
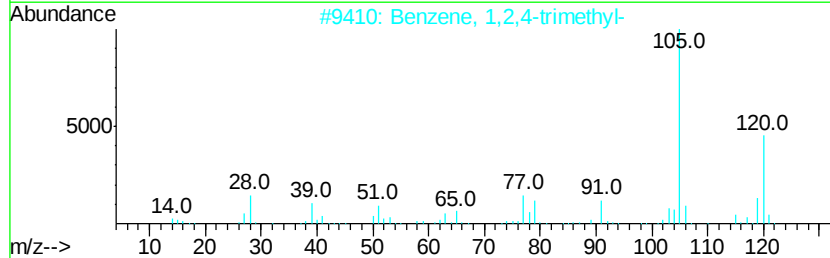
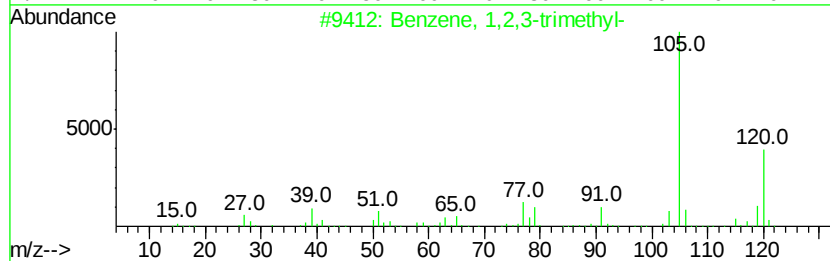
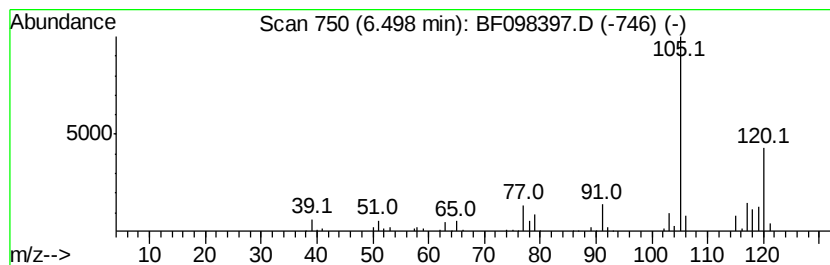
Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF081517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 1 Benzene, 1,2,3-trimethyl- Concentration Rank 16

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 6.50 | 6.67 ng | 197471 | 1,4-Dichlorobenzene-d4 | 6.67 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 93   |
| 2    |    |   | Benzene, 1,2,4-trimethyl-  | 120 | C9H12   | 000095-63-6 | 91   |
| 3    |    |   | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 87   |
| 4    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 87   |
| 5    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 81   |



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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

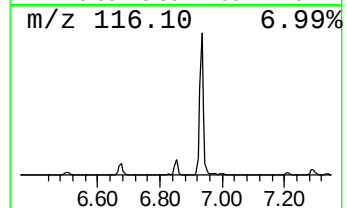
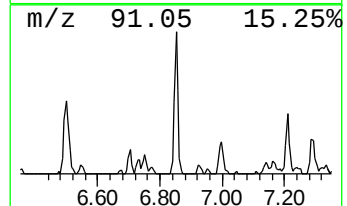
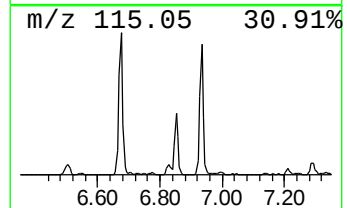
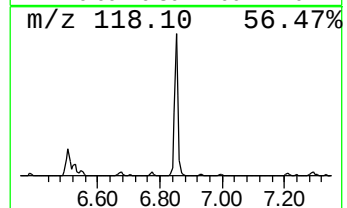
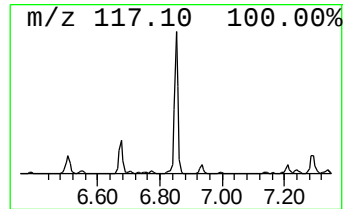
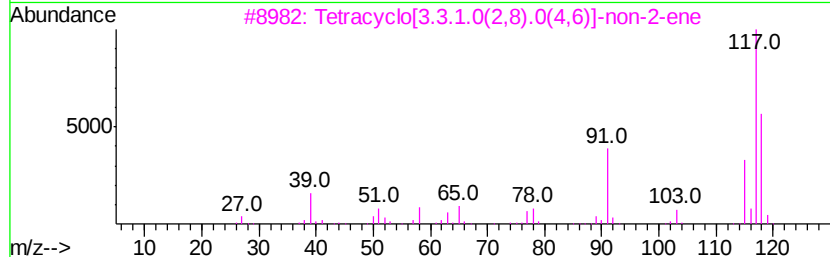
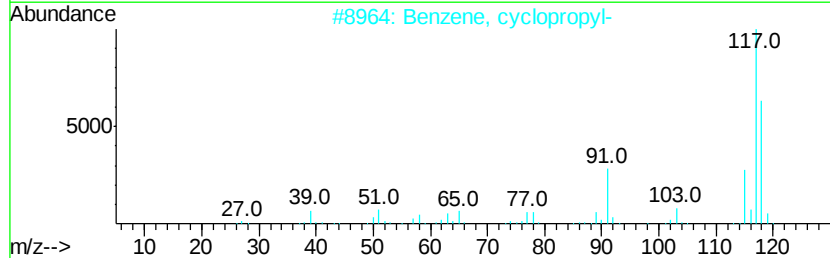
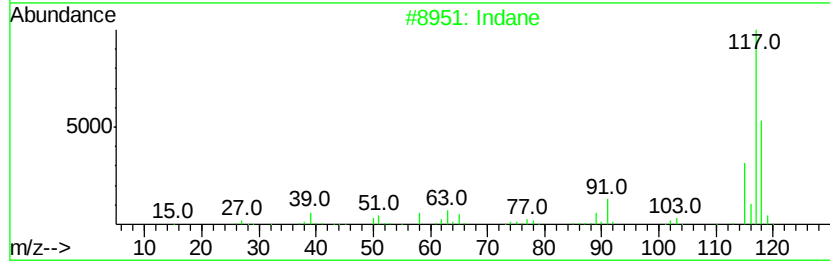
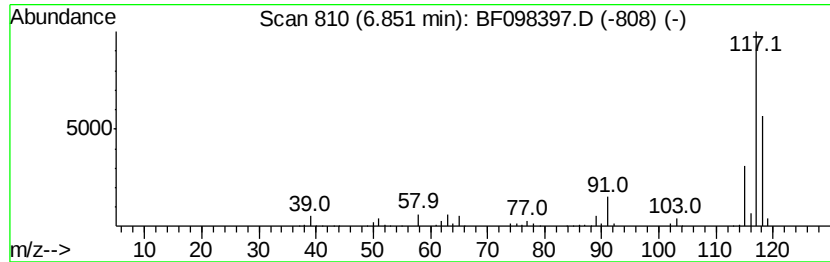
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 Indane Concentration Rank 11

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 6.85 | 7.83 ng | 231814 | 1,4-Dichlorobenzene-d4 | 6.67 |

| Hit# | of | 5 | Tentative ID                     | MW  | MolForm | CAS#         | Qual |
|------|----|---|----------------------------------|-----|---------|--------------|------|
| 1    |    |   | Indane                           | 118 | C9H10   | 000496-11-7  | 91   |
| 2    |    |   | Benzene, cyclopropyl-            | 118 | C9H10   | 000873-49-4  | 87   |
| 3    |    |   | Tetracyclo[3.3.1.0(2,8).0(4,6)]- | 118 | C9H10   | 1000191-13-7 | 80   |
| 4    |    |   | Benzene, 1-propenyl-             | 118 | C9H10   | 000637-50-3  | 72   |
| 5    |    |   | Benzene, 2-propenyl-             | 118 | C9H10   | 000300-57-2  | 64   |



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Data File : BF098397.D  
Acq On : 9 Sep 2017 20:30  
Operator : SJ/JU  
Sample : I5151-03 50X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

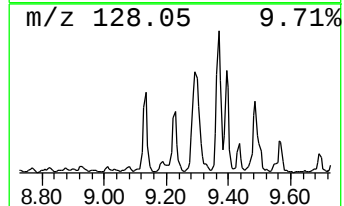
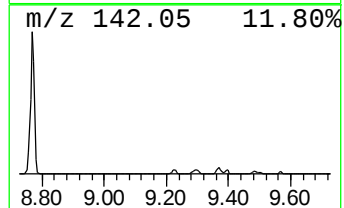
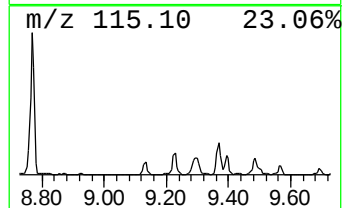
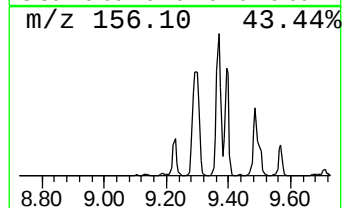
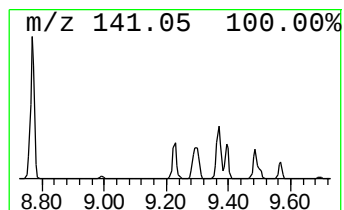
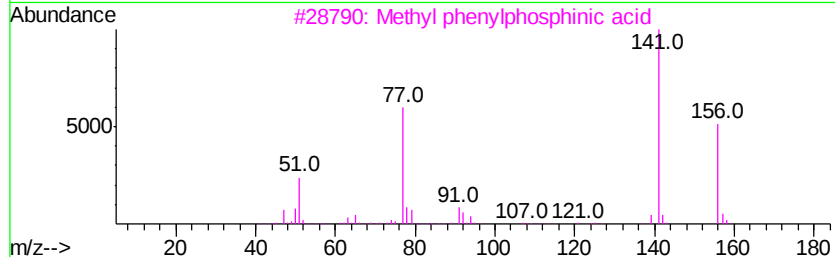
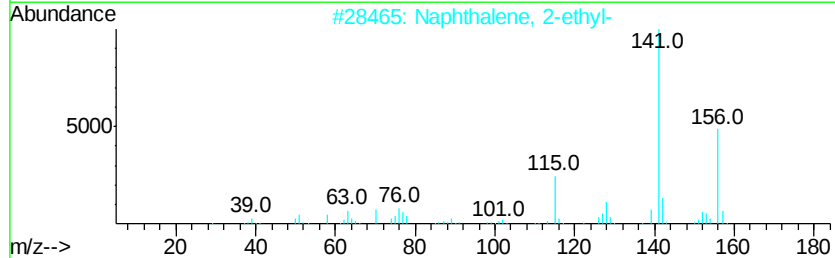
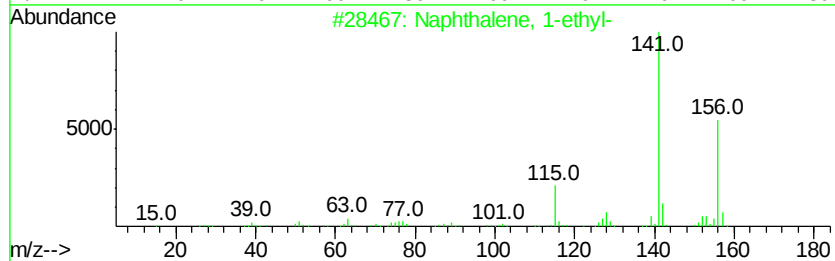
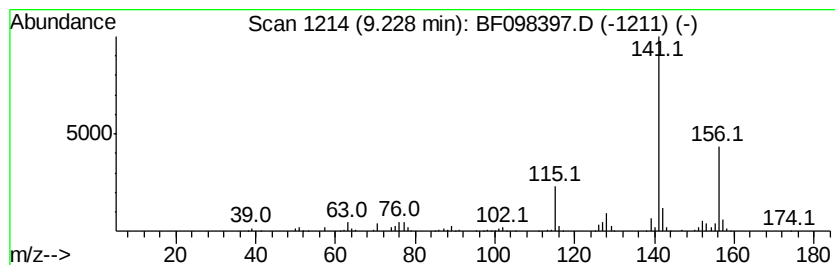
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ClientSampleId :  
NYSEG-PH4-ESMI2-3

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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 4 Naphthalene, 1-ethyl- Concentration Rank 13

| R.T.    | EstConc | Area                         | Relative to ISTD | R.T.           |
|---------|---------|------------------------------|------------------|----------------|
| 9.23    | 7.56 ng | 466800                       | Acenaphthene-d10 | 9.71           |
| Hit# of | 5       | Tentative ID                 | MW MolForm       | CAS# Qual      |
| 1       |         | Naphthalene, 1-ethyl-        | 156 C12H12       | 001127-76-0 96 |
| 2       |         | Naphthalene, 2-ethyl-        | 156 C12H12       | 000939-27-5 94 |
| 3       |         | Methyl phenylphosphinic acid | 156 C7H9O2P      | 004271-13-0 53 |
| 4       |         | Naphthalene, 1-propyl-       | 170 C13H14       | 002765-18-6 53 |
| 5       |         | 3',4'-Difluoroacetophenone   | 156 C8H6F2O      | 000369-33-5 53 |



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF090917\  
Data File : BF098397.D  
Acq On : 9 Sep 2017 20:30  
Operator : SJ/JU  
Sample : I5151-03 50X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

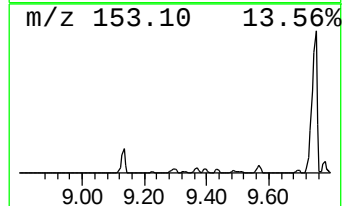
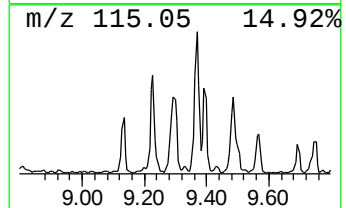
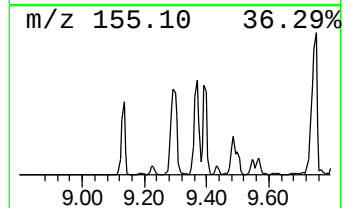
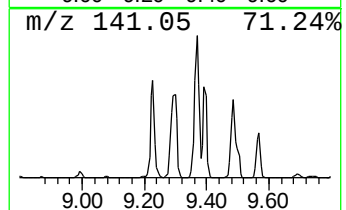
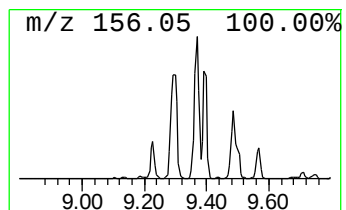
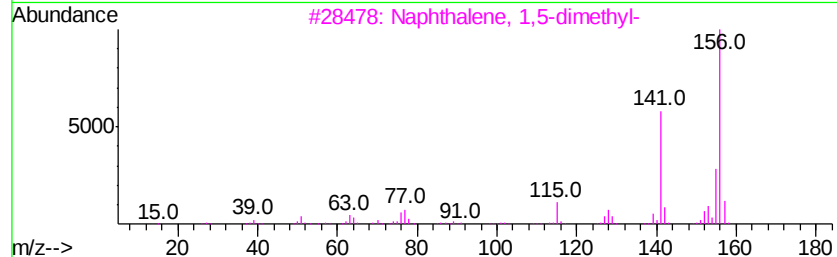
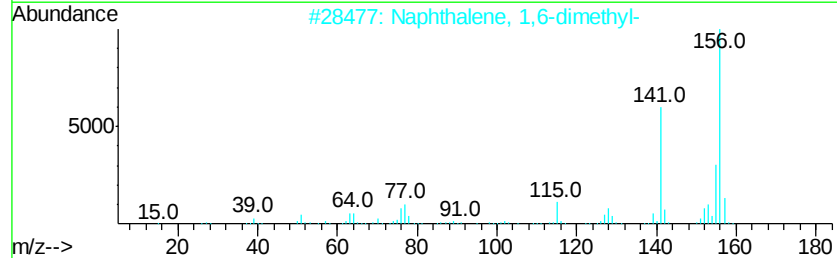
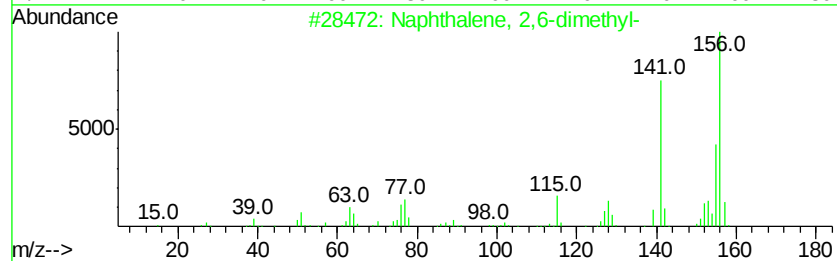
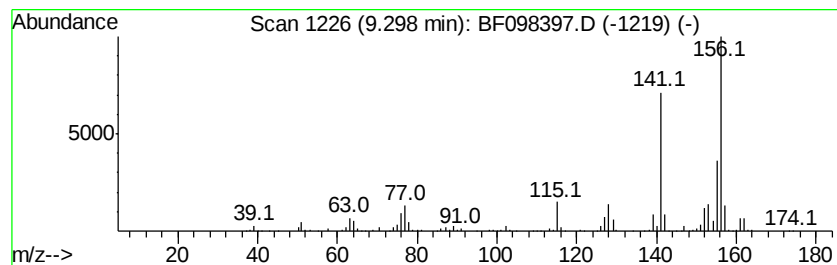
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ClientSampleId :  
NYSEG-PH4-ESMI2-3

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.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 5 Naphthalene, 2,6-dimethyl- Concentration Rank 2

| R.T.    | EstConc                    | Area         | Relative to ISTD |             | R.T. |      |
|---------|----------------------------|--------------|------------------|-------------|------|------|
| 9.30    | 18.57 ng                   | 1146470      | Acenaphthene-d10 |             | 9.71 |      |
| Hit# of | 5                          | Tentative ID | MW               | MolForm     | CAS# | Qual |
| 1       | Naphthalene, 2,6-dimethyl- | 156          | C12H12           | 000581-42-0 | 98   |      |
| 2       | Naphthalene, 1,6-dimethyl- | 156          | C12H12           | 000575-43-9 | 97   |      |
| 3       | Naphthalene, 1,5-dimethyl- | 156          | C12H12           | 000571-61-9 | 97   |      |
| 4       | Naphthalene, 1,7-dimethyl- | 156          | C12H12           | 000575-37-1 | 97   |      |
| 5       | Naphthalene, 2,7-dimethyl- | 156          | C12H12           | 000582-16-1 | 97   |      |



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF090917\  
Data File : BF098397.D  
Acq On : 9 Sep 2017 20:30  
Operator : SJ/JU  
Sample : I5151-03 50X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

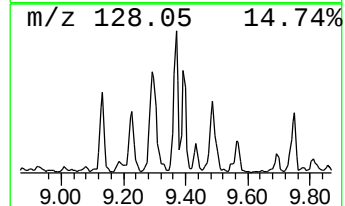
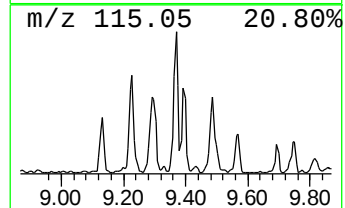
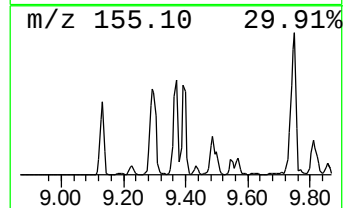
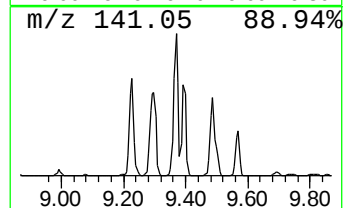
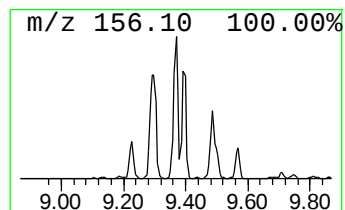
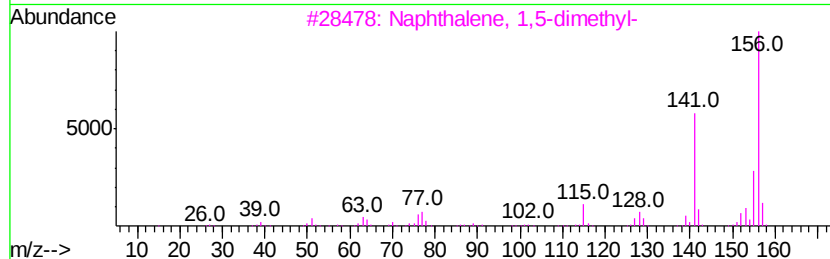
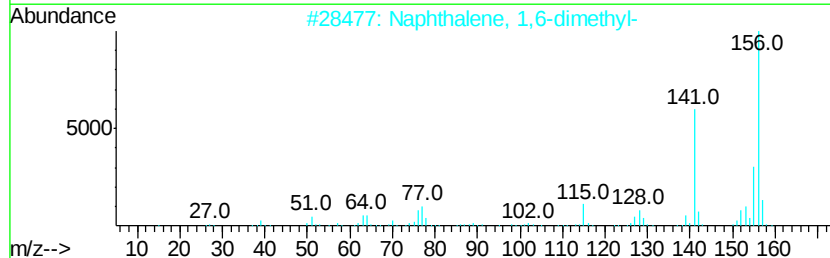
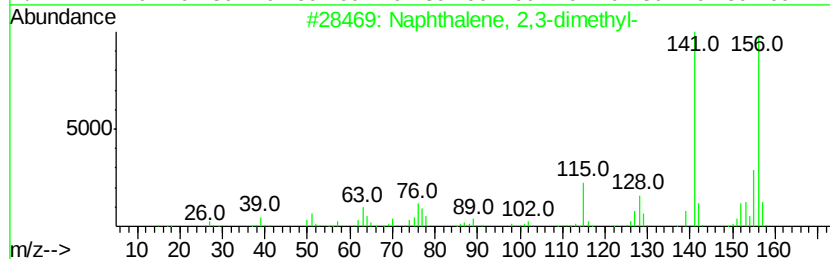
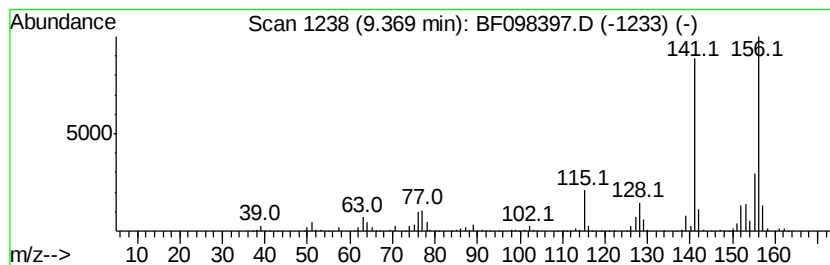
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ClientSampleId :  
NYSEG-PH4-ESMI2-3

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 6 Naphthalene, 2,3-dimethyl- Concentration Rank 4

| R.T.      | EstConc                    | Area    | Relative to ISTD | R.T.        |      |
|-----------|----------------------------|---------|------------------|-------------|------|
| 9.37      | 18.28 ng                   | 1128600 | Acenaphthene-d10 | 9.71        |      |
| Hit# of 5 | Tentative ID               | MW      | MolForm          | CAS#        | Qual |
| 1         | Naphthalene, 2,3-dimethyl- | 156     | C12H12           | 000581-40-8 | 98   |
| 2         | Naphthalene, 1,6-dimethyl- | 156     | C12H12           | 000575-43-9 | 97   |
| 3         | Naphthalene, 1,5-dimethyl- | 156     | C12H12           | 000571-61-9 | 97   |
| 4         | Naphthalene, 1,4-dimethyl- | 156     | C12H12           | 000571-58-4 | 97   |
| 5         | Naphthalene, 2,7-dimethyl- | 156     | C12H12           | 000582-16-1 | 97   |



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF090917\  
Data File : BF098397.D  
Acq On : 9 Sep 2017 20:30  
Operator : SJ/JU  
Sample : I5151-03 50X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
NYSEG-PH4-ESMI2-3

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

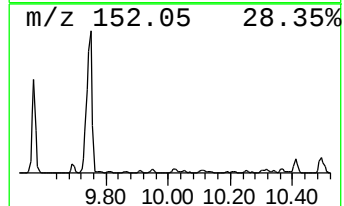
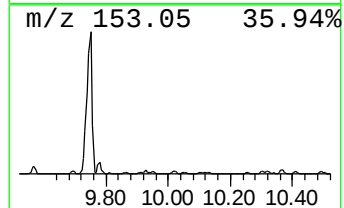
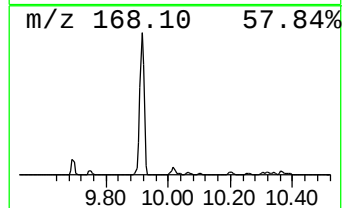
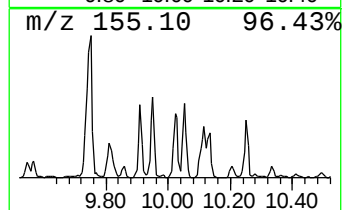
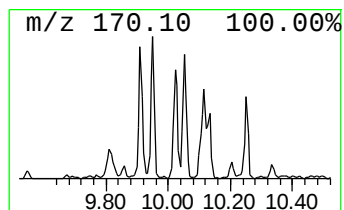
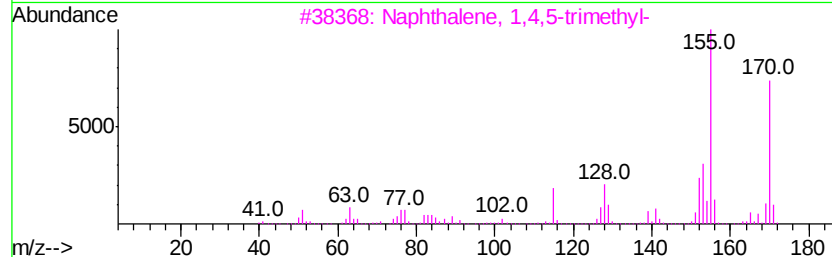
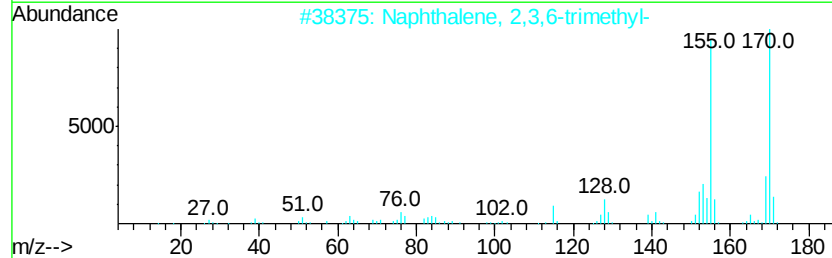
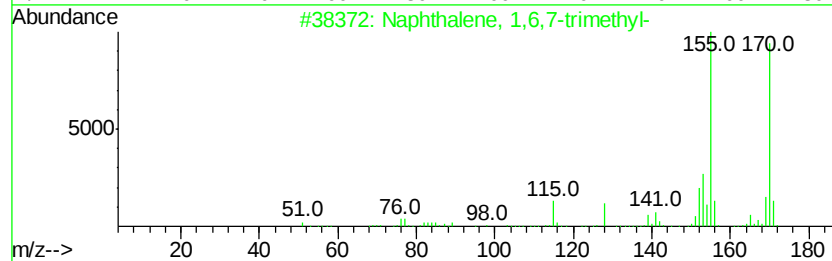
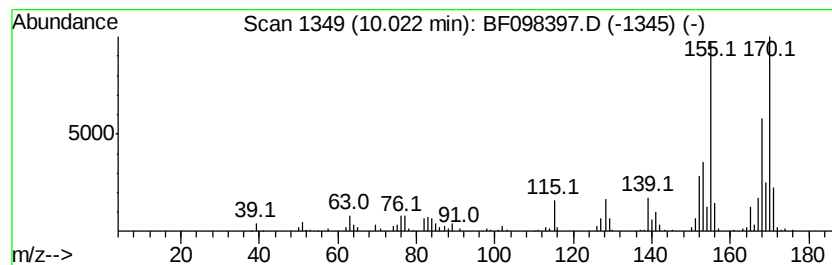
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 Naphthalene, 1,6,7-trimethyl- Concentration Rank 14

| R.T.  | EstConc | Area   | Relative to ISTD | R.T. |
|-------|---------|--------|------------------|------|
| 10.02 | 7.03 ng | 434053 | Acenaphthene-d10 | 9.71 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------------------|-----|---------|--------------|------|
| 1    | 5  | Naphthalene, 1,6,7-trimethyl-       | 170 | C13H14  | 002245-38-7  | 95   |
| 2    |    | Naphthalene, 2,3,6-trimethyl-       | 170 | C13H14  | 000829-26-5  | 94   |
| 3    |    | Naphthalene, 1,4,5-trimethyl-       | 170 | C13H14  | 002131-41-1  | 92   |
| 4    |    | Naphthalene, 1,4,6-trimethyl-       | 170 | C13H14  | 002131-42-2  | 76   |
| 5    |    | 1-(2,2-Dimethylcyclopropyl)-2-ph... | 170 | C13H14  | 1000294-91-1 | 70   |



Data Path : \\74.0.250.170\SVOASRV\HPCHEM1\BNA F\DATA\BF090917\  
Data File : BF098397.D  
Acq On : 9 Sep 2017 20:30  
Operator : SJ/JU  
Sample : I5151-03 50X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
NYSEG-PH4-ESMI2-3

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

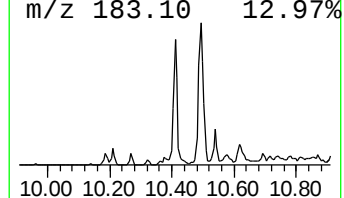
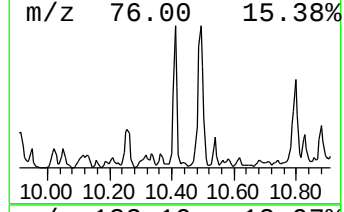
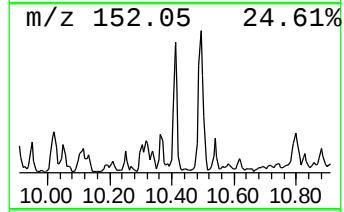
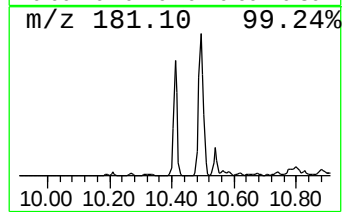
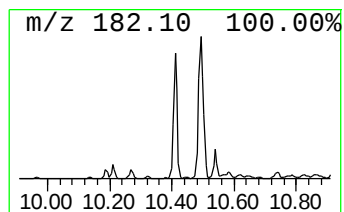
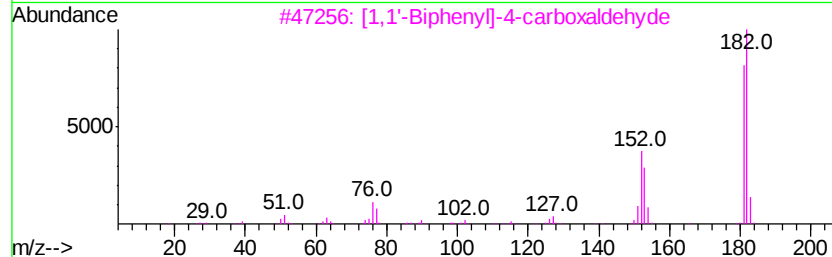
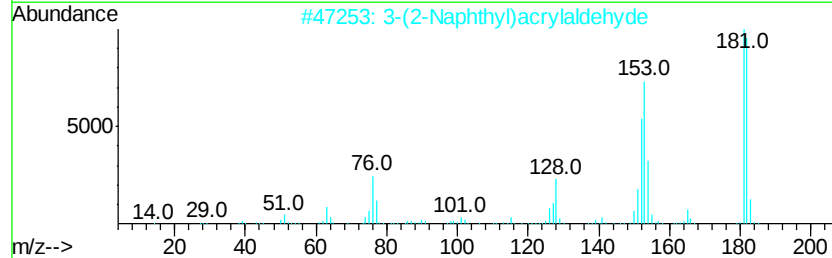
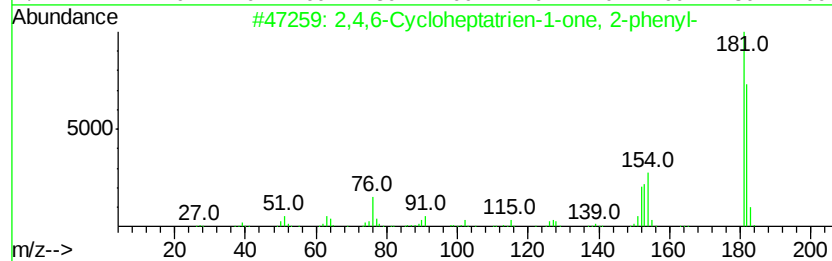
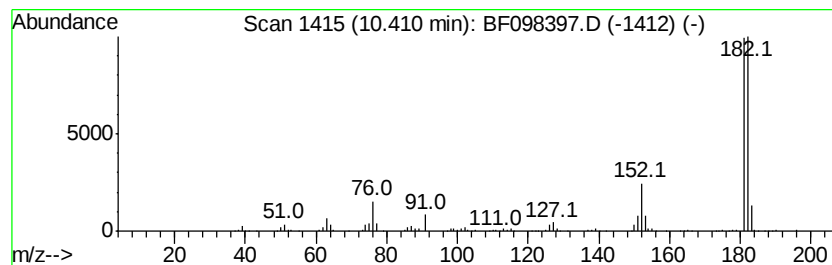
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 2,4,6-Cycloheptatrien-1-one... Concentration Rank 7

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T. |
|-------|----------|--------|------------------|------|
| 10.41 | 11.55 ng | 713247 | Acenaphthene-d10 | 9.71 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 2,4,6-Cycloheptatrien-1-one, 2-p... | 182 | C13H10O | 014562-09-5 | 91   |
| 2    |    |   | 3-(2-Naphthyl)acrylaldehyde         | 182 | C13H10O | 020884-05-3 | 90   |
| 3    |    |   | [1,1'-Biphenyl]-4-carboxaldehyde    | 182 | C13H10O | 003218-36-8 | 87   |
| 4    |    |   | 9H-Fluoren-9-ol                     | 182 | C13H10O | 001689-64-1 | 83   |
| 5    |    |   | Dibenzofuran, 4-methyl-             | 182 | C13H10O | 007320-53-8 | 68   |



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF090917\  
Data File : BF098397.D  
Acq On : 9 Sep 2017 20:30  
Operator : SJ/JU  
Sample : I5151-03 50X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

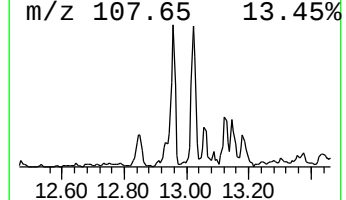
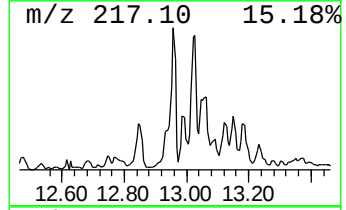
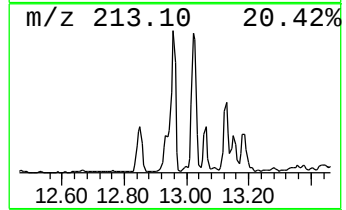
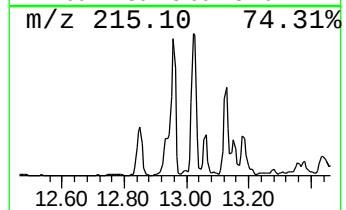
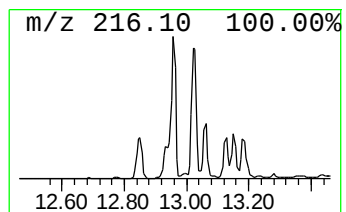
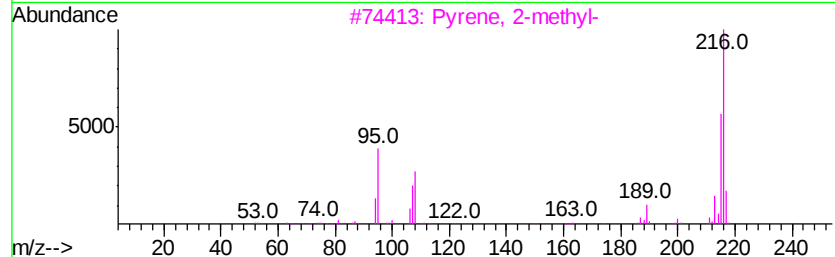
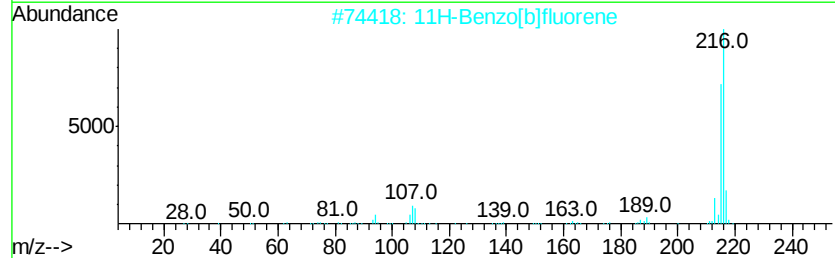
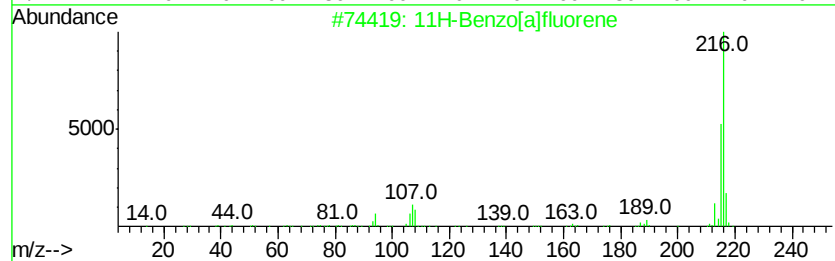
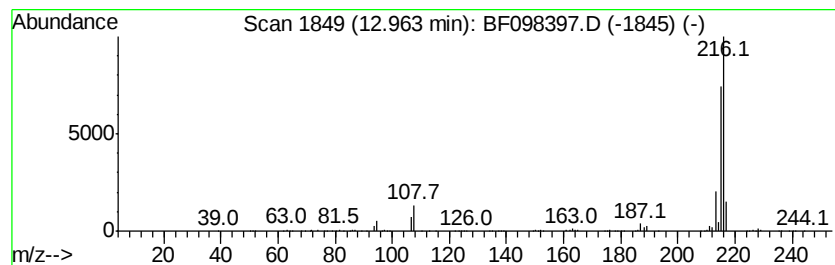
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ClientSampleId :  
NYSEG-PH4-ESMI2-3

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.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 12 11H-Benzo[a]fluorene Concentration Rank 20

| R.T.      | EstConc              | Area   | Relative to ISTD | R.T.        |      |
|-----------|----------------------|--------|------------------|-------------|------|
| 12.96     | 5.14 ng              | 892087 | Chrysene-d12     | 13.83       |      |
| Hit# of 5 | Tentative ID         | MW     | MolForm          | CAS#        | Qual |
| 1         | 11H-Benzo[a]fluorene | 216    | C17H12           | 000238-84-6 | 90   |
| 2         | 11H-Benzo[b]fluorene | 216    | C17H12           | 000243-17-4 | 90   |
| 3         | Pvrene, 2-methyl-    | 216    | C17H12           | 003442-78-2 | 76   |
| 4         | Pvrene, 1-methyl-    | 216    | C17H12           | 002381-21-7 | 74   |
| 5         | 7H-Benzo[c]fluorene  | 216    | C17H12           | 000205-12-9 | 74   |



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF090917\  
Data File : BF098397.D  
Acq On : 9 Sep 2017 20:30  
Operator : SJ/JU  
Sample : I5151-03 50X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

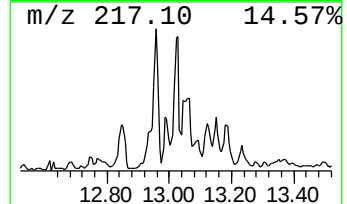
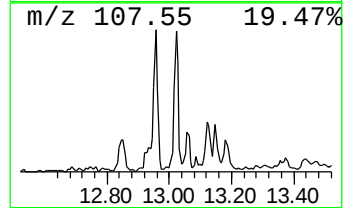
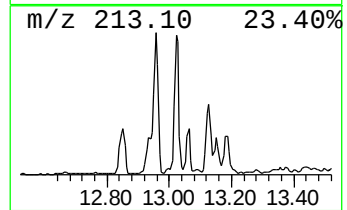
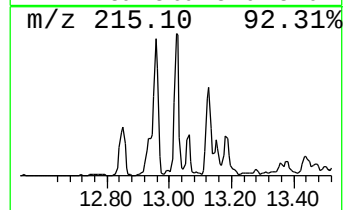
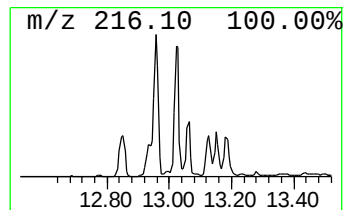
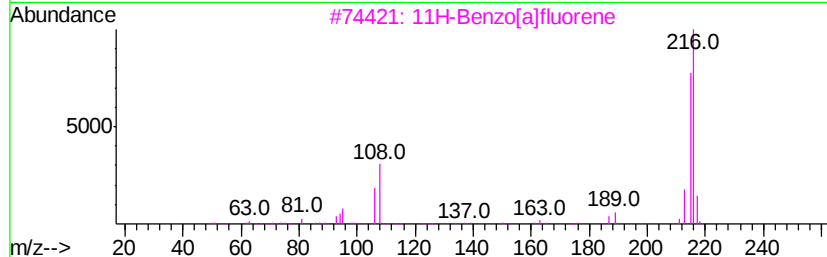
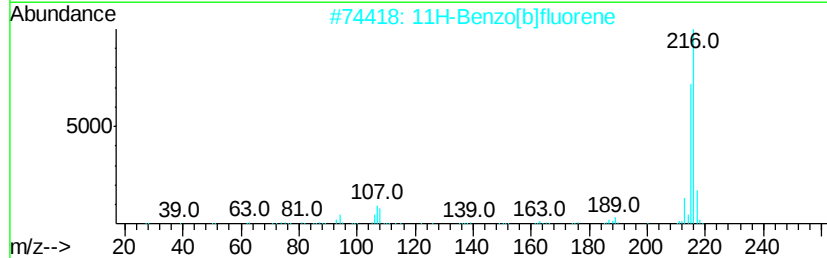
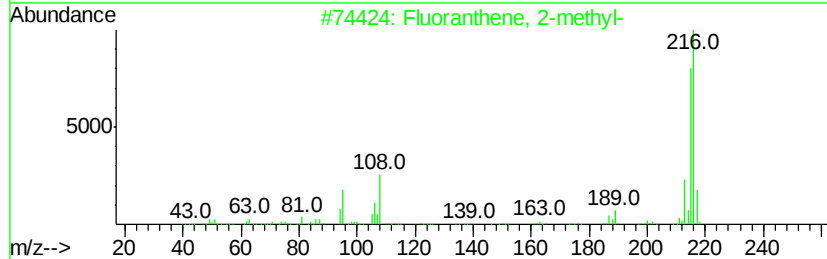
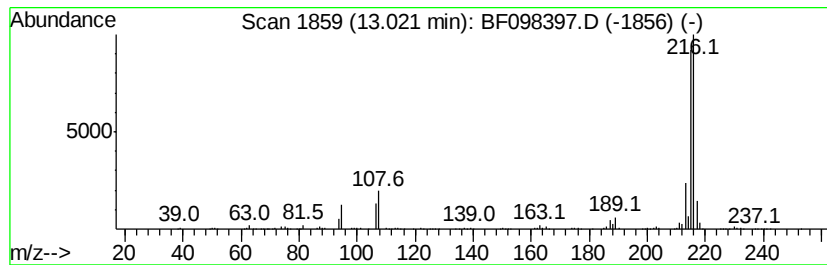
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ClientSampleId :  
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 13 Fluoranthene, 2-methyl- Concentration Rank 19

| R.T.      | EstConc                 | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------|--------|------------------|-------------|------|
| 13.02     | 5.19 ng                 | 899995 | Chrysene-d12     | 13.83       |      |
| Hit# of 5 | Tentative ID            | MW     | MolForm          | CAS#        | Qual |
| 1         | Fluoranthene, 2-methyl- | 216    | C17H12           | 033543-31-6 | 93   |
| 2         | 11H-Benzo[b]fluorene    | 216    | C17H12           | 000243-17-4 | 87   |
| 3         | 11H-Benzo[a]fluorene    | 216    | C17H12           | 000238-84-6 | 81   |
| 4         | Pyrene, 1-methyl-       | 216    | C17H12           | 002381-21-7 | 64   |
| 5         | Pyrene, 2-methyl-       | 216    | C17H12           | 003442-78-2 | 62   |



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF090917\  
Data File : BF098397.D  
Acq On : 9 Sep 2017 20:30  
Operator : SJ/JU  
Sample : I5151-03 50X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
NYSEG-PH4-ESMI2-3

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF081517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

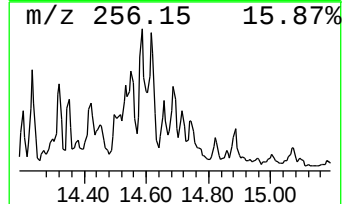
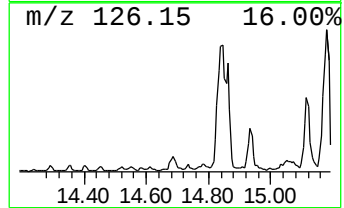
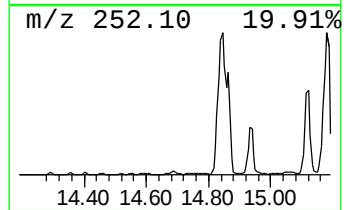
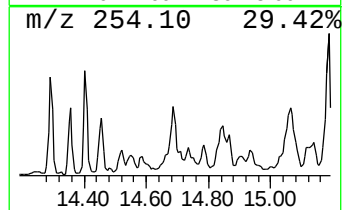
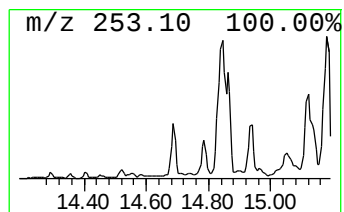
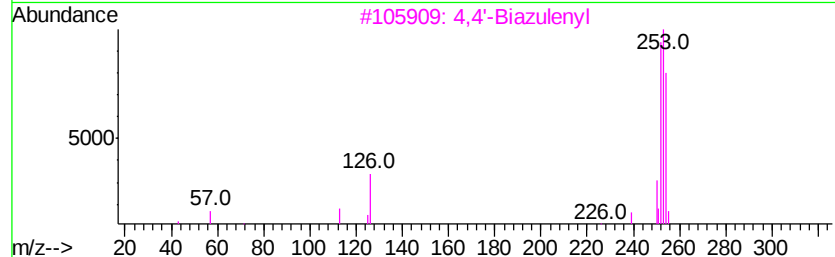
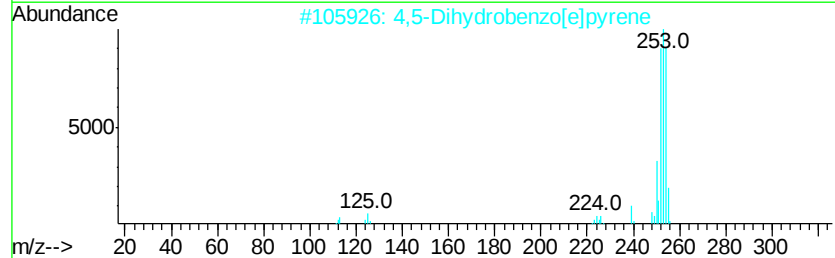
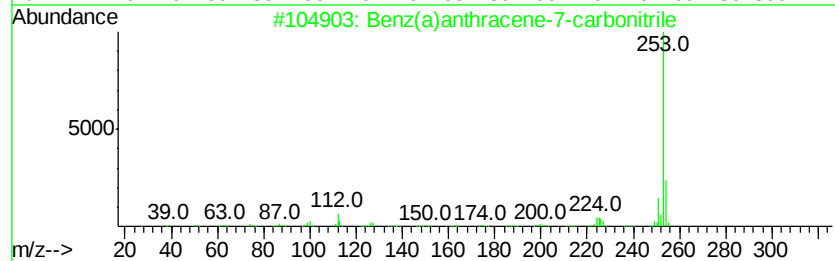
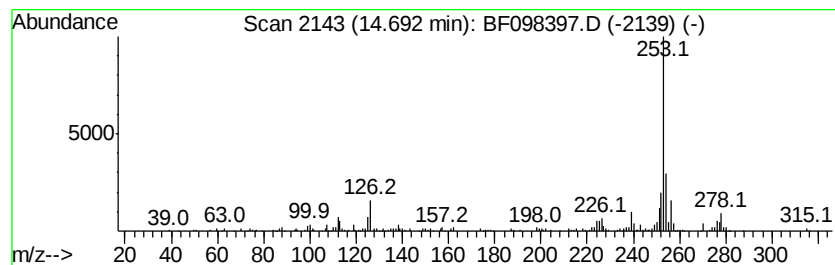
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TIC Integration Parameters: LSCINT.P

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Peak Number 14 Benz(a)anthracene-7-carboni... Concentration Rank 15

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 14.69 | 6.92 ng | 189127 | Perylene-d12     | 15.23 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-------------|--------------|------|
| 1    |    |   | Benz(a)anthracene-7-carbonitrile    | 253 | C19H11N     | 007476-08-6  | 83   |
| 2    |    |   | 4,5-Dihydrobenzo[el]pyrene          | 254 | C20H14      | 095676-42-9  | 50   |
| 3    |    |   | 4,4'-Biazulenvl                     | 254 | C20H14      | 094053-54-0  | 38   |
| 4    |    |   | 1,1'-Binaphthalene                  | 254 | C20H14      | 000604-53-5  | 38   |
| 5    |    |   | Propanephosphonic acid, bis(trim... | 268 | C9H25O3PSi2 | 1000193-07-4 | 28   |



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF090917\  
Data File : BF098397.D  
Acq On : 9 Sep 2017 20:30  
Operator : SJ/JU  
Sample : I5151-03 50X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

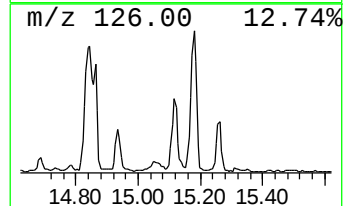
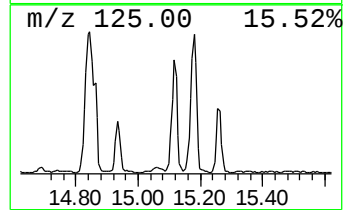
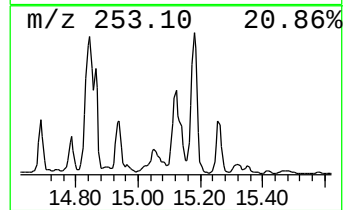
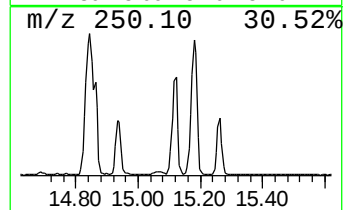
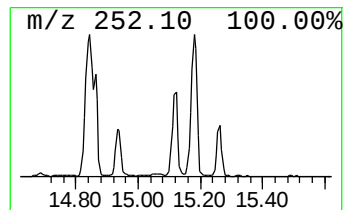
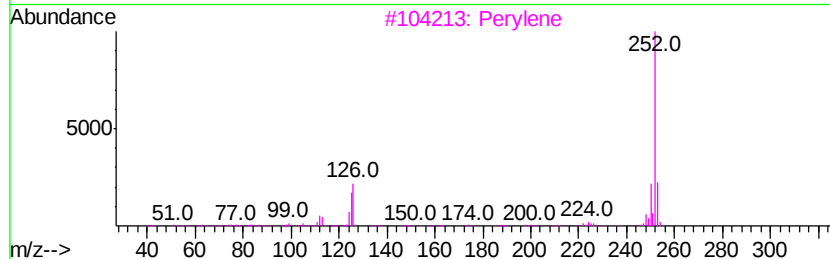
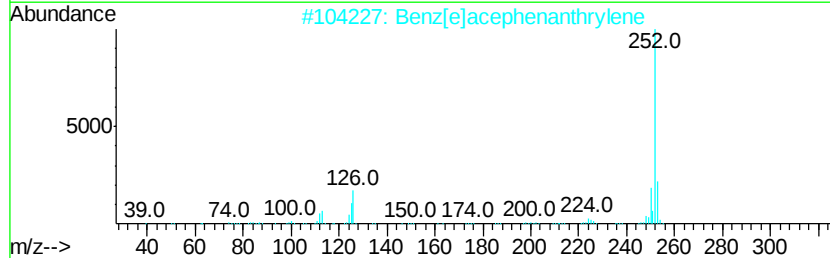
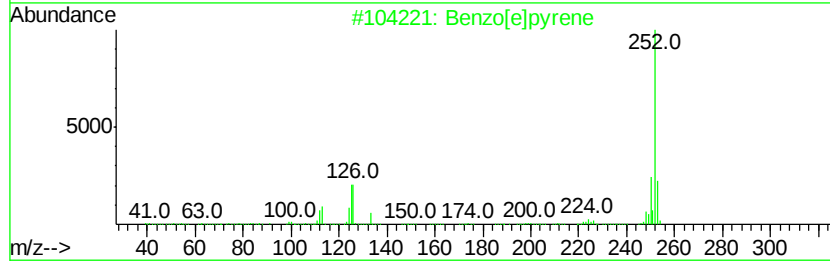
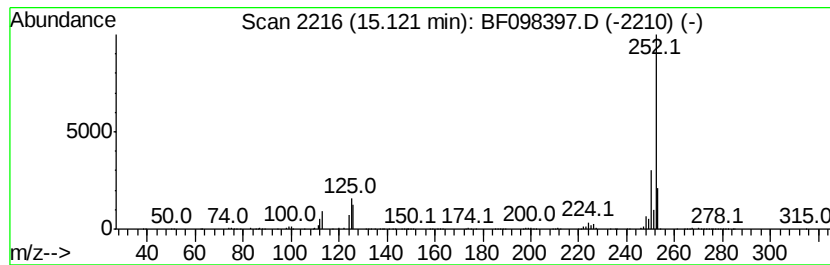
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ClientSampleId :  
NYSEG-PH4-ESMI2-3

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 16 Benzo[e]pyrene Concentration Rank 1

| R.T.    | EstConc  | Area                      | Relative to ISTD | R.T.           |
|---------|----------|---------------------------|------------------|----------------|
| 15.12   | 43.00 ng | 1176030                   | Perylene-d12     | 15.23          |
| Hit# of | 5        | Tentative ID              | MW MolForm       | CAS# Qual      |
| 1       |          | Benzo[e]pyrene            | 252 C20H12       | 000192-97-2 98 |
| 2       |          | Benzo[e]acephenanthrylene | 252 C20H12       | 000205-99-2 97 |
| 3       |          | Perylene                  | 252 C20H12       | 000198-55-0 95 |
| 4       |          | Benzo[a]pyrene            | 252 C20H12       | 000050-32-8 94 |
| 5       |          | Benzo[k]fluoranthene      | 252 C20H12       | 000207-08-9 93 |



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF090917\  
Data File : BF098397.D  
Acq On : 9 Sep 2017 20:30  
Operator : SJ/JU  
Sample : I5151-03 50X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
NYSEG-PH4-ESMI2-3

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF081517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

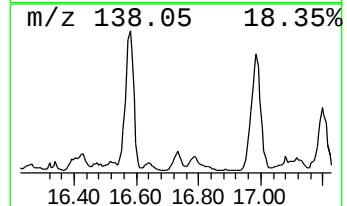
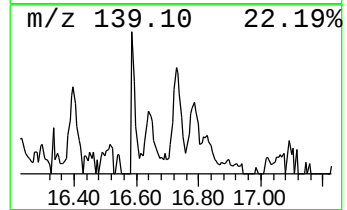
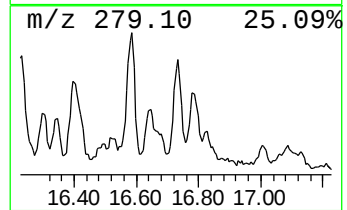
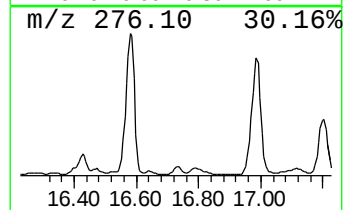
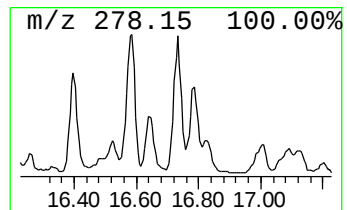
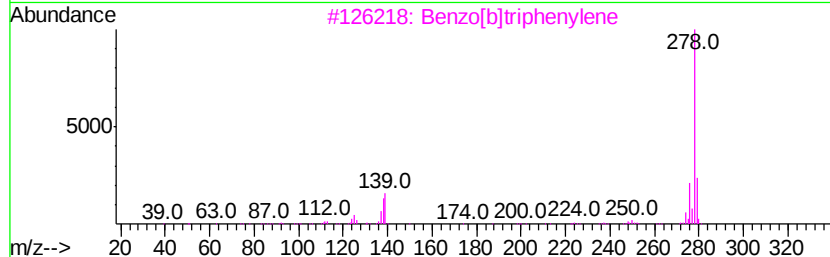
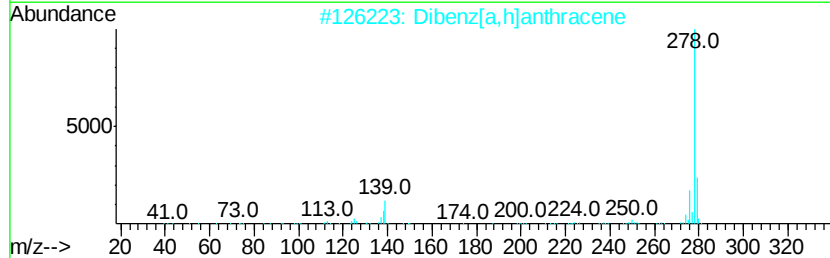
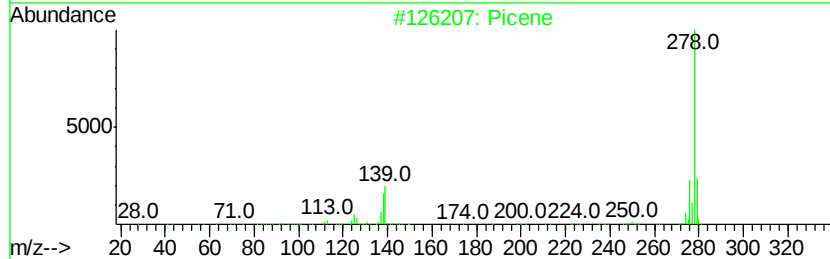
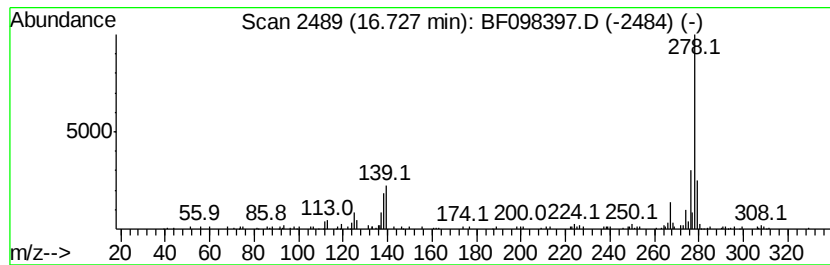
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 17 Picene Concentration Rank 10

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 16.73 | 8.58 ng | 234655 | Perylene-d12     | 15.23 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------|-----|---------|-------------|------|
| 1    |    |   | Picene                | 278 | C22H14  | 000213-46-7 | 97   |
| 2    |    |   | Dibenz[a,h]anthracene | 278 | C22H14  | 000053-70-3 | 95   |
| 3    |    |   | Benzo[b]triphenylene  | 278 | C22H14  | 000215-58-7 | 94   |
| 4    |    |   | Benzo[a]naphthacene   | 278 | C22H14  | 000226-88-0 | 93   |
| 5    |    |   | Pentacene             | 278 | C22H14  | 000135-48-8 | 89   |



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF090917\  
Data File : BF098397.D  
Acq On : 9 Sep 2017 20:30  
Operator : SJ/JU  
Sample : I5151-03 50X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
NYSEG-PH4-ESMI2-3

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

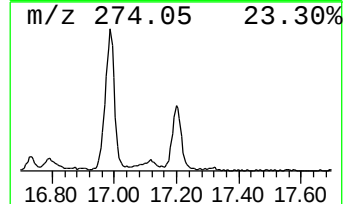
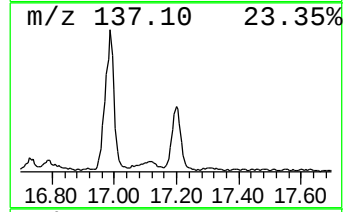
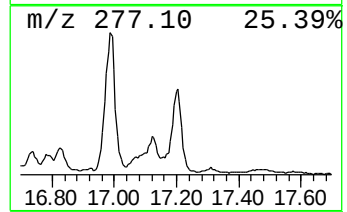
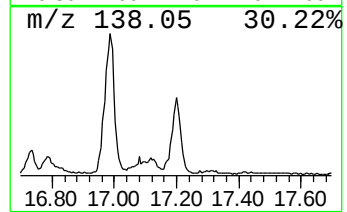
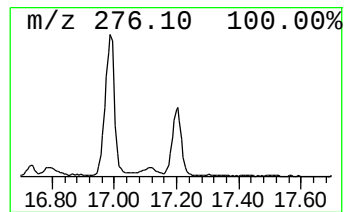
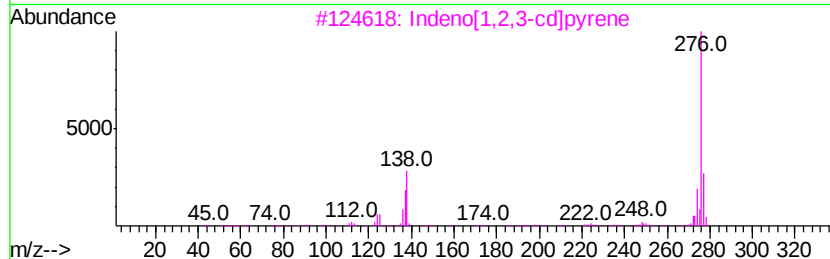
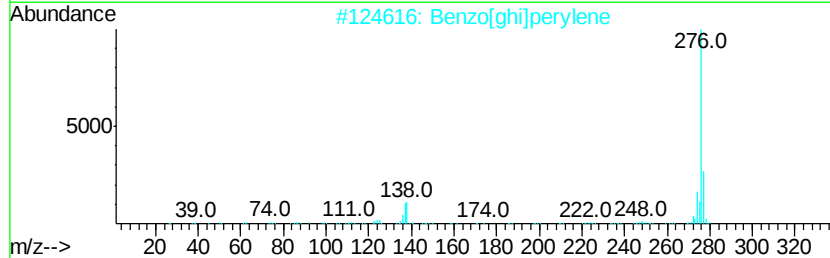
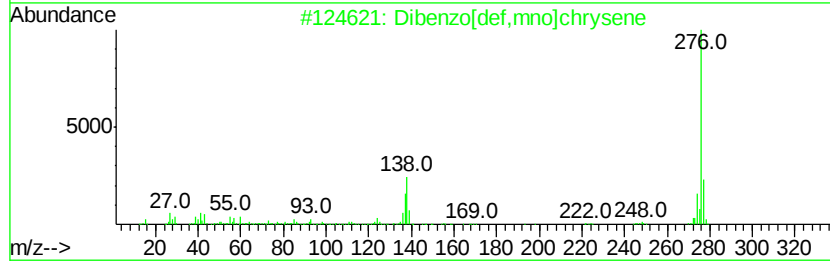
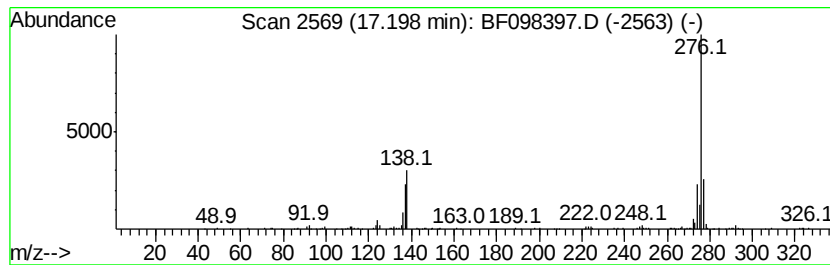
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TIC Integration Parameters: LSCINT.P

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Peak Number 18 Dibenzo[def,mno]chrysene Concentration Rank 5

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 17.20 | 13.96 ng | 381668 | Perylene-d12     | 15.23 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|------|----|-------------------------------------|-----|------------|-------------|------|
| 1    | 5  | Dibenzo[def,mno]chrysene            | 276 | C22H12     | 000191-26-4 | 98   |
| 2    |    | Benzo[ghi]perylene                  | 276 | C22H12     | 000191-24-2 | 96   |
| 3    |    | Indeno[1,2,3-cd]pyrene              | 276 | C22H12     | 000193-39-5 | 81   |
| 4    |    | Indeno[1,2,3-cd]fluoranthene        | 276 | C22H12     | 000193-43-1 | 74   |
| 5    |    | Naphthalen-1-yl-(3-nitro-benzyl)... | 276 | C17H12N2O2 | 200421-53-0 | 47   |



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF090917\  
Data File : BF098397.D  
Acq On : 9 Sep 2017 20:30  
Operator : SJ/JU  
Sample : I5151-03 50X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

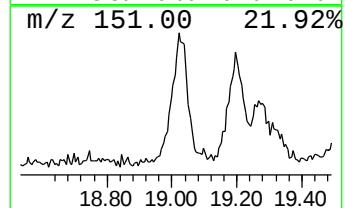
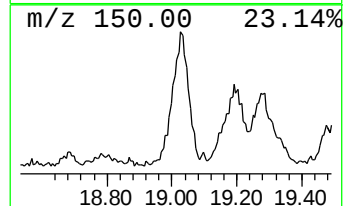
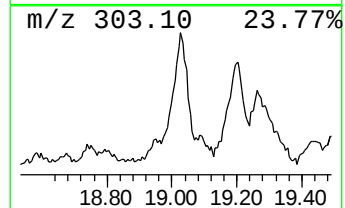
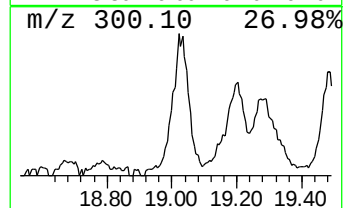
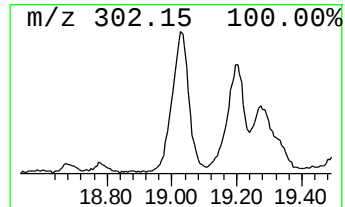
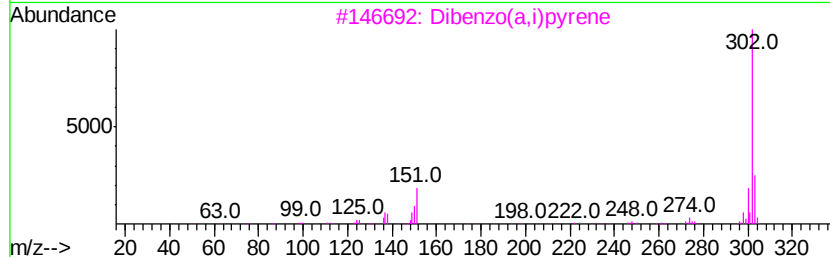
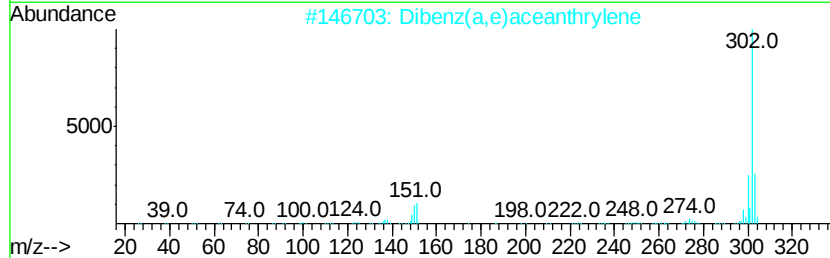
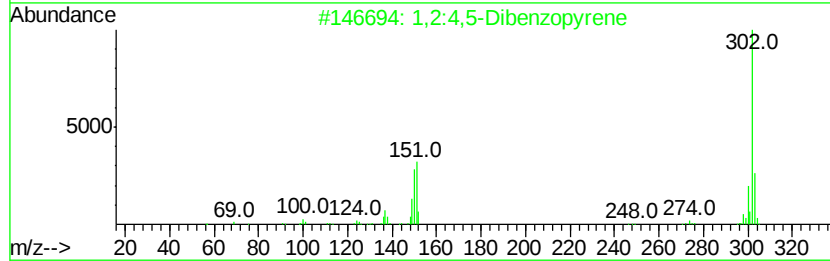
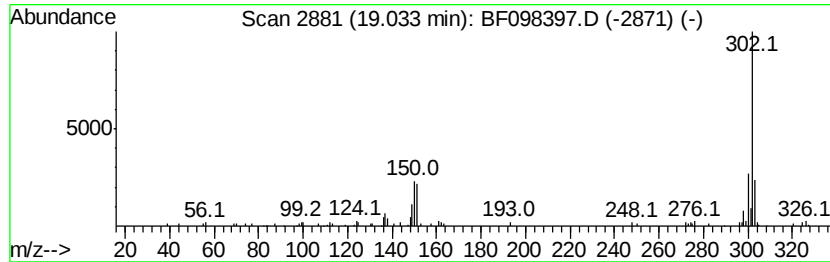
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ClientSampleId :  
NYSEG-PH4-ESMI2-3

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF081517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 19 1,2:4,5-Dibenzopyrene Concentration Rank 9

| R.T.      | EstConc                     | Area   | Relative to ISTD | R.T.        |      |
|-----------|-----------------------------|--------|------------------|-------------|------|
| 19.03     | 9.61 ng                     | 262907 | Perylene-d12     | 15.23       |      |
| Hit# of 5 | Tentative ID                | MW     | MolForm          | CAS#        | Qual |
| 1         | 1,2:4,5-Dibenzopvrene       | 302    | C24H14           | 000192-65-4 | 98   |
| 2         | Dibenz(a,e)aceanthrylene    | 302    | C24H14           | 005385-75-1 | 95   |
| 3         | Dibenzo(a,i)pvrene          | 302    | C24H14           | 000189-55-9 | 94   |
| 4         | 1,2,9,10-Dibenzopvrene      | 302    | C24H14           | 000191-30-0 | 93   |
| 5         | Indeno[1,2,3-fg]naphthacene | 302    | C24H14           | 000203-11-2 | 89   |



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF090917\  
Data File : BF098397.D  
Acq On : 9 Sep 2017 20:30  
Operator : SJ/JU  
Sample : I5151-03 50X  
Misc :  
ALS Vial : 22 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
NYSEG-PH4-ESMI2-3

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

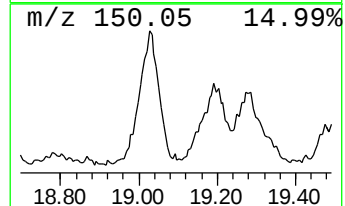
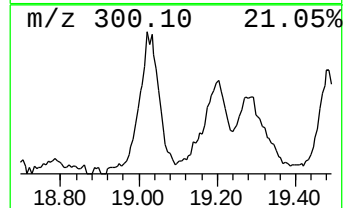
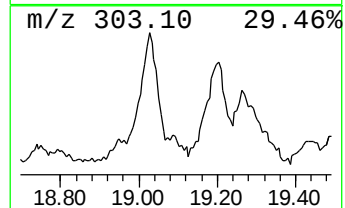
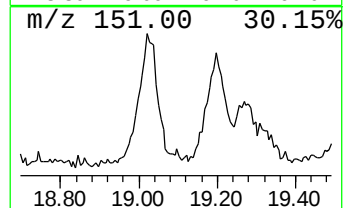
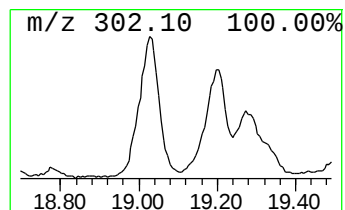
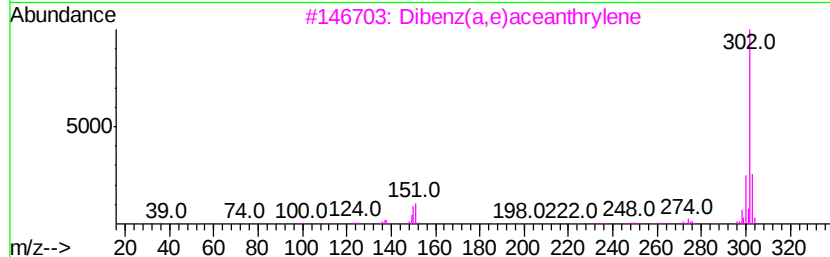
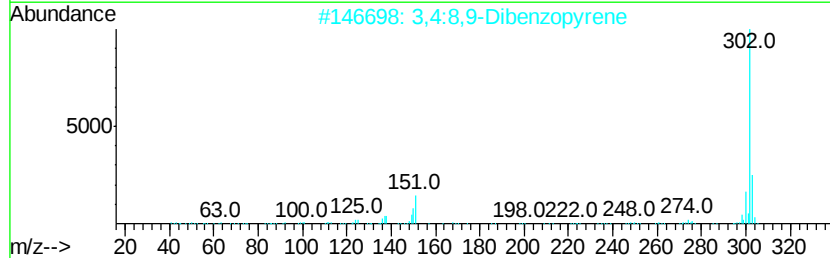
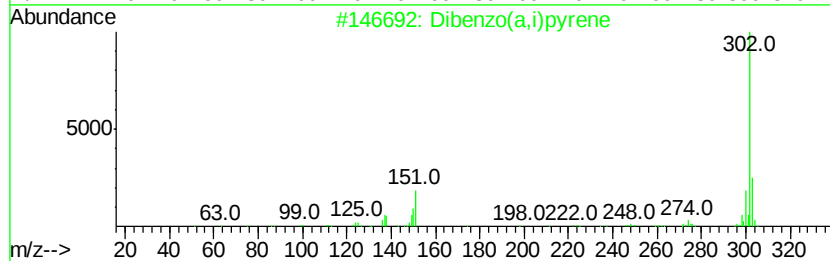
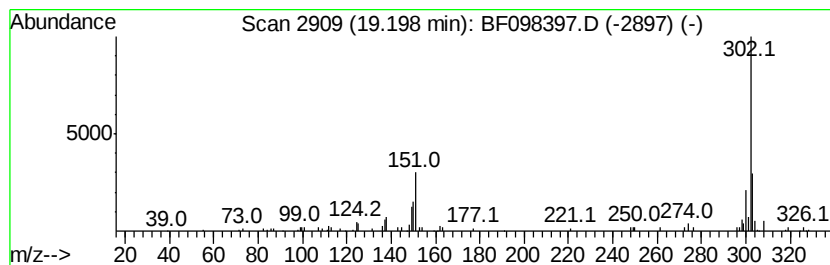
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 Dibenzo(a,i)pyrene Concentration Rank 12

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 19.20 | 7.73 ng | 211443 | Perylene-d12     | 15.23 |

| Hit# | of | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------|-----|---------|-------------|------|
| 1    | 5  | Dibenzo(a,i)pyrene       | 302 | C24H14  | 000189-55-9 | 98   |
| 2    |    | 3,4:8,9-Dibenzopyrene    | 302 | C24H14  | 000189-64-0 | 96   |
| 3    |    | Dibenz(a,e)aceanthrylene | 302 | C24H14  | 005385-75-1 | 96   |
| 4    |    | 1,2:4,5-Dibenzopyrene    | 302 | C24H14  | 000192-65-4 | 95   |
| 5    |    | 1,2,9,10-Dibenzopyrene   | 302 | C24H14  | 000191-30-0 | 93   |



Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA\_F\DATA\BF090917\  
 Data File : BF098397.D  
 Acq On : 9 Sep 2017 20:30  
 Operator : SJ/JU  
 Sample : I5151-03 50X  
 Misc :  
 ALS Vial : 22 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 NYSEG-PH4-ESMI2-3

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF081517.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 |
| Benzene, 1,2,3-tr... | 6.50  | 6.7     | ng    | 197471   | 1                     | 6.67  | 592412  | 20.0 |
| Indane               | 6.85  | 7.8     | ng    | 231814   | 1                     | 6.67  | 592412  | 20.0 |
| Naphthalene, 1-et... | 9.23  | 7.6     | ng    | 466800   | 3                     | 9.71  | 1234950 | 20.0 |
| Naphthalene, 2,6-... | 9.30  | 18.6    | ng    | 1146470  | 3                     | 9.71  | 1234950 | 20.0 |
| Naphthalene, 2,3-... | 9.37  | 18.3    | ng    | 1128600  | 3                     | 9.71  | 1234950 | 20.0 |
| Naphthalene, 1,6,... | 10.02 | 7.0     | ng    | 434053   | 3                     | 9.71  | 1234950 | 20.0 |
| 2,4,6-Cycloheptat... | 10.41 | 11.6    | ng    | 713247   | 3                     | 9.71  | 1234950 | 20.0 |
| 11H-Benzo[a]fluorene | 12.96 | 5.1     | ng    | 892087   | 5                     | 13.83 | 3468060 | 20.0 |
| Fluoranthene, 2-m... | 13.02 | 5.2     | ng    | 899995   | 5                     | 13.83 | 3468060 | 20.0 |
| Benz(a)anthracene... | 14.69 | 6.9     | ng    | 189127   | 6                     | 15.23 | 546966  | 20.0 |
| Benzo[e]pyrene       | 15.12 | 43.0    | ng    | 1176030  | 6                     | 15.23 | 546966  | 20.0 |
| Picene               | 16.73 | 8.6     | ng    | 234655   | 6                     | 15.23 | 546966  | 20.0 |
| Dibenzo[def,mno]c... | 17.20 | 14.0    | ng    | 381668   | 6                     | 15.23 | 546966  | 20.0 |
| 1,2:4,5-Dibenzopy... | 19.03 | 9.6     | ng    | 262907   | 6                     | 15.23 | 546966  | 20.0 |
| Dibenzo(a,i)pyrene   | 19.20 | 7.7     | ng    | 211443   | 6                     | 15.23 | 546966  | 20.0 |