

Data Path : Z:\HPCHEM1\BNA G\DATA\BG080515\  
 Data File : BG018288.D  
 Acq On : 6 Aug 2015 1:32  
 Operator : UM/TP  
 Sample : G3109-14RE  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 C01W2RE

## Integration Parameters: LSCINT.P

Integrator: RTE  
 Smoothing : OFF  
 Sampling : 1  
 Start Thrs: 0.2  
 Stop Thrs : 0

Filtering: 5  
 Min Area: 0.1 % 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\_G\METHODS\SOM02.2-EPA-BG080515.M  
 Title : SVOA 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         | 3.249       | 91            | 95          | 99           | rBV2     | 16540          | 19065         | 0.96%           | 0.102%        |
| 2         | 3.425       | 121           | 125         | 130          | rBV4     | 8308           | 9649          | 0.49%           | 0.051%        |
| 3         | 4.947       | 382           | 384         | 389          | rBV3     | 5898           | 7860          | 0.40%           | 0.042%        |
| 4         | 5.082       | 401           | 407         | 409          | rBV4     | 7185           | 10210         | 0.51%           | 0.054%        |
| 5         | 5.446       | 465           | 469         | 476          | rVB4     | 11470          | 18904         | 0.95%           | 0.101%        |
| 6         | 6.527       | 649           | 653         | 657          | rVV3     | 7443           | 10950         | 0.55%           | 0.058%        |
| 7         | 6.586       | 657           | 663         | 668          | rVB3     | 7142           | 12795         | 0.64%           | 0.068%        |
| 8         | 6.645       | 668           | 673         | 686          | rVB      | 148668         | 247512        | 12.47%          | 1.321%        |
| 9         | 6.780       | 690           | 696         | 701          | rVV      | 103884         | 153145        | 7.71%           | 0.817%        |
| 10        | 6.974       | 722           | 729         | 737          | rVB      | 155558         | 250159        | 12.60%          | 1.335%        |
| 11        | 7.086       | 740           | 748         | 756          | rVB2     | 18114          | 28631         | 1.44%           | 0.153%        |
| 12        | 7.438       | 800           | 808         | 817          | rBV      | 166511         | 268457        | 13.52%          | 1.432%        |
| 13        | 7.544       | 822           | 826         | 835          | rVB2     | 12287          | 21356         | 1.08%           | 0.114%        |
| 14        | 7.979       | 896           | 900         | 903          | rBV3     | 5335           | 8305          | 0.42%           | 0.044%        |
| 15        | 8.073       | 912           | 916         | 923          | rVB3     | 9099           | 15235         | 0.77%           | 0.081%        |
| 16        | 8.178       | 928           | 934         | 941          | rBV      | 168630         | 273412        | 13.77%          | 1.459%        |
| 17        | 8.261       | 941           | 948         | 955          | rVB5     | 7724           | 18108         | 0.91%           | 0.097%        |
| 18        | 8.537       | 989           | 995         | 999          | rBV3     | 12014          | 19655         | 0.99%           | 0.105%        |
| 19        | 8.596       | 999           | 1005        | 1013         | rVV      | 148788         | 245930        | 12.39%          | 1.312%        |
| 20        | 8.919       | 1055          | 1060        | 1063         | rBV3     | 9221           | 14024         | 0.71%           | 0.075%        |
| 21        | 9.013       | 1071          | 1076        | 1080         | rBV4     | 5883           | 8501          | 0.43%           | 0.045%        |
| 22        | 9.124       | 1091          | 1095        | 1099         | rBV5     | 6920           | 9831          | 0.50%           | 0.052%        |
| 23        | 9.312       | 1118          | 1127        | 1135         | rBV      | 153077         | 258323        | 13.01%          | 1.378%        |
| 24        | 9.630       | 1176          | 1181        | 1185         | rVB4     | 9654           | 13626         | 0.69%           | 0.073%        |
| 25        | 9.865       | 1215          | 1221        | 1229         | rVV      | 214131         | 350069        | 17.63%          | 1.868%        |
| 26        | 10.006      | 1241          | 1245        | 1251         | rVB2     | 7496           | 12998         | 0.65%           | 0.069%        |
| 27        | 10.223      | 1276          | 1282        | 1287         | rVV      | 252141         | 417233        | 21.01%          | 2.226%        |
| 28        | 10.264      | 1287          | 1289        | 1296         | rVB3     | 16777          | 28139         | 1.42%           | 0.150%        |
| 29        | 10.370      | 1300          | 1307        | 1317         | rVV      | 73744          | 132818        | 6.69%           | 0.709%        |
| 30        | 10.975      | 1406          | 1410        | 1415         | rBV4     | 5577           | 9031          | 0.45%           | 0.048%        |
| 31        | 11.063      | 1423          | 1425        | 1431         | rVV5     | 5208           | 9056          | 0.46%           | 0.048%        |
| 32        | 11.145      | 1434          | 1439        | 1445         | rVB4     | 5844           | 12286         | 0.62%           | 0.066%        |
| 33        | 11.604      | 1511          | 1517        | 1519         | rBV4     | 6552           | 12413         | 0.63%           | 0.066%        |
| 34        | 11.903      | 1563          | 1568        | 1575         | rVB2     | 19151          | 34061         | 1.72%           | 0.182%        |

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 Max Peaks: 100  
 Peak Location: TOP

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

Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG080515.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |        |        |        |        |
|----|--------|------|------|------|------|--------|--------|--------|--------|
| 35 | 12.127 | 1602 | 1606 | 1614 | rVB2 | 12493  | 21944  | 1.11%  | 0.117% |
| 36 | 12.644 | 1691 | 1694 | 1697 | rVV5 | 8560   | 9938   | 0.50%  | 0.053% |
| 37 | 12.685 | 1697 | 1701 | 1704 | rVV5 | 12628  | 22064  | 1.11%  | 0.118% |
| 38 | 12.714 | 1704 | 1706 | 1709 | rVB4 | 8857   | 8671   | 0.44%  | 0.046% |
| 39 | 12.826 | 1721 | 1725 | 1727 | rBV5 | 4881   | 7408   | 0.37%  | 0.040% |
| 40 | 12.873 | 1729 | 1733 | 1737 | rVV7 | 10154  | 15849  | 0.80%  | 0.085% |
| 41 | 12.943 | 1741 | 1745 | 1751 | rVB8 | 10227  | 17894  | 0.90%  | 0.095% |
| 42 | 13.084 | 1766 | 1769 | 1772 | rBV5 | 4191   | 7492   | 0.38%  | 0.040% |
| 43 | 13.296 | 1803 | 1805 | 1808 | rVB4 | 10659  | 9791   | 0.49%  | 0.052% |
| 44 | 13.431 | 1824 | 1828 | 1833 | rBV7 | 20452  | 34311  | 1.73%  | 0.183% |
| 45 | 13.490 | 1835 | 1838 | 1840 | rVV3 | 8226   | 9646   | 0.49%  | 0.051% |
| 46 | 13.537 | 1840 | 1846 | 1850 | rVV  | 374147 | 541539 | 27.27% | 2.889% |
| 47 | 13.584 | 1850 | 1854 | 1864 | rVV  | 167037 | 247142 | 12.45% | 1.319% |
| 48 | 13.666 | 1864 | 1868 | 1871 | rVB6 | 10805  | 17085  | 0.86%  | 0.091% |
| 49 | 13.807 | 1886 | 1892 | 1903 | rVV  | 328492 | 518284 | 26.10% | 2.765% |
| 50 | 13.936 | 1911 | 1914 | 1920 | rVB6 | 27426  | 42681  | 2.15%  | 0.228% |
| 51 | 14.071 | 1935 | 1937 | 1940 | rBV3 | 8054   | 9832   | 0.50%  | 0.052% |
| 52 | 14.118 | 1940 | 1945 | 1952 | rVV2 | 360941 | 562105 | 28.31% | 2.999% |
| 53 | 14.189 | 1952 | 1957 | 1964 | rVB4 | 34094  | 57012  | 2.87%  | 0.304% |
| 54 | 14.389 | 1986 | 1991 | 1996 | rVB2 | 65221  | 100826 | 5.08%  | 0.538% |
| 55 | 14.524 | 2007 | 2014 | 2018 | rBV3 | 26344  | 58437  | 2.94%  | 0.312% |
| 56 | 14.659 | 2032 | 2037 | 2040 | rBV6 | 11376  | 19486  | 0.98%  | 0.104% |
| 57 | 14.747 | 2050 | 2052 | 2055 | rBV3 | 8507   | 10608  | 0.53%  | 0.057% |
| 58 | 14.835 | 2065 | 2067 | 2068 | rBV2 | 10308  | 7454   | 0.38%  | 0.040% |
| 59 | 14.912 | 2075 | 2080 | 2086 | rBV7 | 33655  | 63114  | 3.18%  | 0.337% |
| 60 | 14.994 | 2091 | 2094 | 2096 | rBV4 | 20576  | 22757  | 1.15%  | 0.121% |
| 61 | 15.123 | 2111 | 2116 | 2121 | rVB  | 376194 | 534912 | 26.94% | 2.854% |
| 62 | 15.182 | 2121 | 2126 | 2129 | rBV2 | 33227  | 51315  | 2.58%  | 0.274% |
| 63 | 15.276 | 2138 | 2142 | 2145 | rBV  | 42526  | 51285  | 2.58%  | 0.274% |
| 64 | 15.358 | 2151 | 2156 | 2157 | rBV5 | 37012  | 48730  | 2.45%  | 0.260% |
| 65 | 15.863 | 2239 | 2242 | 2243 | rBV3 | 30790  | 32504  | 1.64%  | 0.173% |
| 66 | 16.404 | 2330 | 2334 | 2338 | rBV5 | 52625  | 76789  | 3.87%  | 0.410% |
| 67 | 16.657 | 2372 | 2377 | 2381 | rBV2 | 310735 | 370160 | 18.64% | 1.975% |
| 68 | 16.762 | 2391 | 2395 | 2399 | rBV2 | 132236 | 148245 | 7.47%  | 0.791% |
| 69 | 16.880 | 2411 | 2415 | 2419 | rBV  | 317899 | 456644 | 23.00% | 2.436% |
| 70 | 16.927 | 2419 | 2423 | 2427 | rVB  | 278723 | 388045 | 19.54% | 2.070% |
| 71 | 16.980 | 2427 | 2432 | 2436 | rBV2 | 308891 | 439562 | 22.14% | 2.345% |

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ALS Vial : 17 Sample Multiplier: 1

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ClientSampleId :  
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Peak Location: TOP

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Peak separation: 5

Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG080515.M  
Title : SVOA CALIBRATION

|     |        |      |      |      |      |         |         |         |         |
|-----|--------|------|------|------|------|---------|---------|---------|---------|
| 72  | 17.491 | 2514 | 2519 | 2523 | rVB3 | 95390   | 170685  | 8.60%   | 0.911%  |
| 73  | 17.802 | 2569 | 2572 | 2578 | rVB3 | 110389  | 163427  | 8.23%   | 0.872%  |
| 74  | 17.867 | 2578 | 2583 | 2587 | rBV  | 407163  | 520962  | 26.24%  | 2.780%  |
| 75  | 17.955 | 2594 | 2598 | 2603 | rBV2 | 238562  | 333061  | 16.77%  | 1.777%  |
| 76  | 18.020 | 2606 | 2609 | 2613 | rVB3 | 101559  | 122962  | 6.19%   | 0.656%  |
| 77  | 18.061 | 2613 | 2616 | 2619 | rBV5 | 61265   | 79885   | 4.02%   | 0.426%  |
| 78  | 18.249 | 2644 | 2648 | 2654 | rBV5 | 109157  | 183908  | 9.26%   | 0.981%  |
| 79  | 18.425 | 2675 | 2678 | 2682 | rBV6 | 48454   | 68596   | 3.45%   | 0.366%  |
| 80  | 18.795 | 2738 | 2741 | 2742 | rBV2 | 99015   | 103591  | 5.22%   | 0.553%  |
| 81  | 18.819 | 2742 | 2745 | 2755 | rVB4 | 205327  | 344960  | 17.37%  | 1.841%  |
| 82  | 18.960 | 2764 | 2769 | 2773 | rBV  | 543547  | 692665  | 34.88%  | 3.696%  |
| 83  | 19.107 | 2790 | 2794 | 2799 | rVB2 | 289871  | 404442  | 20.37%  | 2.158%  |
| 84  | 19.171 | 2802 | 2805 | 2809 | rVV4 | 96614   | 108980  | 5.49%   | 0.581%  |
| 85  | 19.295 | 2822 | 2826 | 2829 | rBV2 | 310557  | 465768  | 23.46%  | 2.485%  |
| 86  | 19.324 | 2829 | 2831 | 2835 | rVB  | 373569  | 419407  | 21.12%  | 2.238%  |
| 87  | 19.448 | 2849 | 2852 | 2856 | rVB4 | 101610  | 120072  | 6.05%   | 0.641%  |
| 88  | 19.553 | 2867 | 2870 | 2877 | rVB2 | 239432  | 341179  | 17.18%  | 1.820%  |
| 89  | 19.818 | 2912 | 2915 | 2921 | rVB3 | 214538  | 292147  | 14.71%  | 1.559%  |
| 90  | 19.876 | 2922 | 2925 | 2929 | rVB  | 175755  | 201248  | 10.14%  | 1.074%  |
| 91  | 20.100 | 2960 | 2963 | 2966 | rBV  | 173884  | 217390  | 10.95%  | 1.160%  |
| 92  | 20.241 | 2984 | 2987 | 2993 | rVB  | 256917  | 295236  | 14.87%  | 1.575%  |
| 93  | 20.417 | 3015 | 3017 | 3021 | rVB4 | 128105  | 138732  | 6.99%   | 0.740%  |
| 94  | 20.728 | 3067 | 3070 | 3075 | rBV2 | 102894  | 154575  | 7.78%   | 0.825%  |
| 95  | 21.022 | 3116 | 3120 | 3127 | rBV  | 1756690 | 1985649 | 100.00% | 10.595% |
| 96  | 21.099 | 3127 | 3133 | 3136 | rBV2 | 465572  | 921616  | 46.41%  | 4.917%  |
| 97  | 21.134 | 3136 | 3139 | 3149 | rVB  | 341745  | 461192  | 23.23%  | 2.461%  |
| 98  | 21.739 | 3239 | 3242 | 3247 | rBV3 | 260511  | 309904  | 15.61%  | 1.654%  |
| 99  | 21.886 | 3264 | 3267 | 3271 | rBV  | 427716  | 516075  | 25.99%  | 2.754%  |
| 100 | 22.626 | 3389 | 3393 | 3398 | rBV  | 320952  | 600619  | 30.25%  | 3.205%  |

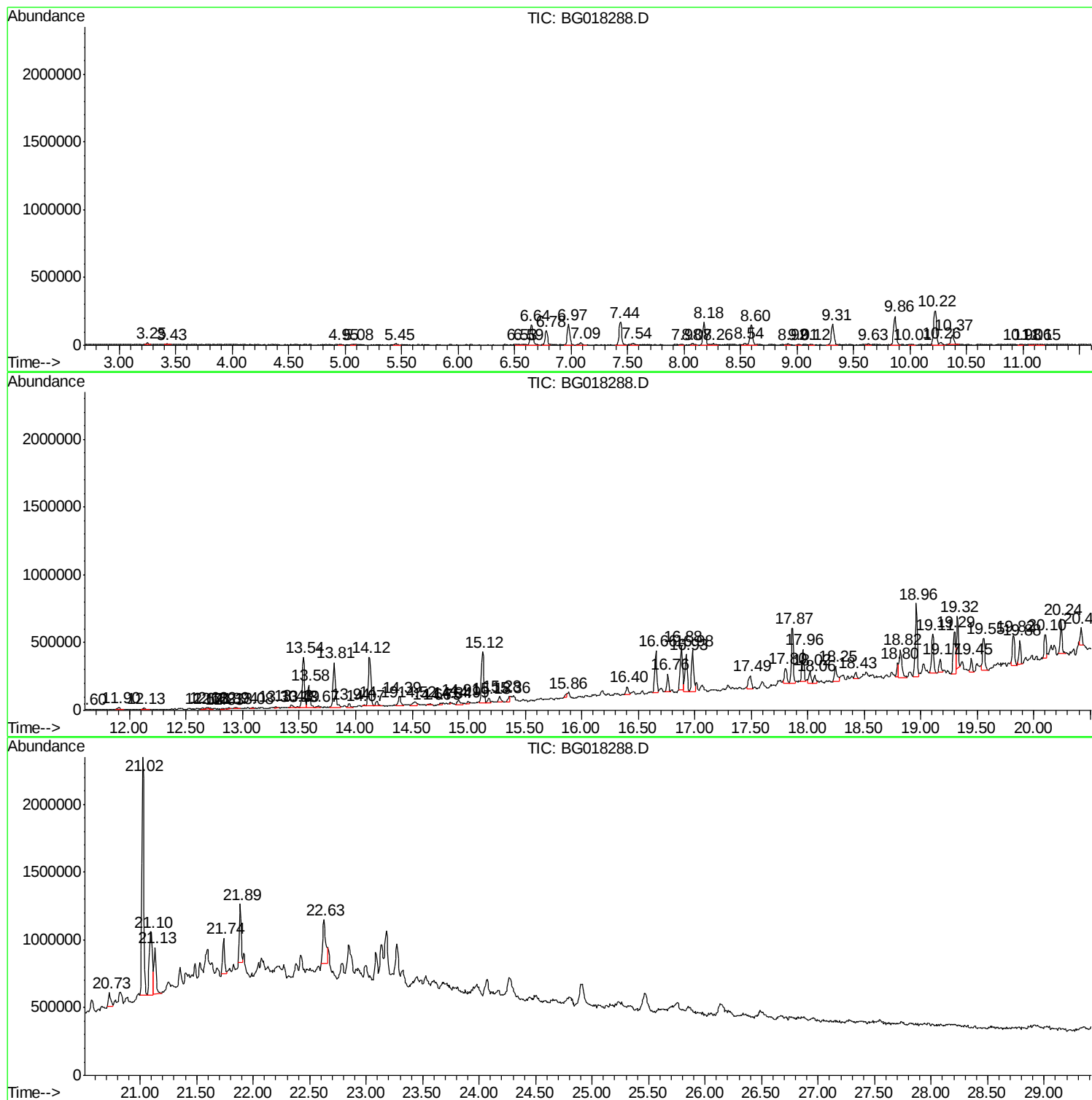
Sum of corrected areas: 18742241

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Instrument :  
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Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P



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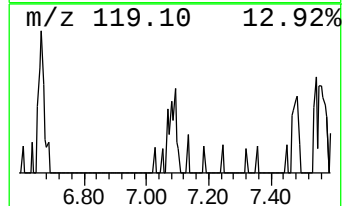
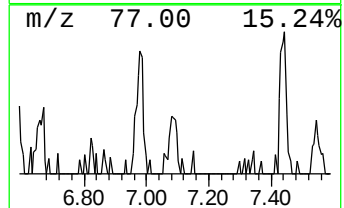
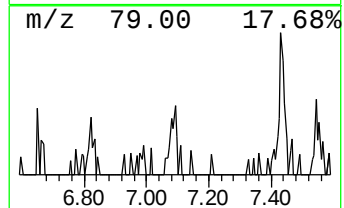
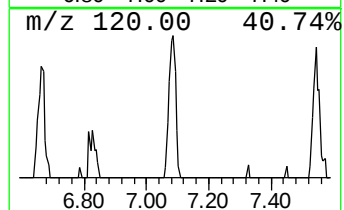
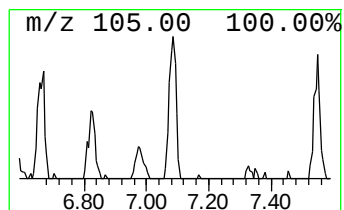
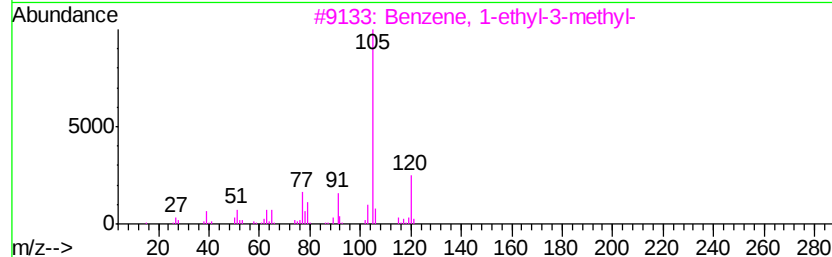
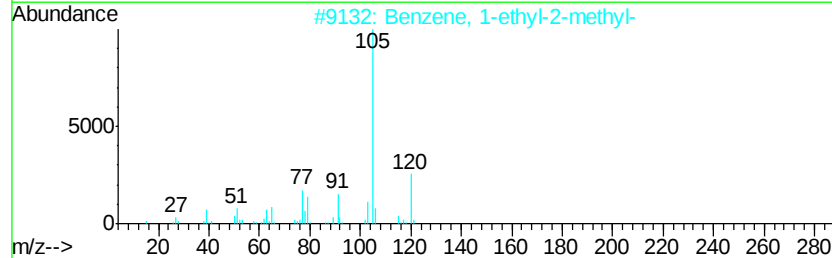
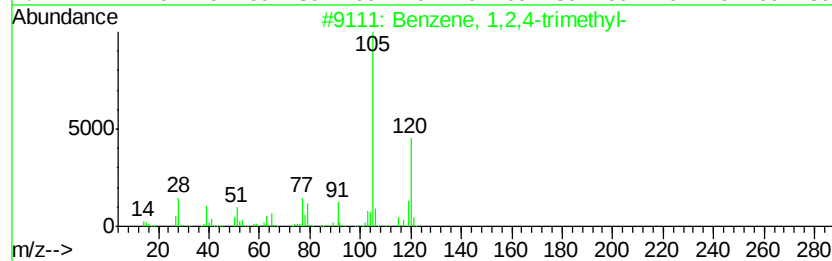
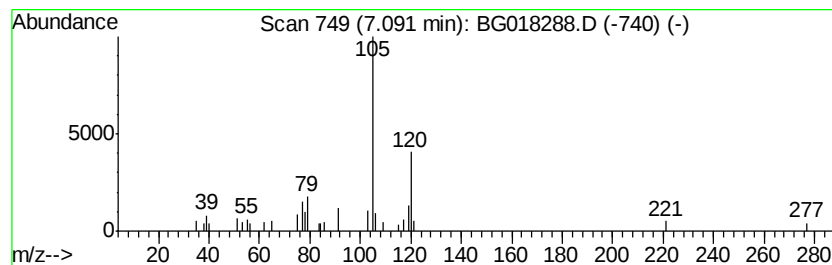
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 1 Benzene, 1,2,4-trimethyl- Concentration Rank 24

| R.T. | EstConc    | Area  | Relative to ISTD       | R.T. |
|------|------------|-------|------------------------|------|
| 7.09 | 2.13 ng/ul | 28631 | 1,4-Dichlorobenzene-d4 | 7.44 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2,4-trimethyl-  | 120 | C9H12   | 000095-63-6 | 81   |
| 2    |    | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 70   |
| 3    |    | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 70   |
| 4    |    | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 70   |
| 5    |    | Benzene, (1-methylethyl)-  | 120 | C9H12   | 000098-82-8 | 59   |



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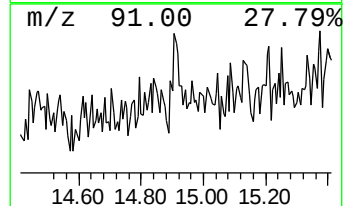
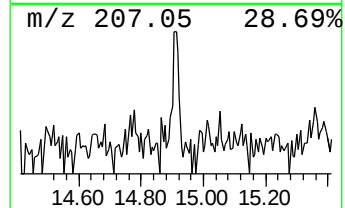
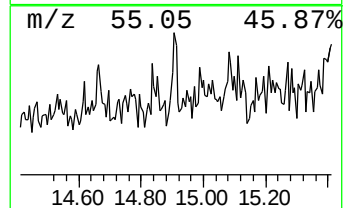
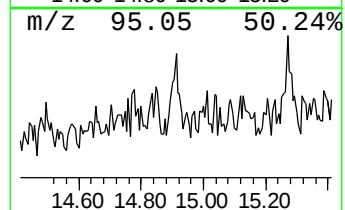
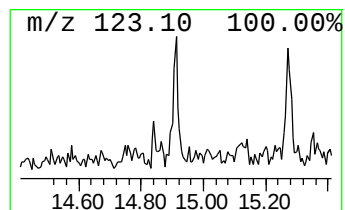
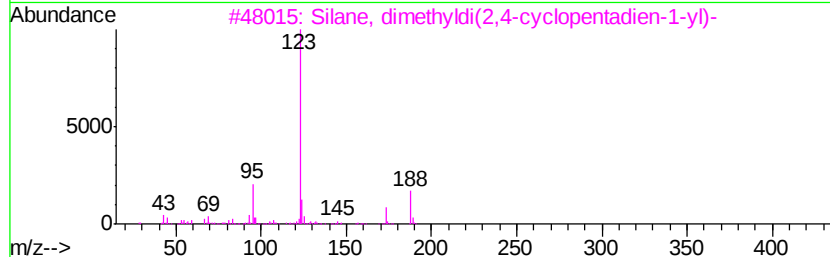
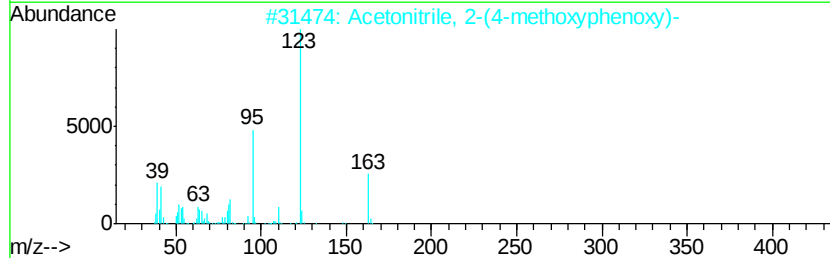
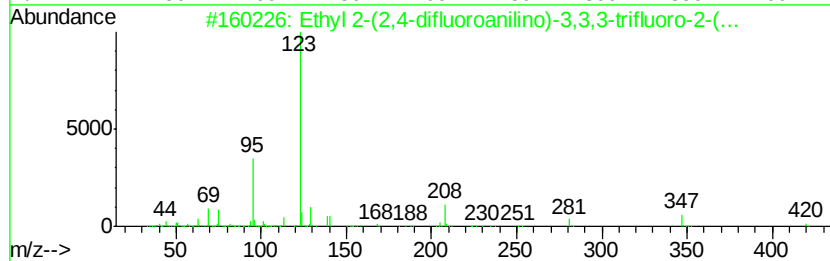
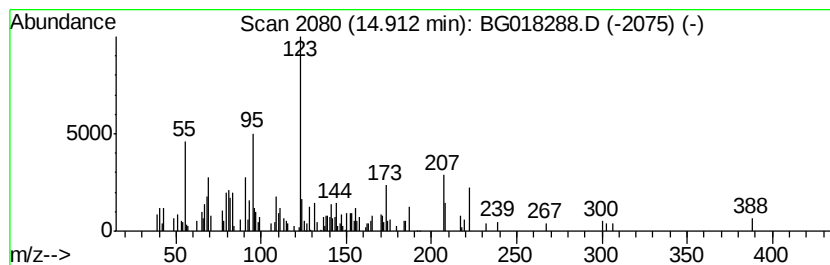
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Quant Title : SVOA CALIBRATION

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\*\*\*\*\*  
Peak Number 2 unknown-01 Concentration Rank 23

| R.T.    | EstConc    | Area                                | Relative to ISTD | R.T.            |
|---------|------------|-------------------------------------|------------------|-----------------|
| 14.91   | 2.25 ng/ul | 63114                               | Acenaphthene-d10 | 14.12           |
| Hit# of | 5          | Tentative ID                        | MW MolForm       | CAS# Qual       |
| 1       |            | Ethyl 2-(2,4-difluoroanilino)-3,... | 420 C18H14F6N2O3 | 339353-07-0 47  |
| 2       |            | Acetonitrile, 2-(4-methoxyphenoxy)- | 163 C9H9NO2      | 022446-12-4 46  |
| 3       |            | Silane, dimethyldi(2,4-cyclopent... | 188 C12H16Si     | 018053-74-2 43  |
| 4       |            | 2-(2-Hydroxycyclohexyl)-furan       | 166 C10H14O2     | 1000145-92-8 43 |
| 5       |            | 3-Fluorobenzoic acid, 2-ethylcyc... | 250 C15H19FO2    | 1000279-04-9 43 |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG080515\  
Data File : BG018288.D  
Acq On : 6 Aug 2015 1:32  
Operator : UM/TP  
Sample : G3109-14RE  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

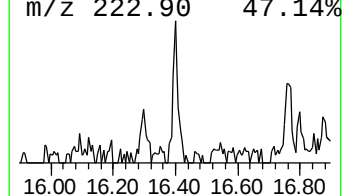
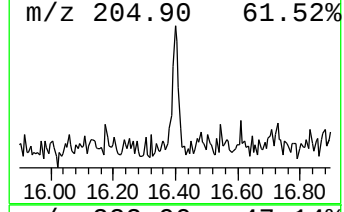
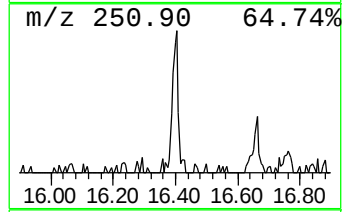
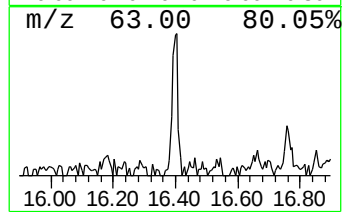
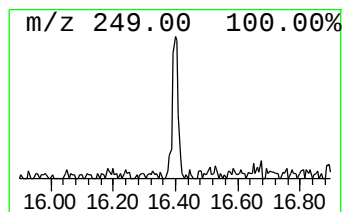
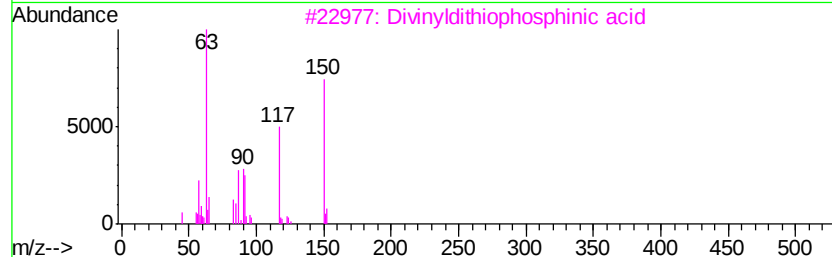
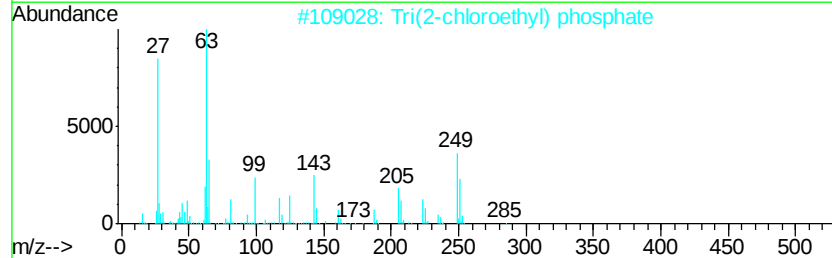
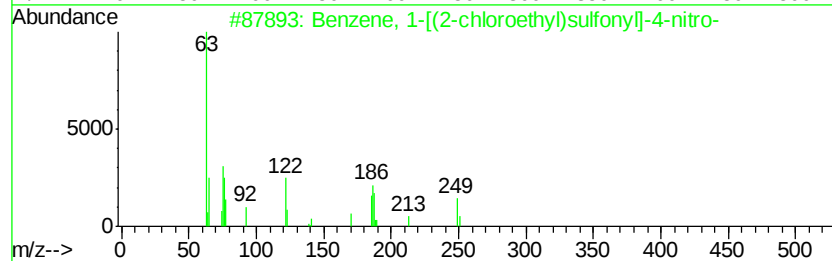
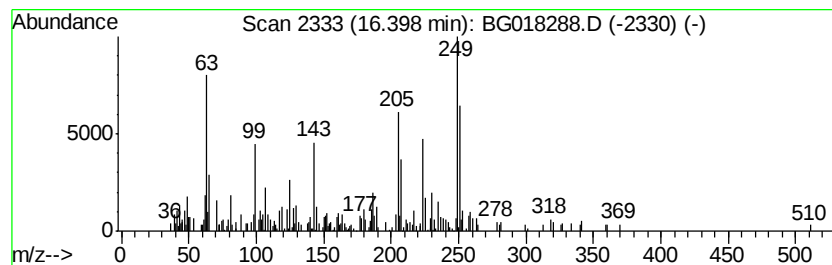
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BNA\_G  
ClientSampleId :  
C01W2RE

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 unknown-02 Concentration Rank 17

| R.T.    | EstConc    | Area                                | Relative to ISTD | R.T.        |             |      |
|---------|------------|-------------------------------------|------------------|-------------|-------------|------|
| 16.40   | 3.36 ng/ul | 76789                               | Phenanthrene-d10 | 16.88       |             |      |
| Hit# of | 5          | Tentative ID                        | MW               | MolForm     | CAS#        | Qual |
| 1       |            | Benzene, 1-[(2-chloroethyl)sulfo... | 249              | C8H8ClN04S  | 006461-63-8 | 47   |
| 2       |            | Tri(2-chloroethyl) phosphate        | 284              | C6H12Cl304P | 000115-96-8 | 42   |
| 3       |            | Divinylidithiophosphinic acid       | 150              | C4H7PS2     | 085417-00-1 | 38   |
| 4       |            | Propanoic acid, 2-chloro-, ethyl... | 136              | C5H9ClO2    | 000535-13-7 | 38   |
| 5       |            | Propane, 1,1,2-trichloro-           | 146              | C3H5Cl3     | 000598-77-6 | 38   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG080515\  
Data File : BG018288.D  
Acq On : 6 Aug 2015 1:32  
Operator : UM/TP  
Sample : G3109-14RE  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01W2RE

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
Quant Title : SVOA CALIBRATION

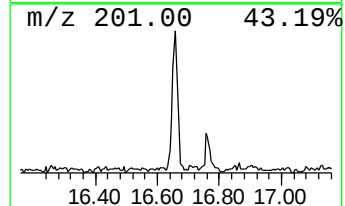
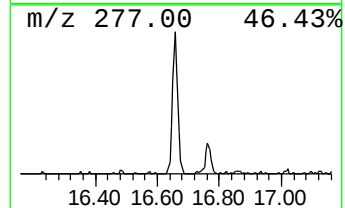
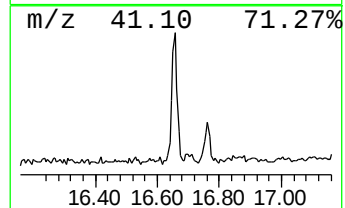
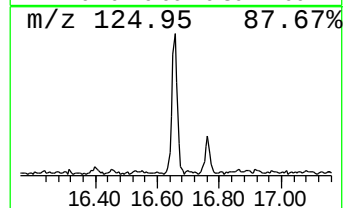
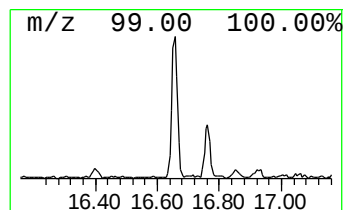
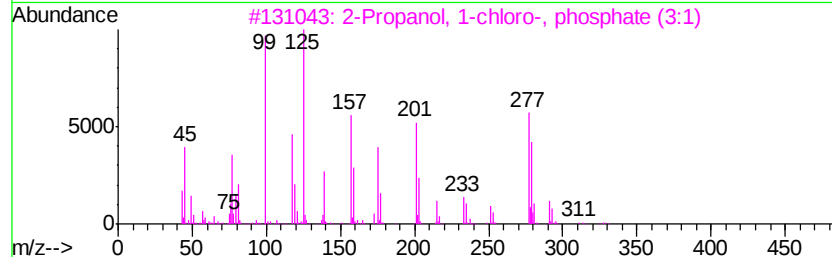
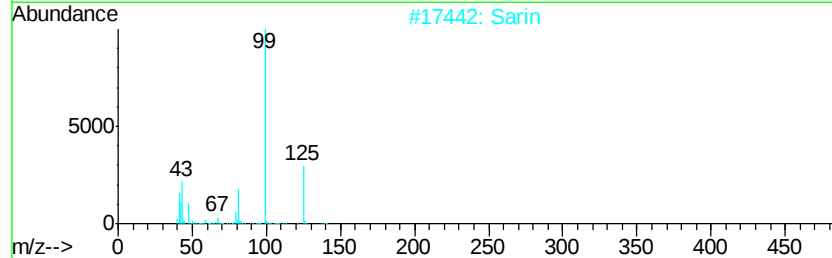
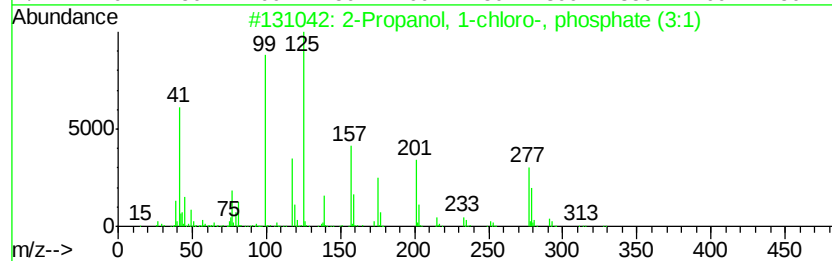
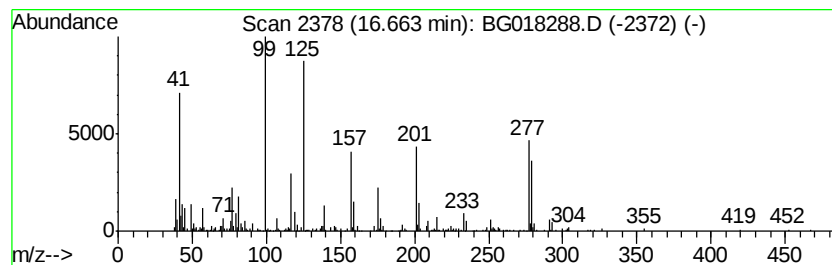
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 4 2-Propanol, 1-chloro-, phos... Concentration Rank 1

| R.T.  | EstConc     | Area   | Relative to ISTD | R.T.  |
|-------|-------------|--------|------------------|-------|
| 16.66 | 16.21 ng/ul | 370160 | Phenanthrene-d10 | 16.88 |

| Hit# | of | Tentative ID                        | MW  | MolForm     | CAS#        | Qual |
|------|----|-------------------------------------|-----|-------------|-------------|------|
| 1    | 5  | 2-Propanol, 1-chloro-, phosphate... | 326 | C9H18Cl3O4P | 013674-84-5 | 93   |
| 2    |    | Sarin                               | 140 | C4H10F02P   | 000107-44-8 | 35   |
| 3    |    | 2-Propanol, 1-chloro-, phosphate... | 326 | C9H18Cl3O4P | 013674-84-5 | 18   |
| 4    |    | Bis(3-chloro-1-propyl)(1-chloro-... | 326 | C9H18Cl3O4P | 137888-35-8 | 16   |
| 5    |    | Bis(1-chloro-2-propyl)(3-chloro-... | 326 | C9H18Cl3O4P | 137909-40-1 | 14   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG080515\  
Data File : BG018288.D  
Acq On : 6 Aug 2015 1:32  
Operator : UM/TP  
Sample : G3109-14RE  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01W2RE

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
Quant Title : SVOA CALIBRATION

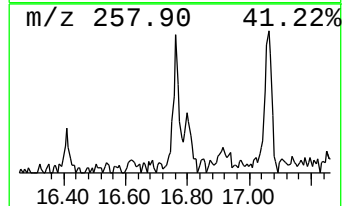
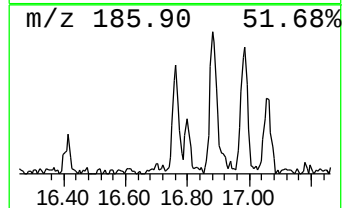
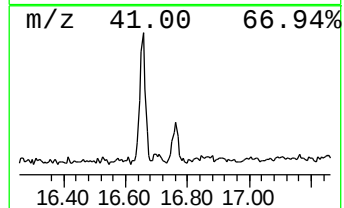
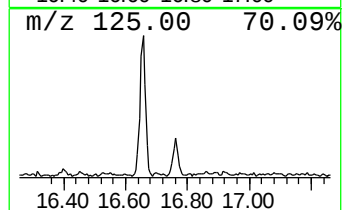
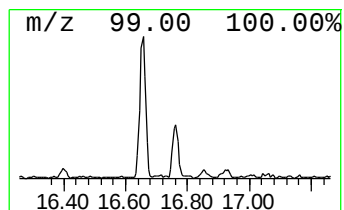
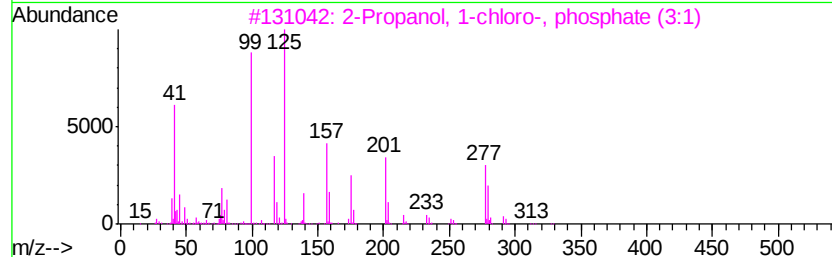
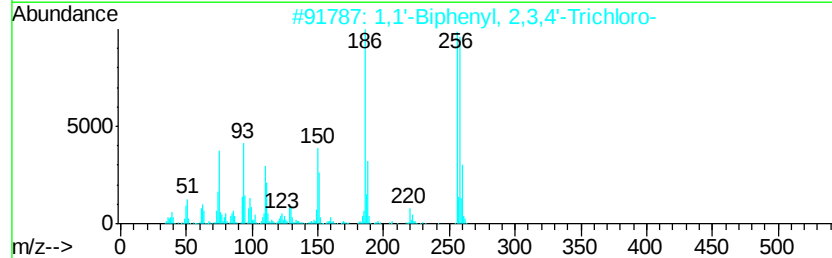
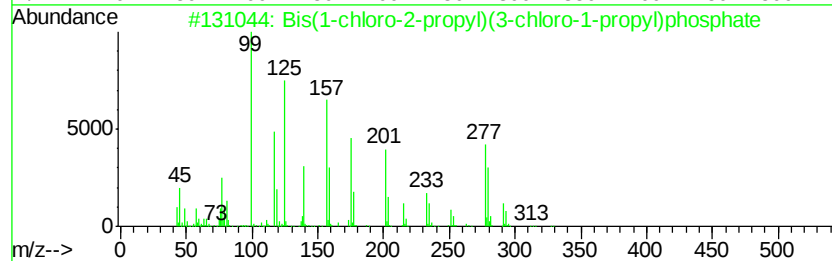
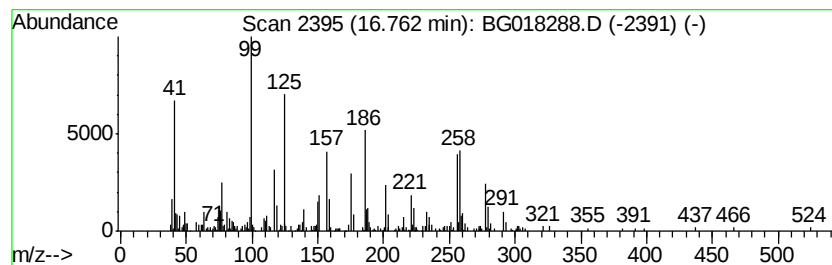
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 Bis(1-chloro-2-propyl)(3-ch... Concentration Rank 10

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 16.76 | 6.49 ng/ul | 148245 | Phenanthrene-d10 | 16.88 |

| Hit# | of | Tentative ID                        | MW  | MolForm     | CAS#        | Qual |
|------|----|-------------------------------------|-----|-------------|-------------|------|
| 1    | 5  | Bis(1-chloro-2-propyl)(3-chloro-... | 326 | C9H18Cl3O4P | 137909-40-1 | 60   |
| 2    |    | 1,1'-Biphenyl, 2,3,4'-Trichloro-    | 256 | C12H7Cl3    | 038444-85-8 | 52   |
| 3    |    | 2-Propanol, 1-chloro-, phosphate... | 326 | C9H18Cl3O4P | 013674-84-5 | 38   |
| 4    |    | 1,1'-Biphenyl, 2,4',5-trichloro-    | 256 | C12H7Cl3    | 016606-02-3 | 30   |
| 5    |    | 1,1'-Biphenyl, 2,4,4'-trichloro-    | 256 | C12H7Cl3    | 007012-37-5 | 25   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG080515\  
Data File : BG018288.D  
Acq On : 6 Aug 2015 1:32  
Operator : UM/TP  
Sample : G3109-14RE  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampled :  
C01W2RE

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
Quant Title : SVOA CALIBRATION

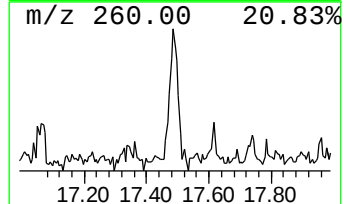
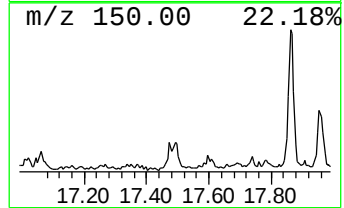
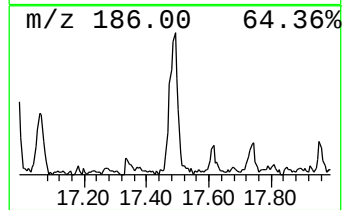
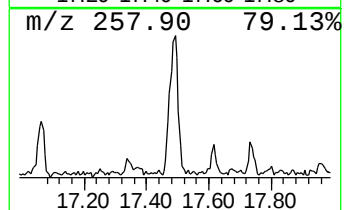
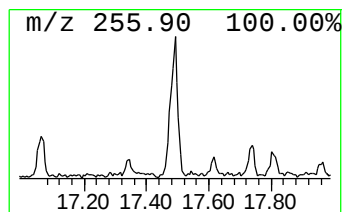
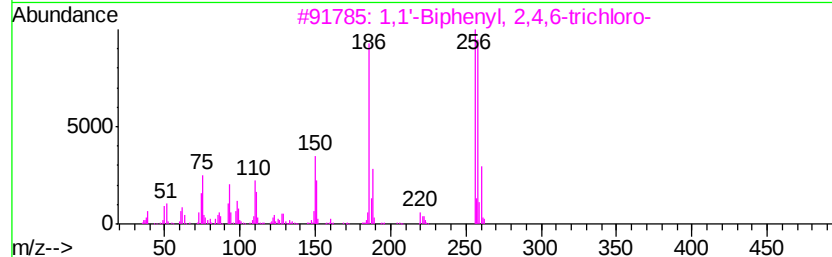
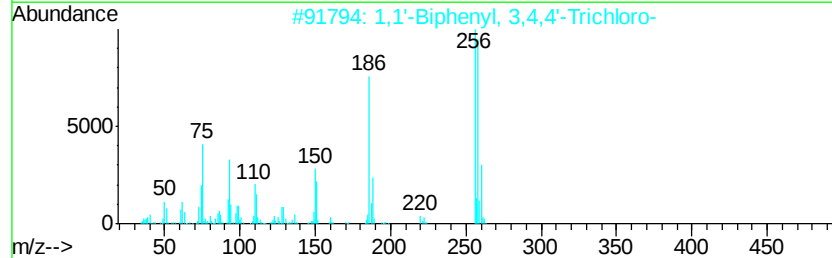
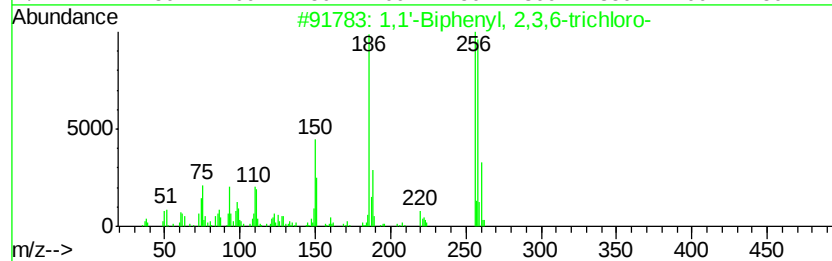
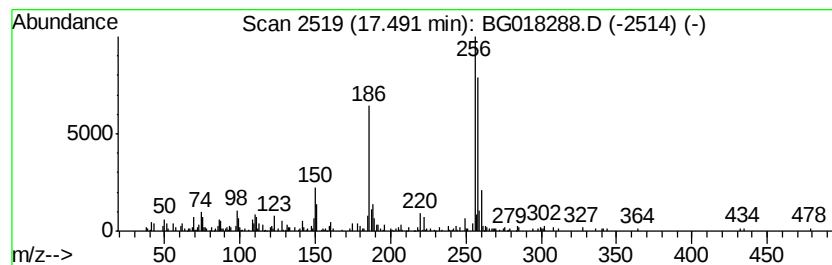
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 6 1,1'-Biphenyl, 2,3,6-trichl... Concentration Rank 6

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 17.49 | 7.48 ng/ul | 170685 | Phenanthrene-d10 | 16.88 |

| Hit# | of | 5 | Tentative ID                     | MW  | MolForm  | CAS#        | Qual |
|------|----|---|----------------------------------|-----|----------|-------------|------|
| 1    |    |   | 1,1'-Biphenyl, 2,3,6-trichloro-  | 256 | C12H7Cl3 | 055702-45-9 | 95   |
| 2    |    |   | 1,1'-Biphenyl, 3,4,4'-Trichloro- | 256 | C12H7Cl3 | 038444-90-5 | 90   |
| 3    |    |   | 1,1'-Biphenyl, 2,4,6-trichloro-  | 256 | C12H7Cl3 | 035693-92-6 | 89   |
| 4    |    |   | 1,1'-Biphenyl, 2,4,4'-trichloro- | 256 | C12H7Cl3 | 007012-37-5 | 89   |
| 5    |    |   | 1,1'-Biphenyl, 2,4',5-trichloro- | 256 | C12H7Cl3 | 016606-02-3 | 89   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG080515\  
Data File : BG018288.D  
Acq On : 6 Aug 2015 1:32  
Operator : UM/TP  
Sample : G3109-14RE  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01W2RE

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
Quant Title : SVOA CALIBRATION

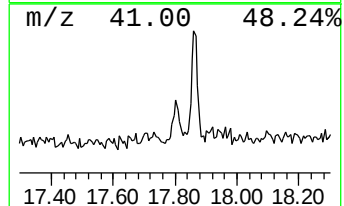
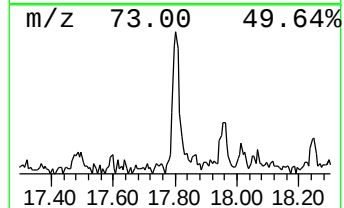
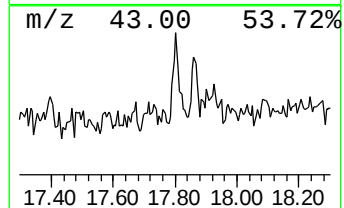
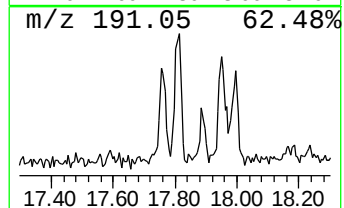
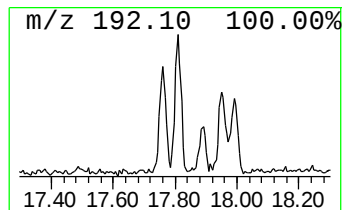
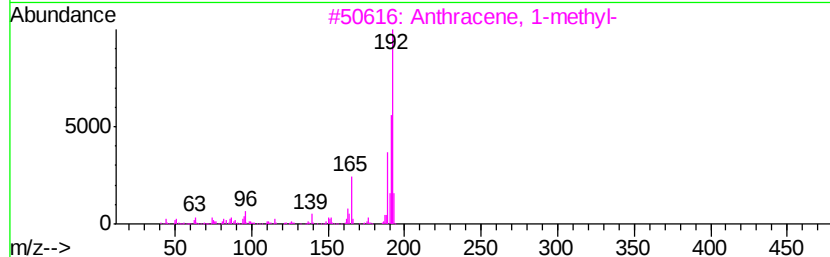
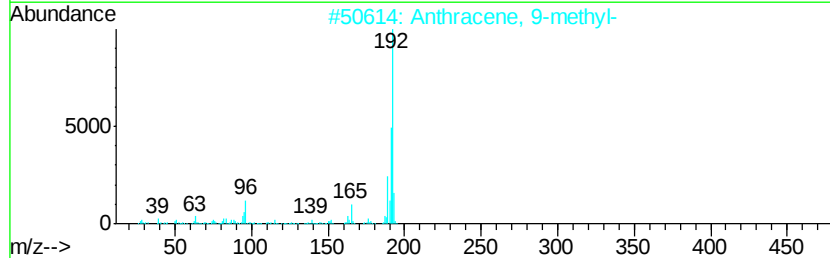
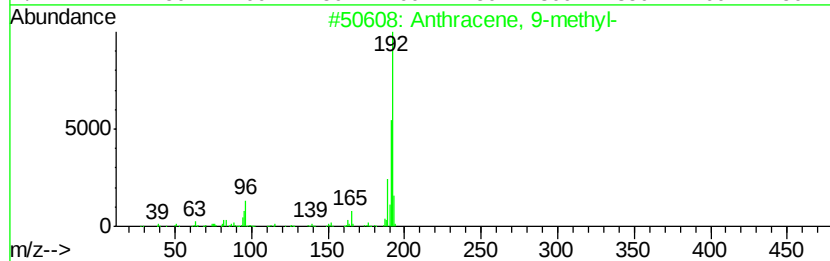
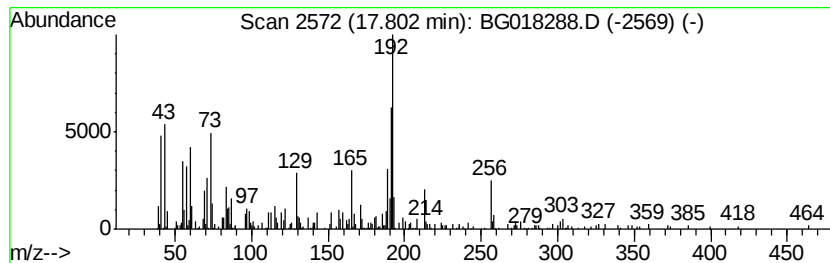
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 7 Anthracene, 9-methyl- Concentration Rank 8

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 17.80 | 7.16 ng/ul | 163427 | Phenanthrene-d10 | 16.88 |

| Hit# | of | Tentative ID          | MW  | MolForm  | CAS#        | Qual |
|------|----|-----------------------|-----|----------|-------------|------|
| 1    | 5  | Anthracene, 9-methyl- | 192 | C15H12   | 000779-02-2 | 55   |
| 2    |    | Anthracene, 9-methyl- | 192 | C15H12   | 000779-02-2 | 55   |
| 3    |    | Anthracene, 1-methyl- | 192 | C15H12   | 000610-48-0 | 55   |
| 4    |    | n-Hexadecanoic acid   | 256 | C16H32O2 | 000057-10-3 | 53   |
| 5    |    | 1H-Indene, 2-phenyl-  | 192 | C15H12   | 004505-48-0 | 45   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG080515\  
Data File : BG018288.D  
Acq On : 6 Aug 2015 1:32  
Operator : UM/TP  
Sample : G3109-14RE  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampled :  
C01W2RE

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
Quant Title : SVOA CALIBRATION

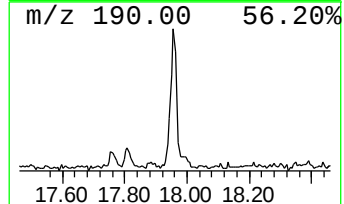
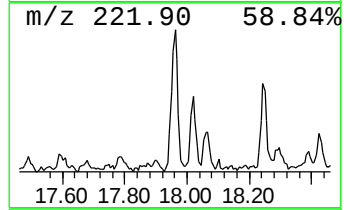
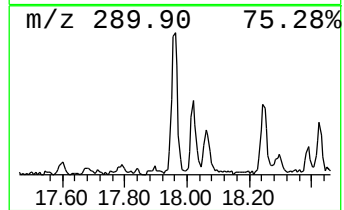
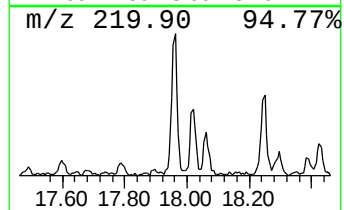
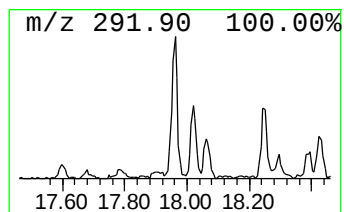
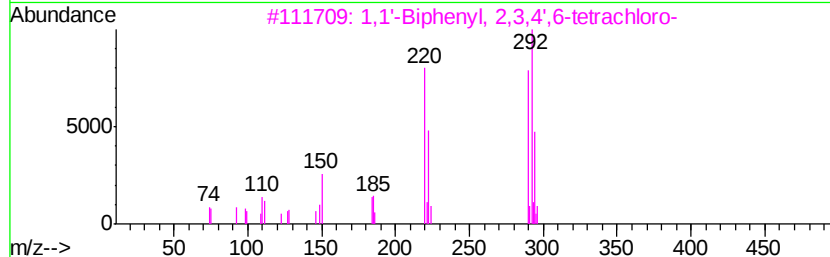
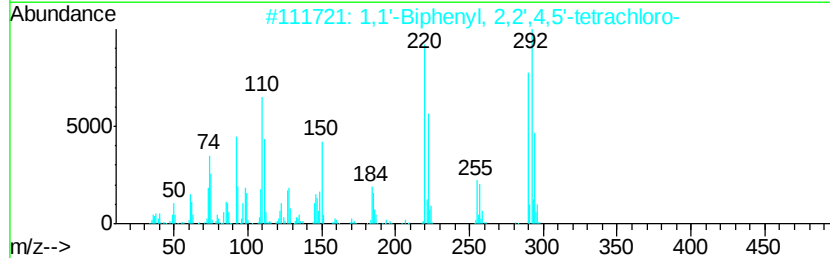
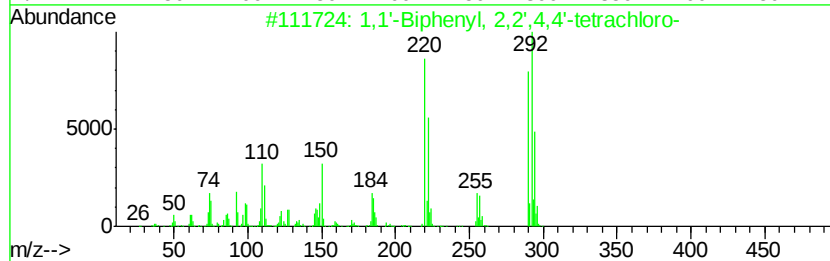
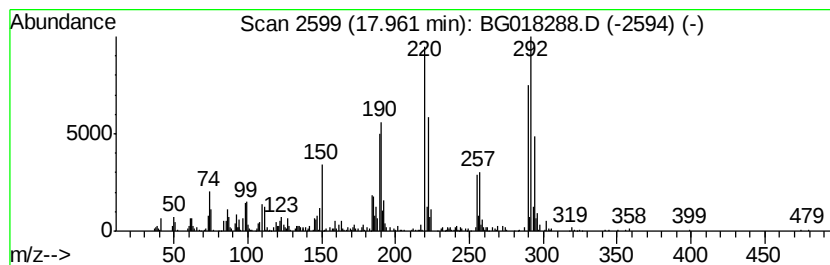
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 1,1'-Biphenyl, 2,2',4,4'-te... Concentration Rank 3

| R.T.  | EstConc     | Area   | Relative to ISTD | R.T.  |
|-------|-------------|--------|------------------|-------|
| 17.96 | 14.59 ng/ul | 333061 | Phenanthrene-d10 | 16.88 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 1,1'-Biphenyl, 2,2',4,4'-tetrach... | 290 | C12H6Cl4 | 002437-79-8 | 98   |
| 2    |    | 1,1'-Biphenyl, 2,2',4,5'-tetrach... | 290 | C12H6Cl4 | 041464-40-8 | 94   |
| 3    |    | 1,1'-Biphenyl, 2,3,4',6-tetrachl... | 290 | C12H6Cl4 | 052663-58-8 | 93   |
| 4    |    | 1,1'-Biphenyl, 3,3',5,5'-tetrach... | 290 | C12H6Cl4 | 033284-52-5 | 93   |
| 5    |    | 1,1'-Biphenyl, 2,4,4',6-tetrachl... | 290 | C12H6Cl4 | 032598-12-2 | 92   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG080515\  
Data File : BG018288.D  
Acq On : 6 Aug 2015 1:32  
Operator : UM/TP  
Sample : G3109-14RE  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01W2RE

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
Quant Title : SVOA CALIBRATION

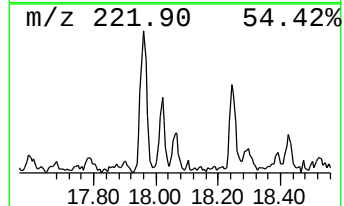
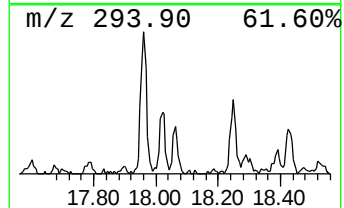
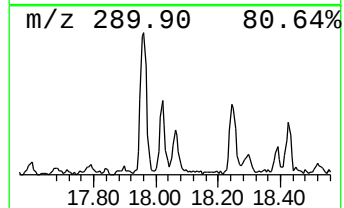
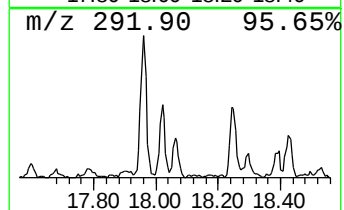
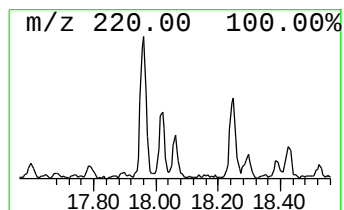
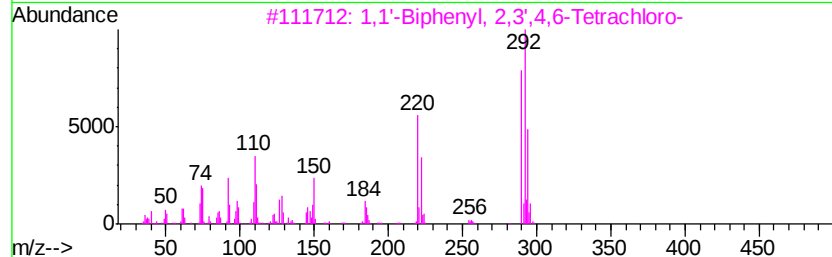
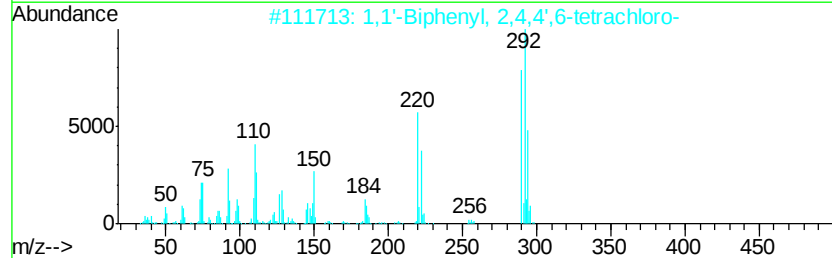
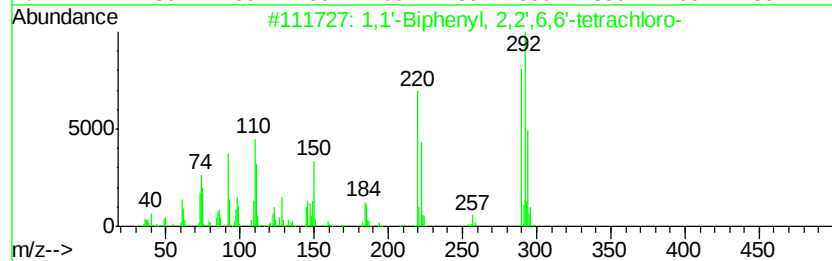
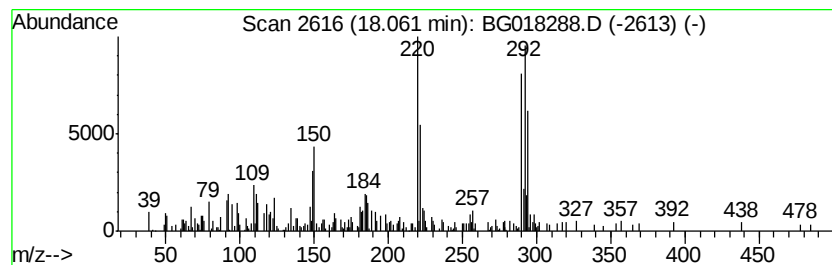
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 1,1'-Biphenyl, 2,2',6,6'-te... Concentration Rank 16

| R.T.  | EstConc    | Area  | Relative to ISTD | R.T.  |
|-------|------------|-------|------------------|-------|
| 18.06 | 3.50 ng/ul | 79885 | Phenanthrene-d10 | 16.88 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 1,1'-Biphenyl, 2,2',6,6'-tetrach... | 290 | C12H6Cl4 | 015968-05-5 | 94   |
| 2    |    |   | 1,1'-Biphenyl, 2,4,4',6-tetrachl... | 290 | C12H6Cl4 | 032598-12-2 | 94   |
| 3    |    |   | 1,1'-Biphenyl, 2,3',4,6-Tetrachl... | 290 | C12H6Cl4 | 060233-24-1 | 94   |
| 4    |    |   | 1,1'-Biphenyl, 2,3,4,4'-tetrachl... | 290 | C12H6Cl4 | 033025-41-1 | 94   |
| 5    |    |   | 1,1'-Biphenyl, 2,2',4,4'-tetrach... | 290 | C12H6Cl4 | 002437-79-8 | 93   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG080515\  
Data File : BG018288.D  
Acq On : 6 Aug 2015 1:32  
Operator : UM/TP  
Sample : G3109-14RE  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01W2RE

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
Quant Title : SVOA CALIBRATION

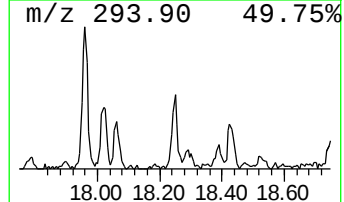
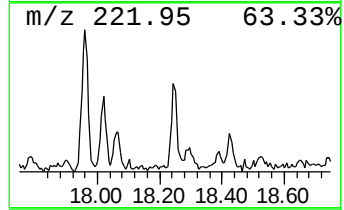
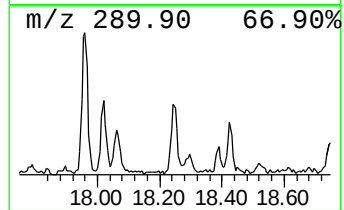
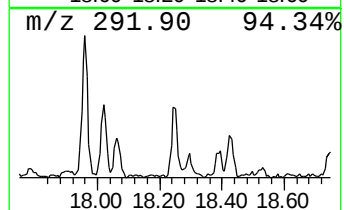
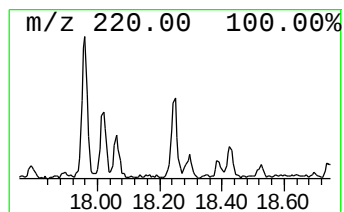
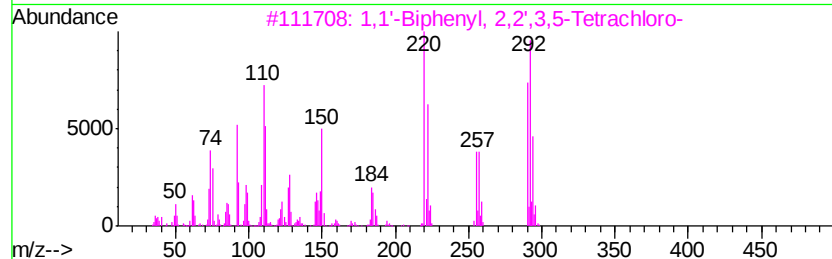
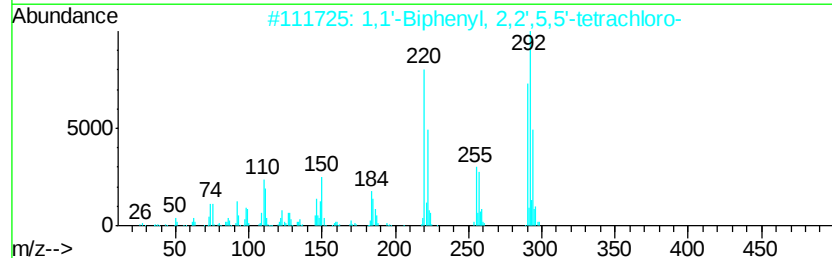
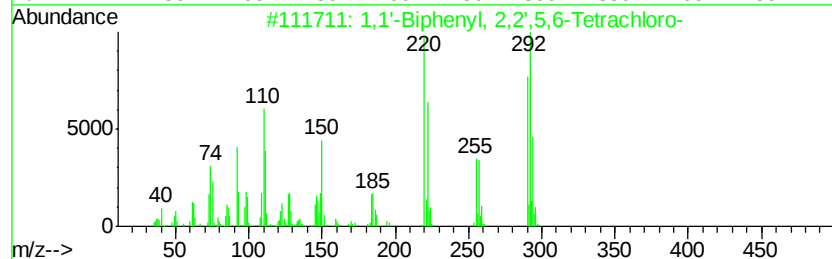
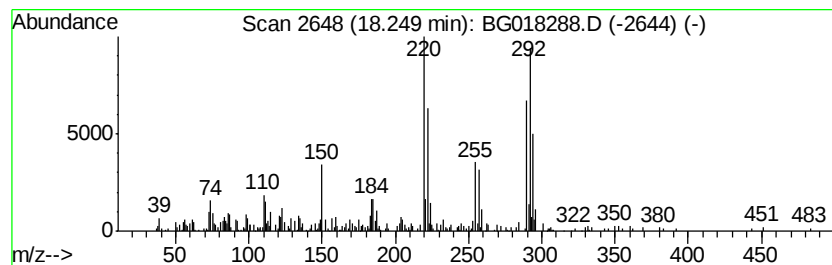
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 1,1'-Biphenyl, 2,2',5,6-Tet... Concentration Rank 5

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.25 | 8.05 ng/ul | 183908 | Phenanthrene-d10 | 16.88 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 1,1'-Biphenyl, 2,2',5,6-Tetrachl... | 290 | C12H6Cl4 | 041464-41-9 | 95   |
| 2    |    | 1,1'-Biphenyl, 2,2',5,5'-tetrach... | 290 | C12H6Cl4 | 035693-99-3 | 94   |
| 3    |    | 1,1'-Biphenyl, 2,2',3,5-Tetrachl... | 290 | C12H6Cl4 | 070362-46-8 | 94   |
| 4    |    | 1,1'-Biphenyl, 2,2',5,5'-tetrach... | 290 | C12H6Cl4 | 035693-99-3 | 94   |
| 5    |    | 1,1'-Biphenyl, 2,2',4,5'-tetrach... | 290 | C12H6Cl4 | 041464-40-8 | 93   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG080515\  
Data File : BG018288.D  
Acq On : 6 Aug 2015 1:32  
Operator : UM/TP  
Sample : G3109-14RE  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01W2RE

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
Quant Title : SVOA CALIBRATION

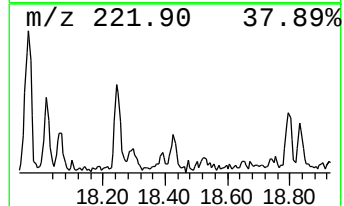
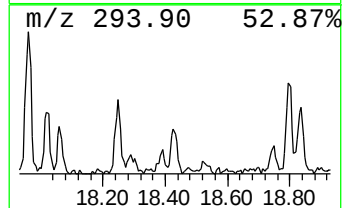
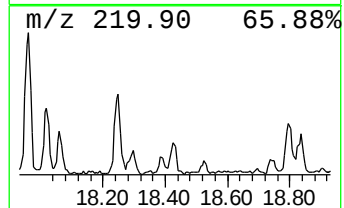
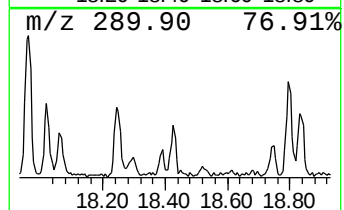
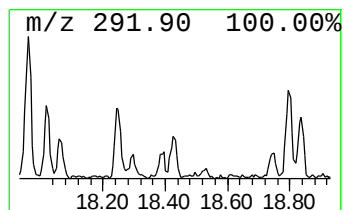
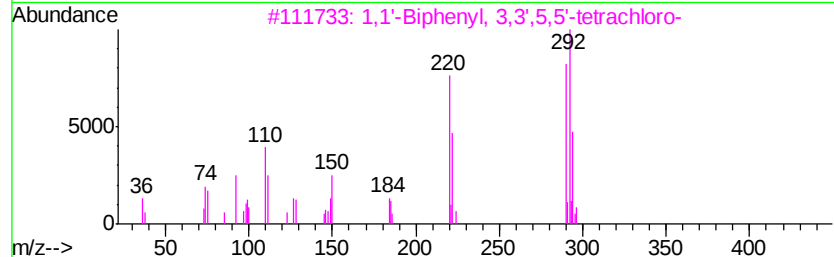
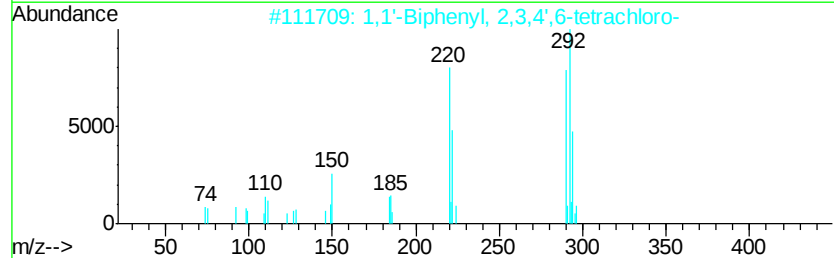
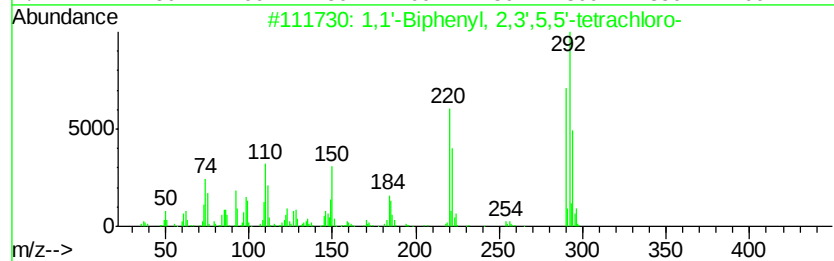
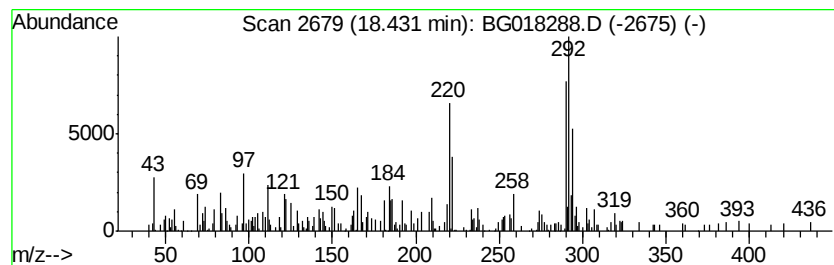
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 1,1'-Biphenyl, 2,3',5,5'-te... Concentration Rank 20

| R.T.  | EstConc    | Area  | Relative to ISTD | R.T.  |
|-------|------------|-------|------------------|-------|
| 18.43 | 3.00 ng/ul | 68596 | Phenanthrene-d10 | 16.88 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 1,1'-Biphenyl, 2,3',5,5'-tetrach... | 290 | C12H6Cl4 | 041464-42-0 | 94   |
| 2    |    |   | 1,1'-Biphenyl, 2,3,4',6-tetrachl... | 290 | C12H6Cl4 | 052663-58-8 | 93   |
| 3    |    |   | 1,1'-Biphenyl, 3,3',5,5'-tetrach... | 290 | C12H6Cl4 | 033284-52-5 | 93   |
| 4    |    |   | 1,1'-Biphenyl, 2,3',4,4'-tetrach... | 290 | C12H6Cl4 | 032598-10-0 | 93   |
| 5    |    |   | 1,1'-Biphenyl, 2,4,4',6-tetrachl... | 290 | C12H6Cl4 | 032598-12-2 | 93   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG080515\  
Data File : BG018288.D  
Acq On : 6 Aug 2015 1:32  
Operator : UM/TP  
Sample : G3109-14RE  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01W2RE

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
Quant Title : SVOA CALIBRATION

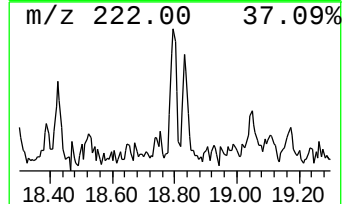
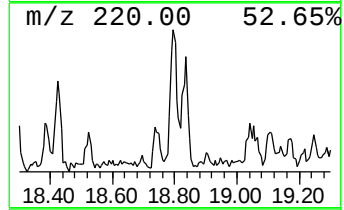
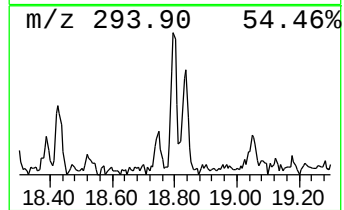
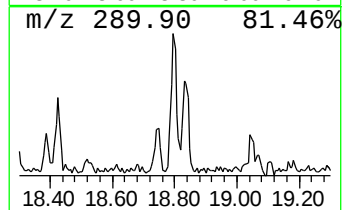
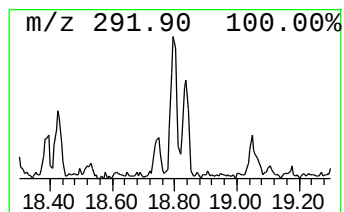
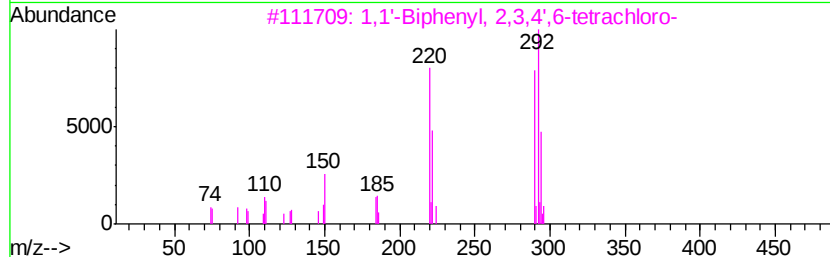
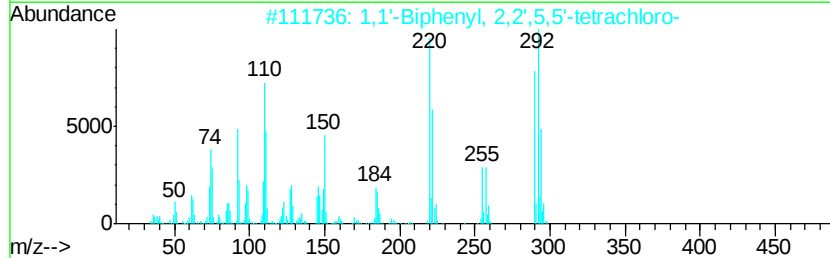
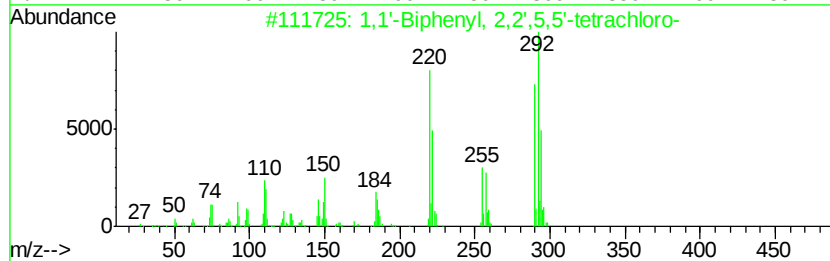
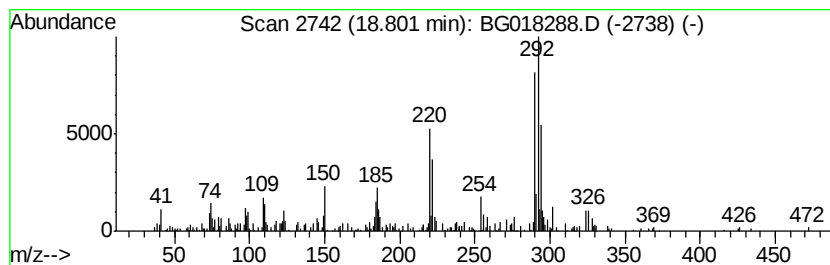
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 1,1'-Biphenyl, 2,2',5,5'-te... Concentration Rank 14

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.80 | 4.54 ng/ul | 103591 | Phenanthrene-d10 | 16.88 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 1,1'-Biphenyl, 2,2',5,5'-tetrach... | 290 | C12H6Cl4 | 035693-99-3 | 94   |
| 2    |    |   | 1,1'-Biphenyl, 2,2',5,5'-tetrach... | 290 | C12H6Cl4 | 035693-99-3 | 94   |
| 3    |    |   | 1,1'-Biphenyl, 2,3,4',6-tetrachl... | 290 | C12H6Cl4 | 052663-58-8 | 94   |
| 4    |    |   | 1,1'-Biphenyl, 2,2',6,6'-tetrach... | 290 | C12H6Cl4 | 015968-05-5 | 94   |
| 5    |    |   | 1,1'-Biphenyl, 2,4,4',6-tetrachl... | 290 | C12H6Cl4 | 032598-12-2 | 94   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG080515\  
Data File : BG018288.D  
Acq On : 6 Aug 2015 1:32  
Operator : UM/TP  
Sample : G3109-14RE  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01W2RE

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
Quant Title : SVOA CALIBRATION

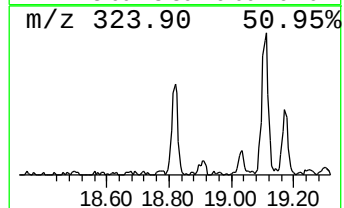
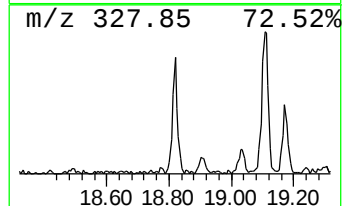
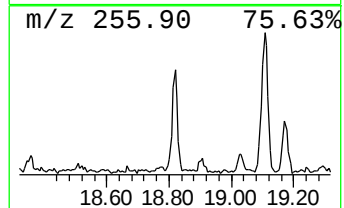
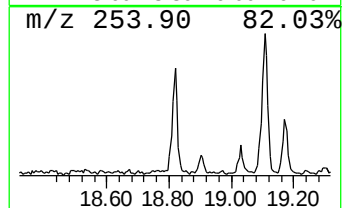
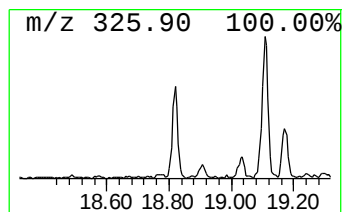
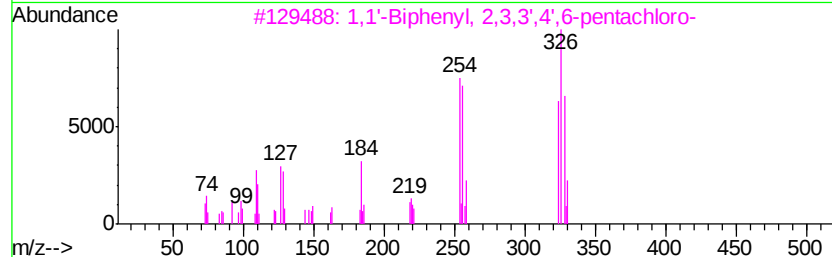
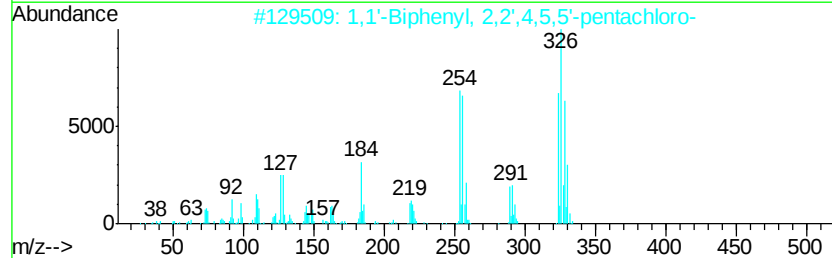
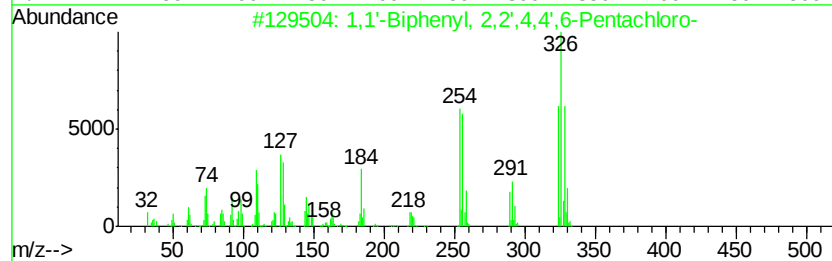
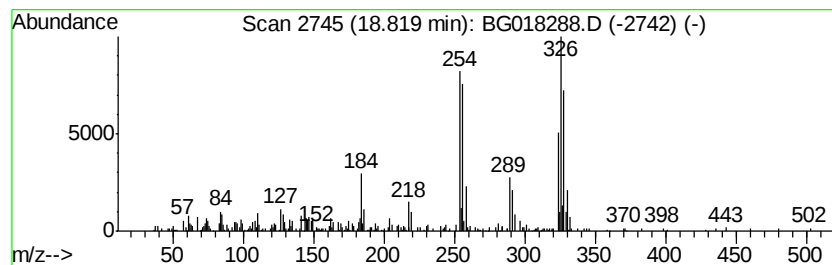
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 1,1'-Biphenyl, 2,2',4,4',6-... Concentration Rank 2

| R.T.  | EstConc     | Area   | Relative to ISTD | R.T.  |
|-------|-------------|--------|------------------|-------|
| 18.82 | 15.11 ng/ul | 344960 | Phenanthrene-d10 | 16.88 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 1,1'-Biphenyl, 2,2',4,4',6-Penta... | 324 | C12H5Cl5 | 039485-83-1 | 96   |
| 2    |    | 1,1'-Biphenyl, 2,2',4,5,5'-penta... | 324 | C12H5Cl5 | 037680-73-2 | 95   |
| 3    |    | 1,1'-Biphenyl, 2,3,3',4',6-penta... | 324 | C12H5Cl5 | 038380-03-9 | 94   |
| 4    |    | 1,1'-Biphenyl, 2,2',4,5,5'-penta... | 324 | C12H5Cl5 | 037680-73-2 | 94   |
| 5    |    | 1,1'-Biphenyl, 2,3',4,4',5-penta... | 324 | C12H5Cl5 | 031508-00-6 | 94   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG080515\  
Data File : BG018288.D  
Acq On : 6 Aug 2015 1:32  
Operator : UM/TP  
Sample : G3109-14RE  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01W2RE

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
Quant Title : SVOA CALIBRATION

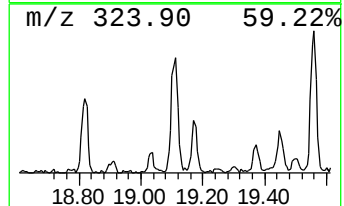
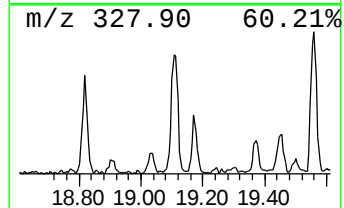
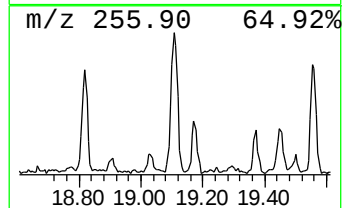
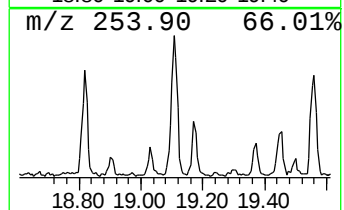
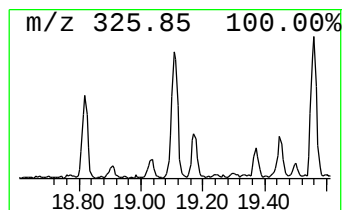
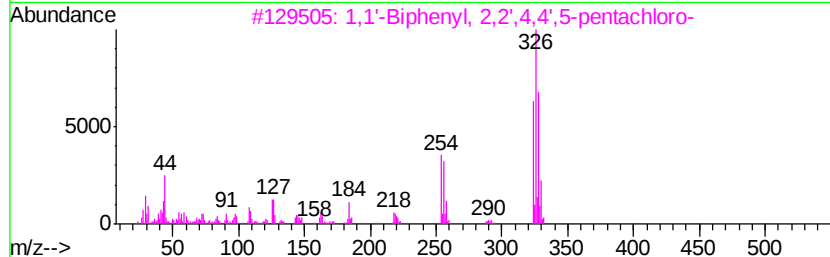
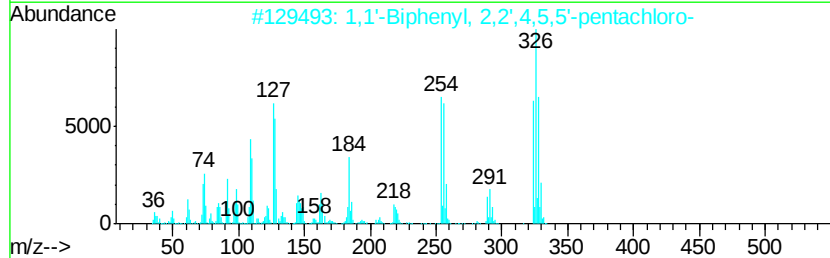
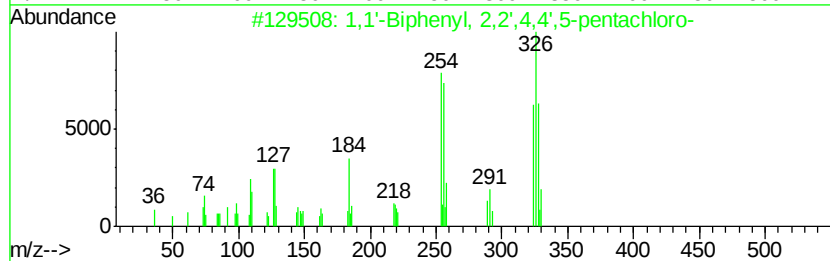
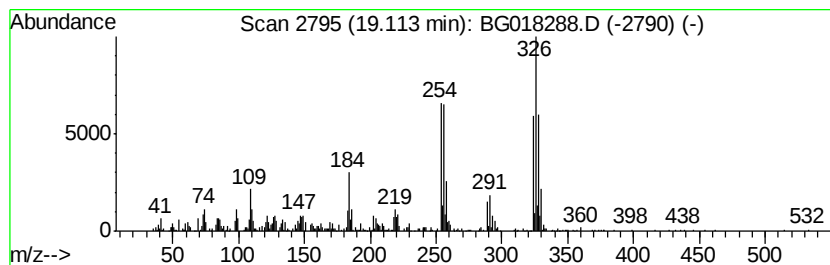
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 1,1'-Biphenyl, 2,2',4,4',5-... Concentration Rank 4

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 19.11 | 8.78 ng/ul | 404442 | Chrysene-d12     | 21.10 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 1,1'-Biphenyl, 2,2',4,4',5-penta... | 324 | C12H5Cl5 | 038380-01-7 | 95   |
| 2    |    | 1,1'-Biphenyl, 2,2',4,5,5'-penta... | 324 | C12H5Cl5 | 037680-73-2 | 95   |
| 3    |    | 1,1'-Biphenyl, 2,2',4,4',5-penta... | 324 | C12H5Cl5 | 038380-01-7 | 95   |
| 4    |    | 1,1'-Biphenyl, 2,3',4,4',5-penta... | 324 | C12H5Cl5 | 031508-00-6 | 95   |
| 5    |    | 1,1'-Biphenyl, 2,3',4,4',6-Penta... | 324 | C12H5Cl5 | 056558-17-9 | 95   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG080515\  
Data File : BG018288.D  
Acq On : 6 Aug 2015 1:32  
Operator : UM/TP  
Sample : G3109-14RE  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01W2RE

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
Quant Title : SVOA CALIBRATION

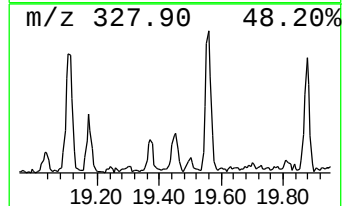
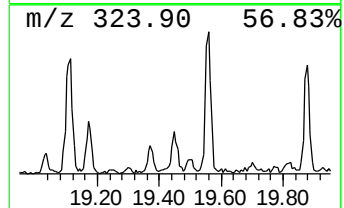
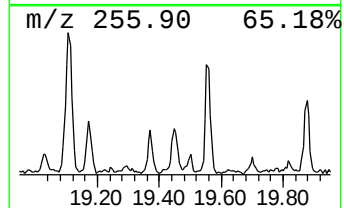
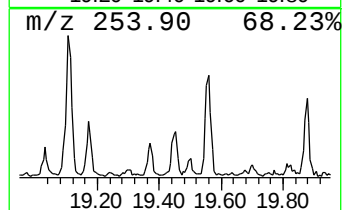
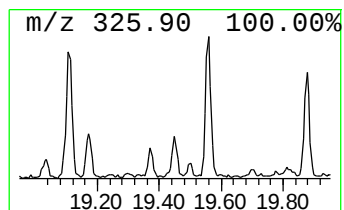
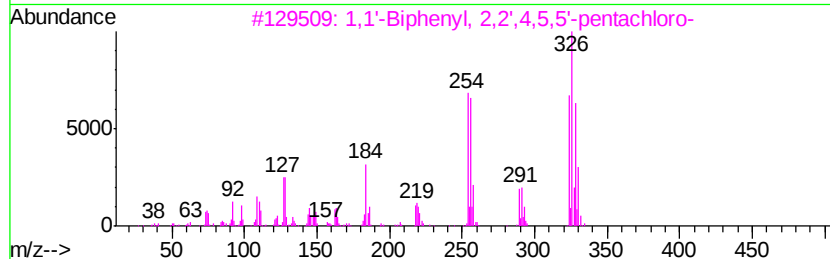
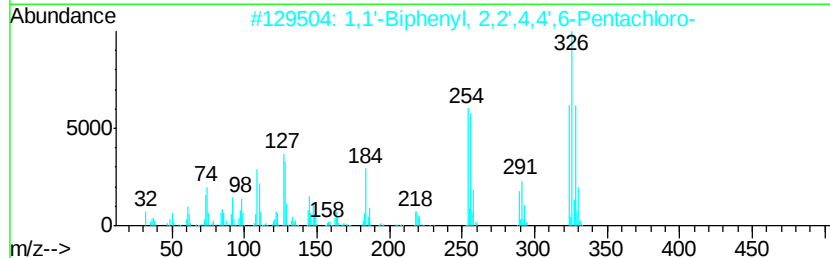
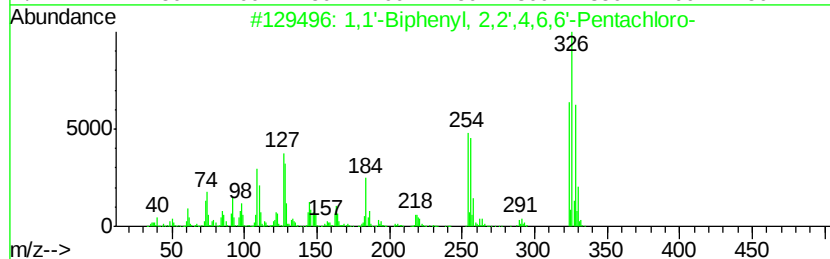
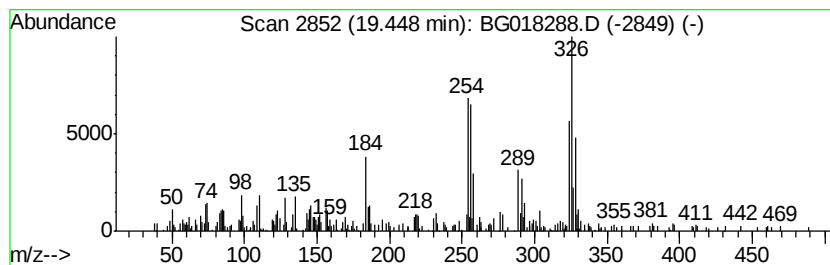
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 17 1,1'-Biphenyl, 2,2',4,6,6'-... Concentration Rank 21

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 19.45 | 2.61 ng/ul | 120072 | Chrysene-d12     | 21.10 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 1,1'-Biphenyl, 2,2',4,6,6'-Penta... | 324 | C12H5Cl5 | 056558-16-8 | 91   |
| 2    |    |   | 1,1'-Biphenyl, 2,2',4,4',6-Penta... | 324 | C12H5Cl5 | 039485-83-1 | 91   |
| 3    |    |   | 1,1'-Biphenyl, 2,2',4,5,5'-penta... | 324 | C12H5Cl5 | 037680-73-2 | 87   |
| 4    |    |   | 1,1'-Biphenyl, 2,2',4,5,6'-penta... | 324 | C12H5Cl5 | 068194-06-9 | 83   |
| 5    |    |   | 1,1'-Biphenyl, 2,2',3,3',5-penta... | 324 | C12H5Cl5 | 060145-20-2 | 81   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG080515\  
Data File : BG018288.D  
Acq On : 6 Aug 2015 1:32  
Operator : UM/TP  
Sample : G3109-14RE  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01W2RE

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
Quant Title : SVOA CALIBRATION

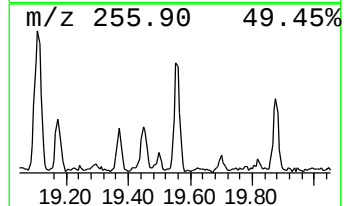
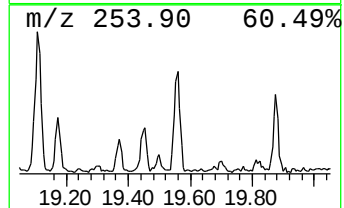
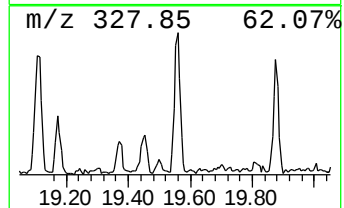
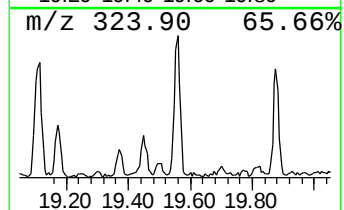
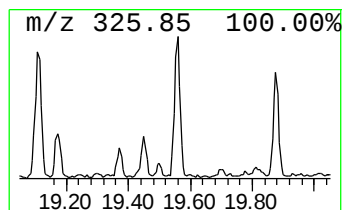
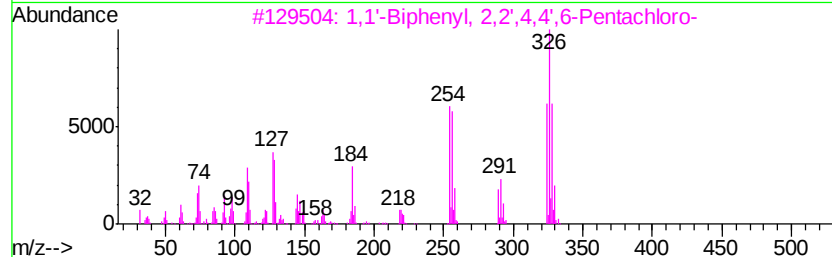
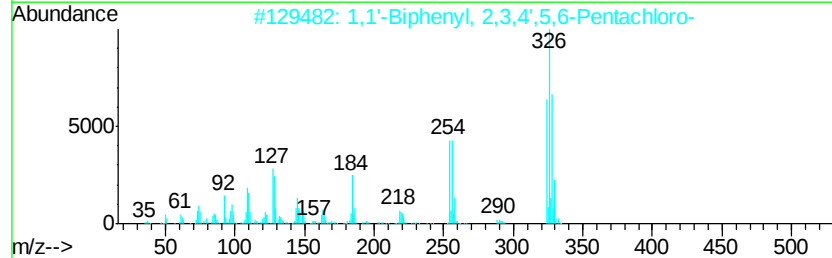
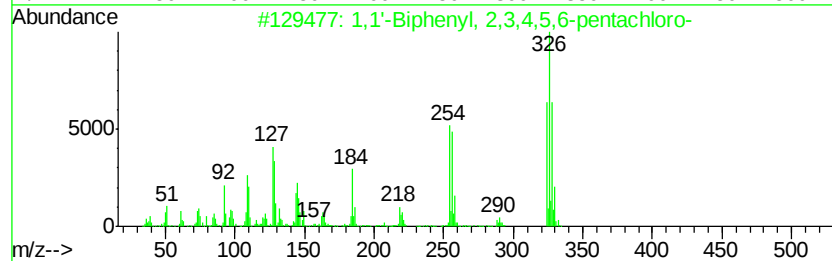
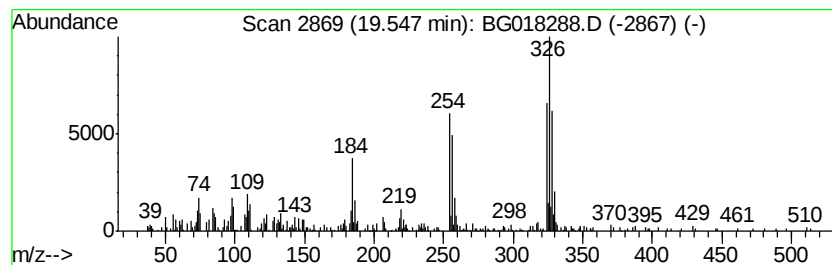
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 18 1,1'-Biphenyl, 2,3,4,5,6-pe... Concentration Rank 7

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 19.55 | 7.40 ng/ul | 341179 | Chrysene-d12     | 21.10 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 1,1'-Biphenyl, 2,3,4,5,6-pentach... | 324 | C12H5Cl5 | 018259-05-7 | 95   |
| 2    |    | 1,1'-Biphenyl, 2,3,4',5,6-Pentac... | 324 | C12H5Cl5 | 068194-11-6 | 95   |
| 3    |    | 1,1'-Biphenyl, 2,2',4,4',6-Penta... | 324 | C12H5Cl5 | 039485-83-1 | 95   |
| 4    |    | 1,1'-Biphenyl, 2,3,4,4',6-Pentac... | 324 | C12H5Cl5 | 074472-38-1 | 94   |
| 5    |    | 1,1'-Biphenyl, 2',3,4,5,5'-Penta... | 324 | C12H5Cl5 | 070424-70-3 | 94   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG080515\  
Data File : BG018288.D  
Acq On : 6 Aug 2015 1:32  
Operator : UM/TP  
Sample : G3109-14RE  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampled :  
C01W2RE

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
Quant Title : SVOA CALIBRATION

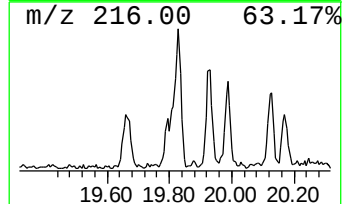
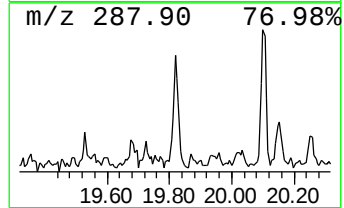
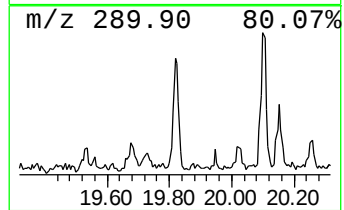
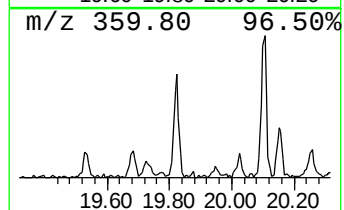
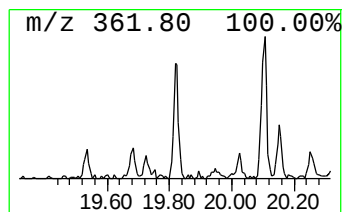
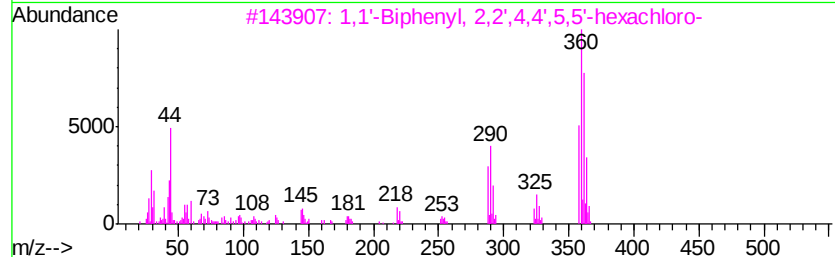
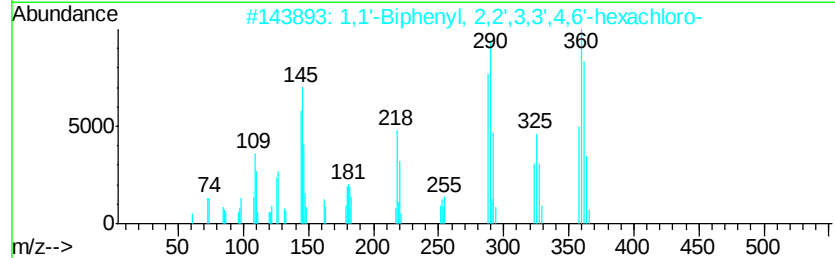
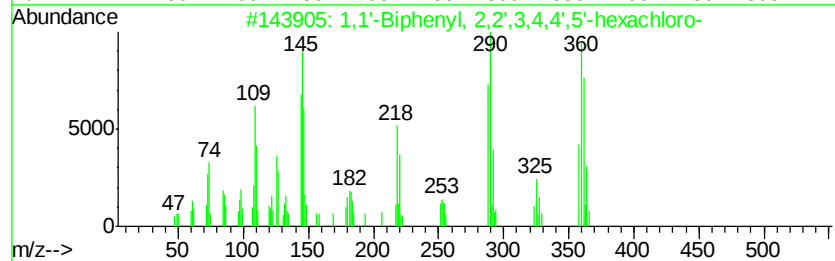
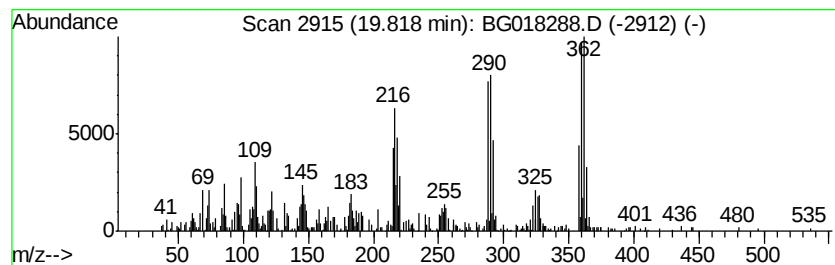
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 19 1,1'-Biphenyl, 2,2',3,4,4',... Concentration Rank 11

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 19.82 | 6.34 ng/ul | 292147 | Chrysene-d12     | 21.10 |

| Hit# | of   | 5         | Tentative ID         | MW  | MolForm  | CAS#        | Qual |
|------|------|-----------|----------------------|-----|----------|-------------|------|
| 1    | 1,1' | -Biphenyl | 2,2',3,4,4',5'-he... | 358 | C12H4Cl6 | 035065-28-2 | 94   |
| 2    | 1,1' | -Biphenyl | 2,2',3,3',4,6'-he... | 358 | C12H4Cl6 | 038380-05-1 | 91   |
| 3    | 1,1' | -Biphenyl | 2,2',4,4',5,5'-he... | 358 | C12H4Cl6 | 035065-27-1 | 86   |
| 4    | 1,1' | -Biphenyl | 2,2',3,4,4',6-Hex... | 358 | C12H4Cl6 | 056030-56-9 | 86   |
| 5    | 1,1' | -Biphenyl | 3,3',4,4',5,5'-he... | 358 | C12H4Cl6 | 032774-16-6 | 83   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG080515\  
Data File : BG018288.D  
Acq On : 6 Aug 2015 1:32  
Operator : UM/TP  
Sample : G3109-14RE  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01W2RE

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
Quant Title : SVOA CALIBRATION

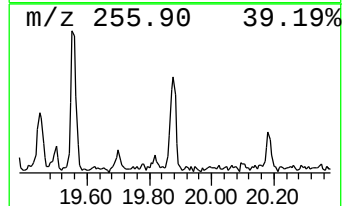
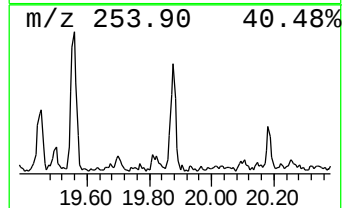
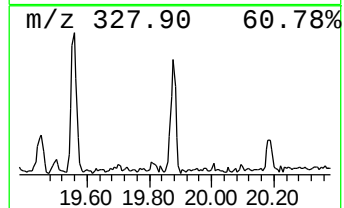
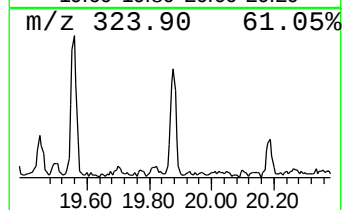
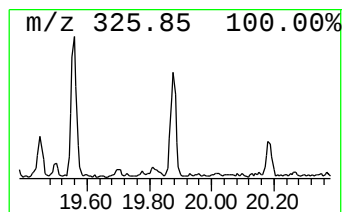
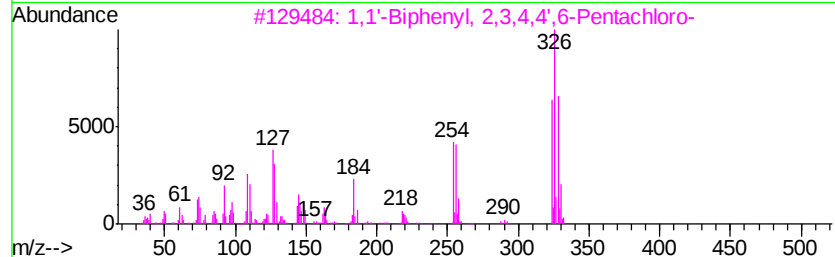
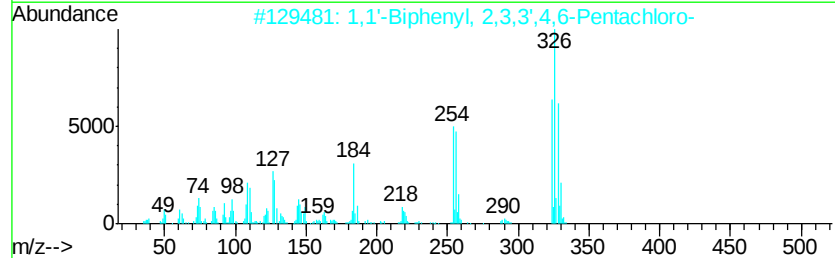
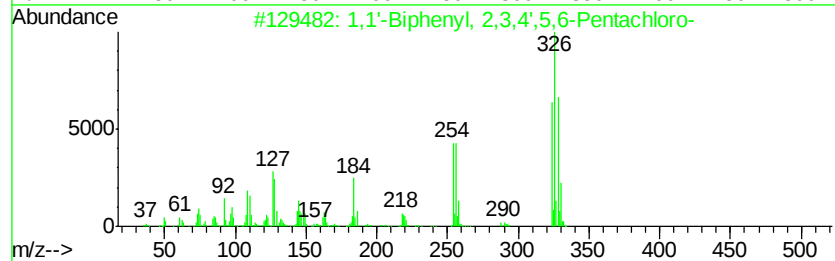
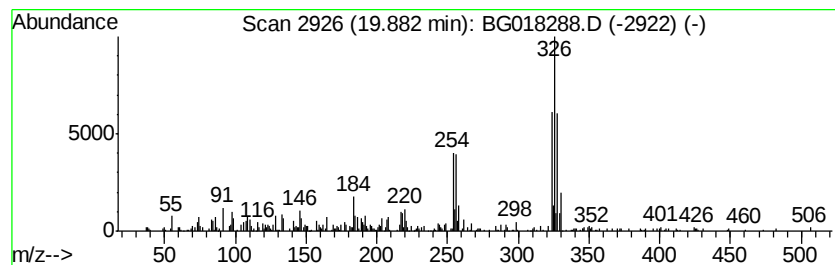
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 20 1,1'-Biphenyl, 2,3,4',5,6-P... Concentration Rank 15

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 19.88 | 4.37 ng/ul | 201248 | Chrysene-d12     | 21.10 |

| Hit# | of   | 5          | Tentative ID         | MW  | MolForm  | CAS#        | Qual |
|------|------|------------|----------------------|-----|----------|-------------|------|
| 1    | 1,1' | -Biphenyl, | 2,3,4',5,6-Pentac... | 324 | C12H5Cl5 | 068194-11-6 | 96   |
| 2    | 1,1' | -Biphenyl, | 2,3,3',4,6-Pentac... | 324 | C12H5Cl5 | 074472-35-8 | 96   |
| 3    | 1,1' | -Biphenyl, | 2,3,4,4',6-Pentac... | 324 | C12H5Cl5 | 074472-38-1 | 95   |
| 4    | 1,1' | -Biphenyl, | 2,3',4,4',6-Penta... | 324 | C12H5Cl5 | 056558-17-9 | 95   |
| 5    | 1,1' | -Biphenyl, | 2,2',3,5,5'-penta... | 324 | C12H5Cl5 | 052663-61-3 | 95   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG080515\  
Data File : BG018288.D  
Acq On : 6 Aug 2015 1:32  
Operator : UM/TP  
Sample : G3109-14RE  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01W2RE

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
Quant Title : SVOA CALIBRATION

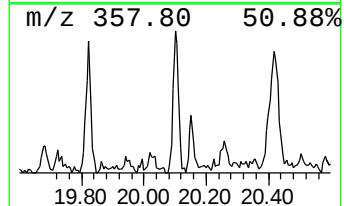
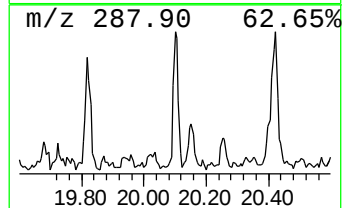
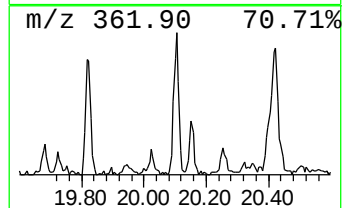
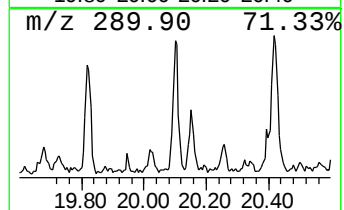
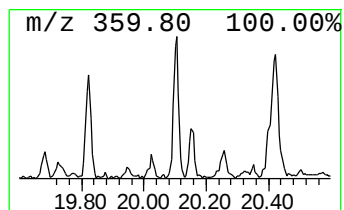
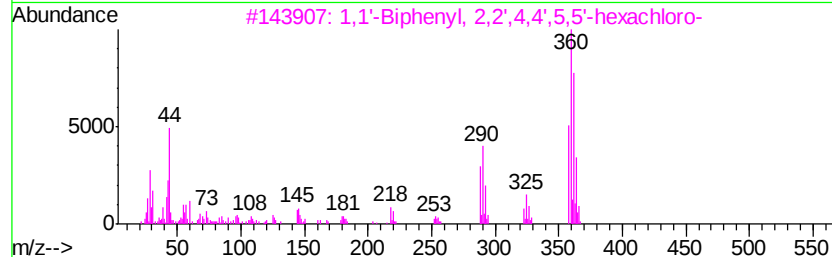
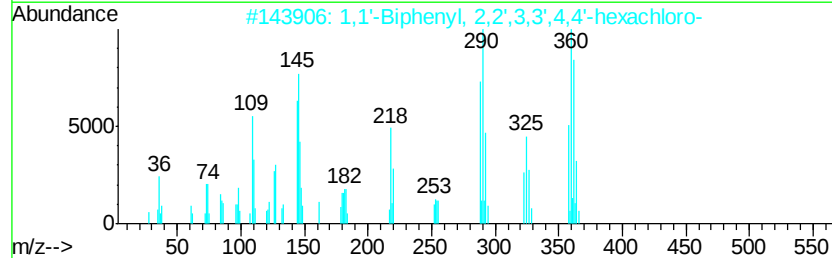
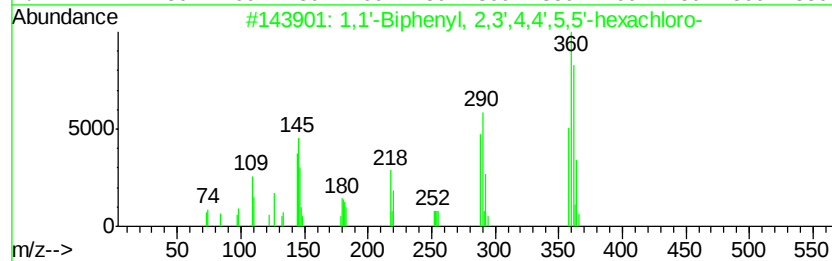
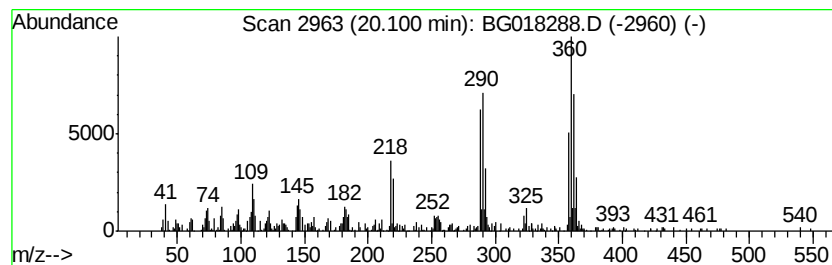
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 21 1,1'-Biphenyl, 2,3',4,4',5,... Concentration Rank 13

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 20.10 | 4.72 ng/ul | 217390 | Chrysene-d12     | 21.10 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 1,1'-Biphenyl, 2,3',4,4',5,5'-he... | 358 | C12H4Cl6 | 052663-72-6 | 96   |
| 2    |    |   | 1,1'-Biphenyl, 2,2',3,3',4,4'-he... | 358 | C12H4Cl6 | 038380-07-3 | 95   |
| 3    |    |   | 1,1'-Biphenyl, 2,2',4,4',5,5'-he... | 358 | C12H4Cl6 | 035065-27-1 | 94   |
| 4    |    |   | 1,1'-Biphenyl, 2,2',3,4,5,5'-hex... | 358 | C12H4Cl6 | 052712-04-6 | 94   |
| 5    |    |   | 1,1'-Biphenyl, 2,2',3,3',5,6'-he... | 358 | C12H4Cl6 | 052744-13-5 | 94   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG080515\  
Data File : BG018288.D  
Acq On : 6 Aug 2015 1:32  
Operator : UM/TP  
Sample : G3109-14RE  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01W2RE

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
Quant Title : SVOA CALIBRATION

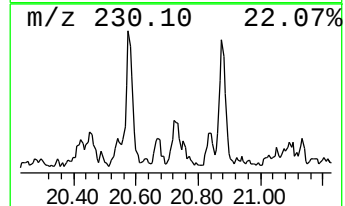
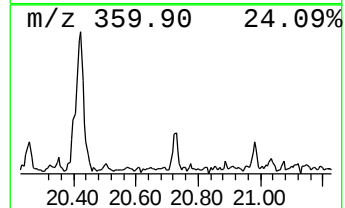
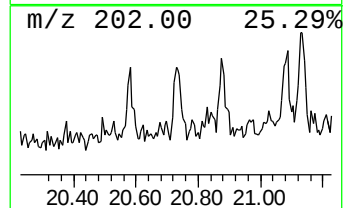
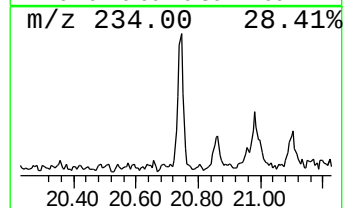
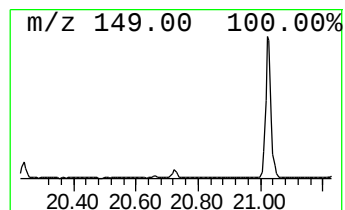
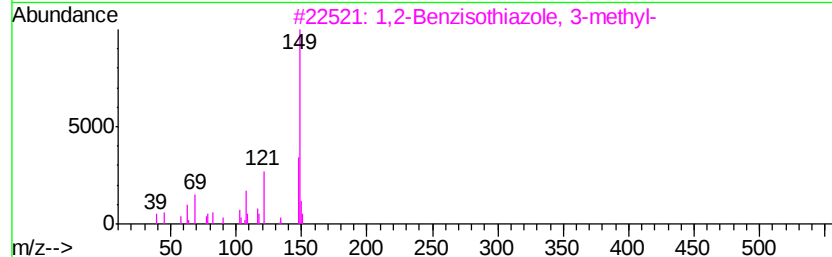
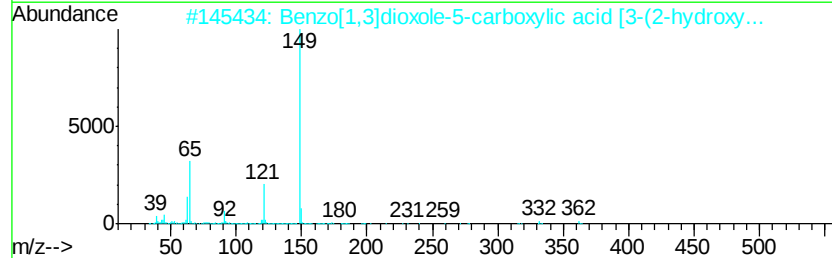
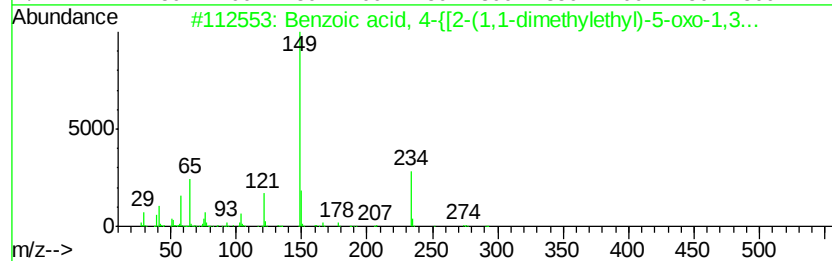
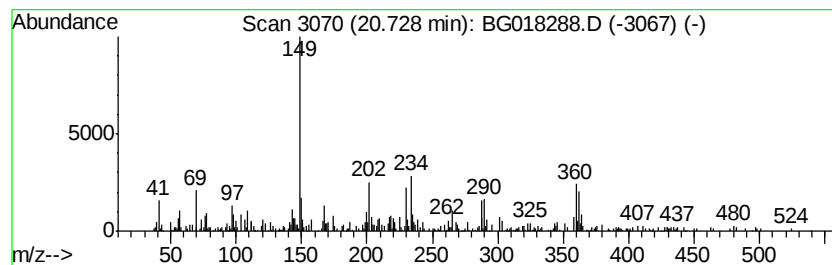
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 23 unknown-03 Concentration Rank 18

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 20.73 | 3.35 ng/ul | 154575 | Chrysene-d12     | 21.10 |

| Hit# | of | 5 | Tentative ID                         | MW  | MolForm     | CAS#         | Qual |
|------|----|---|--------------------------------------|-----|-------------|--------------|------|
| 1    |    |   | Benzoic acid, 4-{[2-(1,1-dimethyl... | 291 | C15H17N05   | 1000197-16-9 | 43   |
| 2    |    |   | Benzo[1,3]dioxole-5-carboxylic a...  | 362 | C16H14N206S | 1000296-89-7 | 43   |
| 3    |    |   | 1,2-Benzisothiazole, 3-methyl-       | 149 | C8H7NS      | 006187-89-9  | 30   |
| 4    |    |   | 1,2-Benzenedicarboxylic acid, bu...  | 278 | C16H2204    | 017851-53-5  | 30   |
| 5    |    |   | Benzothiazole, 2-methyl-             | 149 | C8H7NS      | 000120-75-2  | 30   |



Data Path : Z:\HPCHEM1\BNA G\DATA\BG080515\  
Data File : BG018288.D  
Acq On : 6 Aug 2015 1:32  
Operator : UM/TP  
Sample : G3109-14RE  
Misc :  
ALS Vial : 17 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
C01W2RE

Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG080515.M  
Quant Title : SVOA CALIBRATION

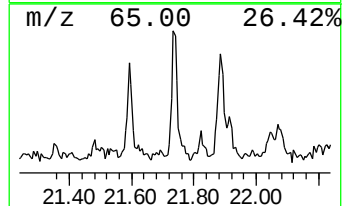
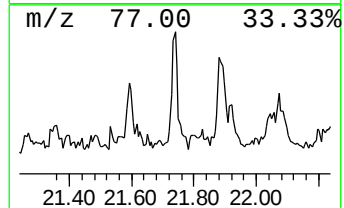
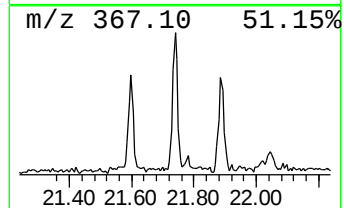
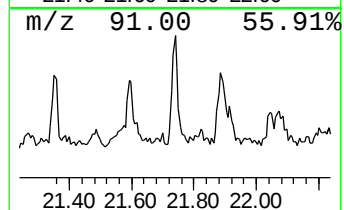
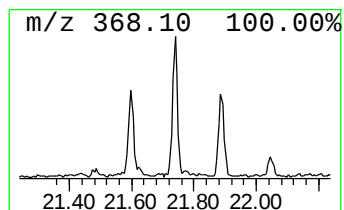
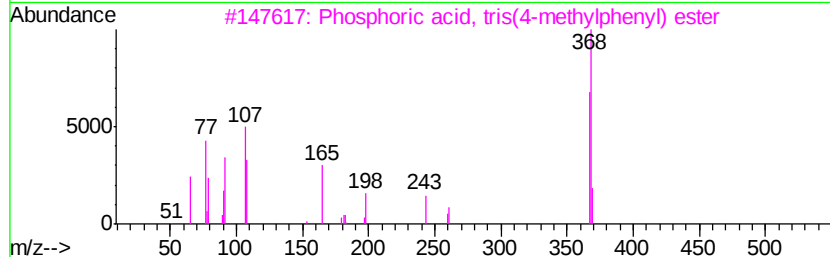
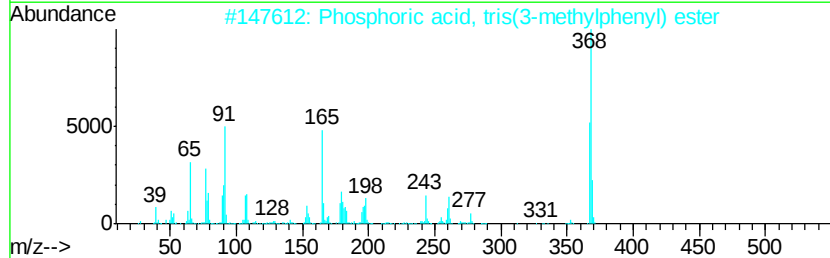
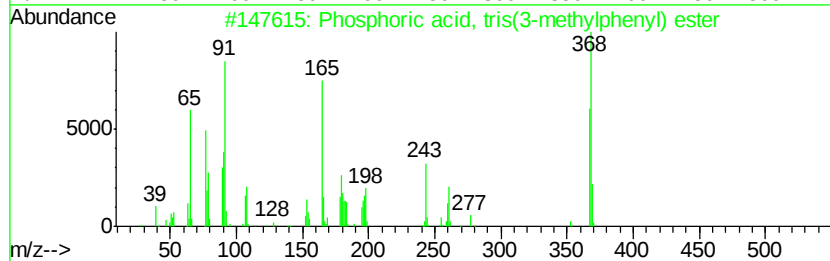
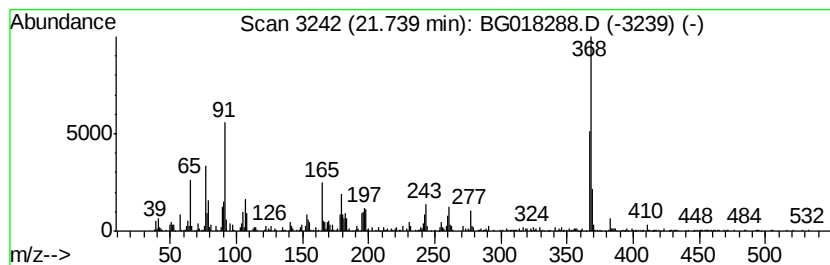
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 24 Phosphoric acid, tris(3-met... Concentration Rank 9

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 21.74 | 6.73 ng/ul | 309904 | Chrysene-d12     | 21.10 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|-------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | Phosphoric acid, tris(3-methylph... | 368 | C21H21O4P | 000563-04-2 | 99   |
| 2    |    | Phosphoric acid, tris(3-methylph... | 368 | C21H21O4P | 000563-04-2 | 97   |
| 3    |    | Phosphoric acid, tris(4-methylph... | 368 | C21H21O4P | 000078-32-0 | 89   |
| 4    |    | Phosphoric acid, tris(4-methylph... | 368 | C21H21O4P | 000078-32-0 | 53   |
| 5    |    | Phosphoric acid, tris(2-methylph... | 368 | C21H21O4P | 000078-30-8 | 41   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG080515\  
 Data File : BG018288.D  
 Acq On : 6 Aug 2015 1:32  
 Operator : UM/TP  
 Sample : G3109-14RE  
 Misc :  
 ALS Vial : 17 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 C01W2RE

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\SOM02.2-EPA-BG080515.M  
 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L  
 TIC Integration Parameters: LSCINT.P

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |        |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|--------|------|
|                      |       |         |       |          | #                     | RT    | Resp   | Conc |
| Benzene, 1,2,4-tr... | 7.09  | 2.1     | ng/ul | 28631    | 1                     | 7.44  | 268457 | 20.0 |
| unknown-01           | 14.91 | 2.3     | ng/ul | 63114    | 3                     | 14.12 | 562105 | 20.0 |
| unknown-02           | 16.40 | 3.4     | ng/ul | 76789    | 4                     | 16.88 | 456644 | 20.0 |
| 2-Propanol, 1-chl... | 16.66 | 16.2    | ng/ul | 370160   | 4                     | 16.88 | 456644 | 20.0 |
| Bis(1-chloro-2-pr... | 16.76 | 6.5     | ng/ul | 148245   | 4                     | 16.88 | 456644 | 20.0 |
| 1,1'-Biphenyl, 2,... | 17.49 | 7.5     | ng/ul | 170685   | 4                     | 16.88 | 456644 | 20.0 |
| Anthracene, 9-met... | 17.80 | 7.2     | ng/ul | 163427   | 4                     | 16.88 | 456644 | 20.0 |
| 1,1'-Biphenyl, 2,... | 17.96 | 14.6    | ng/ul | 333061   | 4                     | 16.88 | 456644 | 20.0 |
| 1,1'-Biphenyl, 2,... | 18.06 | 3.5     | ng/ul | 79885    | 4                     | 16.88 | 456644 | 20.0 |
| 1,1'-Biphenyl, 2,... | 18.25 | 8.1     | ng/ul | 183908   | 4                     | 16.88 | 456644 | 20.0 |
| 1,1'-Biphenyl, 2,... | 18.43 | 3.0     | ng/ul | 68596    | 4                     | 16.88 | 456644 | 20.0 |
| 1,1'-Biphenyl, 2,... | 18.80 | 4.5     | ng/ul | 103591   | 4                     | 16.88 | 456644 | 20.0 |
| 1,1'-Biphenyl, 2,... | 18.82 | 15.1    | ng/ul | 344960   | 4                     | 16.88 | 456644 | 20.0 |
| 1,1'-Biphenyl, 2,... | 19.11 | 8.8     | ng/ul | 404442   | 5                     | 21.10 | 921616 | 20.0 |
| 1,1'-Biphenyl, 2,... | 19.45 | 2.6     | ng/ul | 120072   | 5                     | 21.10 | 921616 | 20.0 |
| 1,1'-Biphenyl, 2,... | 19.55 | 7.4     | ng/ul | 341179   | 5                     | 21.10 | 921616 | 20.0 |
| 1,1'-Biphenyl, 2,... | 19.82 | 6.3     | ng/ul | 292147   | 5                     | 21.10 | 921616 | 20.0 |
| 1,1'-Biphenyl, 2,... | 19.88 | 4.4     | ng/ul | 201248   | 5                     | 21.10 | 921616 | 20.0 |
| 1,1'-Biphenyl, 2,... | 20.10 | 4.7     | ng/ul | 217390   | 5                     | 21.10 | 921616 | 20.0 |
| unknown-03           | 20.73 | 3.4     | ng/ul | 154575   | 5                     | 21.10 | 921616 | 20.0 |
| Phosphoric acid, ... | 21.74 | 6.7     | ng/ul | 309904   | 5                     | 21.10 | 921616 | 20.0 |