

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052720\  
 Data File : BM026138.D  
 Acq On : 27 May 2020 10:35  
 Operator : MA/SJ  
 Sample : L2756-04DL 10X  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 ETKW8DL

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:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM051320MA.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         | 3.028       | 19            | 24          | 30           | rVB      | 172016         | 249149        | 0.71%           | 0.234%        |
| 2         | 3.304       | 67            | 71          | 81           | rVB2     | 33942          | 53920         | 0.15%           | 0.051%        |
| 3         | 3.387       | 81            | 85          | 97           | rBV      | 1254314        | 1503002       | 4.28%           | 1.410%        |
| 4         | 3.763       | 143           | 149         | 162          | rBV      | 10451254       | 14389229      | 40.96%          | 13.498%       |
| 5         | 3.963       | 179           | 183         | 193          | rVB      | 90062          | 131231        | 0.37%           | 0.123%        |
| 6         | 5.045       | 361           | 367         | 376          | rBV      | 2013269        | 2869027       | 8.17%           | 2.691%        |
| 7         | 5.175       | 382           | 389         | 403          | rBV      | 2718092        | 5968984       | 16.99%          | 5.599%        |
| 8         | 5.387       | 420           | 425         | 432          | rBV      | 49010          | 77974         | 0.22%           | 0.073%        |
| 9         | 5.534       | 439           | 450         | 460          | rBV2     | 1662484        | 2883612       | 8.21%           | 2.705%        |
| 10        | 6.004       | 523           | 530         | 536          | rBV      | 47220          | 77570         | 0.22%           | 0.073%        |
| 11        | 6.357       | 582           | 590         | 596          | rBV      | 54551          | 86548         | 0.25%           | 0.081%        |
| 12        | 6.487       | 606           | 612         | 617          | rBV      | 115415         | 170383        | 0.49%           | 0.160%        |
| 13        | 6.592       | 621           | 630         | 635          | rVV      | 415560         | 587600        | 1.67%           | 0.551%        |
| 14        | 6.651       | 635           | 640         | 648          | rVV      | 247715         | 403178        | 1.15%           | 0.378%        |
| 15        | 6.728       | 648           | 653         | 659          | rVB      | 331656         | 476498        | 1.36%           | 0.447%        |
| 16        | 6.834       | 664           | 671         | 675          | rBV      | 53949          | 86934         | 0.25%           | 0.082%        |
| 17        | 6.887       | 675           | 680         | 684          | rVV      | 156483         | 240562        | 0.68%           | 0.226%        |
| 18        | 6.928       | 684           | 687         | 696          | rVB2     | 55982          | 86985         | 0.25%           | 0.082%        |
| 19        | 7.028       | 696           | 704         | 710          | rVB      | 55444          | 89418         | 0.25%           | 0.084%        |
| 20        | 7.092       | 710           | 715         | 718          | rBV      | 28079          | 40451         | 0.12%           | 0.038%        |
| 21        | 7.145       | 718           | 724         | 733          | rVV2     | 1324701        | 2147615       | 6.11%           | 2.015%        |
| 22        | 7.234       | 733           | 739         | 748          | rVB      | 137655         | 211932        | 0.60%           | 0.199%        |
| 23        | 7.481       | 775           | 781         | 787          | rBV      | 597601         | 924107        | 2.63%           | 0.867%        |
| 24        | 7.598       | 794           | 801         | 808          | rVB      | 285608         | 440943        | 1.26%           | 0.414%        |
| 25        | 7.669       | 808           | 813         | 819          | rVB      | 224550         | 328567        | 0.94%           | 0.308%        |
| 26        | 7.745       | 822           | 826         | 832          | rBV      | 29878          | 45512         | 0.13%           | 0.043%        |
| 27        | 7.851       | 832           | 844         | 849          | rBV2     | 185226         | 348386        | 0.99%           | 0.327%        |
| 28        | 7.916       | 849           | 855         | 863          | rVV      | 638812         | 1076898       | 3.07%           | 1.010%        |
| 29        | 8.010       | 863           | 871         | 881          | rVV      | 2661883        | 4206226       | 11.97%          | 3.946%        |
| 30        | 8.122       | 881           | 890         | 896          | rVV3     | 109969         | 224723        | 0.64%           | 0.211%        |
| 31        | 8.204       | 899           | 904         | 908          | rVV      | 49291          | 81413         | 0.23%           | 0.076%        |
| 32        | 8.263       | 908           | 914         | 918          | rVV2     | 44653          | 99654         | 0.28%           | 0.093%        |
| 33        | 8.304       | 919           | 921         | 928          | rVB2     | 38987          | 52567         | 0.15%           | 0.049%        |
| 34        | 8.433       | 936           | 943         | 948          | rBV3     | 41163          | 72694         | 0.21%           | 0.068%        |

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 Sampling : 1  
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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:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM051320MA.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |          |          |         |         |
|----|--------|------|------|------|------|----------|----------|---------|---------|
| 35 | 8.486  | 948  | 952  | 955  | rBV  | 37393    | 56413    | 0.16%   | 0.053%  |
| 36 | 8.581  | 964  | 968  | 973  | rBV  | 107355   | 178116   | 0.51%   | 0.167%  |
| 37 | 8.628  | 973  | 976  | 989  | rVB4 | 71010    | 175538   | 0.50%   | 0.165%  |
| 38 | 8.751  | 992  | 997  | 1004 | rBV  | 63279    | 129976   | 0.37%   | 0.122%  |
| 39 | 8.910  | 1019 | 1024 | 1028 | rVV3 | 25638    | 43231    | 0.12%   | 0.041%  |
| 40 | 9.057  | 1045 | 1049 | 1053 | rBV2 | 29033    | 47696    | 0.14%   | 0.045%  |
| 41 | 9.098  | 1053 | 1056 | 1061 | rVV  | 49678    | 72269    | 0.21%   | 0.068%  |
| 42 | 9.163  | 1061 | 1067 | 1068 | rVV  | 115180   | 162846   | 0.46%   | 0.153%  |
| 43 | 9.198  | 1068 | 1073 | 1078 | rVV  | 834494   | 1303670  | 3.71%   | 1.223%  |
| 44 | 9.239  | 1078 | 1080 | 1087 | rVV2 | 64628    | 88233    | 0.25%   | 0.083%  |
| 45 | 9.345  | 1089 | 1098 | 1105 | rVB2 | 53191    | 142219   | 0.40%   | 0.133%  |
| 46 | 9.486  | 1118 | 1122 | 1123 | rVV2 | 38942    | 51382    | 0.15%   | 0.048%  |
| 47 | 9.510  | 1123 | 1126 | 1135 | rVB  | 53639    | 90407    | 0.26%   | 0.085%  |
| 48 | 9.686  | 1145 | 1156 | 1162 | rBV3 | 472882   | 1156664  | 3.29%   | 1.085%  |
| 49 | 9.757  | 1162 | 1168 | 1173 | rBV  | 356587   | 798434   | 2.27%   | 0.749%  |
| 50 | 9.880  | 1183 | 1189 | 1197 | rBV2 | 94120    | 171079   | 0.49%   | 0.160%  |
| 51 | 9.951  | 1197 | 1201 | 1211 | rVB2 | 96999    | 168760   | 0.48%   | 0.158%  |
| 52 | 10.251 | 1244 | 1252 | 1254 | rBV  | 735176   | 1216041  | 3.46%   | 1.141%  |
| 53 | 10.310 | 1254 | 1262 | 1270 | rVV  | 19422797 | 35129060 | 100.00% | 32.953% |
| 54 | 10.410 | 1274 | 1279 | 1287 | rVB  | 350263   | 580021   | 1.65%   | 0.544%  |
| 55 | 10.998 | 1371 | 1379 | 1383 | rBV3 | 31658    | 63825    | 0.18%   | 0.060%  |
| 56 | 11.292 | 1425 | 1429 | 1432 | rVV2 | 32823    | 55047    | 0.16%   | 0.052%  |
| 57 | 11.351 | 1432 | 1439 | 1444 | rVV3 | 35693    | 96286    | 0.27%   | 0.090%  |
| 58 | 11.439 | 1450 | 1454 | 1463 | rVB  | 40451    | 76205    | 0.22%   | 0.071%  |
| 59 | 11.639 | 1483 | 1488 | 1497 | rVB4 | 42838    | 76354    | 0.22%   | 0.072%  |
| 60 | 11.769 | 1504 | 1510 | 1513 | rBV4 | 22760    | 44190    | 0.13%   | 0.041%  |
| 61 | 11.810 | 1513 | 1517 | 1525 | rVB2 | 67605    | 125151   | 0.36%   | 0.117%  |
| 62 | 11.916 | 1525 | 1535 | 1545 | rBV  | 2643529  | 4135422  | 11.77%  | 3.879%  |
| 63 | 12.033 | 1550 | 1555 | 1562 | rVB  | 47352    | 85570    | 0.24%   | 0.080%  |
| 64 | 12.139 | 1562 | 1573 | 1582 | rBV  | 1778684  | 2817423  | 8.02%   | 2.643%  |
| 65 | 12.521 | 1633 | 1638 | 1644 | rBV2 | 29816    | 52896    | 0.15%   | 0.050%  |
| 66 | 12.951 | 1703 | 1711 | 1721 | rBV  | 440940   | 642196   | 1.83%   | 0.602%  |
| 67 | 13.151 | 1740 | 1745 | 1748 | rBV  | 50259    | 83116    | 0.24%   | 0.078%  |
| 68 | 13.292 | 1758 | 1769 | 1777 | rBV3 | 88056    | 243262   | 0.69%   | 0.228%  |
| 69 | 13.445 | 1788 | 1795 | 1801 | rBV2 | 187085   | 336174   | 0.96%   | 0.315%  |
| 70 | 13.498 | 1801 | 1804 | 1809 | rVV  | 98439    | 153055   | 0.44%   | 0.144%  |
| 71 | 13.551 | 1809 | 1813 | 1816 | rVV  | 166643   | 261888   | 0.75%   | 0.246%  |

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## Integration Parameters: LSCINT.P

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 Stop Thrs : 0

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Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM051320MA.M  
 Title : SVOA CALIBRATION

|     |        |      |      |      |      |         |         |       |        |
|-----|--------|------|------|------|------|---------|---------|-------|--------|
| 72  | 13.580 | 1816 | 1818 | 1826 | rVB2 | 108259  | 189695  | 0.54% | 0.178% |
| 73  | 13.686 | 1830 | 1836 | 1839 | rBV  | 96465   | 138898  | 0.40% | 0.130% |
| 74  | 13.845 | 1852 | 1863 | 1871 | rBV2 | 938530  | 1610077 | 4.58% | 1.510% |
| 75  | 14.127 | 1899 | 1911 | 1917 | rBV  | 1165576 | 1754558 | 4.99% | 1.646% |
| 76  | 14.192 | 1917 | 1922 | 1926 | rVV  | 129087  | 166261  | 0.47% | 0.156% |
| 77  | 14.539 | 1976 | 1981 | 1987 | rVB4 | 28841   | 57784   | 0.16% | 0.054% |
| 78  | 14.633 | 1987 | 1997 | 2003 | rBV2 | 33993   | 66642   | 0.19% | 0.063% |
| 79  | 15.004 | 2056 | 2060 | 2068 | rVB  | 30599   | 48077   | 0.14% | 0.045% |
| 80  | 15.127 | 2076 | 2081 | 2086 | rBV  | 285692  | 381338  | 1.09% | 0.358% |
| 81  | 15.180 | 2086 | 2090 | 2098 | rVB2 | 395646  | 588513  | 1.68% | 0.552% |
| 82  | 15.262 | 2100 | 2104 | 2109 | rBV3 | 39131   | 64006   | 0.18% | 0.060% |
| 83  | 15.386 | 2121 | 2125 | 2138 | rVB4 | 18135   | 48496   | 0.14% | 0.045% |
| 84  | 15.598 | 2155 | 2161 | 2166 | rBV2 | 57459   | 82923   | 0.24% | 0.078% |
| 85  | 16.698 | 2339 | 2348 | 2352 | rBV  | 59255   | 93963   | 0.27% | 0.088% |
| 86  | 16.880 | 2372 | 2379 | 2382 | rBV  | 1396644 | 1991202 | 5.67% | 1.868% |
| 87  | 16.915 | 2382 | 2385 | 2391 | rVV  | 795270  | 1098426 | 3.13% | 1.030% |
| 88  | 16.974 | 2391 | 2395 | 2399 | rVV2 | 271318  | 403324  | 1.15% | 0.378% |
| 89  | 17.009 | 2399 | 2401 | 2408 | rVV  | 112117  | 134042  | 0.38% | 0.126% |
| 90  | 17.080 | 2409 | 2413 | 2418 | rVB  | 33703   | 44133   | 0.13% | 0.041% |
| 91  | 17.756 | 2520 | 2528 | 2532 | rBV  | 30216   | 49031   | 0.14% | 0.046% |
| 92  | 17.803 | 2532 | 2536 | 2539 | rBV2 | 34787   | 43540   | 0.12% | 0.041% |
| 93  | 17.945 | 2555 | 2560 | 2564 | rBV3 | 63249   | 99308   | 0.28% | 0.093% |
| 94  | 18.945 | 2725 | 2730 | 2735 | rVB  | 80243   | 107663  | 0.31% | 0.101% |
| 95  | 19.274 | 2781 | 2786 | 2789 | rBV  | 324553  | 444685  | 1.27% | 0.417% |
| 96  | 19.303 | 2789 | 2791 | 2795 | rVV  | 122369  | 149779  | 0.43% | 0.141% |
| 97  | 21.027 | 3081 | 3084 | 3087 | rBV2 | 45335   | 49299   | 0.14% | 0.046% |
| 98  | 21.074 | 3087 | 3092 | 3097 | rBV  | 1804738 | 2304459 | 6.56% | 2.162% |
| 99  | 23.062 | 3426 | 3430 | 3435 | rVB  | 183837  | 290285  | 0.83% | 0.272% |
| 100 | 23.197 | 3447 | 3453 | 3466 | rVB2 | 1236573 | 2258622 | 6.43% | 2.119% |

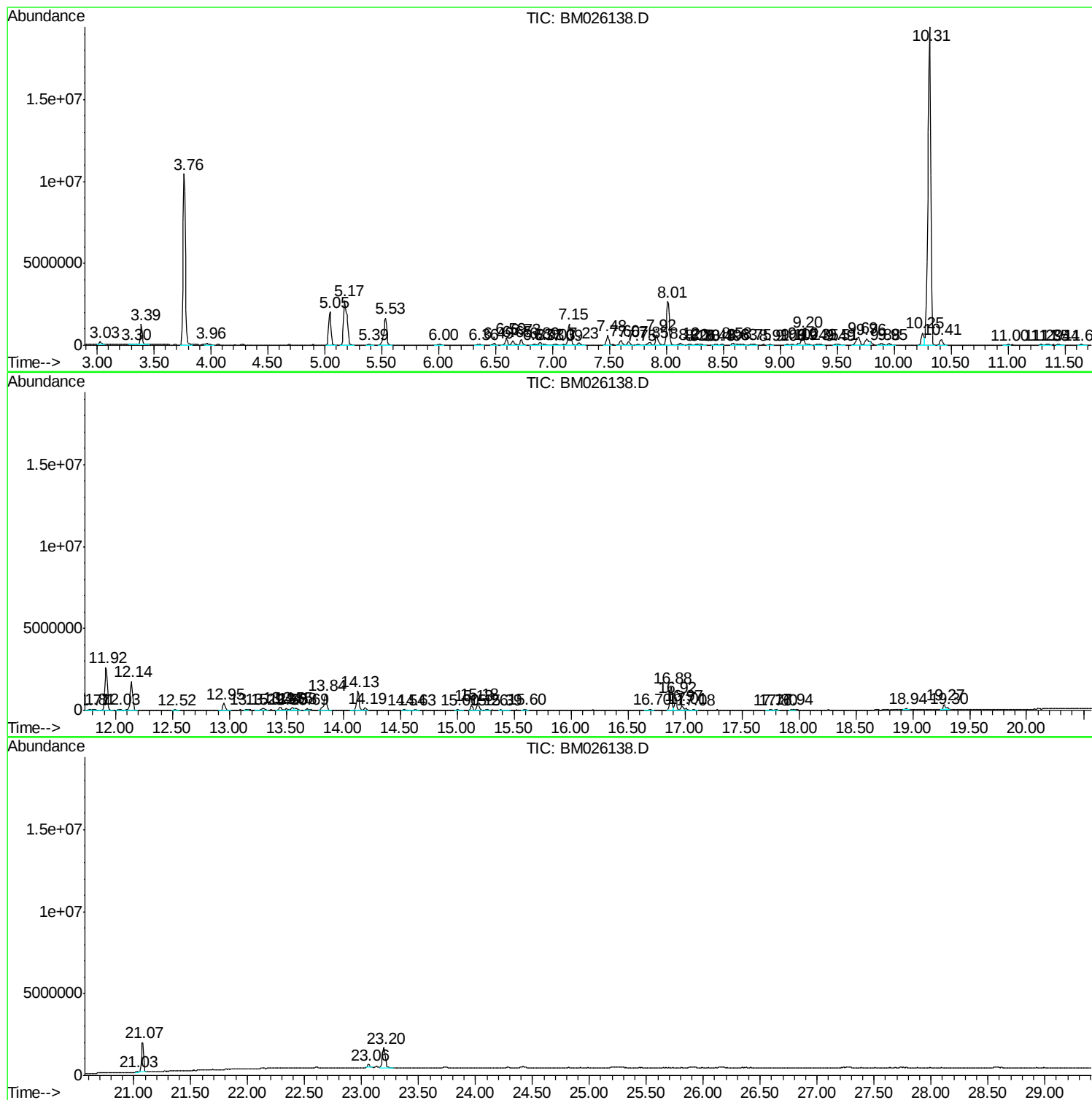
Sum of corrected areas: 106602636

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM051320MA.M  
Quant Title : SVOA CALIBRATION

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



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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM051320MA.M  
Quant Title : SVOA CALIBRATION

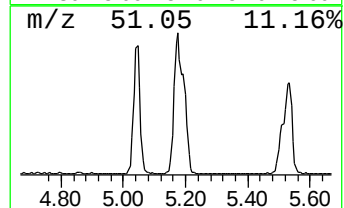
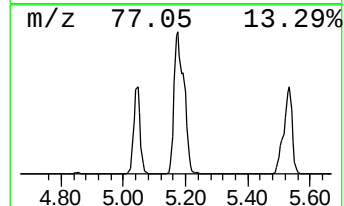
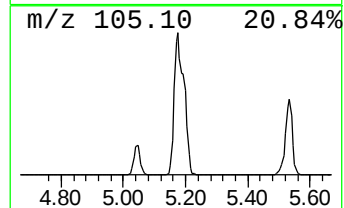
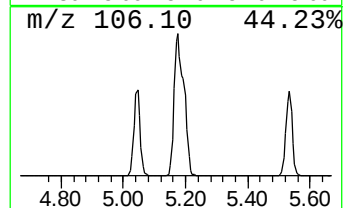
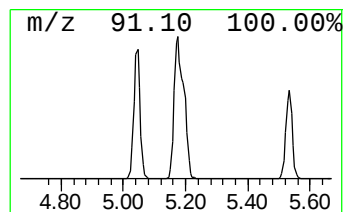
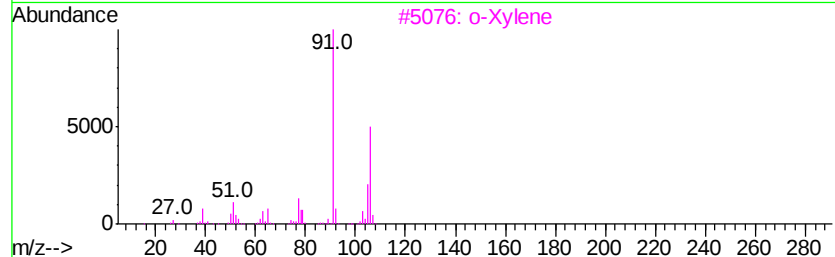
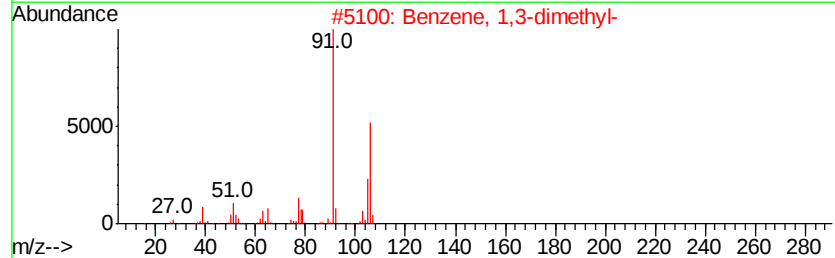
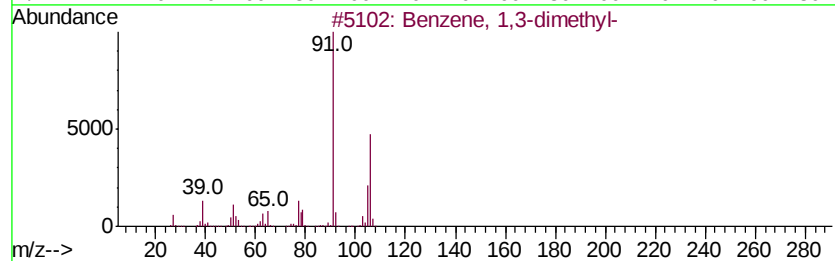
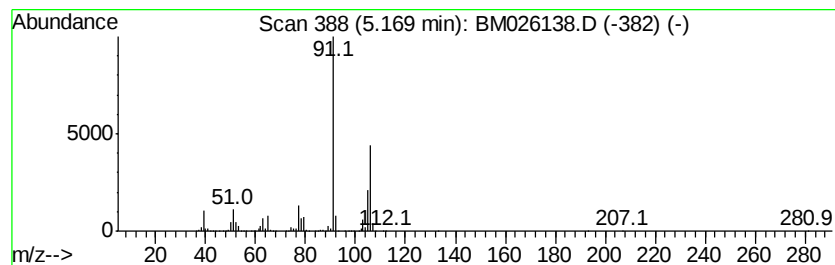
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 Benzene, 1,3-dimethyl- Concentration Rank 2

| R.T. | EstConc      | Area    | Relative to ISTD       | R.T. |
|------|--------------|---------|------------------------|------|
| 5.17 | 129.18 ng/ul | 5968980 | 1,4-Dichlorobenzene-d4 | 7.48 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,3-dimethyl- | 106 | C8H10   | 000108-38-3 | 97   |
| 2    |    | Benzene, 1,3-dimethyl- | 106 | C8H10   | 000108-38-3 | 97   |
| 3    |    | o-Xylene               | 106 | C8H10   | 000095-47-6 | 97   |
| 4    |    | o-Xylene               | 106 | C8H10   | 000095-47-6 | 97   |
| 5    |    | p-Xylene               | 106 | C8H10   | 000106-42-3 | 97   |



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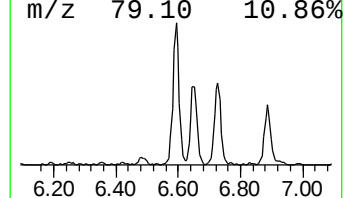
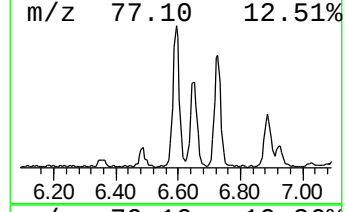
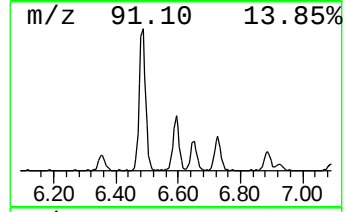
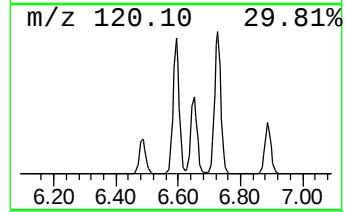
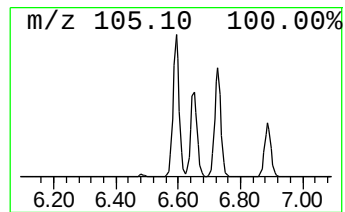
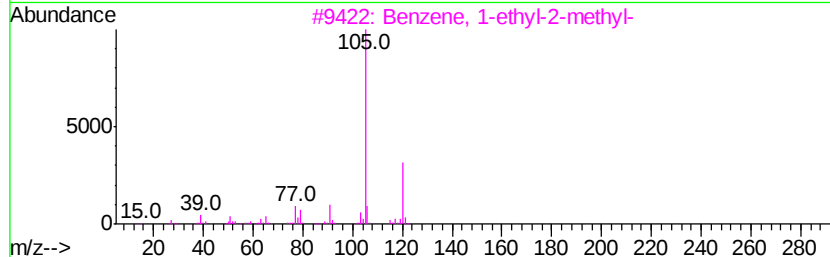
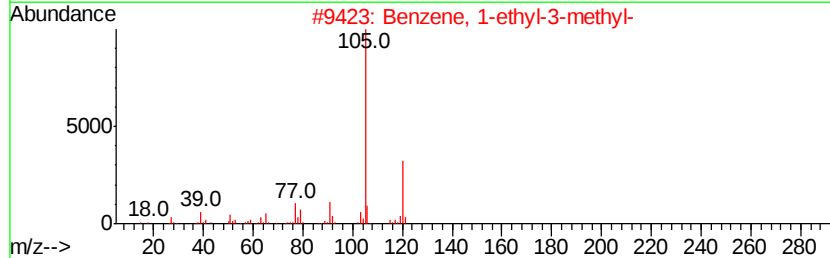
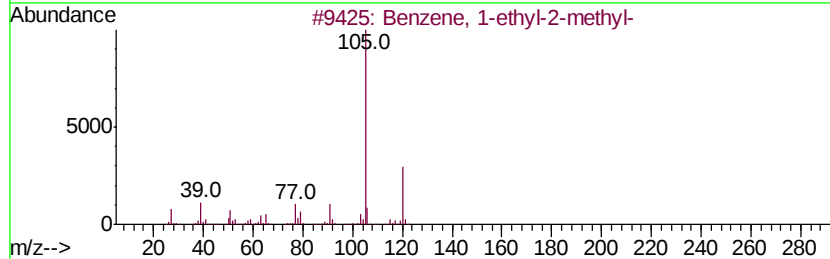
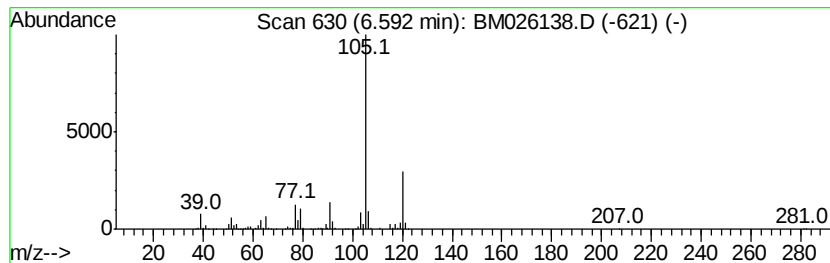
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM051320MA.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 9 Benzene, 1-ethyl-2-methyl- Concentration Rank 11

| R.T. | EstConc     | Area   | Relative to ISTD       | R.T. |
|------|-------------|--------|------------------------|------|
| 6.59 | 12.72 ng/ul | 587600 | 1,4-Dichlorobenzene-d4 | 7.48 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 95   |
| 2    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 95   |
| 3    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 94   |
| 4    |    |   | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 91   |
| 5    |    |   | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052720\  
Data File : BM026138.D  
Acq On : 27 May 2020 10:35  
Operator : MA/SJ  
Sample : L2756-04DL 10X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
ETKW8DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM051320MA.M  
Quant Title : SVOA CALIBRATION

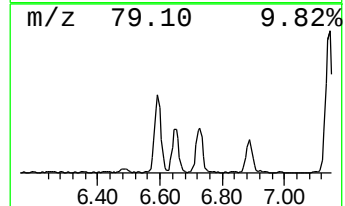
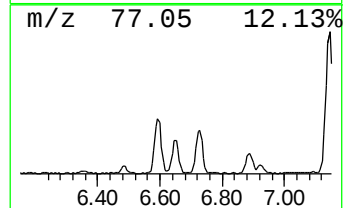
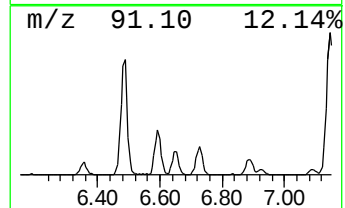
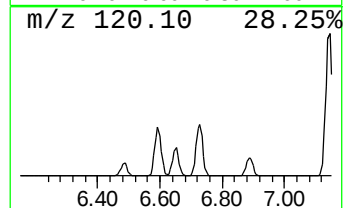
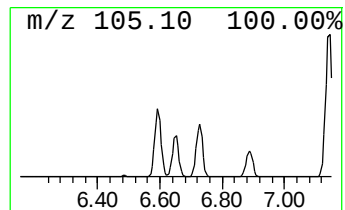
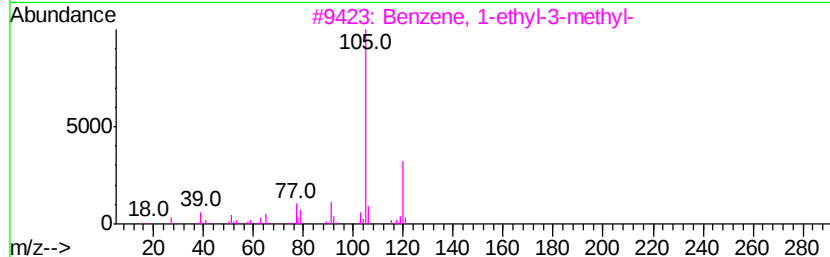
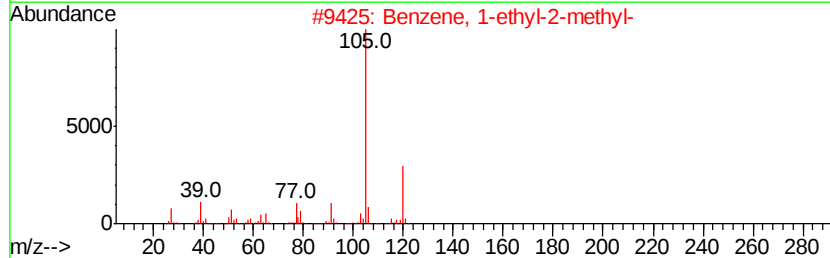
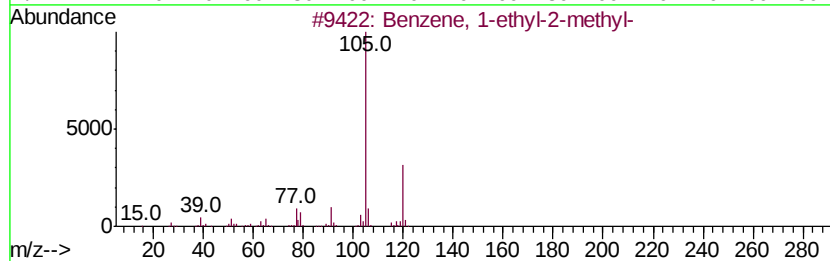
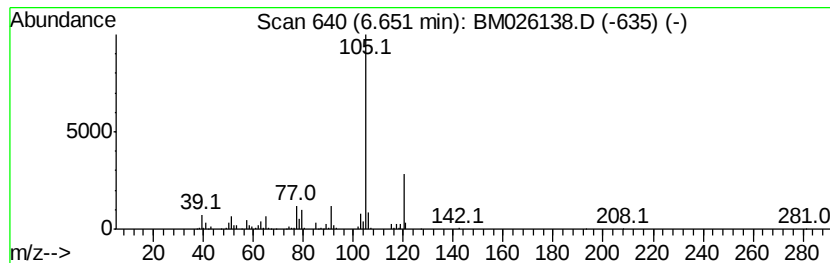
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 Benzene, 1-ethyl-3-methyl- Concentration Rank 15

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 6.65 | 8.73 ng/ul | 403178 | 1,4-Dichlorobenzene-d4 | 7.48 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 94   |
| 2    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 94   |
| 3    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 91   |
| 4    |    |   | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 91   |
| 5    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052720\  
Data File : BM026138.D  
Acq On : 27 May 2020 10:35  
Operator : MA/SJ  
Sample : L2756-04DL 10X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
ETKW8DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM051320MA.M  
Quant Title : SVOA CALIBRATION

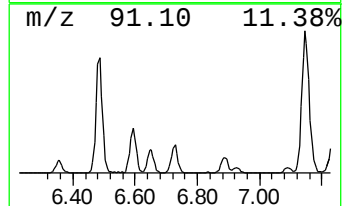
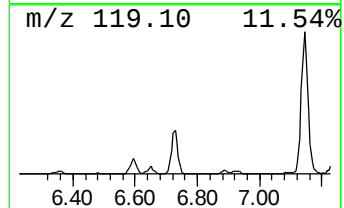
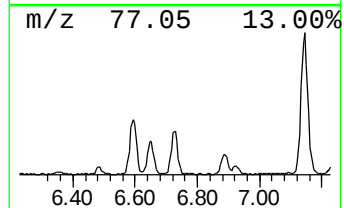
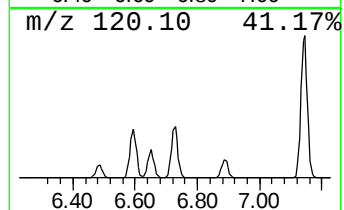
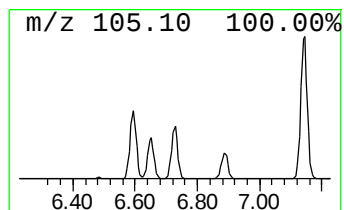
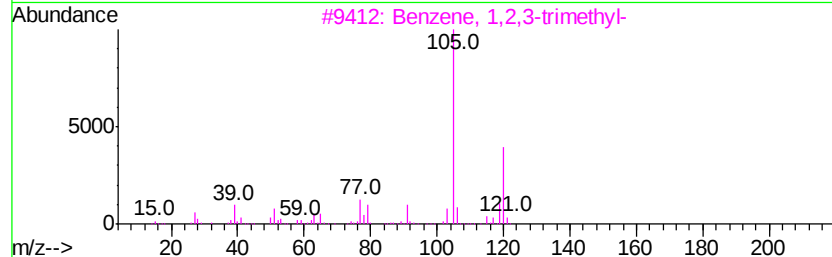
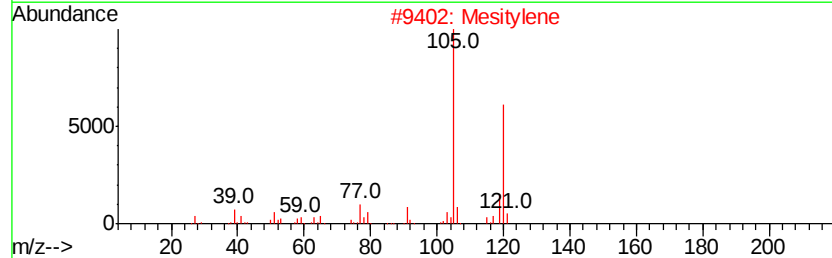
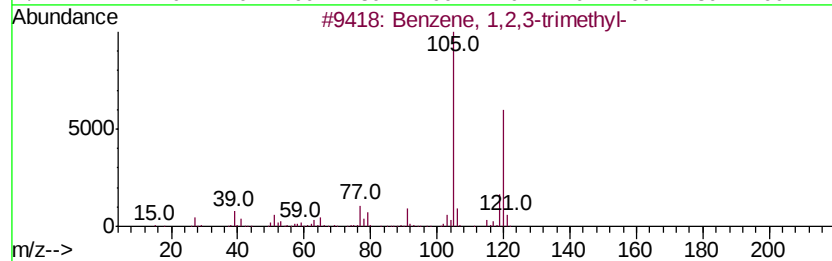
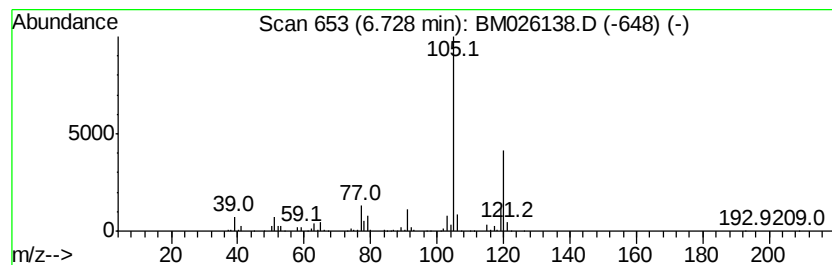
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 Benzene, 1,2,3-trimethyl- Concentration Rank 12

| R.T. | EstConc     | Area   | Relative to ISTD       | R.T. |
|------|-------------|--------|------------------------|------|
| 6.73 | 10.31 ng/ul | 476498 | 1,4-Dichlorobenzene-d4 | 7.48 |

| Hit# | of | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2,3-trimethyl- | 120 | C9H12   | 000526-73-8 | 97   |
| 2    |    | Mesitylene                | 120 | C9H12   | 000108-67-8 | 95   |
| 3    |    | Benzene, 1,2,3-trimethyl- | 120 | C9H12   | 000526-73-8 | 94   |
| 4    |    | Benzene, 1,2,4-trimethyl- | 120 | C9H12   | 000095-63-6 | 94   |
| 5    |    | Mesitylene                | 120 | C9H12   | 000108-67-8 | 94   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052720\  
Data File : BM026138.D  
Acq On : 27 May 2020 10:35  
Operator : MA/SJ  
Sample : L2756-04DL 10X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

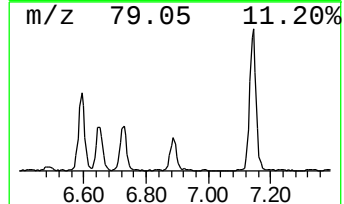
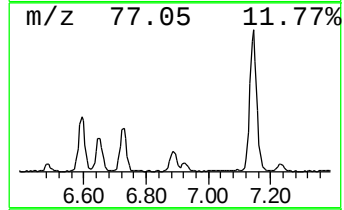
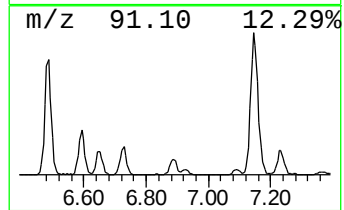
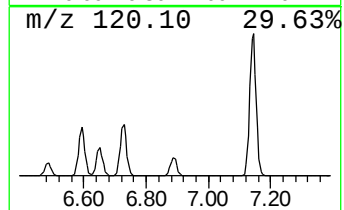
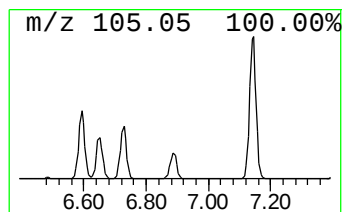
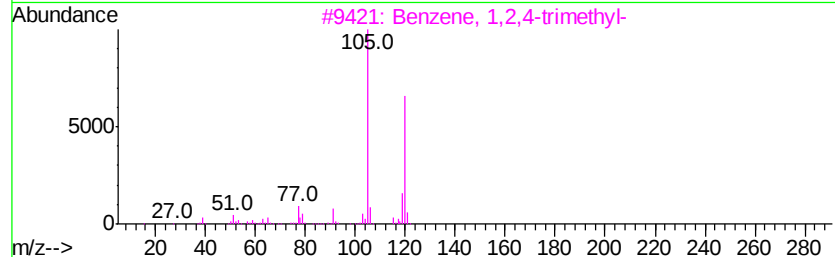
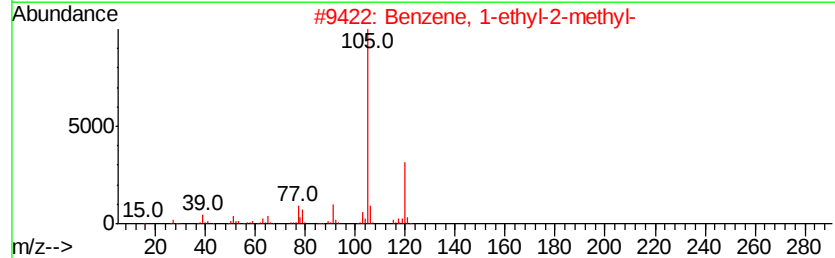
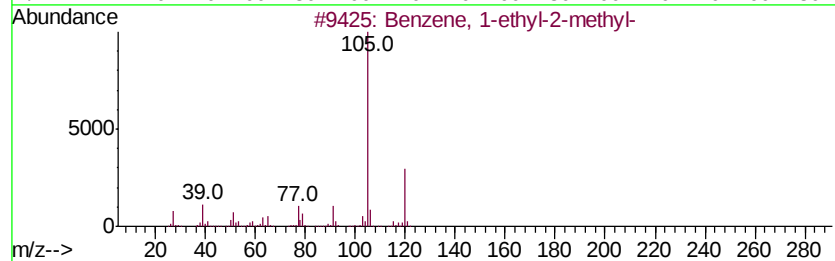
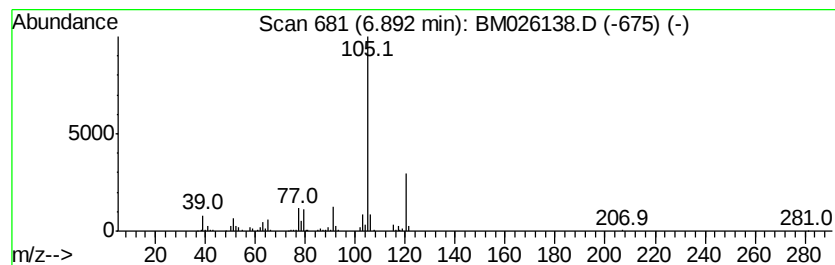
Instrument :  
BNA\_M  
ClientSampleId :  
ETKW8DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM051320MA.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 12 Benzene, 1,2,4-trimethyl- Concentration Rank 19

| R.T.    | EstConc    | Area                       | Relative to ISTD       | R.T.           |
|---------|------------|----------------------------|------------------------|----------------|
| 6.89    | 5.21 ng/ul | 240562                     | 1,4-Dichlorobenzene-d4 | 7.48           |
| Hit# of | 5          | Tentative ID               | MW MolForm             | CAS# Qual      |
| 1       |            | Benzene, 1-ethyl-2-methyl- | 120 C9H12              | 000611-14-3 95 |
| 2       |            | Benzene, 1-ethyl-2-methyl- | 120 C9H12              | 000611-14-3 94 |
| 3       |            | Benzene, 1,2,4-trimethyl-  | 120 C9H12              | 000095-63-6 91 |
| 4       |            | Benzene, 1-ethyl-3-methyl- | 120 C9H12              | 000620-14-4 91 |
| 5       |            | Mesitylene                 | 120 C9H12              | 000108-67-8 91 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052720\  
Data File : BM026138.D  
Acq On : 27 May 2020 10:35  
Operator : MA/SJ  
Sample : L2756-04DL 10X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
ETKW8DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM051320MA.M  
Quant Title : SVOA CALIBRATION

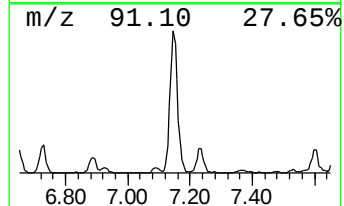
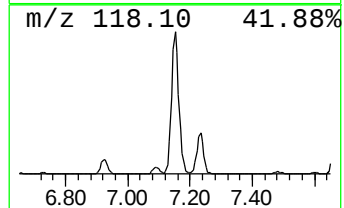
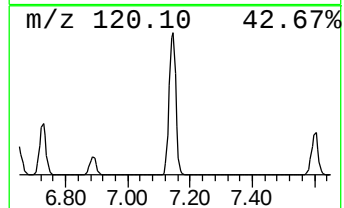
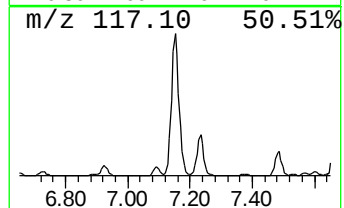
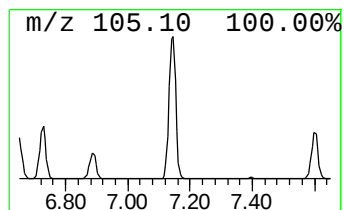
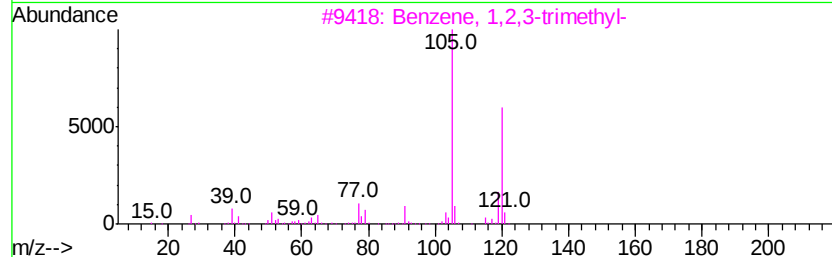
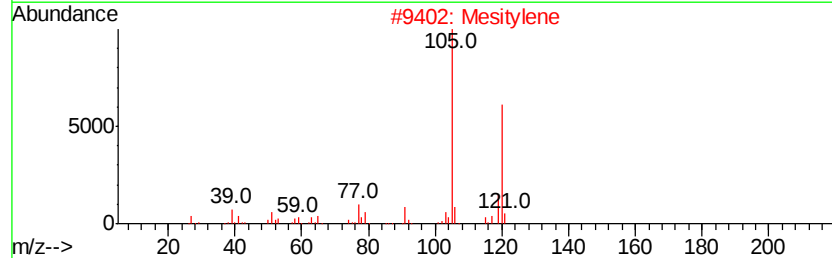
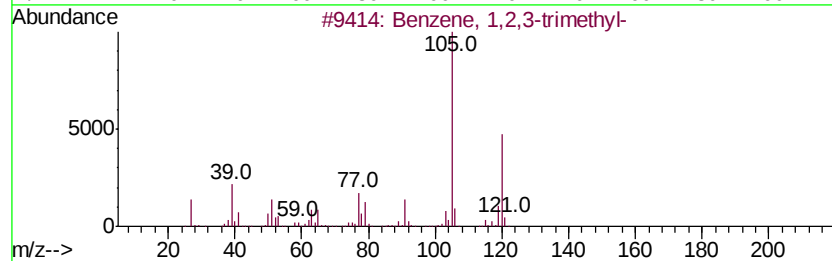
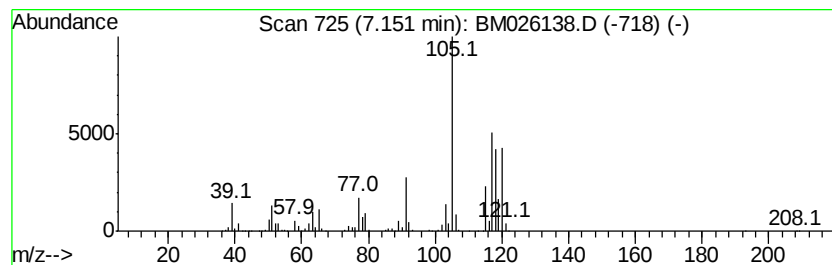
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 13 Mesitylene Concentration Rank 6

| R.T. | EstConc     | Area    | Relative to ISTD       | R.T. |
|------|-------------|---------|------------------------|------|
| 7.15 | 46.48 ng/ul | 2147620 | 1,4-Dichlorobenzene-d4 | 7.48 |

| Hit# | of | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2,3-trimethyl- | 120 | C9H12   | 000526-73-8 | 91   |
| 2    |    | Mesitylene                | 120 | C9H12   | 000108-67-8 | 64   |
| 3    |    | Benzene, 1,2,3-trimethyl- | 120 | C9H12   | 000526-73-8 | 60   |
| 4    |    | 2,4-Nonadiyne             | 120 | C9H12   | 063621-15-8 | 60   |
| 5    |    | Mesitylene                | 120 | C9H12   | 000108-67-8 | 60   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052720\  
Data File : BM026138.D  
Acq On : 27 May 2020 10:35  
Operator : MA/SJ  
Sample : L2756-04DL 10X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
ETKW8DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM051320MA.M  
Quant Title : SVOA CALIBRATION

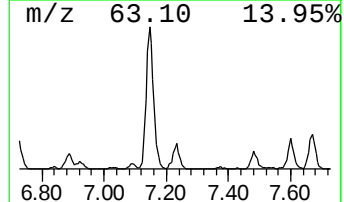
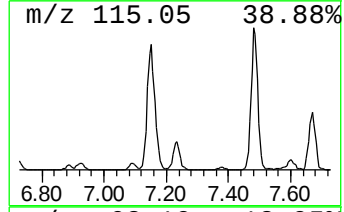
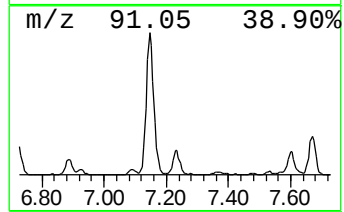
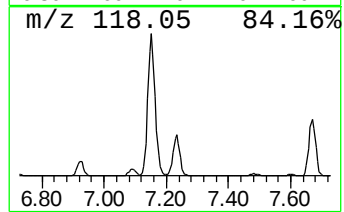
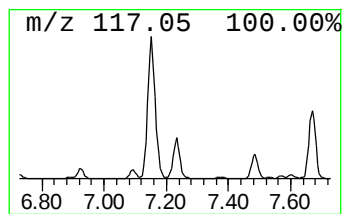
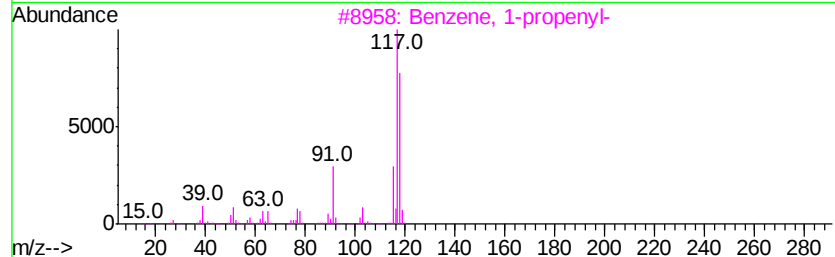
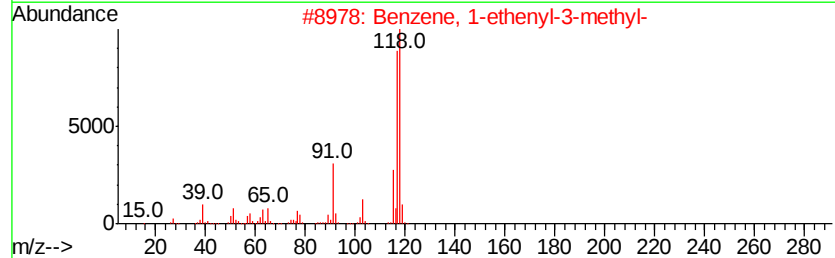
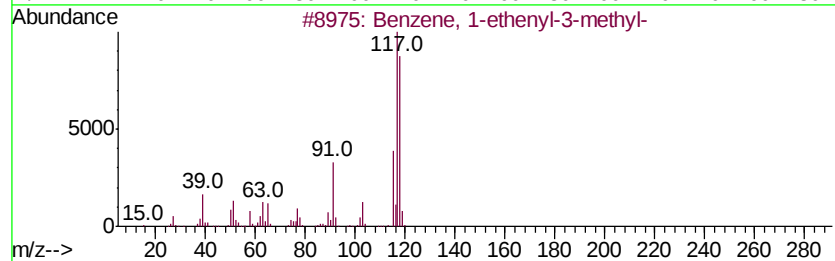
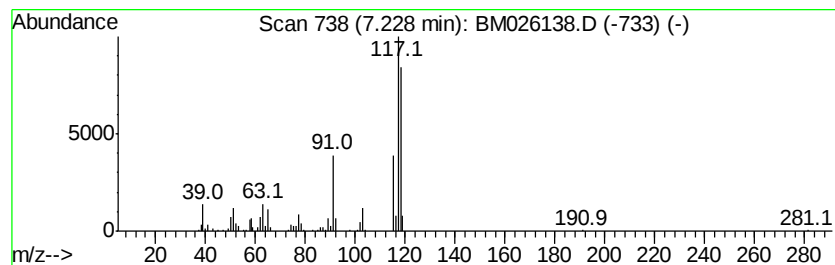
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 Benzene, 1-ethenyl-3-methyl- Concentration Rank 21

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 7.23 | 4.59 ng/ul | 211932 | 1,4-Dichlorobenzene-d4 | 7.48 |

| Hit# | of | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-ethenyl-3-methyl- | 118 | C9H10   | 000100-80-1 | 96   |
| 2    |    | Benzene, 1-ethenyl-3-methyl- | 118 | C9H10   | 000100-80-1 | 96   |
| 3    |    | Benzene, 1-propenyl-         | 118 | C9H10   | 000637-50-3 | 94   |
| 4    |    | Benzene, 1-ethenyl-3-methyl- | 118 | C9H10   | 000100-80-1 | 94   |
| 5    |    | Benzene, 1-ethenyl-2-methyl- | 118 | C9H10   | 000611-15-4 | 94   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052720\  
Data File : BM026138.D  
Acq On : 27 May 2020 10:35  
Operator : MA/SJ  
Sample : L2756-04DL 10X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
ETKW8DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM051320MA.M  
Quant Title : SVOA CALIBRATION

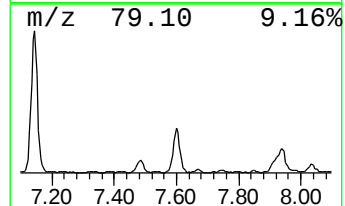
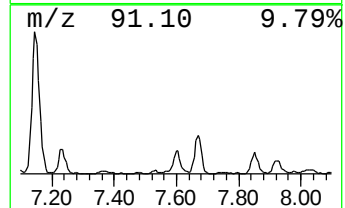
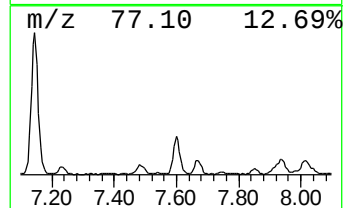
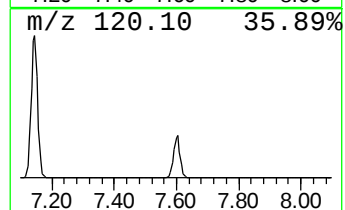
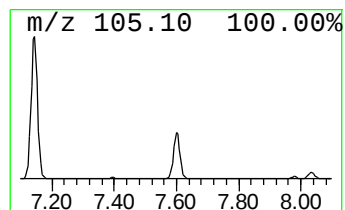
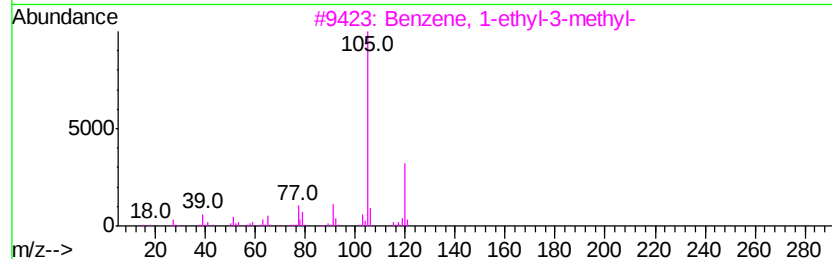
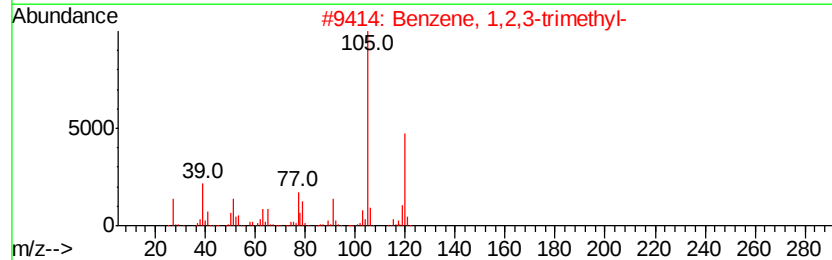
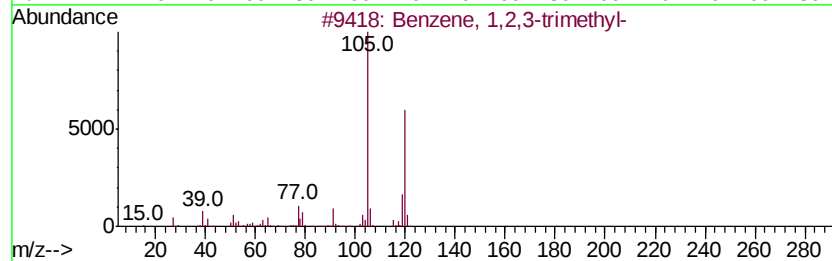
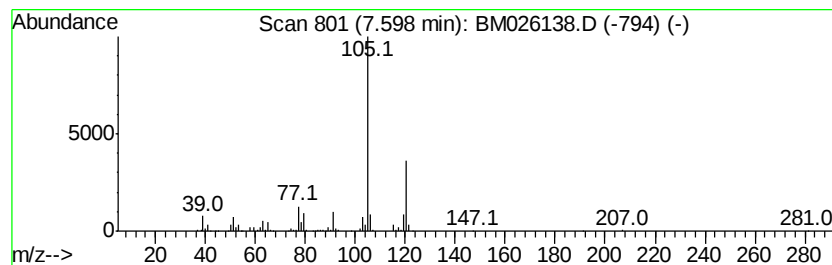
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 unknown-01 Concentration Rank 13

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 7.60 | 9.54 ng/ul | 440943 | 1,4-Dichlorobenzene-d4 | 7.48 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 94   |
| 2    |    |   | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 93   |
| 3    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 91   |
| 4    |    |   | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 91   |
| 5    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052720\  
Data File : BM026138.D  
Acq On : 27 May 2020 10:35  
Operator : MA/SJ  
Sample : L2756-04DL 10X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
ETKW8DL

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Quant Title : SVOA CALIBRATION

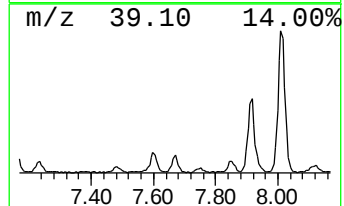
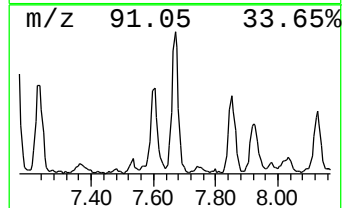
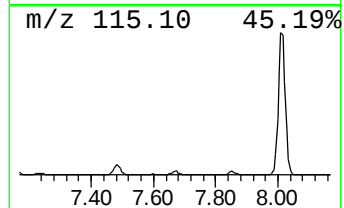
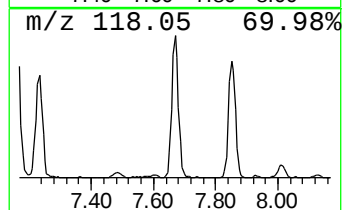
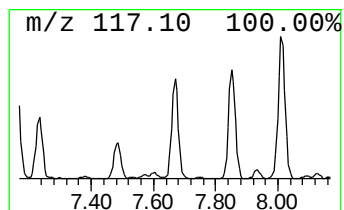
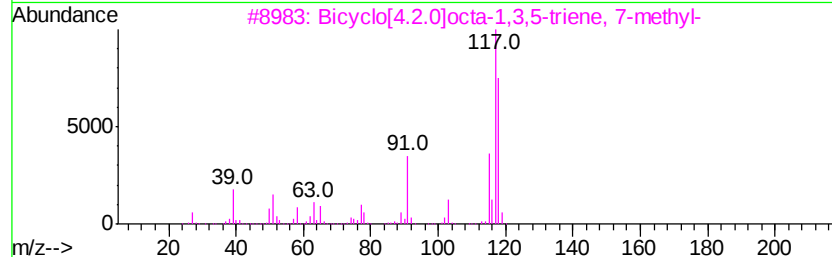
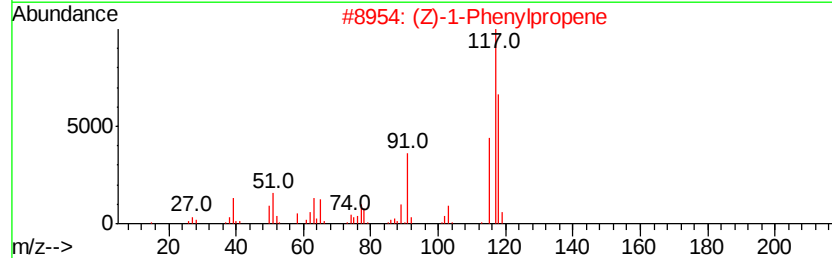
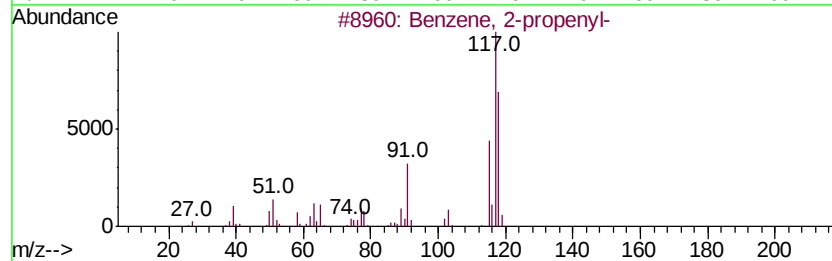
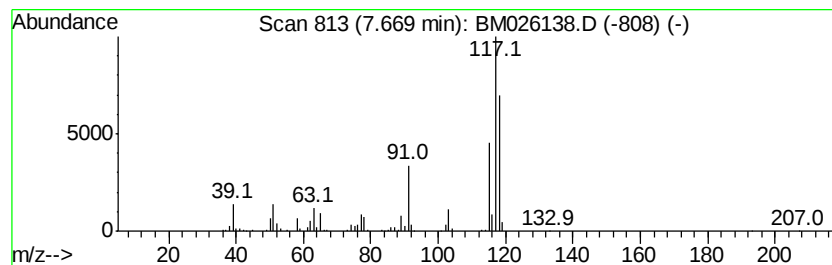
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 Benzene, 2-propenyl- Concentration Rank 17

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 7.67 | 7.11 ng/ul | 328567 | 1,4-Dichlorobenzene-d4 | 7.48 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 2-propenyl-                | 118 | C9H10   | 000300-57-2 | 96   |
| 2    |    |   | (Z)-1-Phenylpropene                 | 118 | C9H10   | 000766-90-5 | 96   |
| 3    |    |   | Bicyclo[4.2.0]octa-1,3,5-triene,... | 118 | C9H10   | 055337-80-9 | 96   |
| 4    |    |   | Benzene, 1-propenyl-                | 118 | C9H10   | 000637-50-3 | 95   |
| 5    |    |   | Benzene, cyclopropyl-               | 118 | C9H10   | 000873-49-4 | 93   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052720\  
Data File : BM026138.D  
Acq On : 27 May 2020 10:35  
Operator : MA/SJ  
Sample : L2756-04DL 10X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
ETKW8DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM051320MA.M  
Quant Title : SVOA CALIBRATION

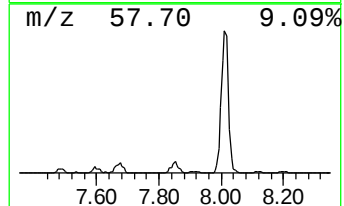
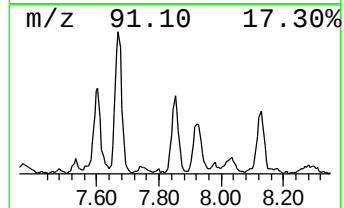
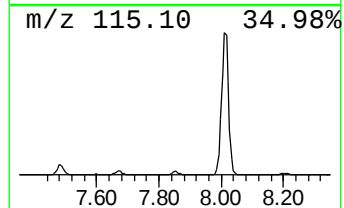
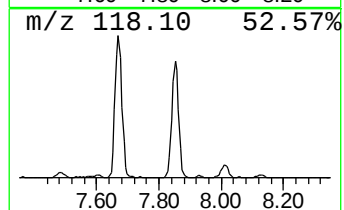
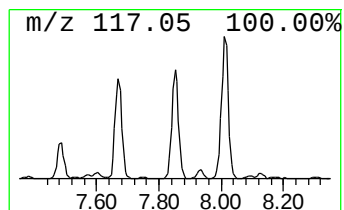
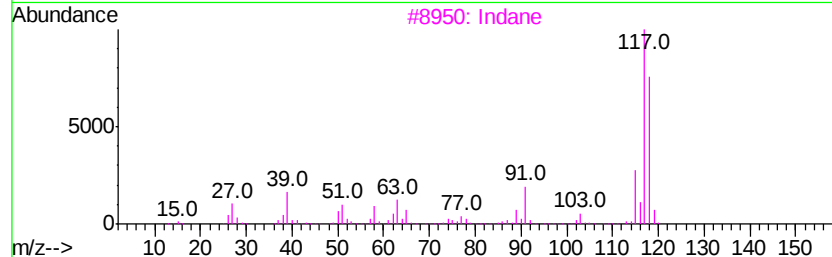
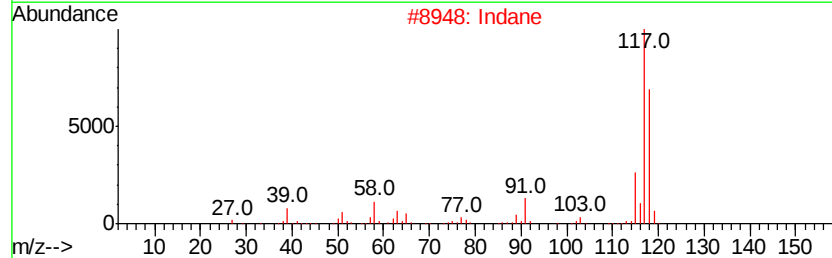
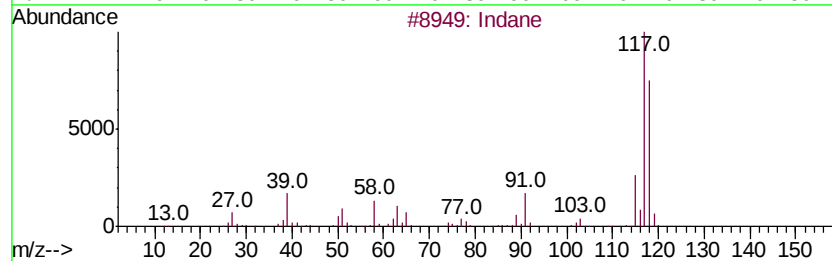
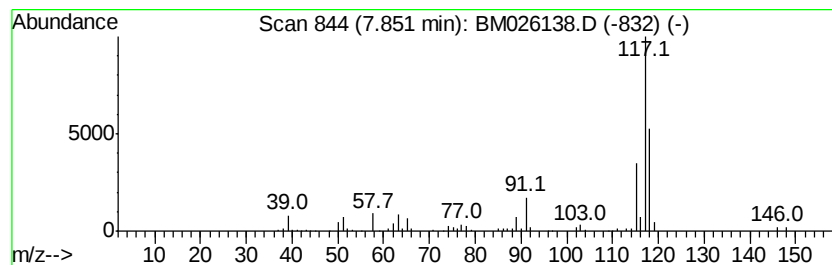
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 17 Indane Concentration Rank 16

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 7.85 | 7.54 ng/ul | 348386 | 1,4-Dichlorobenzene-d4 | 7.48 |

| Hit# | of | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------|-----|---------|-------------|------|
| 1    | 5  | Indane               | 118 | C9H10   | 000496-11-7 | 96   |
| 2    |    | Indane               | 118 | C9H10   | 000496-11-7 | 93   |
| 3    |    | Indane               | 118 | C9H10   | 000496-11-7 | 90   |
| 4    |    | Indane               | 118 | C9H10   | 000496-11-7 | 87   |
| 5    |    | Benzene, 2-propenyl- | 118 | C9H10   | 000300-57-2 | 81   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052720\  
Data File : BM026138.D  
Acq On : 27 May 2020 10:35  
Operator : MA/SJ  
Sample : L2756-04DL 10X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
ETKW8DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM051320MA.M  
Quant Title : SVOA CALIBRATION

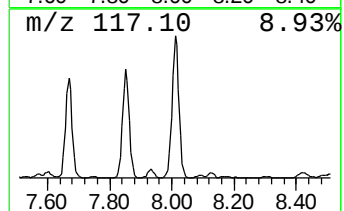
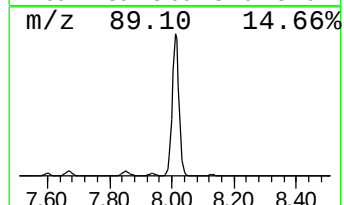
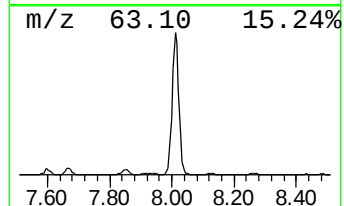
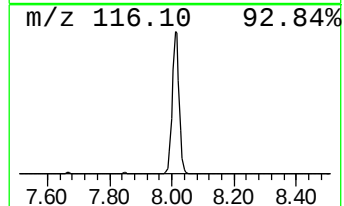
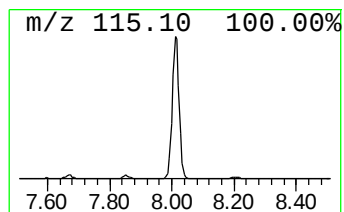
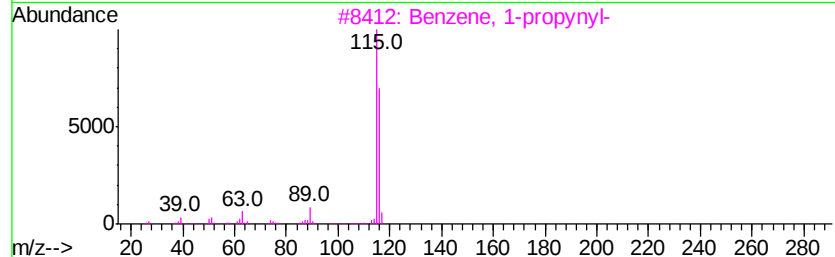
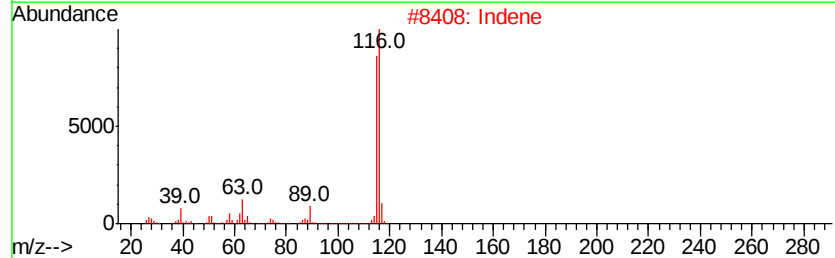
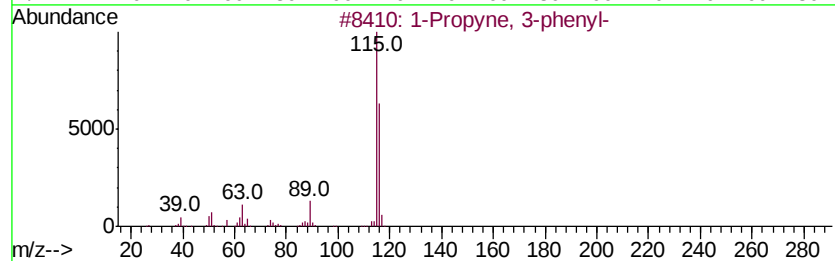
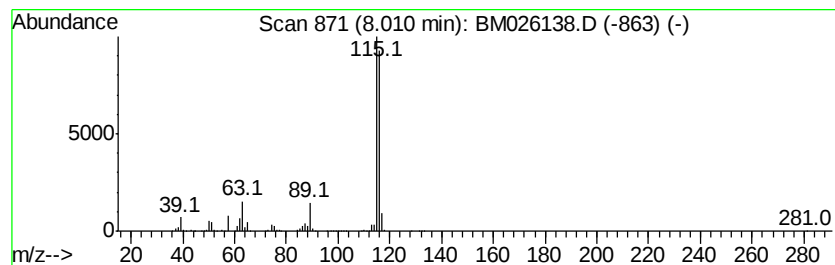
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 18 1-Propyne, 3-phenyl- Concentration Rank 3

| R.T. | EstConc     | Area    | Relative to ISTD       | R.T. |
|------|-------------|---------|------------------------|------|
| 8.01 | 91.03 ng/ul | 4206230 | 1,4-Dichlorobenzene-d4 | 7.48 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | 1-Propyne, 3-phenyl-    | 116 | C9H8    | 010147-11-2 | 95   |
| 2    |    | Indene                  | 116 | C9H8    | 000095-13-6 | 94   |
| 3    |    | Benzene, 1-propynyl-    | 116 | C9H8    | 000673-32-5 | 91   |
| 4    |    | 3-Methylphenylacetylene | 116 | C9H8    | 000766-82-5 | 91   |
| 5    |    | Indene                  | 116 | C9H8    | 000095-13-6 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052720\  
Data File : BM026138.D  
Acq On : 27 May 2020 10:35  
Operator : MA/SJ  
Sample : L2756-04DL 10X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
ETKW8DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM051320MA.M  
Quant Title : SVOA CALIBRATION

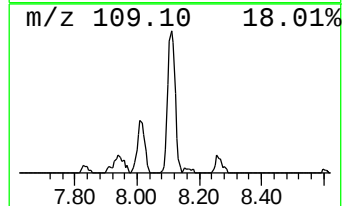
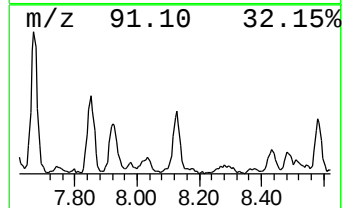
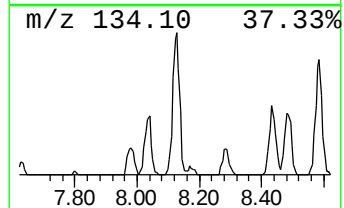
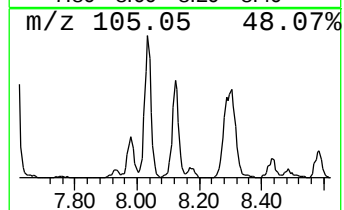
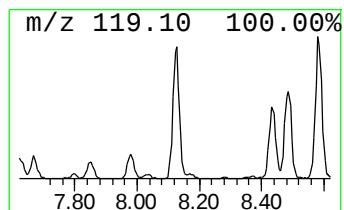
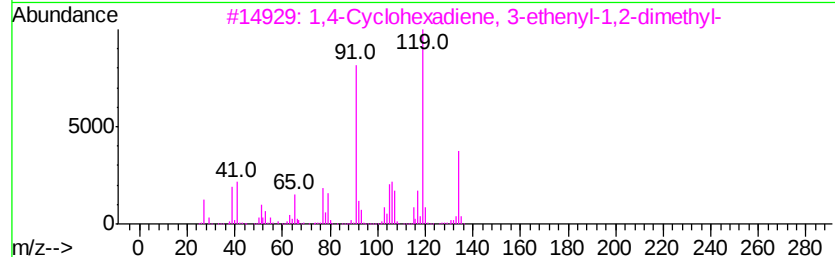
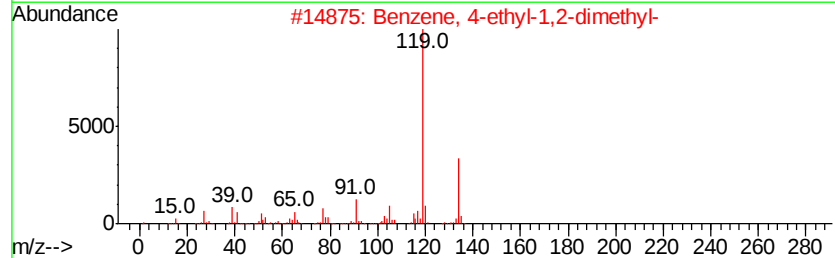
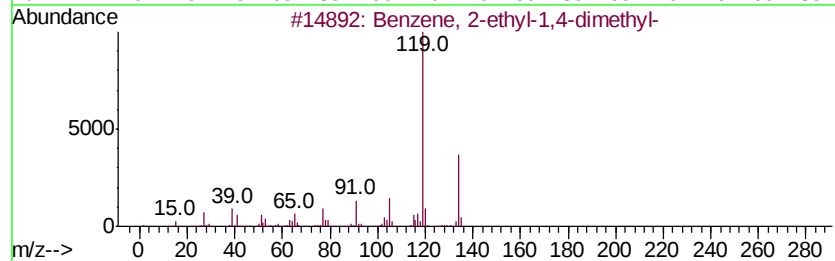
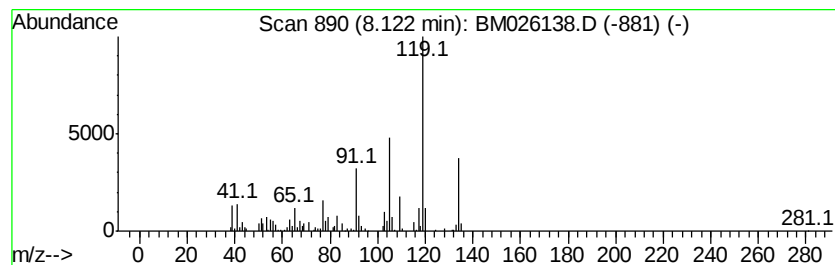
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 20

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 8.12 | 4.86 ng/ul | 224723 | 1,4-Dichlorobenzene-d4 | 7.48 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 2-ethyl-1,4-dimethyl-      | 134 | C10H14  | 001758-88-9 | 91   |
| 2    |    |   | Benzene, 4-ethyl-1,2-dimethyl-      | 134 | C10H14  | 000934-80-5 | 90   |
| 3    |    |   | 1,4-Cyclohexadiene, 3-ethenyl-1,... | 134 | C10H14  | 062338-57-2 | 90   |
| 4    |    |   | Benzene, 1-ethyl-3,5-dimethyl-      | 134 | C10H14  | 000934-74-7 | 90   |
| 5    |    |   | Benzene, 1-ethyl-2,3-dimethyl-      | 134 | C10H14  | 000933-98-2 | 90   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052720\  
Data File : BM026138.D  
Acq On : 27 May 2020 10:35  
Operator : MA/SJ  
Sample : L2756-04DL 10X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
ETKW8DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM051320MA.M  
Quant Title : SVOA CALIBRATION

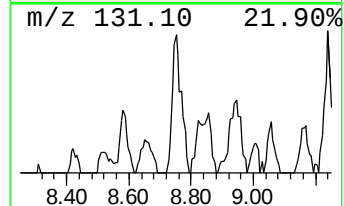
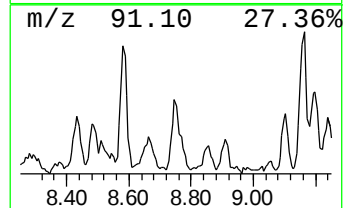
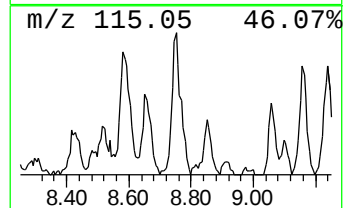
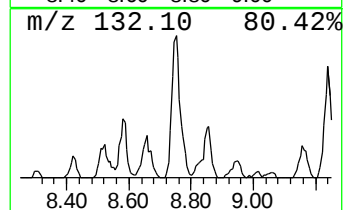
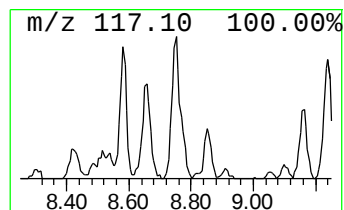
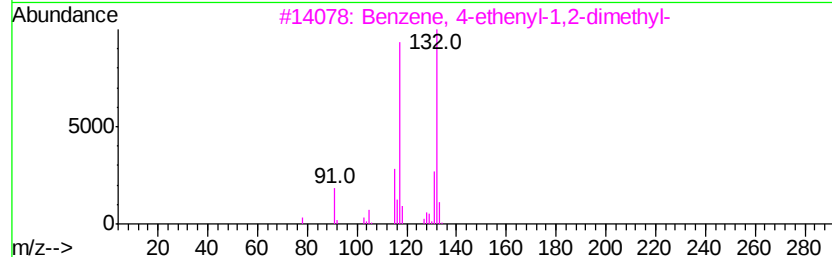
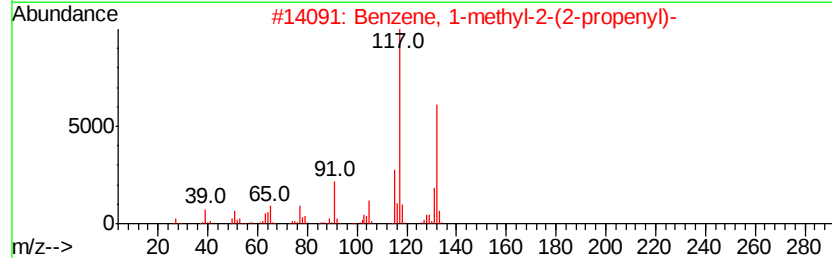
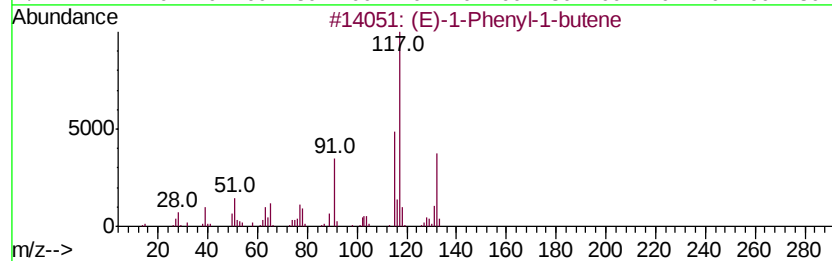
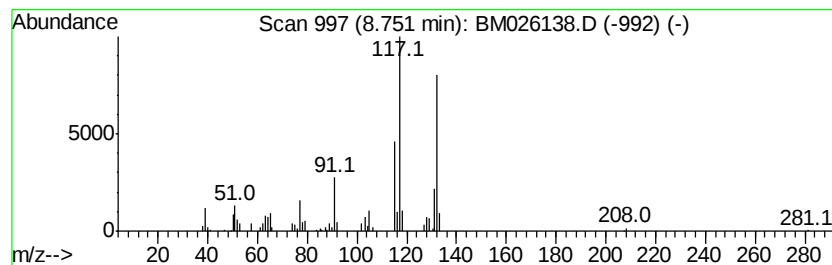
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 20 (E)-1-Phenyl-1-butene Concentration Rank 25

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 8.75 | 2.81 ng/ul | 129976 | 1,4-Dichlorobenzene-d4 | 7.48 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | (E)-1-Phenyl-1-butene               | 132 | C10H12  | 001005-64-7 | 95   |
| 2    |    |   | Benzene, 1-methyl-2-(2-propenyl)-   | 132 | C10H12  | 001587-04-8 | 95   |
| 3    |    |   | Benzene, 4-ethenyl-1,2-dimethyl-    | 132 | C10H12  | 027831-13-6 | 95   |
| 4    |    |   | Benzene, (2-methyl-1-propenyl)-     | 132 | C10H12  | 000768-49-0 | 94   |
| 5    |    |   | Benzene, (1-methyl-1-propenyl)-,... | 132 | C10H12  | 000768-00-3 | 93   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052720\  
Data File : BM026138.D  
Acq On : 27 May 2020 10:35  
Operator : MA/SJ  
Sample : L2756-04DL 10X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
ETKW8DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM051320MA.M  
Quant Title : SVOA CALIBRATION

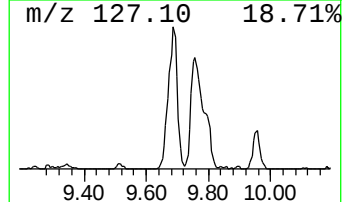
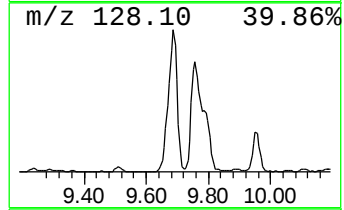
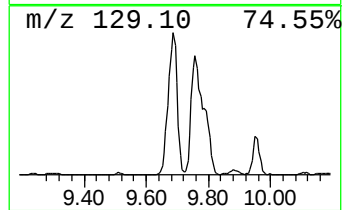
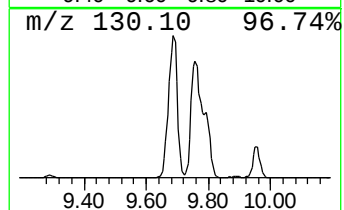
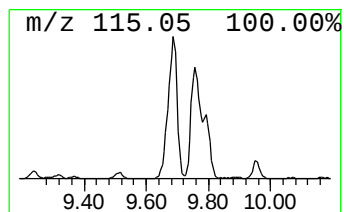
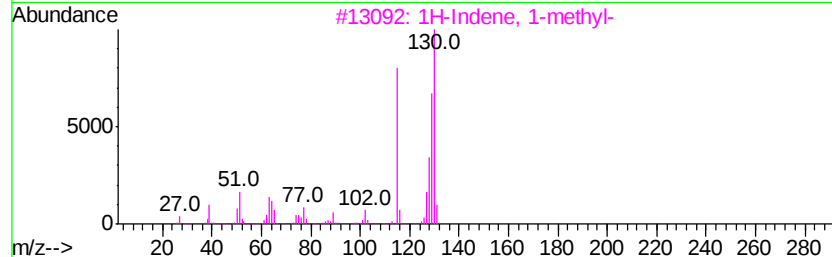
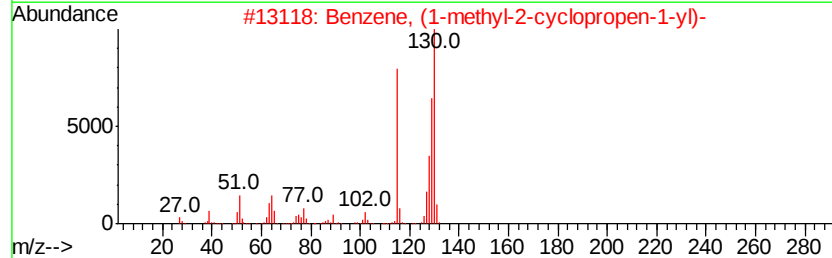
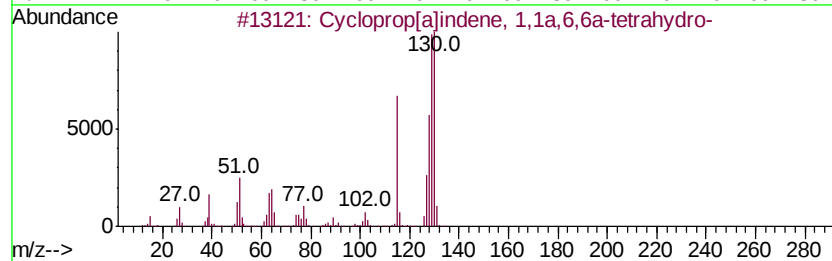
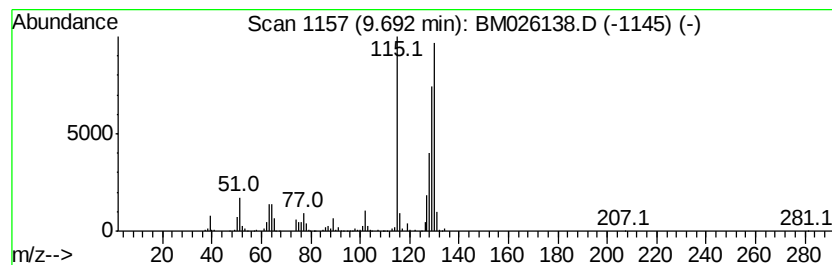
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 21 Cycloprop[a]indene, 1,1a,6,... Concentration Rank 9

| R.T. | EstConc     | Area    | Relative to ISTD | R.T.  |
|------|-------------|---------|------------------|-------|
| 9.69 | 19.02 ng/ul | 1156660 | Naphthalene-d8   | 10.25 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cycloprop[a]indene, 1,1a,6,6a-te... | 130 | C10H10  | 015677-15-3 | 96   |
| 2    |    | Benzene, (1-methyl-2-cyclopropen... | 130 | C10H10  | 065051-83-4 | 96   |
| 3    |    | 1H-Indene, 1-methyl-                | 130 | C10H10  | 000767-59-9 | 96   |
| 4    |    | 2-Methylindene                      | 130 | C10H10  | 002177-47-1 | 96   |
| 5    |    | Cycloprop[a]indene, 1,1a,6,6a-te... | 130 | C10H10  | 015677-15-3 | 95   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052720\  
Data File : BM026138.D  
Acq On : 27 May 2020 10:35  
Operator : MA/SJ  
Sample : L2756-04DL 10X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
ETKW8DL

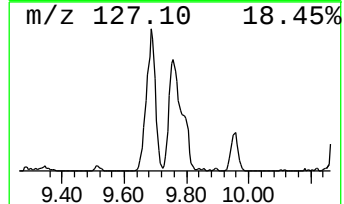
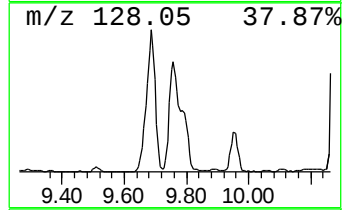
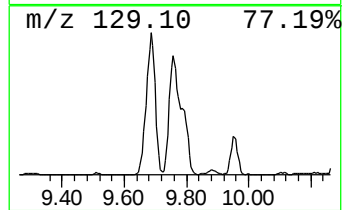
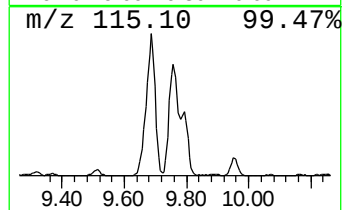
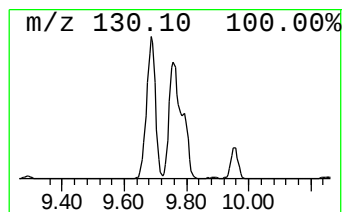
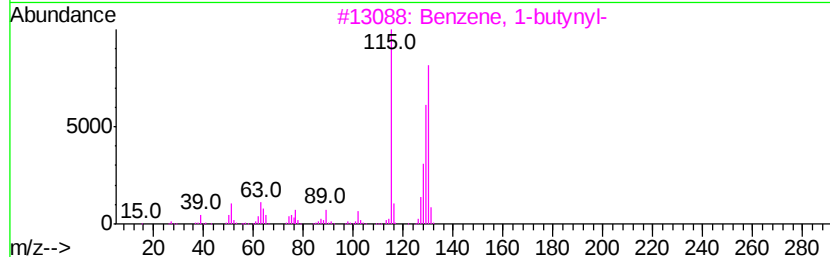
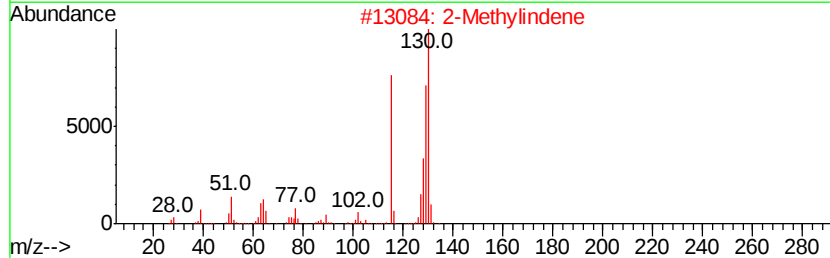
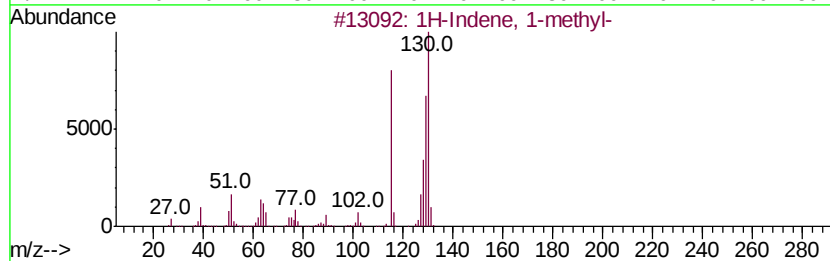
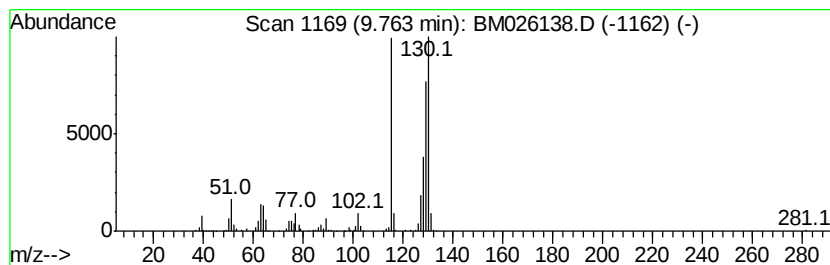
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Quant Title : SVOA CALIBRATION

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

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Peak Number 22 1H-Indene, 1-methyl- Concentration Rank 10

| R.T. | EstConc     | Area   | Relative to ISTD | R.T.  |
|------|-------------|--------|------------------|-------|
| 9.76 | 13.13 ng/ul | 798434 | Naphthalene-d8   | 10.25 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 1H-Indene, 1-methyl-                | 130 | C10H10  | 000767-59-9 | 96   |
| 2    |    |   | 2-Methylindene                      | 130 | C10H10  | 002177-47-1 | 96   |
| 3    |    |   | Benzene, 1-butynyl-                 | 130 | C10H10  | 000622-76-4 | 95   |
| 4    |    |   | Benzene, 1-methyl-1,2-propadienyl-  | 130 | C10H10  | 022433-39-2 | 95   |
| 5    |    |   | Benzene, (1-methyl-2-cyclopropen... | 130 | C10H10  | 065051-83-4 | 95   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052720\  
Data File : BM026138.D  
Acq On : 27 May 2020 10:35  
Operator : MA/SJ  
Sample : L2756-04DL 10X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
ETKW8DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM051320MA.M  
Quant Title : SVOA CALIBRATION

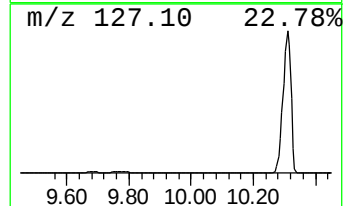
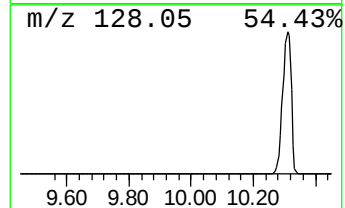
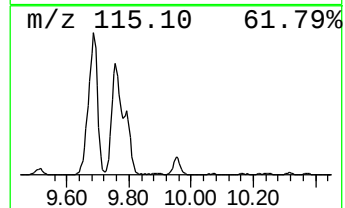
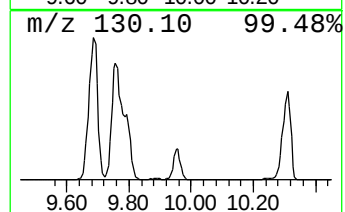
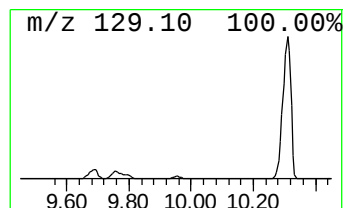
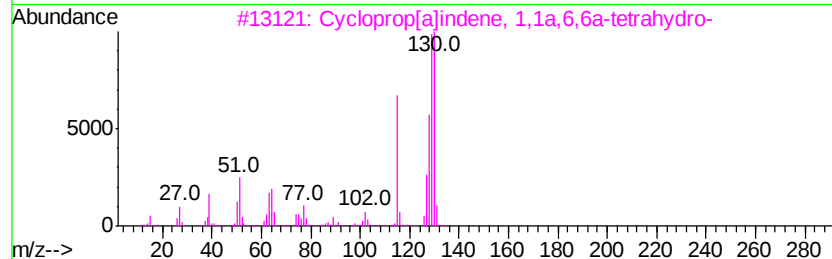
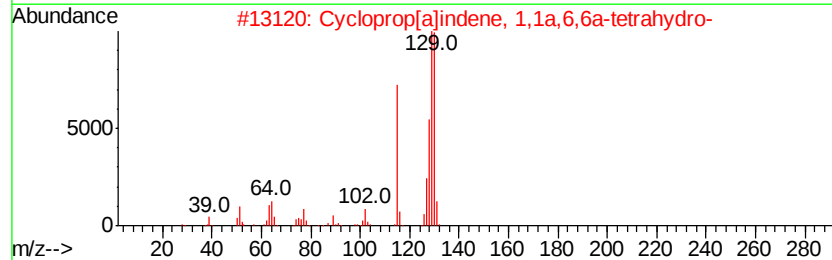
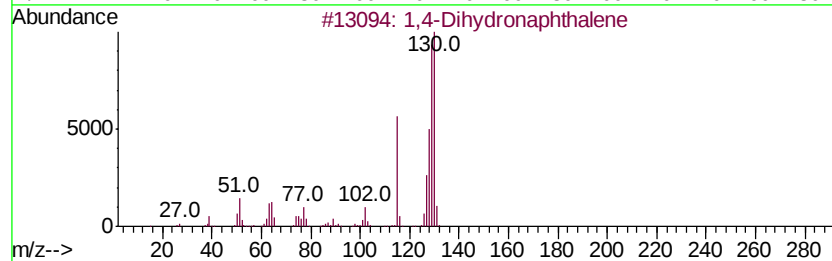
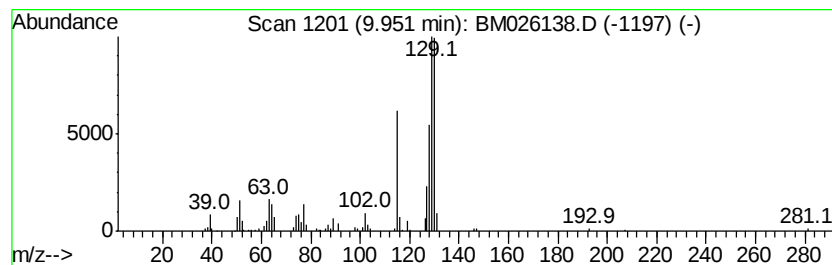
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TIC Integration Parameters: LSCINT.P

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Peak Number 23 1,4-Dihydronaphthalene Concentration Rank 26

| R.T. | EstConc    | Area   | Relative to ISTD | R.T.  |
|------|------------|--------|------------------|-------|
| 9.95 | 2.78 ng/ul | 168760 | Naphthalene-d8   | 10.25 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 1,4-Dihydronaphthalene              | 130 | C10H10  | 000612-17-9 | 96   |
| 2    |    | Cycloprop[alindene, 1,1a,6,6a-te... | 130 | C10H10  | 015677-15-3 | 95   |
| 3    |    | Cycloprop[alindene, 1,1a,6,6a-te... | 130 | C10H10  | 015677-15-3 | 94   |
| 4    |    | 1,4-Dihydronaphthalene              | 130 | C10H10  | 000612-17-9 | 92   |
| 5    |    | 1H-Indene, 3-methyl-                | 130 | C10H10  | 000767-60-2 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052720\  
Data File : BM026138.D  
Acq On : 27 May 2020 10:35  
Operator : MA/SJ  
Sample : L2756-04DL 10X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

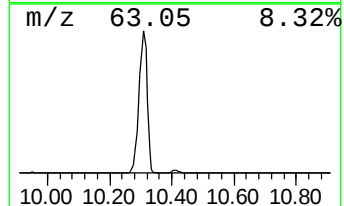
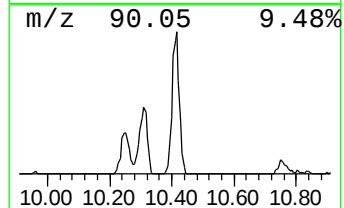
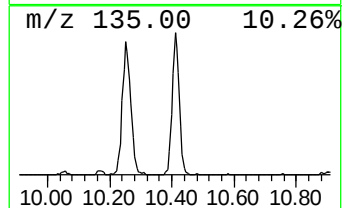
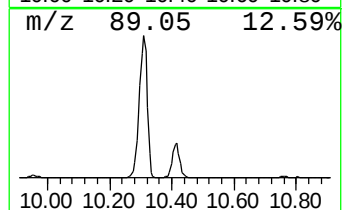
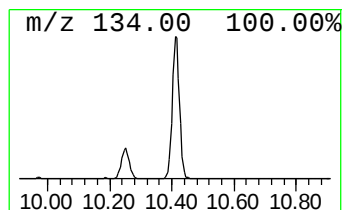
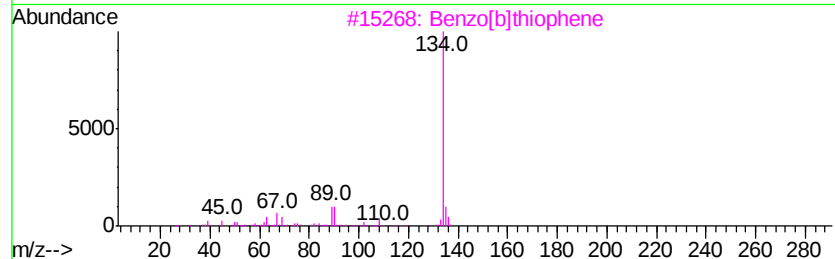
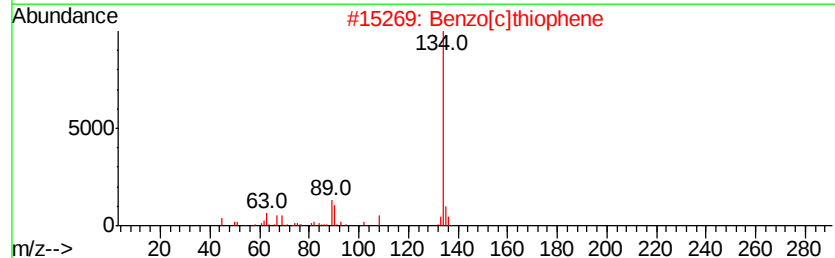
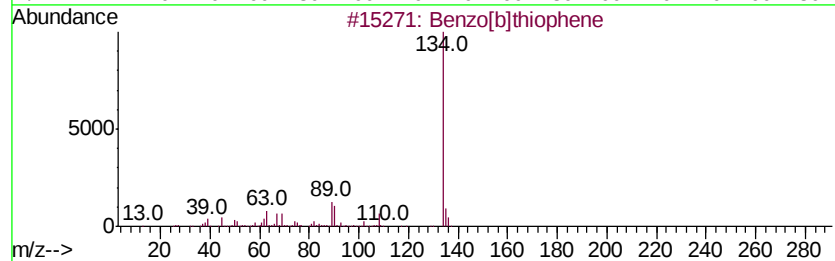
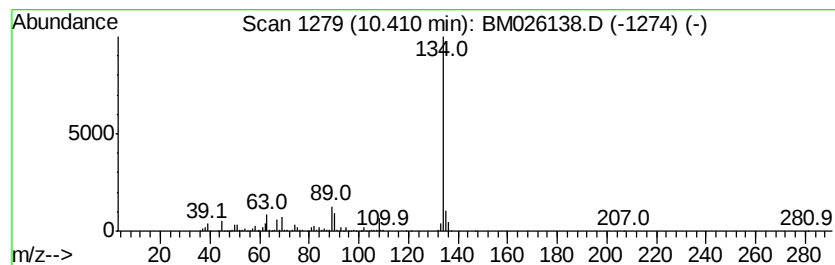
Instrument :  
BNA\_M  
ClientSampleId :  
ETKW8DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM051320MA.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 24 Benzo[b]thiophene Concentration Rank 14

| R.T.    | EstConc    | Area                   | Relative to ISTD | R.T.           |
|---------|------------|------------------------|------------------|----------------|
| 10.41   | 9.54 ng/ul | 580021                 | Naphthalene-d8   | 10.25          |
| Hit# of | 5          | Tentative ID           | MW MolForm       | CAS# Qual      |
| 1       |            | Benzo[b]thiophene      | 134 C8H6S        | 000095-15-8 95 |
| 2       |            | Benzo[c]thiophene      | 134 C8H6S        | 000270-82-6 95 |
| 3       |            | Benzo[b]thiophene      | 134 C8H6S        | 000095-15-8 94 |
| 4       |            | Cyclopenta[c]thiapyran | 134 C8H6S        | 000270-63-3 93 |
| 5       |            | Benzo[b]thiophene      | 134 C8H6S        | 000095-15-8 90 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052720\  
Data File : BM026138.D  
Acq On : 27 May 2020 10:35  
Operator : MA/SJ  
Sample : L2756-04DL 10X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
ETKW8DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM051320MA.M  
Quant Title : SVOA CALIBRATION

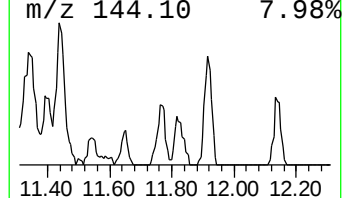
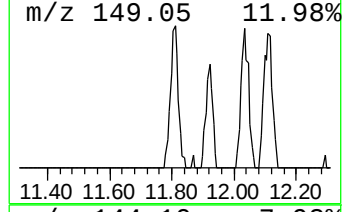
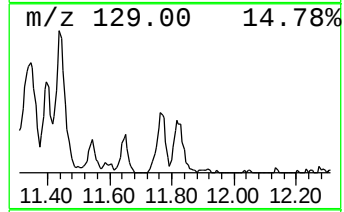
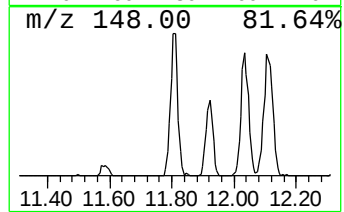
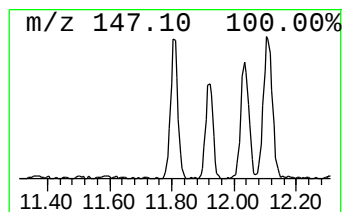
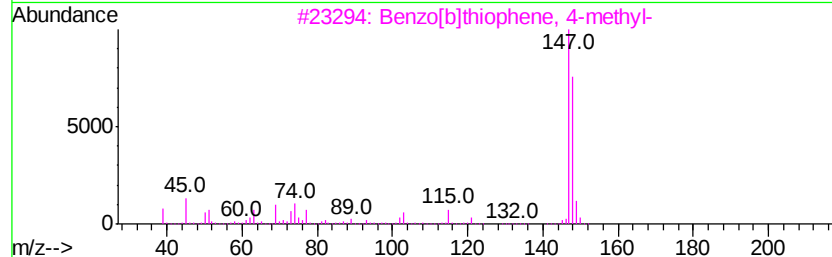
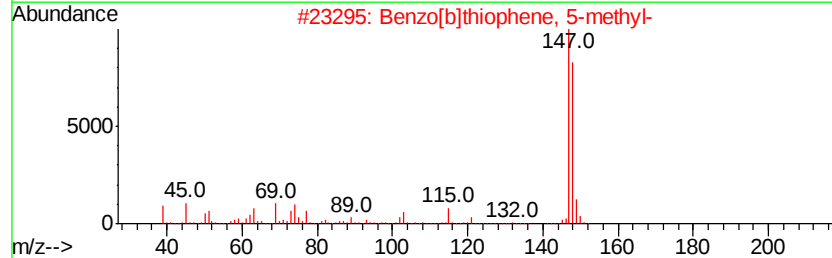
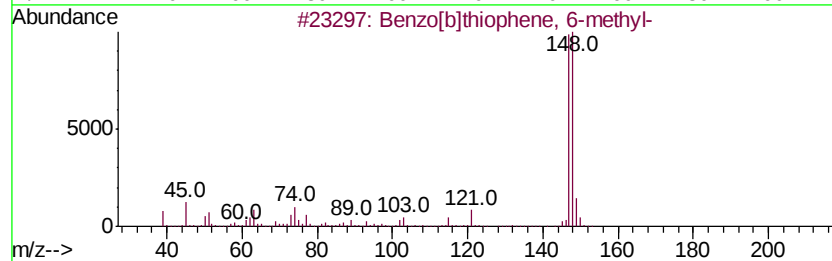
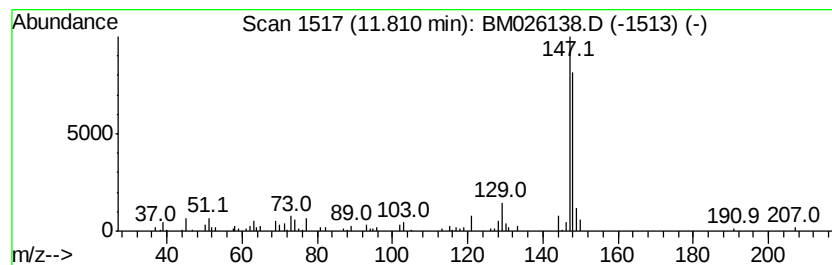
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 25 Benzo[b]thiophene, 6-methyl- Concentration Rank 28

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 11.81 | 2.06 ng/ul | 125151 | Naphthalene-d8   | 10.25 |

| Hit# | of | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzo[b]thiophene, 6-methyl- | 148 | C9H8S   | 016587-47-6 | 87   |
| 2    |    | Benzo[b]thiophene, 5-methyl- | 148 | C9H8S   | 014315-14-1 | 83   |
| 3    |    | Benzo[b]thiophene, 4-methyl- | 148 | C9H8S   | 014315-11-8 | 83   |
| 4    |    | Benzo[b]thiophene, 2-methyl- | 148 | C9H8S   | 001195-14-8 | 80   |
| 5    |    | Benzo[b]thiophene, 2-methyl- | 148 | C9H8S   | 001195-14-8 | 80   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052720\  
Data File : BM026138.D  
Acq On : 27 May 2020 10:35  
Operator : MA/SJ  
Sample : L2756-04DL 10X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
ETKW8DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM051320MA.M  
Quant Title : SVOA CALIBRATION

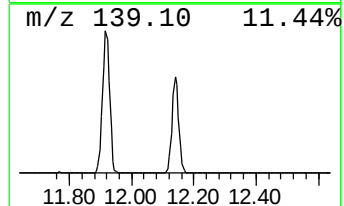
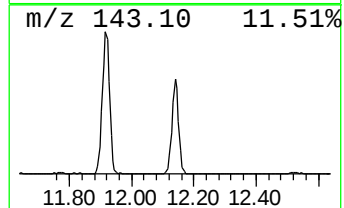
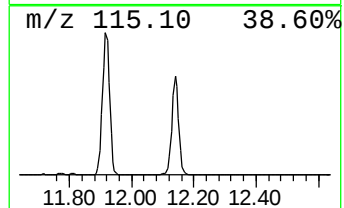
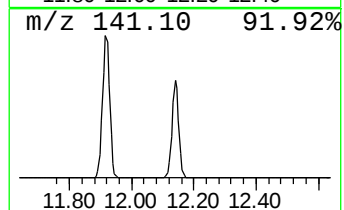
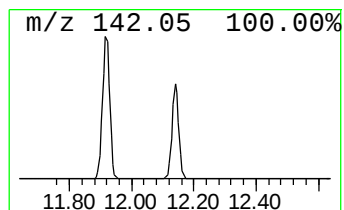
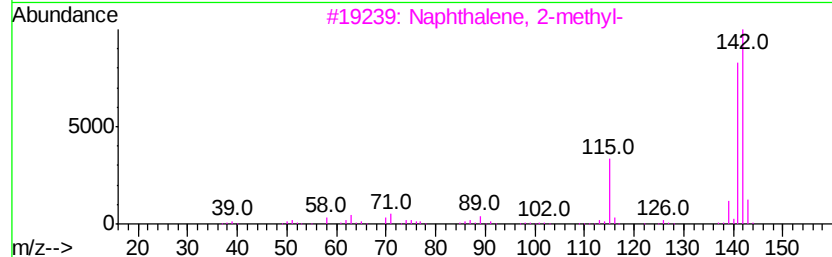
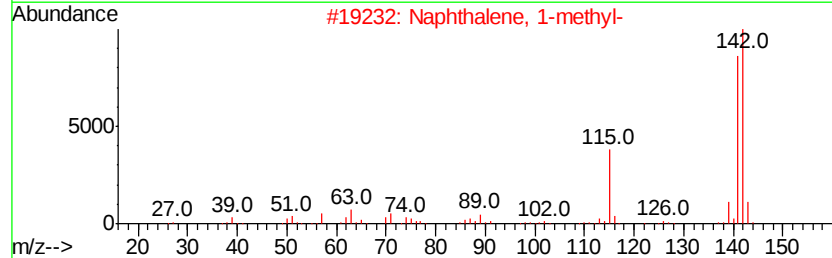
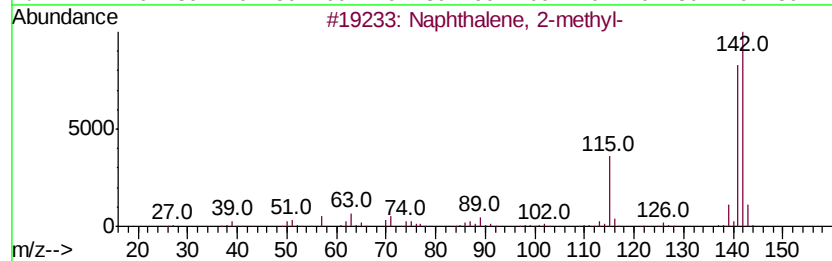
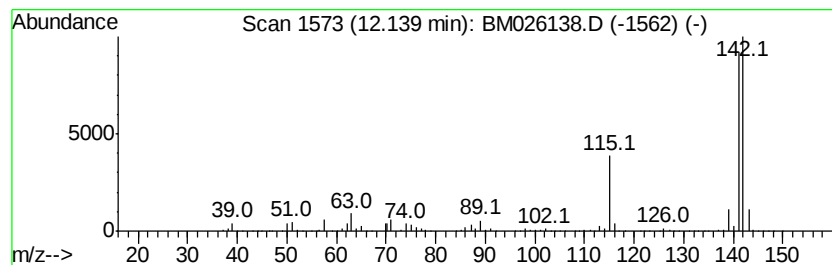
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 26 Naphthalene, 1-methyl- Concentration Rank 7

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 12.14 | 46.34 ng/ul | 2817420 | Naphthalene-d8   | 10.25 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 2-methyl- | 142 | C11H10  | 000091-57-6 | 96   |
| 2    |    | Naphthalene, 1-methyl- | 142 | C11H10  | 000090-12-0 | 96   |
| 3    |    | Naphthalene, 2-methyl- | 142 | C11H10  | 000091-57-6 | 96   |
| 4    |    | Naphthalene, 1-methyl- | 142 | C11H10  | 000090-12-0 | 94   |
| 5    |    | Naphthalene, 2-methyl- | 142 | C11H10  | 000091-57-6 | 93   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052720\  
Data File : BM026138.D  
Acq On : 27 May 2020 10:35  
Operator : MA/SJ  
Sample : L2756-04DL 10X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_M  
ClientSampleId :  
ETKW8DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM051320MA.M  
Quant Title : SVOA CALIBRATION

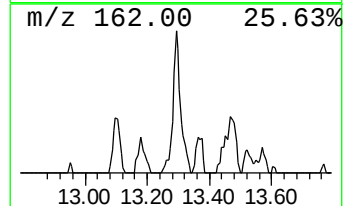
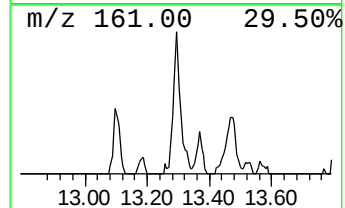
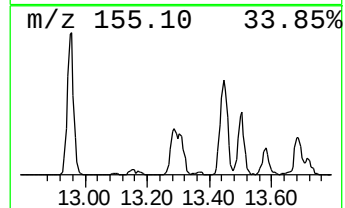
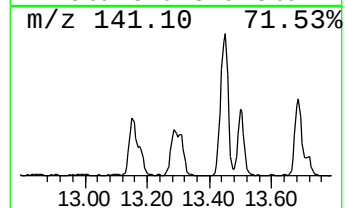
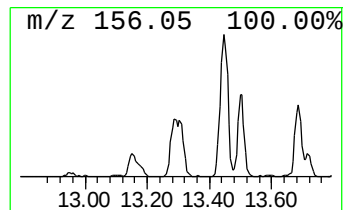
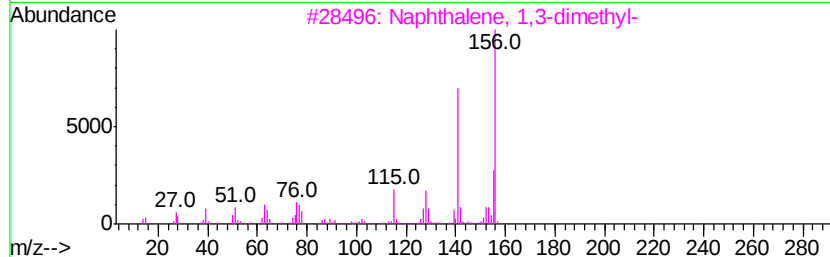
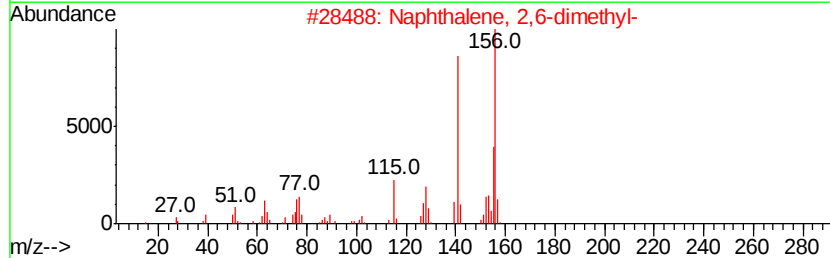
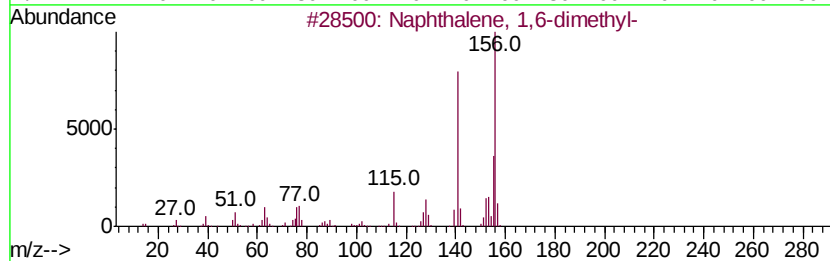
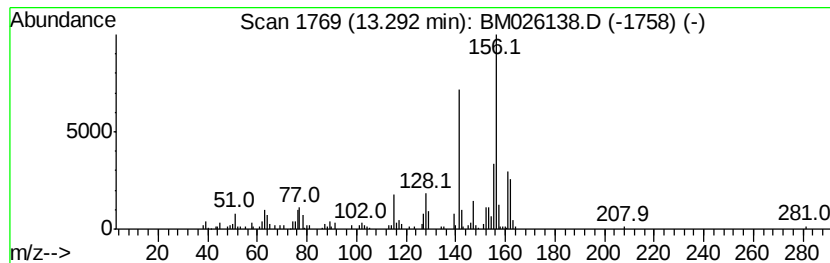
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 27 Naphthalene, 1,6-dimethyl- Concentration Rank 27

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 13.29 | 2.77 ng/ul | 243262 | Acenaphthene-d10 | 14.13 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 98   |
| 2    |    | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 98   |
| 3    |    | Naphthalene, 1,3-dimethyl- | 156 | C12H12  | 000575-41-7 | 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 : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052720\  
Data File : BM026138.D  
Acq On : 27 May 2020 10:35  
Operator : MA/SJ  
Sample : L2756-04DL 10X  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

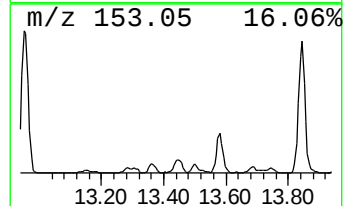
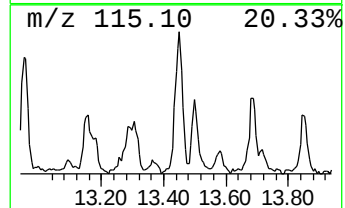
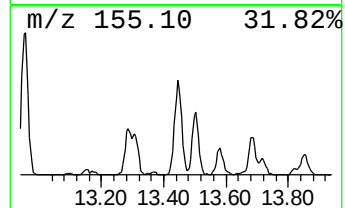
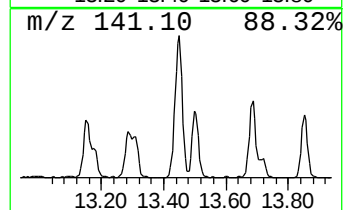
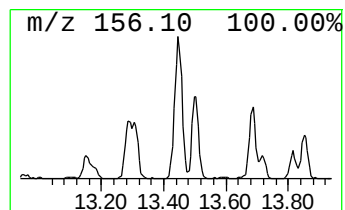
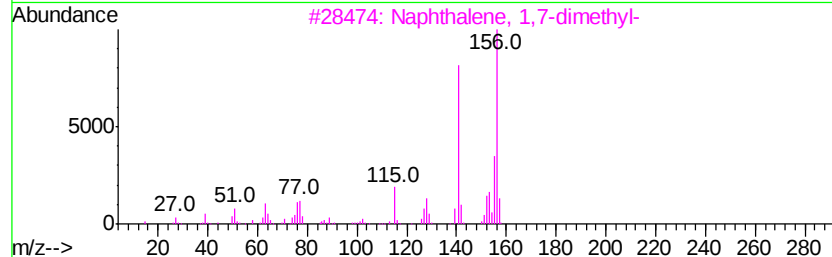
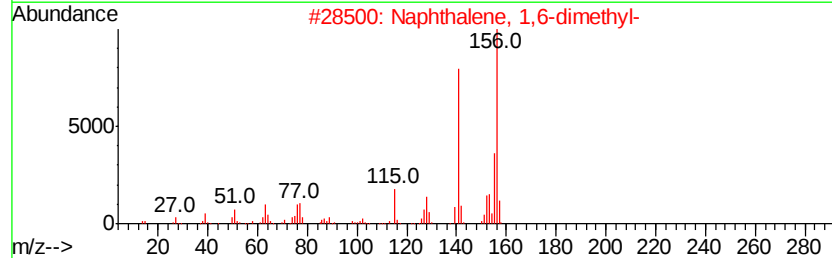
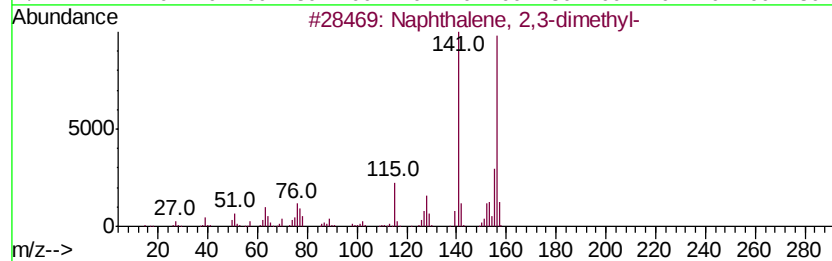
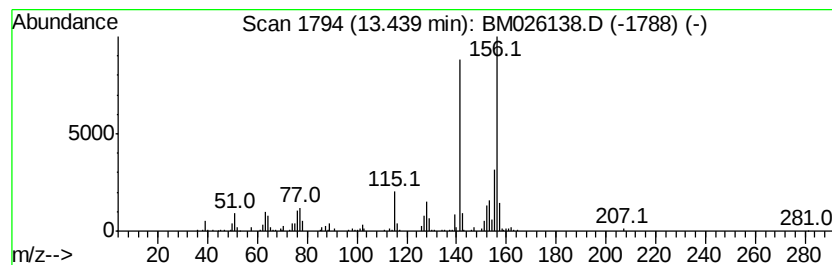
Instrument :  
BNA\_M  
ClientSampleId :  
ETKW8DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM051320MA.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 28 Naphthalene, 2,3-dimethyl- Concentration Rank 22

| R.T.    | EstConc    | Area                       | Relative to ISTD | R.T.           |
|---------|------------|----------------------------|------------------|----------------|
| 13.44   | 3.83 ng/ul | 336174                     | Acenaphthene-d10 | 14.13          |
| 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 98 |
| 3       |            | Naphthalene, 1,7-dimethyl- | 156 C12H12       | 000575-37-1 98 |
| 4       |            | Naphthalene, 1,5-dimethyl- | 156 C12H12       | 000571-61-9 97 |
| 5       |            | Naphthalene, 1,7-dimethyl- | 156 C12H12       | 000575-37-1 97 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_M\DATA\BM052720\  
 Data File : BM026138.D  
 Acq On : 27 May 2020 10:35  
 Operator : MA/SJ  
 Sample : L2756-04DL 10X  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 ETKW8DL

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_M\METHODS\SOM-EPA-BM051320MA.M  
 Quant Title : SVOA 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, 1,3-dime... | 5.17  | 129.2   | ng/ul | 5968980  | 1                     | 7.48  | 924107  | 20.0 |
| Benzene, 1-ethyl-... | 6.59  | 12.7    | ng/ul | 587600   | 1                     | 7.48  | 924107  | 20.0 |
| Benzene, 1-ethyl-... | 6.65  | 8.7     | ng/ul | 403178   | 1                     | 7.48  | 924107  | 20.0 |
| Benzene, 1,2,3-tr... | 6.73  | 10.3    | ng/ul | 476498   | 1                     | 7.48  | 924107  | 20.0 |
| Benzene, 1,2,4-tr... | 6.89  | 5.2     | ng/ul | 240562   | 1                     | 7.48  | 924107  | 20.0 |
| Mesitylene           | 7.15  | 46.5    | ng/ul | 2147620  | 1                     | 7.48  | 924107  | 20.0 |
| Benzene, 1-etheny... | 7.23  | 4.6     | ng/ul | 211932   | 1                     | 7.48  | 924107  | 20.0 |
| unknown-01           | 7.60  | 9.5     | ng/ul | 440943   | 1                     | 7.48  | 924107  | 20.0 |
| Benzene, 2-propenyl- | 7.67  | 7.1     | ng/ul | 328567   | 1                     | 7.48  | 924107  | 20.0 |
| Indane               | 7.85  | 7.5     | ng/ul | 348386   | 1                     | 7.48  | 924107  | 20.0 |
| 1-Propyne, 3-phenyl- | 8.01  | 91.0    | ng/ul | 4206230  | 1                     | 7.48  | 924107  | 20.0 |
| Benzene, 2-ethyl-... | 8.12  | 4.9     | ng/ul | 224723   | 1                     | 7.48  | 924107  | 20.0 |
| (E)-1-Phenyl-1-bu... | 8.75  | 2.8     | ng/ul | 129976   | 1                     | 7.48  | 924107  | 20.0 |
| Cycloprop[a]inden... | 9.69  | 19.0    | ng/ul | 1156660  | 2                     | 10.25 | 1216040 | 20.0 |
| 1H-Indene, 1-methyl- | 9.76  | 13.1    | ng/ul | 798434   | 2                     | 10.25 | 1216040 | 20.0 |
| 1,4-Dihydronaphth... | 9.95  | 2.8     | ng/ul | 168760   | 2                     | 10.25 | 1216040 | 20.0 |
| Benzo[b]thiophene    | 10.41 | 9.5     | ng/ul | 580021   | 2                     | 10.25 | 1216040 | 20.0 |
| Benzo[b]thiophene... | 11.81 | 2.1     | ng/ul | 125151   | 2                     | 10.25 | 1216040 | 20.0 |
| Naphthalene, 1-me... | 12.14 | 46.3    | ng/ul | 2817420  | 2                     | 10.25 | 1216040 | 20.0 |
| Naphthalene, 1,6-... | 13.29 | 2.8     | ng/ul | 243262   | 3                     | 14.13 | 1754560 | 20.0 |
| Naphthalene, 2,3-... | 13.44 | 3.8     | ng/ul | 336174   | 3                     | 14.13 | 1754560 | 20.0 |