

Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
 Data File : BN000563.D  
 Acq On : 23 Mar 2018 04:28  
 Operator : JU/SJ  
 Sample : J1905-03DL2 200X  
 Misc :  
 ALS Vial : 31 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 DAM40DL2

## 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\_N\METHODS\SOM-EPA-BN031918.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         | 6.349       | 489           | 495         | 504          | rVB2     | 258225         | 401921        | 6.42%           | 0.788%        |
| 2         | 7.002       | 601           | 606         | 617          | rVB4     | 77273          | 211343        | 3.38%           | 0.414%        |
| 3         | 7.355       | 662           | 666         | 671          | rVB      | 149073         | 214458        | 3.43%           | 0.420%        |
| 4         | 7.413       | 671           | 676         | 681          | rVB      | 154392         | 212987        | 3.40%           | 0.417%        |
| 5         | 7.466       | 681           | 685         | 690          | rBV      | 140875         | 209383        | 3.34%           | 0.410%        |
| 6         | 7.525       | 690           | 695         | 701          | rBV2     | 245500         | 388101        | 6.20%           | 0.761%        |
| 7         | 7.607       | 701           | 709         | 713          | rBV      | 89118          | 147615        | 2.36%           | 0.289%        |
| 8         | 7.778       | 731           | 738         | 741          | rBV      | 95157          | 170953        | 2.73%           | 0.335%        |
| 9         | 8.002       | 770           | 776         | 780          | rBV      | 796627         | 1141827       | 18.24%          | 2.238%        |
| 10        | 8.043       | 780           | 783         | 792          | rVB      | 216083         | 335540        | 5.36%           | 0.658%        |
| 11        | 8.360       | 832           | 837         | 841          | rBV      | 157580         | 211907        | 3.39%           | 0.415%        |
| 12        | 8.413       | 841           | 846         | 850          | rBV      | 628076         | 1007728       | 16.10%          | 1.975%        |
| 13        | 8.454       | 850           | 853         | 860          | rVB      | 239433         | 366444        | 5.85%           | 0.718%        |
| 14        | 8.525       | 860           | 865         | 870          | rBV3     | 123844         | 235286        | 3.76%           | 0.461%        |
| 15        | 8.690       | 889           | 893         | 900          | rVB2     | 81577          | 131942        | 2.11%           | 0.259%        |
| 16        | 8.919       | 925           | 932         | 935          | rVV2     | 122360         | 249193        | 3.98%           | 0.488%        |
| 17        | 8.978       | 935           | 942         | 948          | rVB2     | 156877         | 391756        | 6.26%           | 0.768%        |
| 18        | 9.049       | 948           | 954         | 961          | rBV2     | 364835         | 616280        | 9.84%           | 1.208%        |
| 19        | 9.160       | 967           | 973         | 977          | rBV      | 177690         | 275076        | 4.39%           | 0.539%        |
| 20        | 9.213       | 977           | 982         | 987          | rVV2     | 91567          | 168681        | 2.69%           | 0.331%        |
| 21        | 9.366       | 1003          | 1008        | 1012         | rBV      | 95853          | 158989        | 2.54%           | 0.312%        |
| 22        | 9.419       | 1013          | 1017        | 1023         | rVB      | 104875         | 184990        | 2.96%           | 0.363%        |
| 23        | 9.519       | 1030          | 1034        | 1044         | rVV2     | 169298         | 302518        | 4.83%           | 0.593%        |
| 24        | 9.625       | 1044          | 1052        | 1058         | rVB      | 933102         | 1418830       | 22.67%          | 2.781%        |
| 25        | 9.766       | 1064          | 1076        | 1084         | rBV6     | 79363          | 284965        | 4.55%           | 0.559%        |
| 26        | 9.860       | 1084          | 1092        | 1097         | rBV3     | 76110          | 218205        | 3.49%           | 0.428%        |
| 27        | 10.049      | 1120          | 1124        | 1130         | rVB3     | 92699          | 152036        | 2.43%           | 0.298%        |
| 28        | 10.113      | 1130          | 1135        | 1139         | rBV      | 82587          | 131162        | 2.10%           | 0.257%        |
| 29        | 10.349      | 1171          | 1175        | 1183         | rVB4     | 127478         | 251834        | 4.02%           | 0.494%        |
| 30        | 10.484      | 1194          | 1198        | 1204         | rVB2     | 115846         | 182997        | 2.92%           | 0.359%        |
| 31        | 10.631      | 1218          | 1223        | 1229         | rBV3     | 142775         | 226338        | 3.62%           | 0.444%        |
| 32        | 10.731      | 1236          | 1240        | 1245         | rVB3     | 76853          | 129662        | 2.07%           | 0.254%        |
| 33        | 10.901      | 1260          | 1269        | 1273         | rBV2     | 68743          | 150079        | 2.40%           | 0.294%        |
| 34        | 11.184      | 1308          | 1317        | 1319         | rBV      | 414713         | 831080        | 13.28%          | 1.629%        |

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Integrator: RTE  
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 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\_N\METHODS\SOM-EPA-BN031918.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 35 | 11.254 | 1324 | 1329 | 1334 | rVV  | 893317  | 1621941 | 25.91% | 3.179% |
| 36 | 11.325 | 1334 | 1341 | 1345 | rVV2 | 221039  | 488625  | 7.81%  | 0.958% |
| 37 | 11.360 | 1345 | 1347 | 1353 | rVB2 | 130079  | 198641  | 3.17%  | 0.389% |
| 38 | 11.590 | 1380 | 1386 | 1393 | rVB  | 382475  | 596699  | 9.53%  | 1.170% |
| 39 | 11.719 | 1399 | 1408 | 1414 | rVB2 | 167073  | 296553  | 4.74%  | 0.581% |
| 40 | 11.925 | 1435 | 1443 | 1454 | rVB7 | 97450   | 289593  | 4.63%  | 0.568% |
| 41 | 12.113 | 1468 | 1475 | 1480 | rBV4 | 97613   | 157298  | 2.51%  | 0.308% |
| 42 | 12.219 | 1487 | 1493 | 1498 | rBV3 | 173469  | 325549  | 5.20%  | 0.638% |
| 43 | 12.307 | 1504 | 1508 | 1523 | rVB  | 329705  | 521376  | 8.33%  | 1.022% |
| 44 | 12.454 | 1528 | 1533 | 1539 | rVB  | 294434  | 417139  | 6.66%  | 0.818% |
| 45 | 12.607 | 1553 | 1559 | 1563 | rBV  | 383288  | 604416  | 9.66%  | 1.185% |
| 46 | 12.648 | 1563 | 1566 | 1572 | rVB  | 258404  | 341589  | 5.46%  | 0.670% |
| 47 | 12.742 | 1573 | 1582 | 1584 | rBV3 | 96607   | 195827  | 3.13%  | 0.384% |
| 48 | 12.848 | 1595 | 1600 | 1603 | rVV  | 395638  | 566655  | 9.05%  | 1.111% |
| 49 | 12.884 | 1603 | 1606 | 1613 | rVB  | 134969  | 197996  | 3.16%  | 0.388% |
| 50 | 13.013 | 1623 | 1628 | 1631 | rBV  | 112313  | 180607  | 2.89%  | 0.354% |
| 51 | 13.095 | 1636 | 1642 | 1650 | rVB2 | 71651   | 172187  | 2.75%  | 0.337% |
| 52 | 13.242 | 1659 | 1667 | 1673 | rBV4 | 80480   | 245385  | 3.92%  | 0.481% |
| 53 | 13.542 | 1713 | 1718 | 1724 | rBV  | 129531  | 213491  | 3.41%  | 0.418% |
| 54 | 13.825 | 1760 | 1766 | 1772 | rBV  | 455870  | 711119  | 11.36% | 1.394% |
| 55 | 14.066 | 1802 | 1807 | 1810 | rBV3 | 103442  | 160089  | 2.56%  | 0.314% |
| 56 | 14.213 | 1821 | 1832 | 1838 | rBV7 | 119025  | 400764  | 6.40%  | 0.785% |
| 57 | 14.348 | 1850 | 1855 | 1862 | rVV4 | 171009  | 450259  | 7.19%  | 0.882% |
| 58 | 14.472 | 1868 | 1876 | 1880 | rVB4 | 164283  | 322713  | 5.16%  | 0.633% |
| 59 | 14.589 | 1892 | 1896 | 1898 | rVV  | 92755   | 130397  | 2.08%  | 0.256% |
| 60 | 14.631 | 1898 | 1903 | 1909 | rVB5 | 196318  | 421468  | 6.73%  | 0.826% |
| 61 | 14.884 | 1941 | 1946 | 1951 | rVB  | 1370452 | 1757572 | 28.08% | 3.445% |
| 62 | 15.031 | 1965 | 1971 | 1980 | rVV2 | 1319205 | 2252136 | 35.98% | 4.414% |
| 63 | 15.142 | 1985 | 1990 | 1998 | rBV  | 500693  | 735551  | 11.75% | 1.442% |
| 64 | 15.219 | 1998 | 2003 | 2005 | rBV3 | 88089   | 156269  | 2.50%  | 0.306% |
| 65 | 15.307 | 2013 | 2018 | 2025 | rBV6 | 99177   | 233954  | 3.74%  | 0.459% |
| 66 | 15.407 | 2031 | 2035 | 2044 | rVB6 | 117522  | 299246  | 4.78%  | 0.587% |
| 67 | 15.489 | 2044 | 2049 | 2053 | rBV2 | 159478  | 248282  | 3.97%  | 0.487% |
| 68 | 15.548 | 2055 | 2059 | 2068 | rVV7 | 155187  | 328060  | 5.24%  | 0.643% |
| 69 | 15.648 | 2068 | 2076 | 2079 | rVV5 | 358525  | 744562  | 11.89% | 1.459% |
| 70 | 15.678 | 2079 | 2081 | 2087 | rVV2 | 188127  | 232154  | 3.71%  | 0.455% |
| 71 | 15.760 | 2090 | 2095 | 2099 | rVV  | 511043  | 753415  | 12.04% | 1.477% |

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Method : Z:\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN031918.M  
 Title : SVOA CALIBRATION

|     |        |      |      |      |      |         |         |         |         |
|-----|--------|------|------|------|------|---------|---------|---------|---------|
| 72  | 15.825 | 2099 | 2106 | 2112 | rVB  | 514289  | 918483  | 14.67%  | 1.800%  |
| 73  | 16.101 | 2149 | 2153 | 2160 | rVB  | 125481  | 194742  | 3.11%   | 0.382%  |
| 74  | 16.225 | 2170 | 2174 | 2178 | rVB  | 157824  | 216151  | 3.45%   | 0.424%  |
| 75  | 16.442 | 2208 | 2211 | 2223 | rVB6 | 85130   | 171912  | 2.75%   | 0.337%  |
| 76  | 16.613 | 2235 | 2240 | 2246 | rBV3 | 222270  | 339208  | 5.42%   | 0.665%  |
| 77  | 16.678 | 2246 | 2251 | 2254 | rBV  | 507152  | 736802  | 11.77%  | 1.444%  |
| 78  | 17.042 | 2305 | 2313 | 2317 | rBV4 | 110909  | 242778  | 3.88%   | 0.476%  |
| 79  | 17.413 | 2369 | 2376 | 2381 | rBV  | 4206473 | 6259920 | 100.00% | 12.269% |
| 80  | 17.460 | 2381 | 2384 | 2389 | rVV  | 477857  | 637091  | 10.18%  | 1.249%  |
| 81  | 17.513 | 2389 | 2393 | 2402 | rVB  | 183457  | 344733  | 5.51%   | 0.676%  |
| 82  | 17.760 | 2427 | 2435 | 2440 | rBV2 | 1670401 | 2477091 | 39.57%  | 4.855%  |
| 83  | 17.872 | 2450 | 2454 | 2460 | rVB4 | 106689  | 149764  | 2.39%   | 0.294%  |
| 84  | 18.007 | 2471 | 2477 | 2479 | rBV4 | 95046   | 197921  | 3.16%   | 0.388%  |
| 85  | 18.195 | 2505 | 2509 | 2518 | rVB  | 381548  | 541342  | 8.65%   | 1.061%  |
| 86  | 18.372 | 2535 | 2539 | 2545 | rVB  | 229806  | 297920  | 4.76%   | 0.584%  |
| 87  | 18.872 | 2621 | 2624 | 2629 | rVB  | 279570  | 336721  | 5.38%   | 0.660%  |
| 88  | 19.495 | 2725 | 2730 | 2732 | rBV  | 264128  | 346437  | 5.53%   | 0.679%  |
| 89  | 19.519 | 2732 | 2734 | 2738 | rVB2 | 239156  | 252421  | 4.03%   | 0.495%  |
| 90  | 19.648 | 2750 | 2756 | 2761 | rVB3 | 98751   | 179899  | 2.87%   | 0.353%  |
| 91  | 20.060 | 2822 | 2826 | 2830 | rVB  | 162559  | 192202  | 3.07%   | 0.377%  |
| 92  | 20.589 | 2912 | 2916 | 2920 | rBV2 | 156971  | 190947  | 3.05%   | 0.374%  |
| 93  | 21.077 | 2995 | 2999 | 3003 | rBV  | 134745  | 147821  | 2.36%   | 0.290%  |
| 94  | 21.536 | 3073 | 3077 | 3084 | rVB  | 290436  | 343216  | 5.48%   | 0.673%  |
| 95  | 21.877 | 3130 | 3135 | 3146 | rVB2 | 1952229 | 2664597 | 42.57%  | 5.223%  |
| 96  | 21.977 | 3149 | 3152 | 3160 | rVB  | 124043  | 170704  | 2.73%   | 0.335%  |
| 97  | 22.360 | 3214 | 3217 | 3224 | rVB2 | 107725  | 148717  | 2.38%   | 0.291%  |
| 98  | 22.454 | 3229 | 3233 | 3239 | rVB2 | 190683  | 271473  | 4.34%   | 0.532%  |
| 99  | 22.960 | 3315 | 3319 | 3322 | rBV  | 86963   | 134521  | 2.15%   | 0.264%  |
| 100 | 24.360 | 3549 | 3557 | 3566 | rBV2 | 1073694 | 2233742 | 35.68%  | 4.378%  |

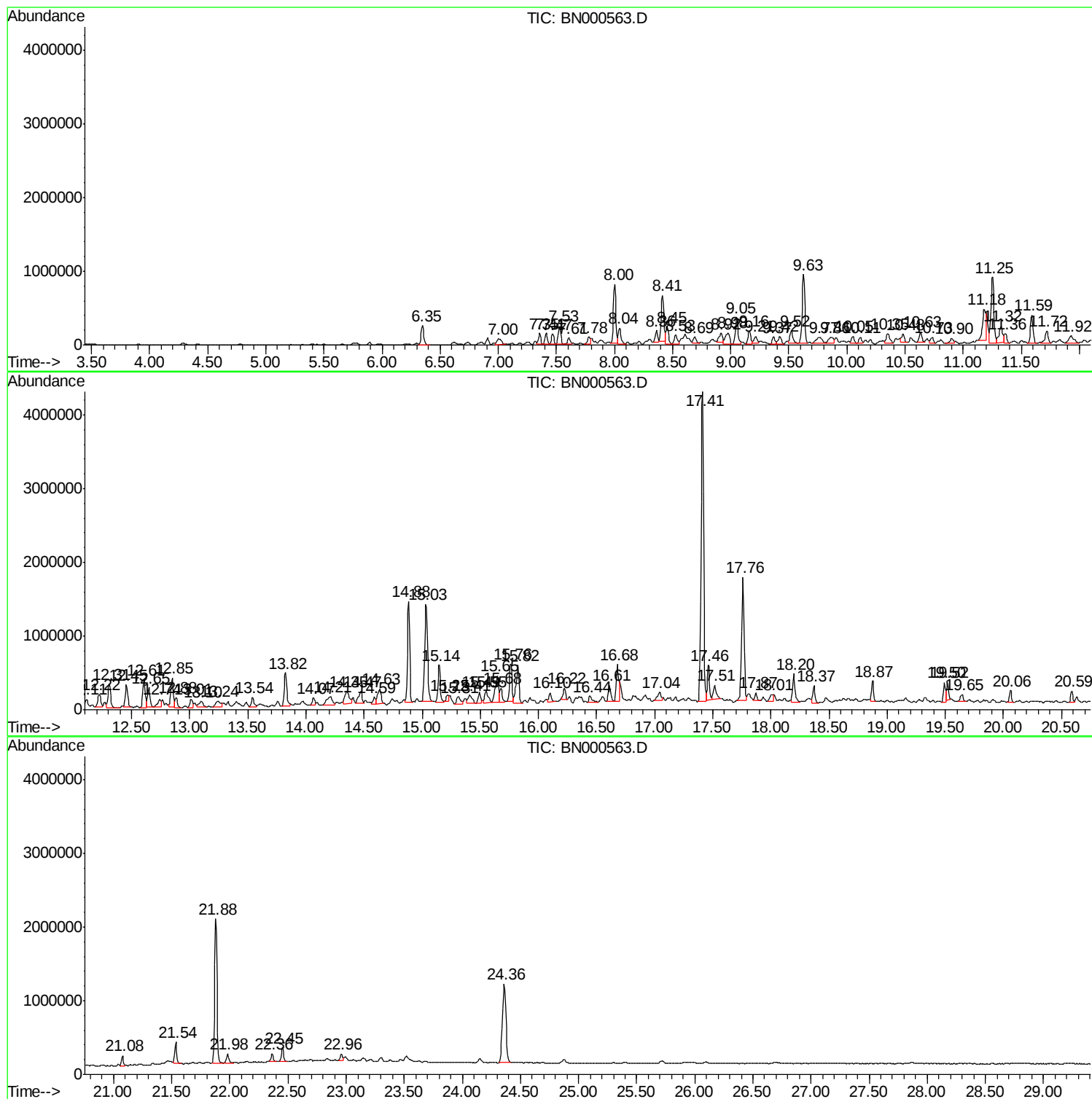
Sum of corrected areas: 51020959

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Quant Method : Z:\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN031918.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 N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

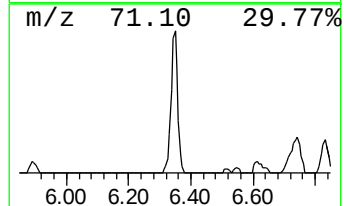
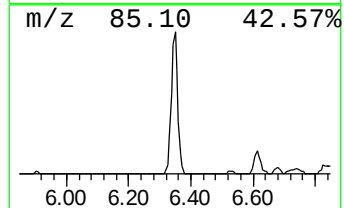
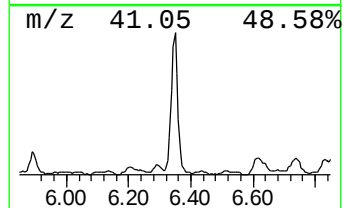
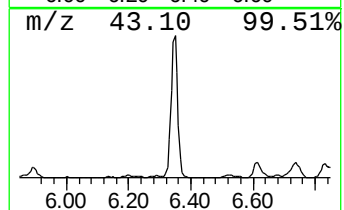
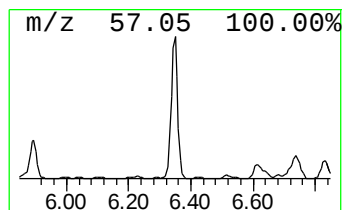
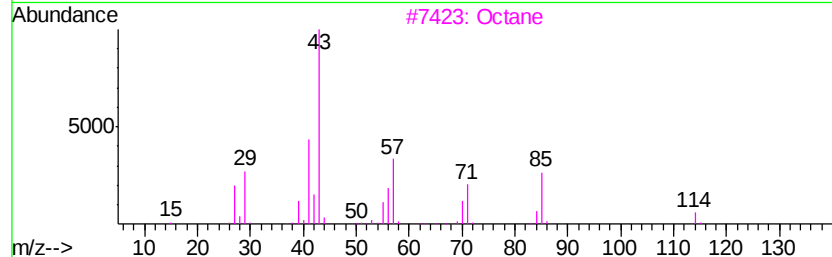
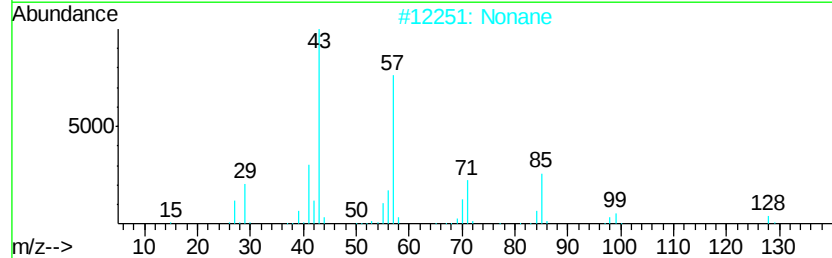
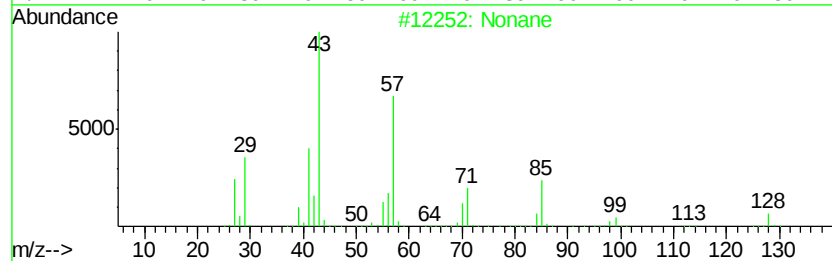
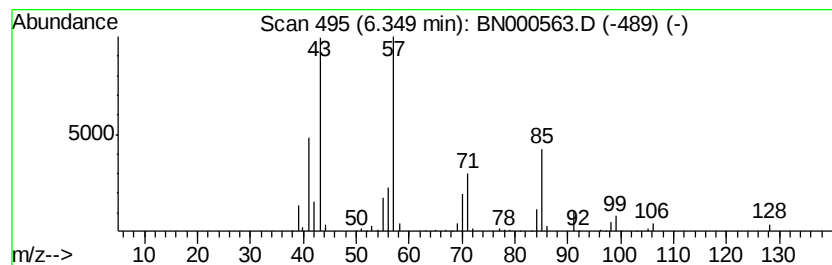
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 7

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 6.35 | 7.98 ng/ul | 401921 | 1,4-Dichlorobenzene-d4 | 8.41 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#         | Qual |
|------|----|-------------------------|-----|---------|--------------|------|
| 1    | 5  | Nonane                  | 128 | C9H20   | 000111-84-2  | 95   |
| 2    |    | Nonane                  | 128 | C9H20   | 000111-84-2  | 91   |
| 3    |    | Octane                  | 114 | C8H18   | 000111-65-9  | 72   |
| 4    |    | 4,4-Dimethylcyclooctene | 138 | C10H18  | 1000113-56-7 | 59   |
| 5    |    | Hexane, 2,4-dimethyl-   | 114 | C8H18   | 000589-43-5  | 50   |



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Quant Method : Z:\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

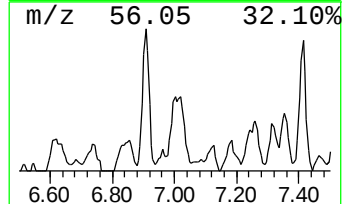
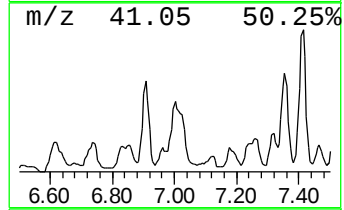
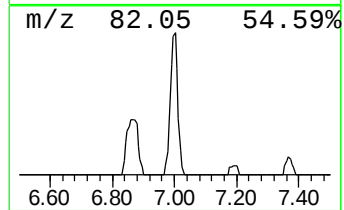
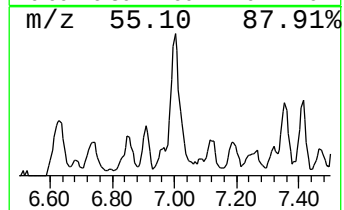
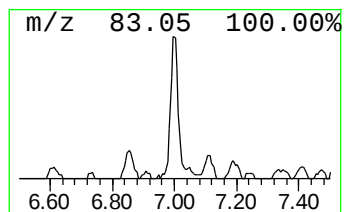
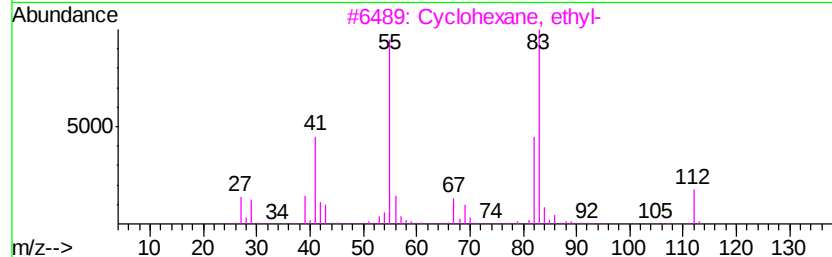
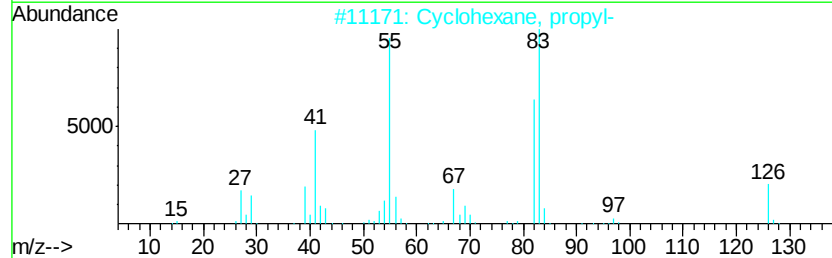
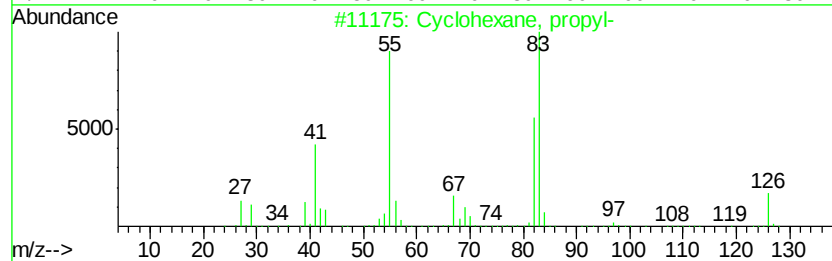
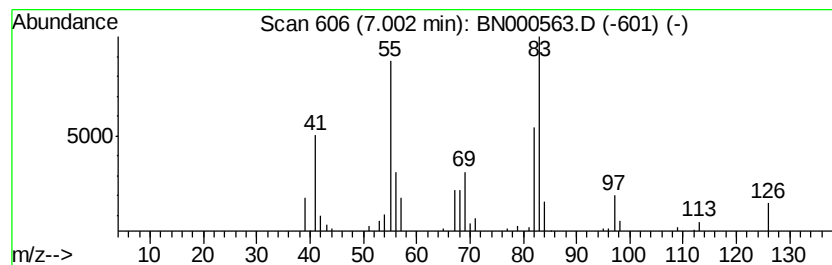
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Peak Number 2 (DEL) Alkane: Cyclic7.00 Concentration Rank 30

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 7.00 | 4.19 ng/ul | 211343 | 1,4-Dichlorobenzene-d4 | 8.41 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, propyl-          | 126 | C9H18   | 001678-92-8 | 64   |
| 2    |    |   | Cyclohexane, propyl-          | 126 | C9H18   | 001678-92-8 | 64   |
| 3    |    |   | Cyclohexane, ethyl-           | 112 | C8H16   | 001678-91-7 | 53   |
| 4    |    |   | Cyclohexane, (1-methylethyl)- | 126 | C9H18   | 000696-29-7 | 53   |
| 5    |    |   | Cyclohexane, propyl-          | 126 | C9H18   | 001678-92-8 | 53   |



Data Path : Z:\HPCHEM1\BNA\_N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampled :  
DAM40DL2

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Quant Title : SVOA CALIBRATION

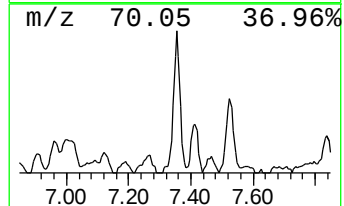
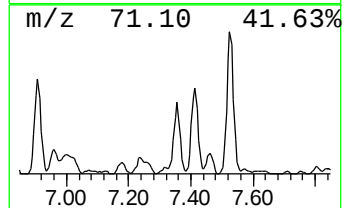
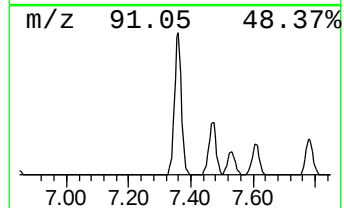
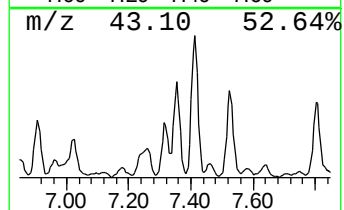
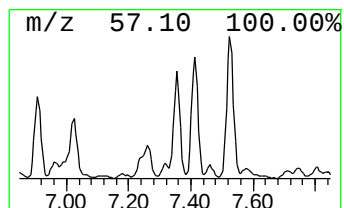
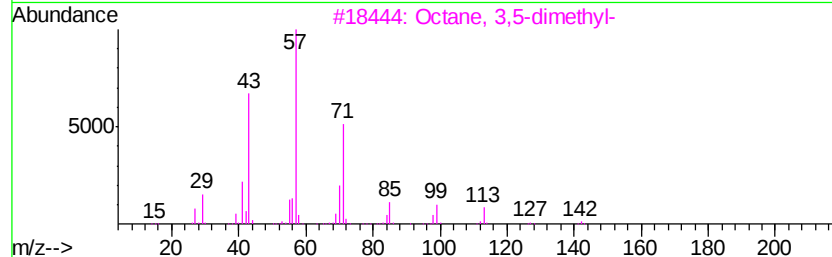
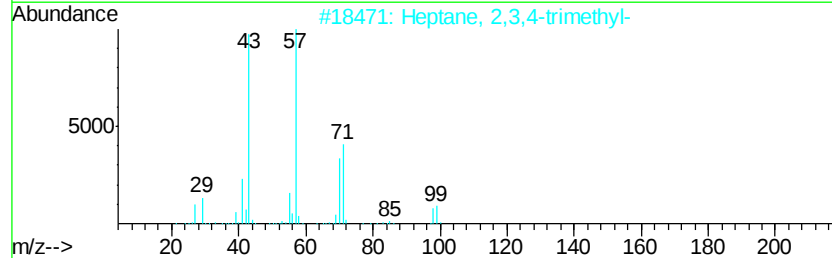
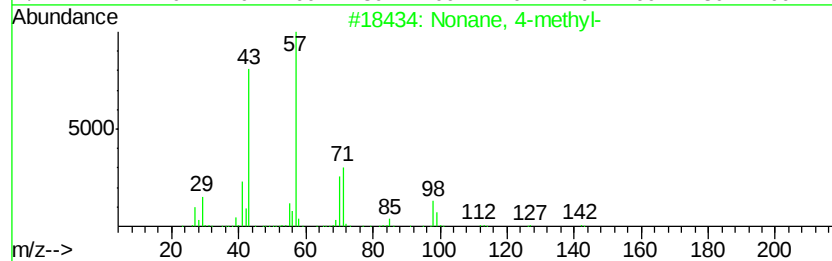
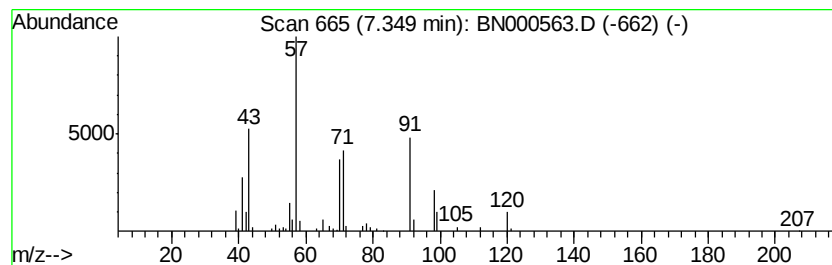
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 26

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 7.35 | 4.26 ng/ul | 214458 | 1,4-Dichlorobenzene-d4 | 8.41 |

| Hit# | of | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------|-----|---------|-------------|------|
| 1    | 5  | Nonane, 4-methyl-             | 142 | C10H22  | 017301-94-9 | 74   |
| 2    |    | Heptane, 2,3,4-trimethyl-     | 142 | C10H22  | 052896-95-4 | 50   |
| 3    |    | Octane, 3,5-dimethyl-         | 142 | C10H22  | 015869-93-9 | 47   |
| 4    |    | Pentane, 2,2,3,3-tetramethyl- | 128 | C9H20   | 007154-79-2 | 47   |
| 5    |    | Nonane, 2-methyl-             | 142 | C10H22  | 000871-83-0 | 47   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

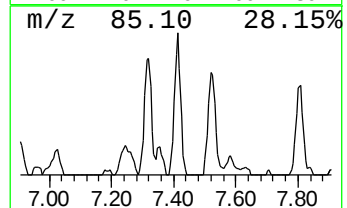
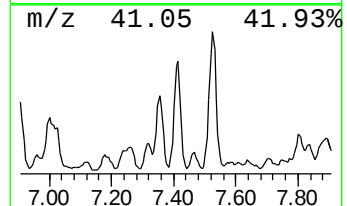
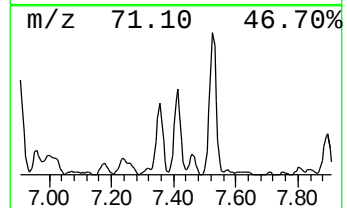
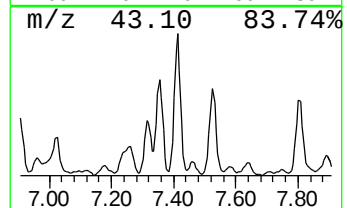
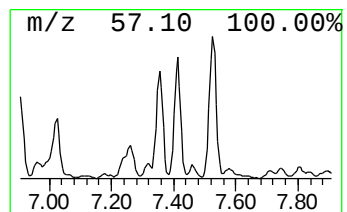
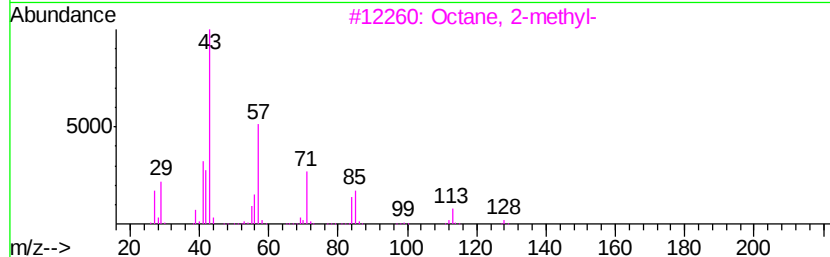
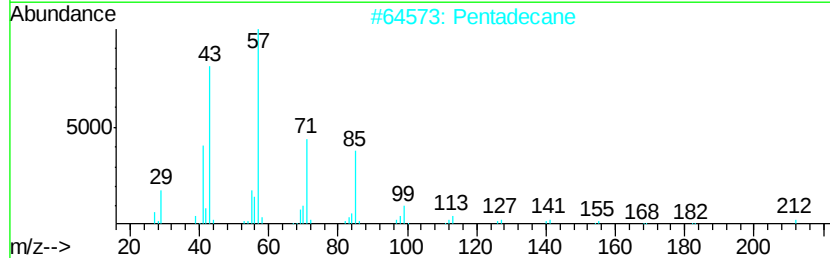
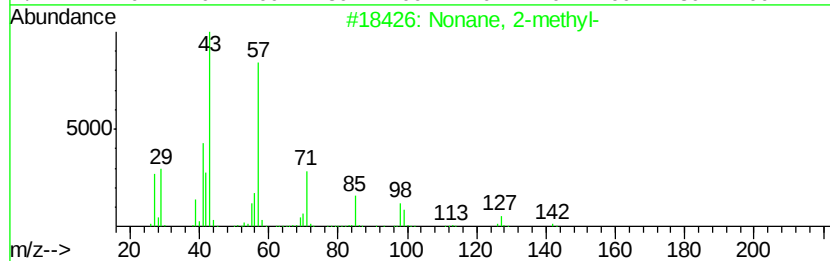
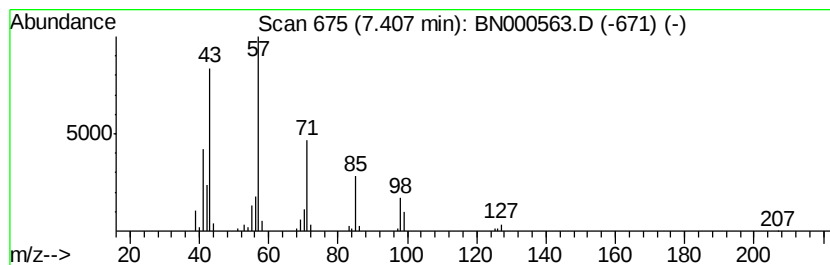
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TIC Integration Parameters: LSCINT.P

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Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 27

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 7.41 | 4.23 ng/ul | 212987 | 1,4-Dichlorobenzene-d4 | 8.41 |

| Hit# | of | 5 | Tentative ID      | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------|-----|---------|-------------|------|
| 1    |    |   | Nonane, 2-methyl- | 142 | C10H22  | 000871-83-0 | 83   |
| 2    |    |   | Pentadecane       | 212 | C15H32  | 000629-62-9 | 72   |
| 3    |    |   | Octane, 2-methyl- | 128 | C9H20   | 003221-61-2 | 72   |
| 4    |    |   | Nonane, 2-methyl- | 142 | C10H22  | 000871-83-0 | 64   |
| 5    |    |   | Eicosane          | 282 | C20H42  | 000112-95-8 | 64   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampled :  
DAM40DL2

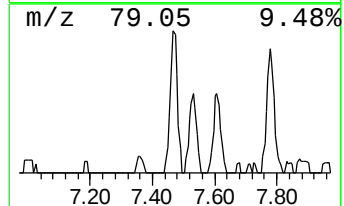
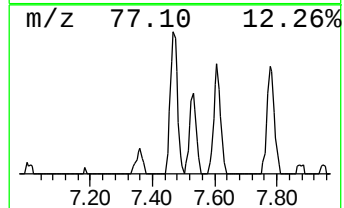
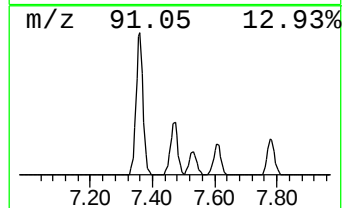
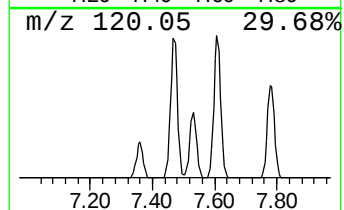
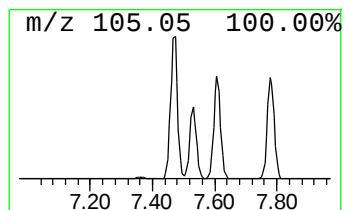
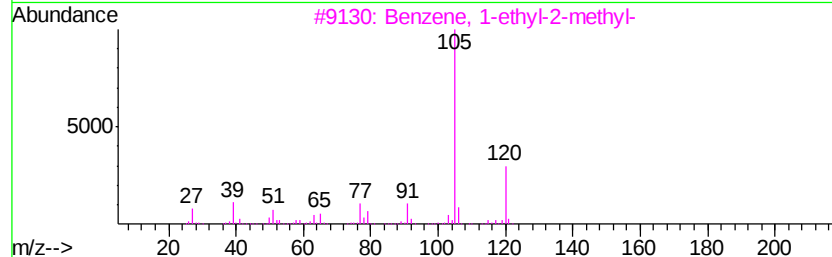
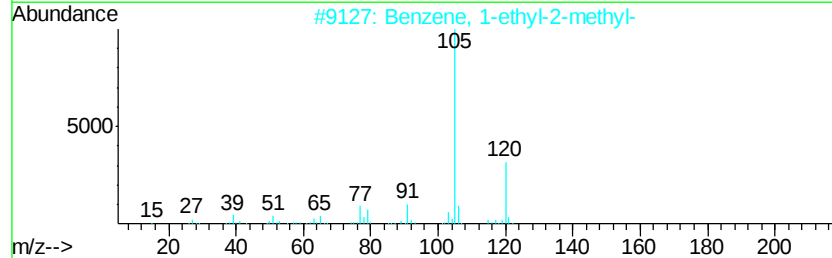
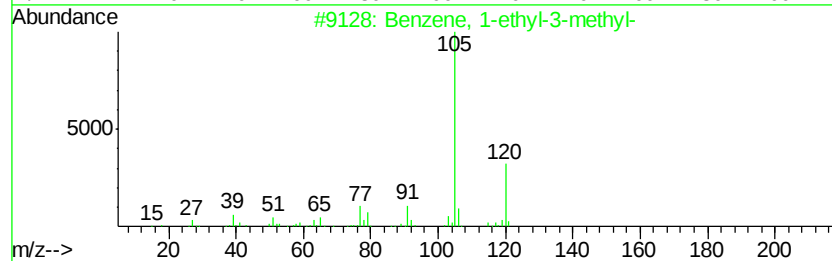
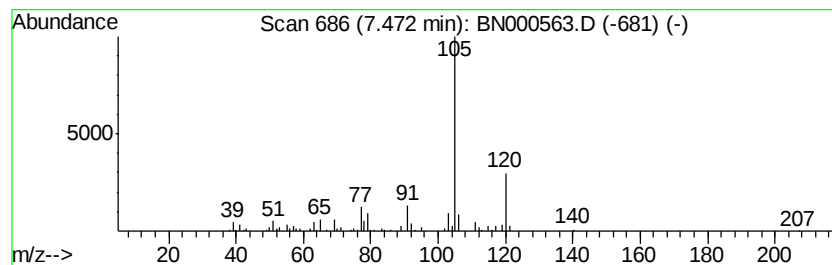
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Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 5 Benzene, 1-ethyl-3-methyl- Concentration Rank 31

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 7.47 | 4.16 ng/ul | 209383 | 1,4-Dichlorobenzene-d4 | 8.41 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 97   |
| 2    |    | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 95   |
| 3    |    | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 95   |
| 4    |    | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 95   |
| 5    |    | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 93   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampled :  
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Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

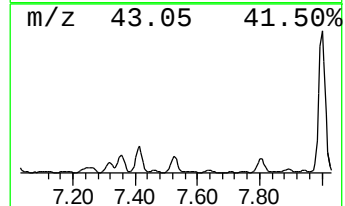
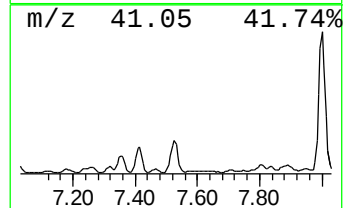
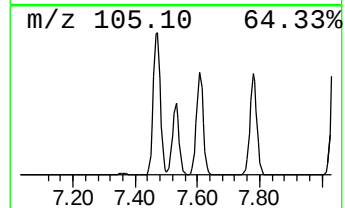
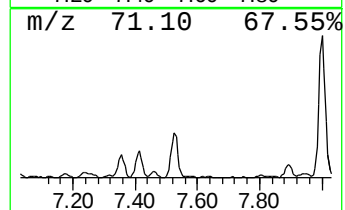
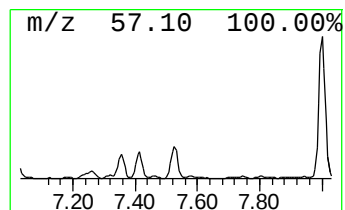
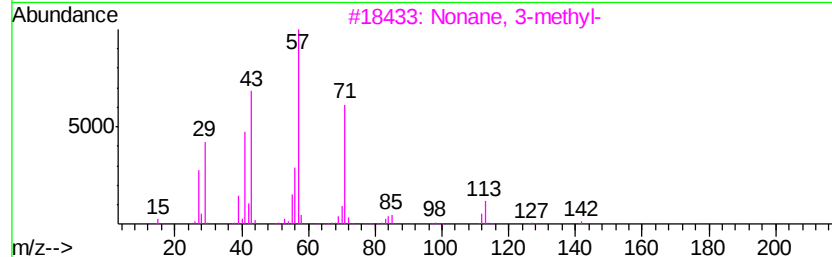
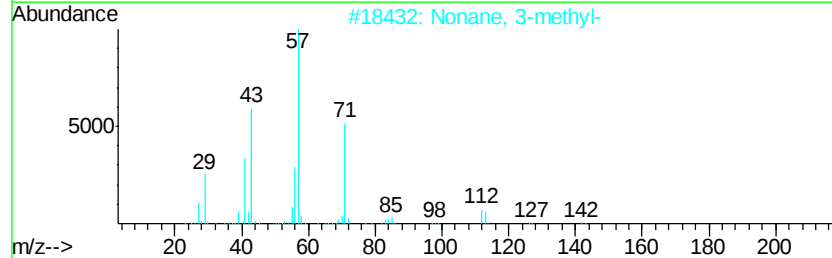
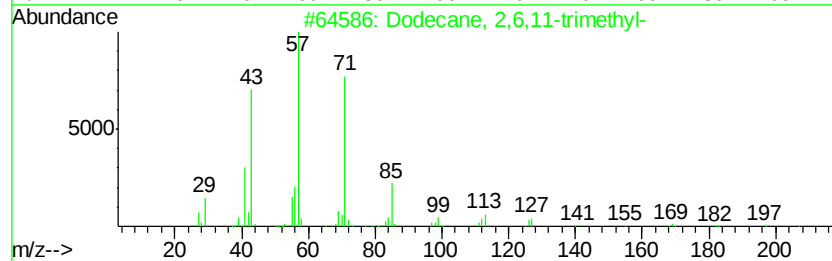
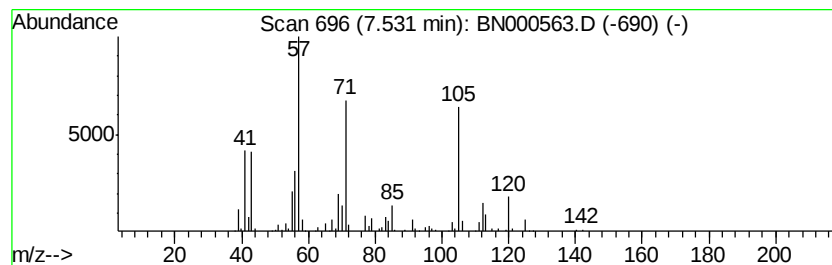
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 9

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 7.53 | 7.70 ng/ul | 388101 | 1,4-Dichlorobenzene-d4 | 8.41 |

| Hit# | of | 5 | Tentative ID                    | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------------|-----|---------|-------------|------|
| 1    |    |   | Dodecane, 2,6,11-trimethyl-     | 212 | C15H32  | 031295-56-4 | 64   |
| 2    |    |   | Nonane, 3-methyl-               | 142 | C10H22  | 005911-04-6 | 59   |
| 3    |    |   | Nonane, 3-methyl-               | 142 | C10H22  | 005911-04-6 | 59   |
| 4    |    |   | Undecane, 2,6-dimethyl-         | 184 | C13H28  | 017301-23-4 | 59   |
| 5    |    |   | 2-Isopropyl-5-methyl-1-heptanol | 172 | C11H24O | 091337-07-4 | 53   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

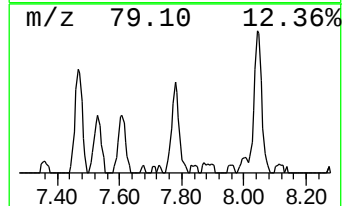
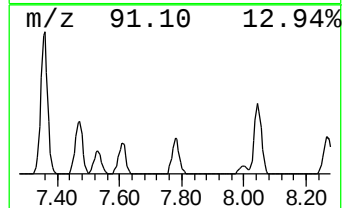
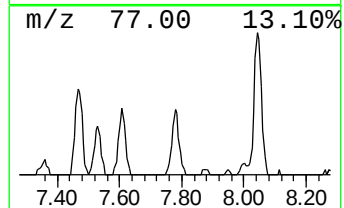
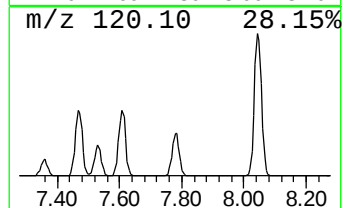
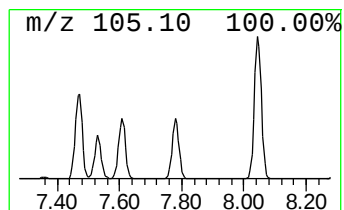
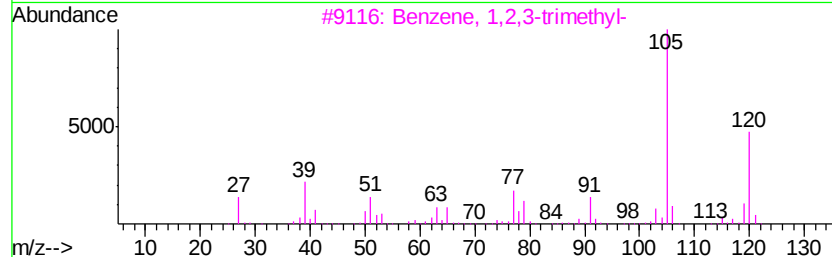
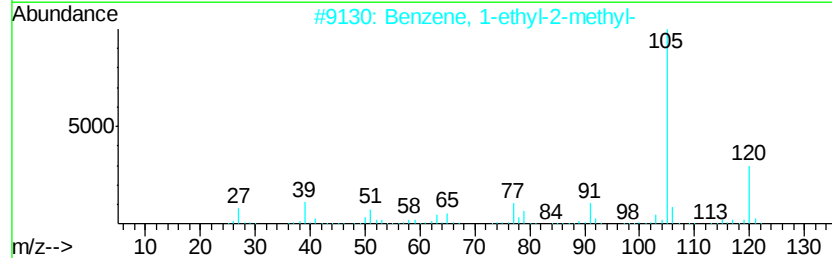
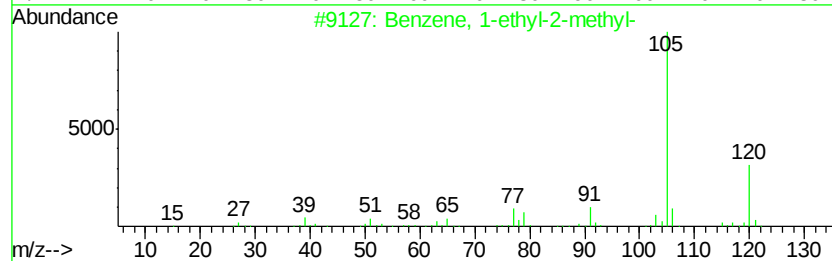
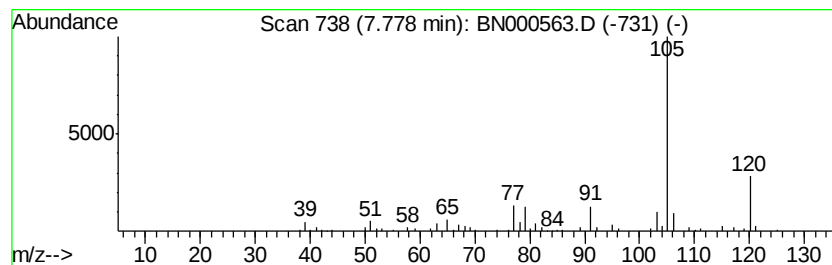
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Quant Title : SVOA CALIBRATION

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

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Peak Number 8 Benzene, 1-ethyl-2-methyl- Concentration Rank 39

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 7.78 | 3.39 ng/ul | 170953 | 1,4-Dichlorobenzene-d4 | 8.41 |

| 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, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 91   |
| 4    |    |   | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 91   |
| 5    |    |   | Benzene, 1,3,5-trimethyl-  | 120 | C9H12   | 000108-67-8 | 91   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

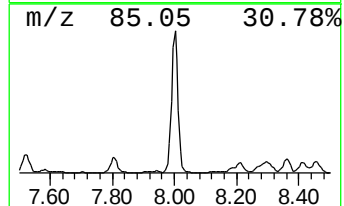
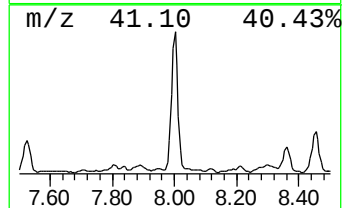
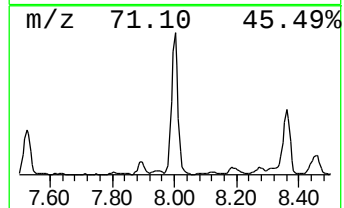
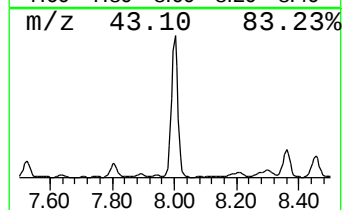
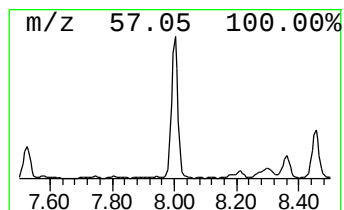
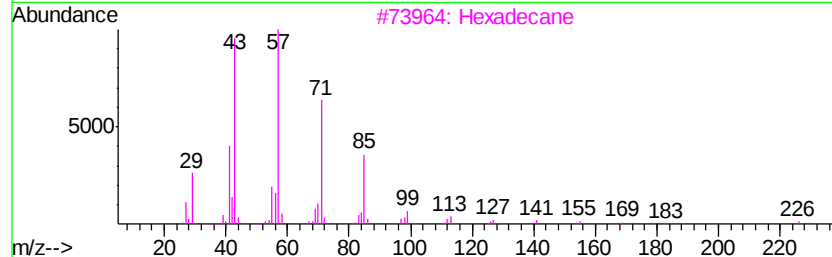
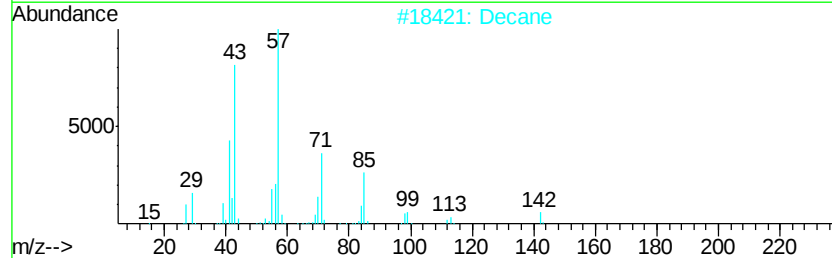
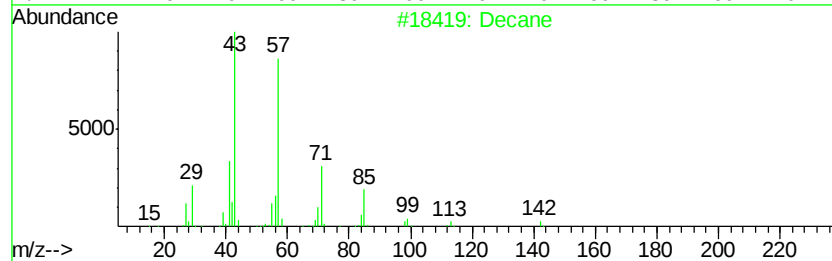
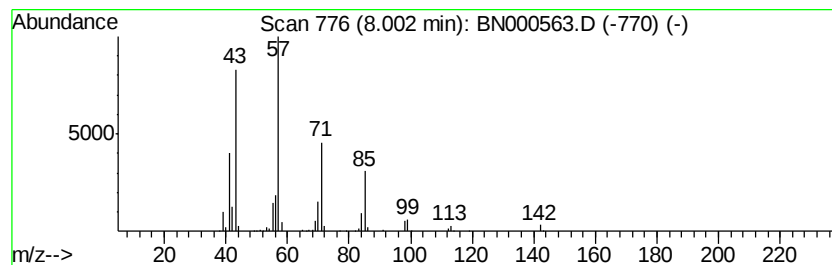
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TIC Integration Parameters: LSCINT.P

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Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 2

| R.T. | EstConc     | Area    | Relative to ISTD       | R.T. |
|------|-------------|---------|------------------------|------|
| 8.00 | 22.66 ng/ul | 1141830 | 1,4-Dichlorobenzene-d4 | 8.41 |

| Hit# | of | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|--------------|-----|---------|-------------|------|
| 1    | 5  | Decane       | 142 | C10H22  | 000124-18-5 | 94   |
| 2    |    | Decane       | 142 | C10H22  | 000124-18-5 | 91   |
| 3    |    | Hexadecane   | 226 | C16H34  | 000544-76-3 | 90   |
| 4    |    | Decane       | 142 | C10H22  | 000124-18-5 | 90   |
| 5    |    | Tridecane    | 184 | C13H28  | 000629-50-5 | 86   |



Data Path : Z:\HPCHEM1\BNA\_N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampled :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

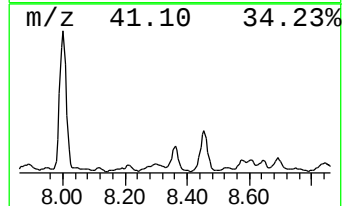
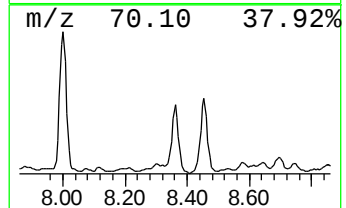
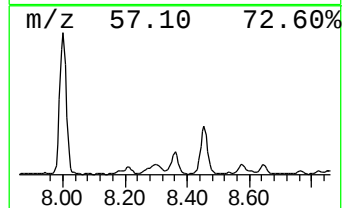
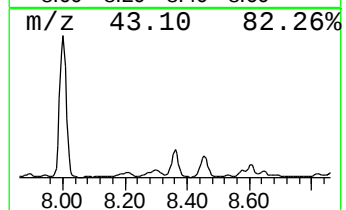
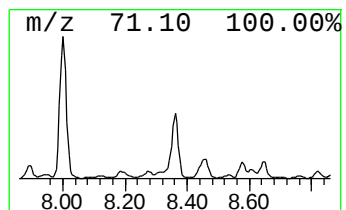
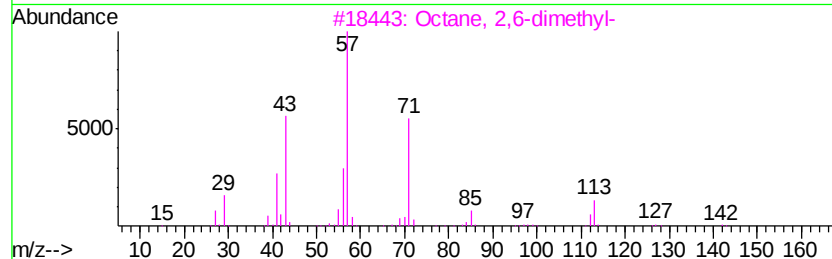
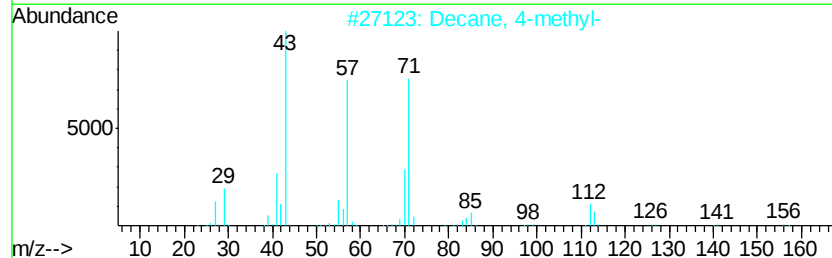
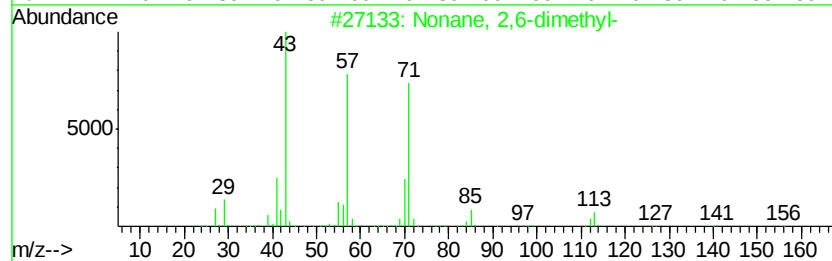
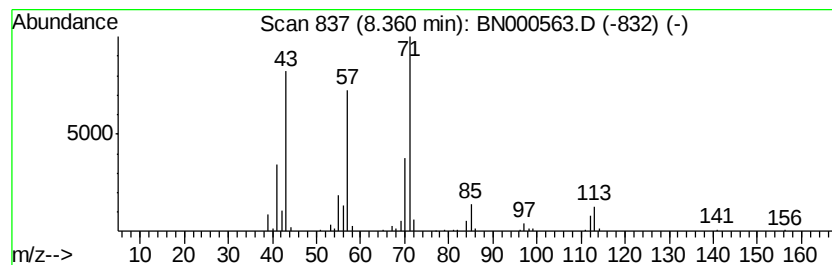
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 29

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 8.36 | 4.21 ng/ul | 211907 | 1,4-Dichlorobenzene-d4 | 8.41 |

| Hit# | of | 5 | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------|-----|---------|-------------|------|
| 1    |    |   | Nonane, 2,6-dimethyl-     | 156 | C11H24  | 017302-28-2 | 91   |
| 2    |    |   | Decane, 4-methyl-         | 156 | C11H24  | 002847-72-5 | 83   |
| 3    |    |   | Octane, 2,6-dimethyl-     | 142 | C10H22  | 002051-30-1 | 80   |
| 4    |    |   | Heptane, 3,3,5-trimethyl- | 142 | C10H22  | 007154-80-5 | 78   |
| 5    |    |   | Octane, 3,3-dimethyl-     | 142 | C10H22  | 004110-44-5 | 78   |



Data Path : Z:\HPCHEM1\BNA\_N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

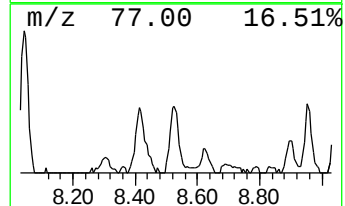
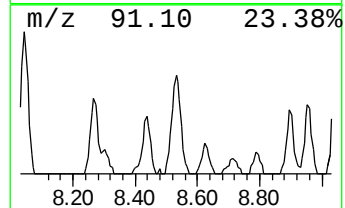
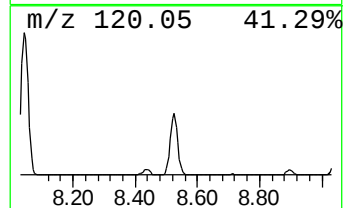
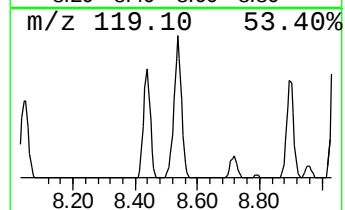
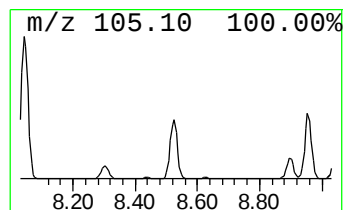
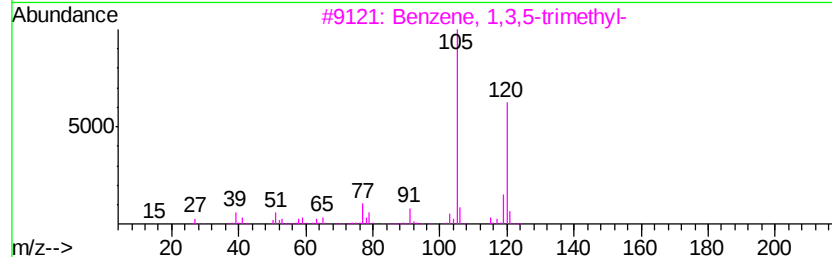
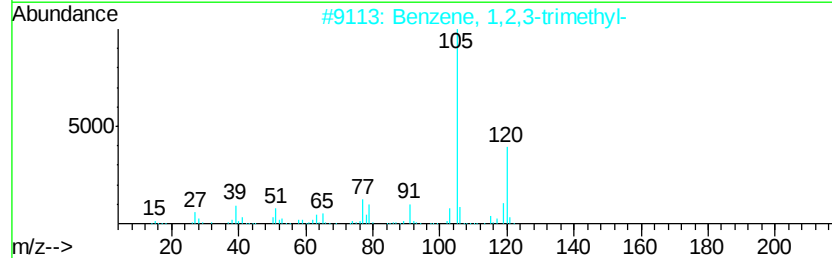
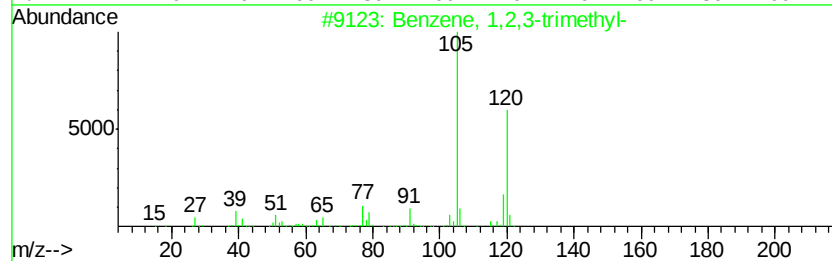
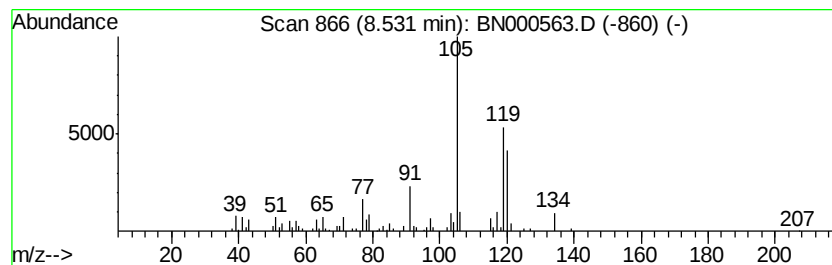
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 Benzene, 1,2,3-trimethyl- Concentration Rank 24

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 8.53 | 4.67 ng/ul | 235286 | 1,4-Dichlorobenzene-d4 | 8.41 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 81   |
| 2    |    | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 81   |
| 3    |    | Benzene, 1,3,5-trimethyl-  | 120 | C9H12   | 000108-67-8 | 81   |
| 4    |    | Benzene, 1,2,4-trimethyl-  | 120 | C9H12   | 000095-63-6 | 81   |
| 5    |    | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 76   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

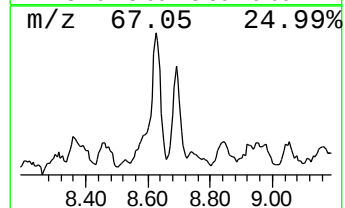
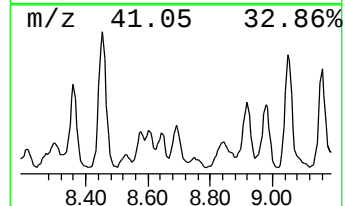
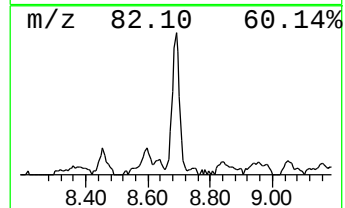
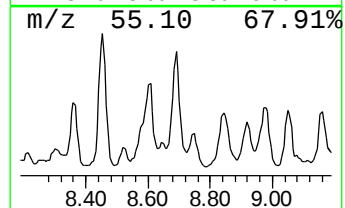
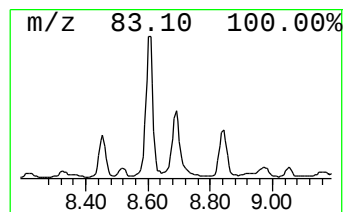
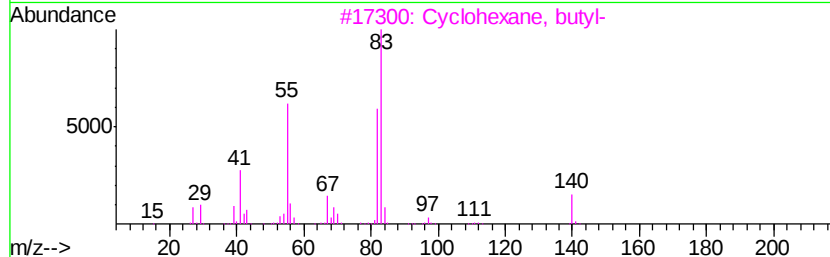
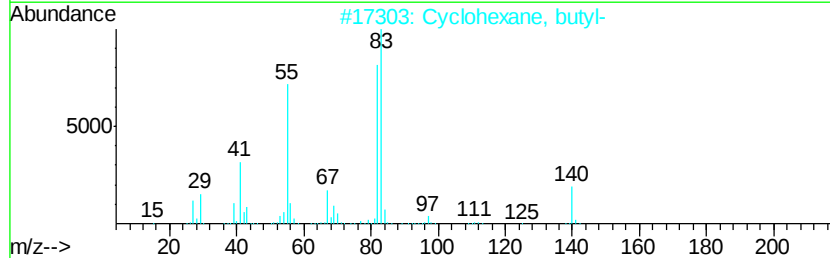
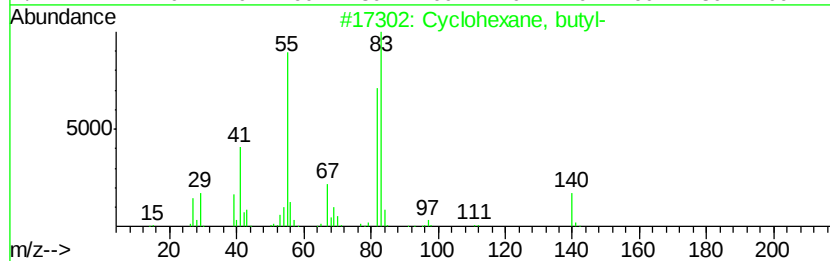
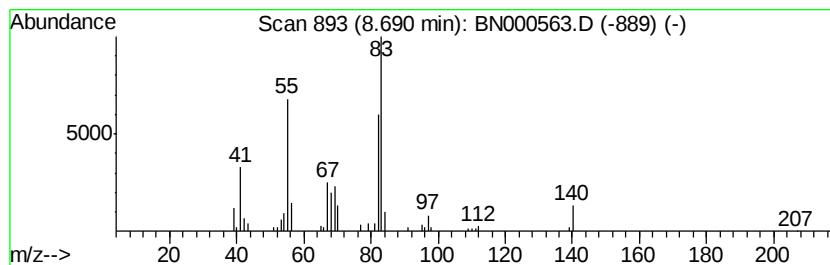
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 13 (DEL) Alkane: Cyclic8.69 Concentration Rank 53

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 8.69 | 2.62 ng/ul | 131942 | 1,4-Dichlorobenzene-d4 | 8.41 |

| Hit# | of | 5 | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, butyl-                | 140 | C10H20  | 001678-93-9 | 72   |
| 2    |    |   | Cyclohexane, butyl-                | 140 | C10H20  | 001678-93-9 | 72   |
| 3    |    |   | Cyclohexane, butyl-                | 140 | C10H20  | 001678-93-9 | 72   |
| 4    |    |   | Cyclohexane, octyl-                | 196 | C14H28  | 001795-15-9 | 64   |
| 5    |    |   | Cyclohexane, 1,1'-(1,4-butanedi... | 222 | C16H30  | 006165-44-2 | 64   |



Data Path : Z:\HPCHEM1\BNA\_N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

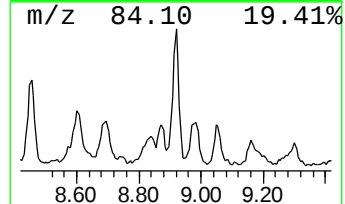
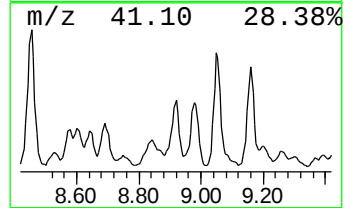
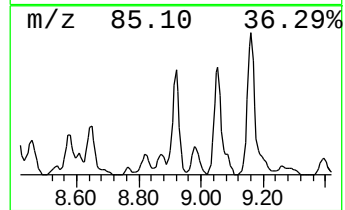
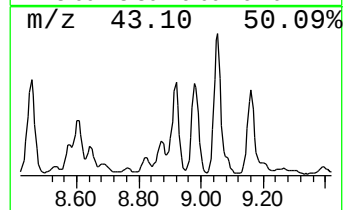
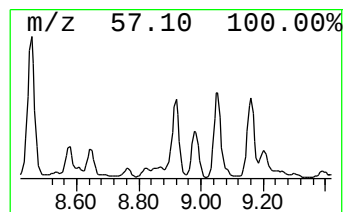
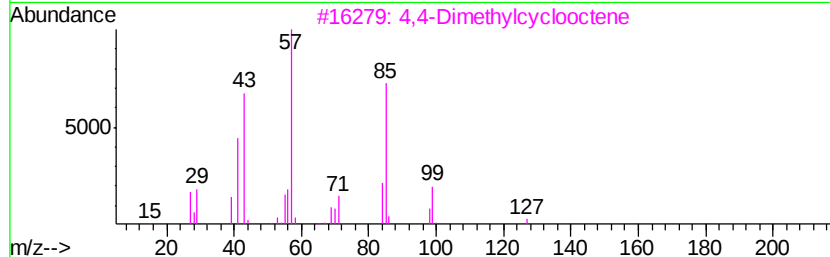
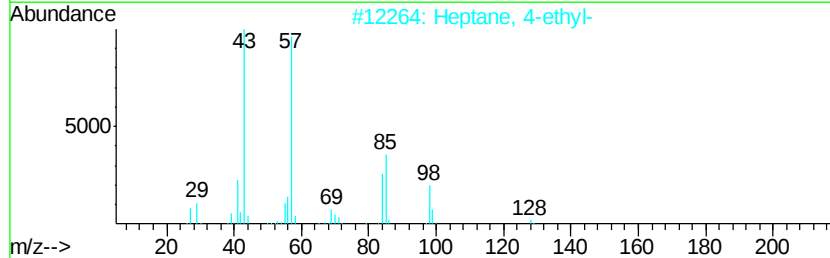
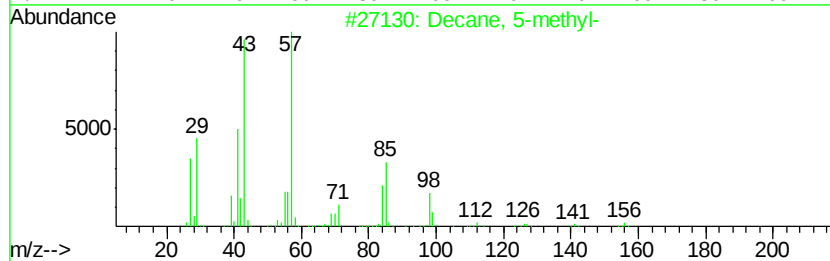
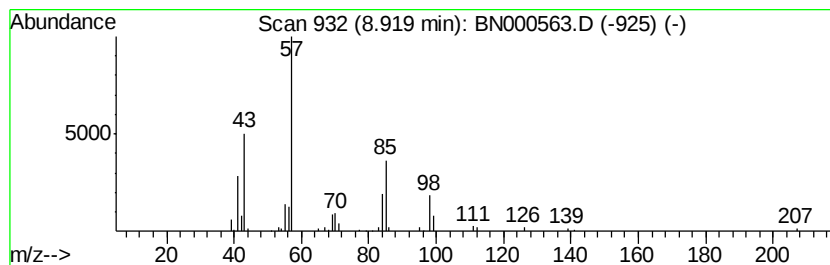
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 23

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 8.92 | 4.95 ng/ul | 249193 | 1,4-Dichlorobenzene-d4 | 8.41 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------|-----|---------|--------------|------|
| 1    |    |   | Decane, 5-methyl-       | 156 | C11H24  | 013151-35-4  | 64   |
| 2    |    |   | Heptane, 4-ethyl-       | 128 | C9H20   | 002216-32-2  | 59   |
| 3    |    |   | 4,4-Dimethylcyclooctene | 138 | C10H18  | 1000113-56-7 | 53   |
| 4    |    |   | Pentadecane             | 212 | C15H32  | 000629-62-9  | 53   |
| 5    |    |   | Heptane, 4-ethyl-       | 128 | C9H20   | 002216-32-2  | 53   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

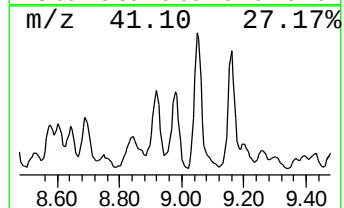
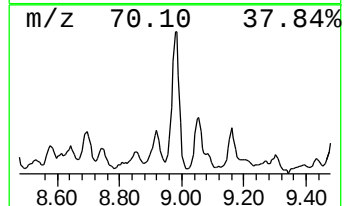
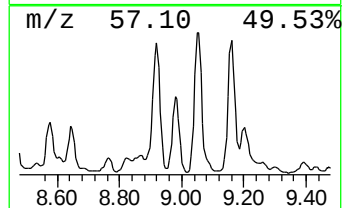
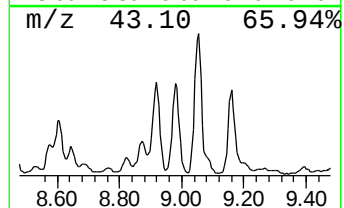
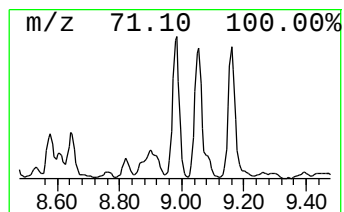
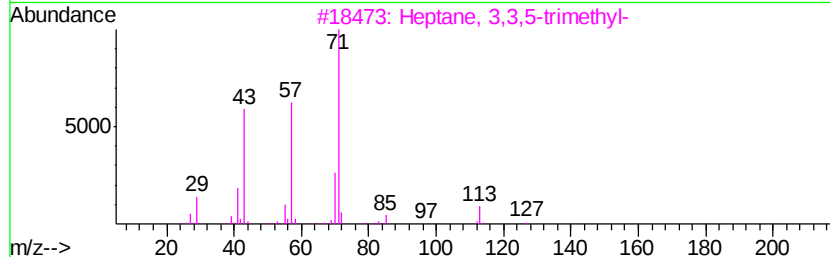
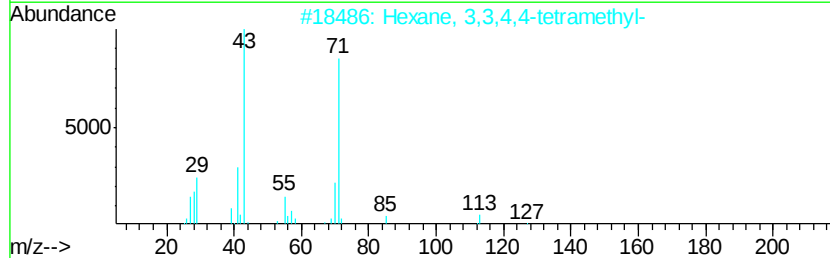
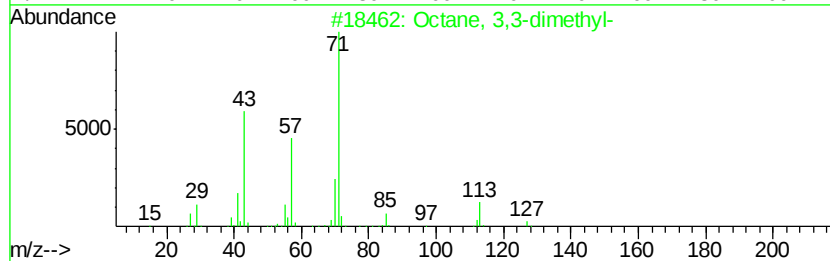
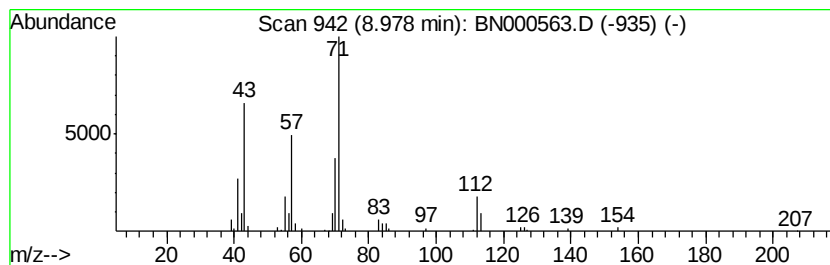
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 8

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 8.98 | 7.78 ng/ul | 391756 | 1,4-Dichlorobenzene-d4 | 8.41 |

| Hit# | of | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------|-----|---------|-------------|------|
| 1    | 5  | Octane, 3,3-dimethyl-        | 142 | C10H22  | 004110-44-5 | 72   |
| 2    |    | Hexane, 3,3,4,4-tetramethyl- | 142 | C10H22  | 005171-84-6 | 64   |
| 3    |    | Heptane, 3,3,5-trimethyl-    | 142 | C10H22  | 007154-80-5 | 64   |
| 4    |    | Octane, 2-bromo-             | 192 | C8H17Br | 000557-35-7 | 59   |
| 5    |    | Octane, 2,6,6-trimethyl-     | 156 | C11H24  | 054166-32-4 | 59   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleID :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

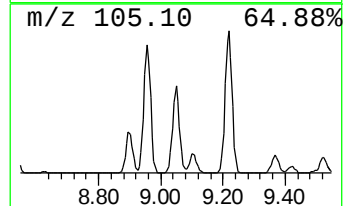
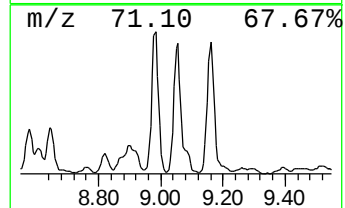
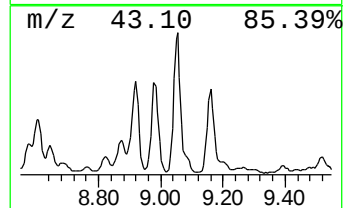
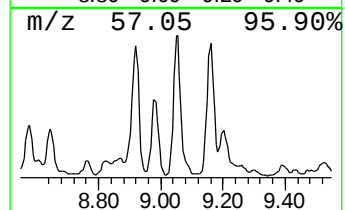
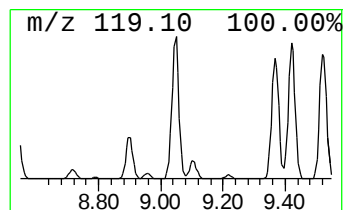
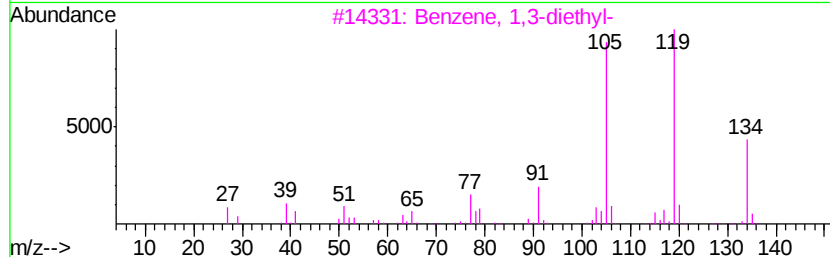
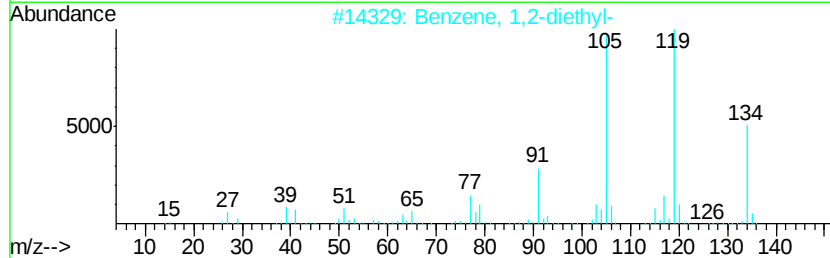
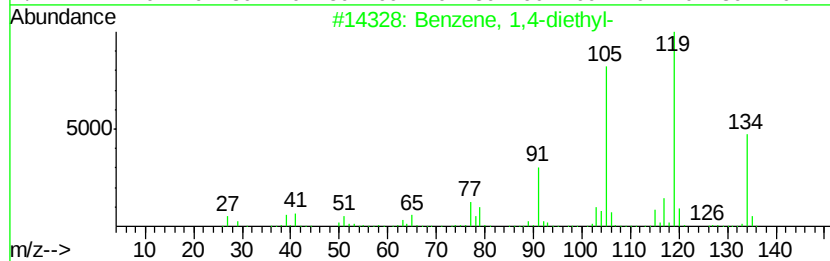
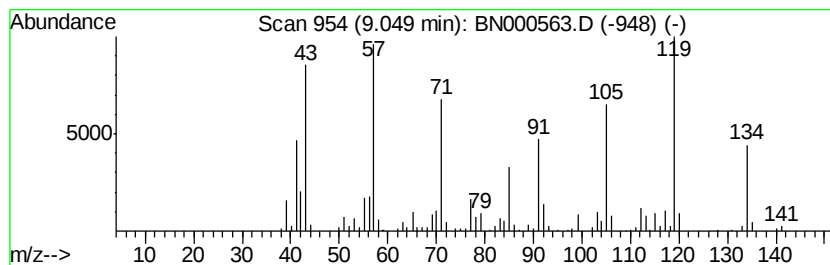
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 16 Benzene, 1,4-diethyl- Concentration Rank 4

| R.T. | EstConc     | Area   | Relative to ISTD       | R.T. |
|------|-------------|--------|------------------------|------|
| 9.05 | 12.23 ng/ul | 616280 | 1,4-Dichlorobenzene-d4 | 8.41 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,4-diethyl-               | 134 | C10H14  | 000105-05-5 | 95   |
| 2    |    | Benzene, 1,2-diethyl-               | 134 | C10H14  | 000135-01-3 | 92   |
| 3    |    | Benzene, 1,3-diethyl-               | 134 | C10H14  | 000141-93-5 | 90   |
| 4    |    | Benzene, 2-ethyl-1,4-dimethyl-      | 134 | C10H14  | 001758-88-9 | 90   |
| 5    |    | Benzene, 1-methyl-2-(1-methyleth... | 134 | C10H14  | 000527-84-4 | 78   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

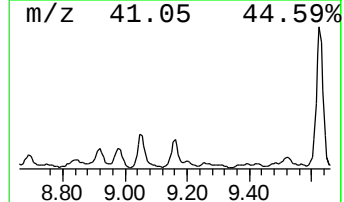
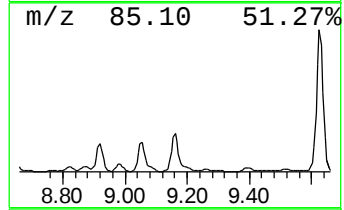
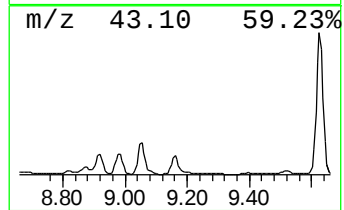
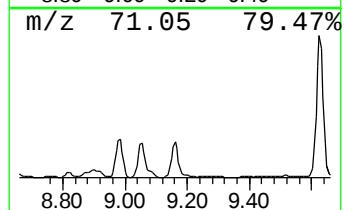
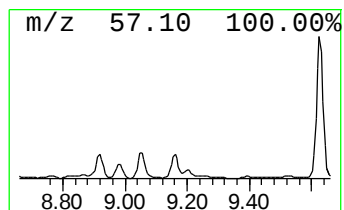
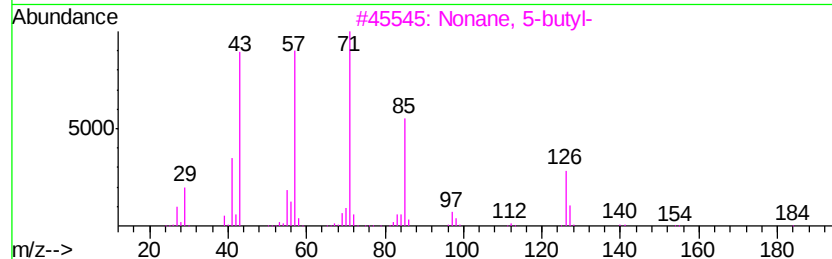
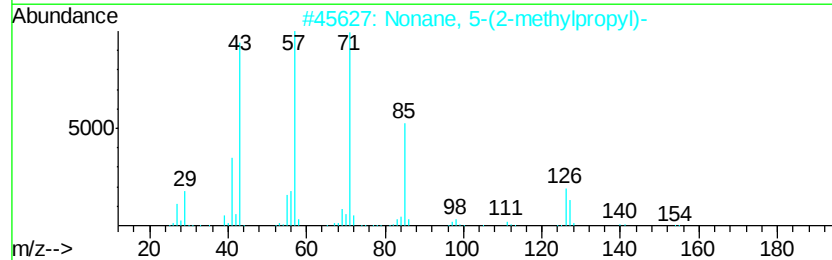
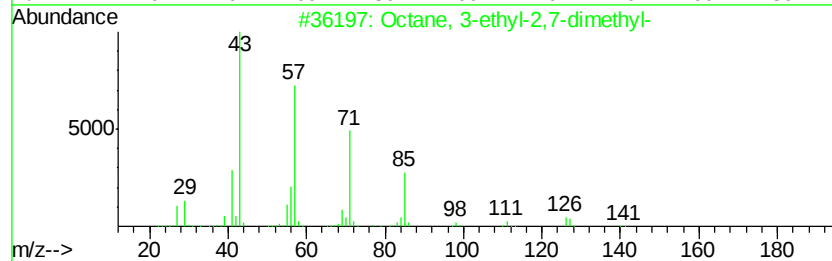
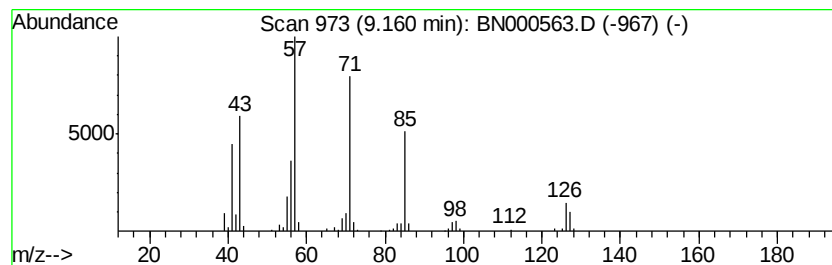
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 21

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 9.16 | 5.46 ng/ul | 275076 | 1,4-Dichlorobenzene-d4 | 8.41 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | Octane, 3-ethyl-2,7-dimethyl- | 170 | C12H26  | 062183-55-5 | 83   |
| 2    |    |   | Nonane, 5-(2-methylpropyl)-   | 184 | C13H28  | 062185-53-9 | 78   |
| 3    |    |   | Nonane, 5-butyl-              | 184 | C13H28  | 017312-63-9 | 72   |
| 4    |    |   | Undecane, 2,7-dimethyl-       | 184 | C13H28  | 017301-24-5 | 72   |
| 5    |    |   | Nonane, 3,7-dimethyl-         | 156 | C11H24  | 017302-32-8 | 64   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

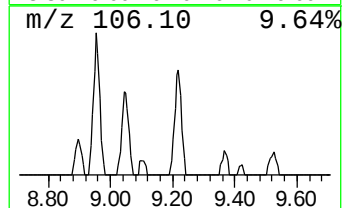
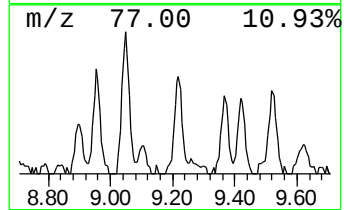
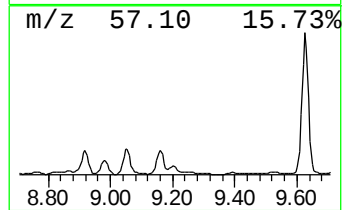
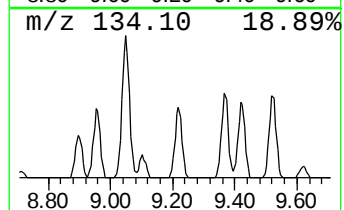
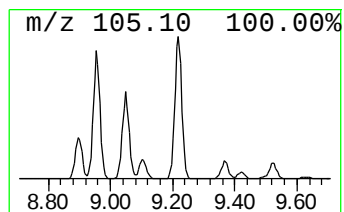
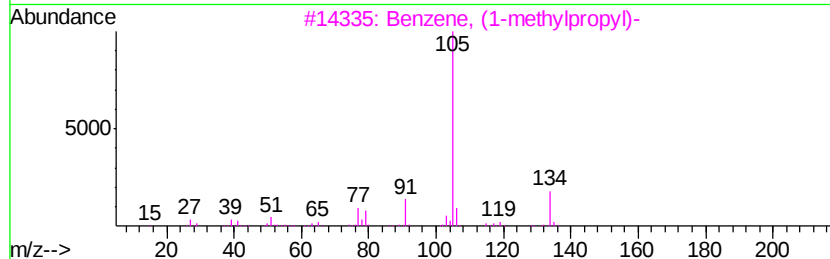
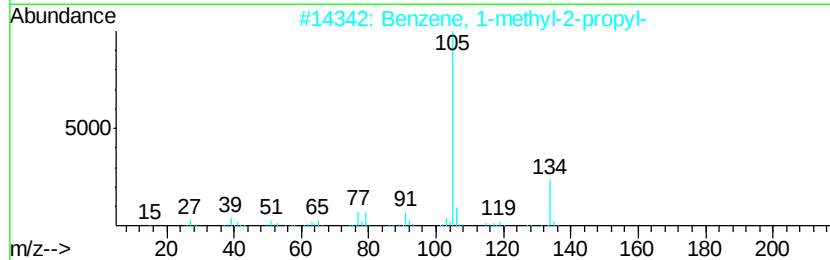
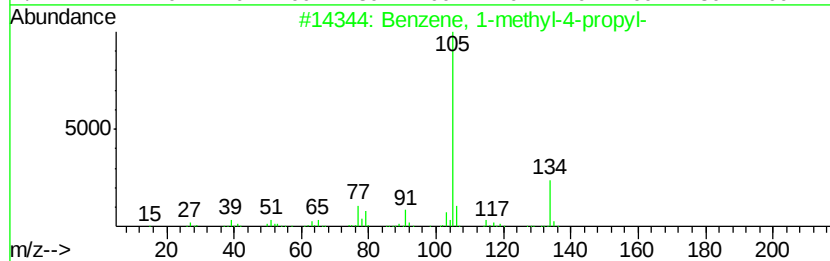
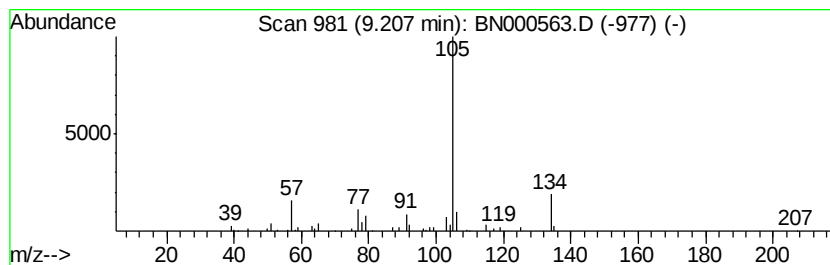
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 18 Benzene, 1-methyl-4-propyl- Concentration Rank 40

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 9.21 | 3.35 ng/ul | 168681 | 1,4-Dichlorobenzene-d4 | 8.41 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-methyl-4-propyl- | 134 | C10H14  | 001074-55-1 | 94   |
| 2    |    | Benzene, 1-methyl-2-propyl- | 134 | C10H14  | 001074-17-5 | 91   |
| 3    |    | Benzene, (1-methylpropyl)-  | 134 | C10H14  | 000135-98-8 | 91   |
| 4    |    | Benzene, 1-methyl-4-propyl- | 134 | C10H14  | 001074-55-1 | 91   |
| 5    |    | Benzene, 1-methyl-2-propyl- | 134 | C10H14  | 001074-17-5 | 90   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

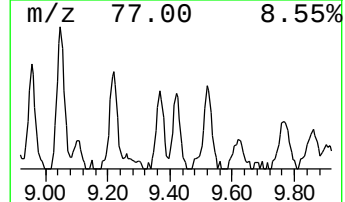
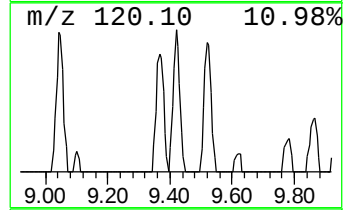
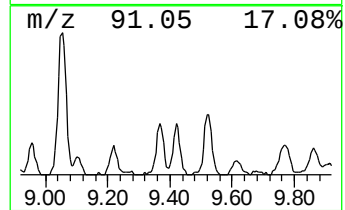
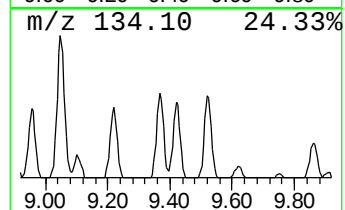
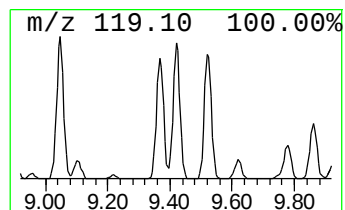
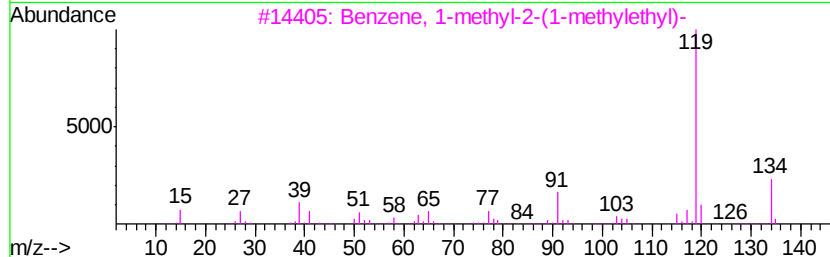
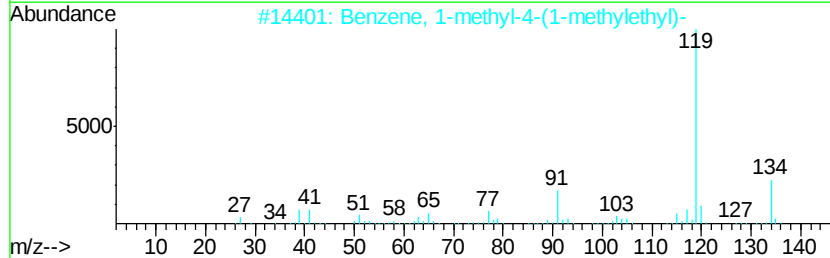
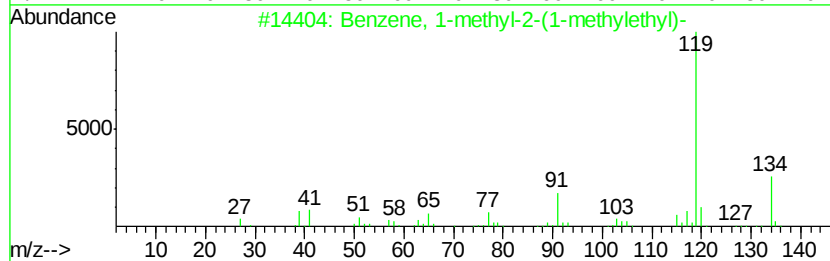
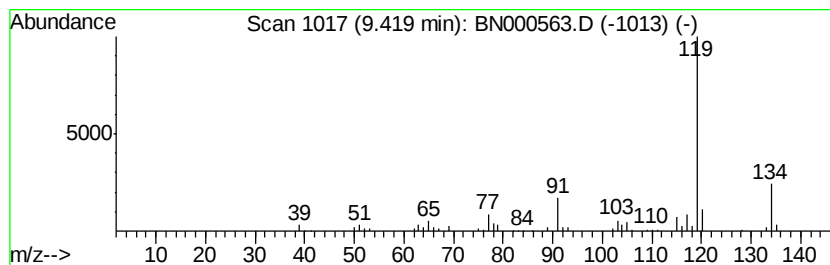
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 20 Benzene, 1-methyl-2-(1-meth... Concentration Rank 35

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 9.42 | 3.67 ng/ul | 184990 | 1,4-Dichlorobenzene-d4 | 8.41 |

| Hit# | of | Tentative ID                         | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1-methyl-2-(1-methylethyl)- | 134 | C10H14  | 000527-84-4 | 97   |
| 2    |    | Benzene, 1-methyl-4-(1-methylethyl)- | 134 | C10H14  | 000099-87-6 | 97   |
| 3    |    | Benzene, 1-methyl-2-(1-methylethyl)- | 134 | C10H14  | 000527-84-4 | 97   |
| 4    |    | Benzene, 1-methyl-2-(1-methylethyl)- | 134 | C10H14  | 000527-84-4 | 95   |
| 5    |    | Benzene, 1-methyl-3-(1-methylethyl)- | 134 | C10H14  | 000535-77-3 | 95   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

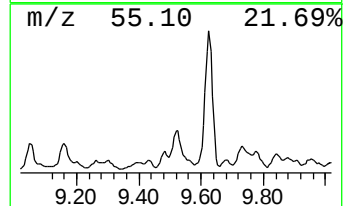
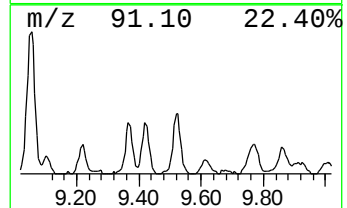
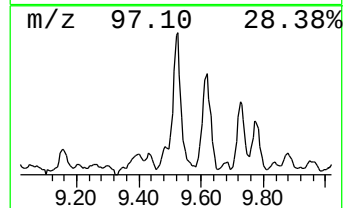
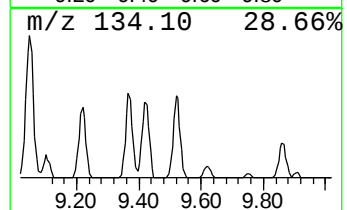
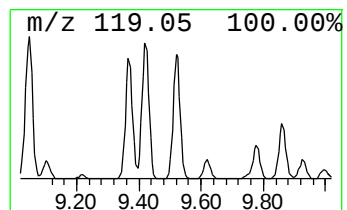
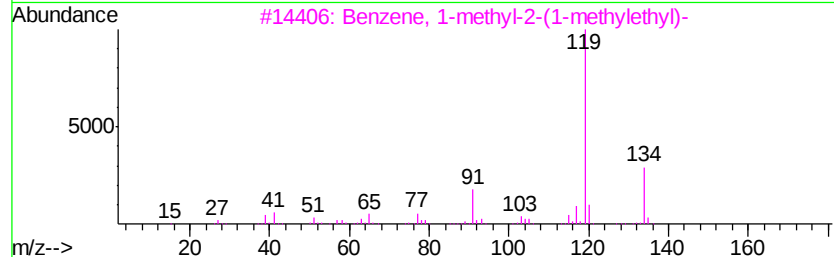
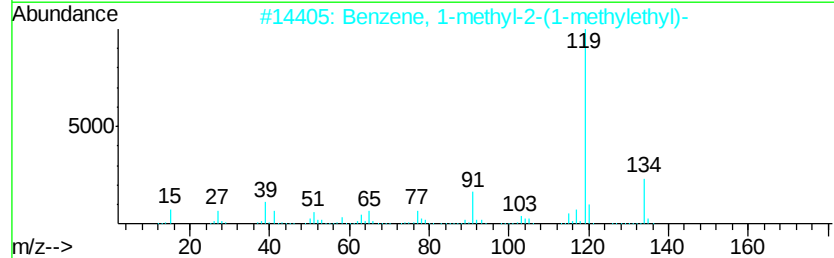
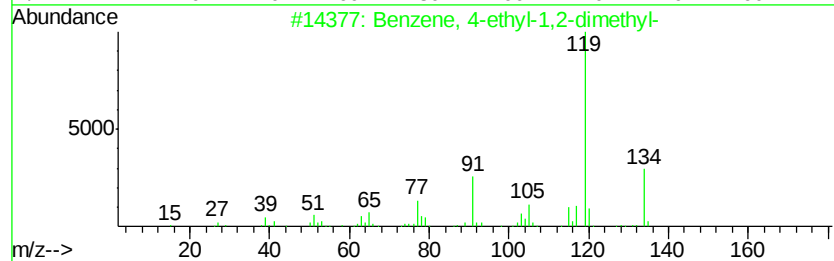
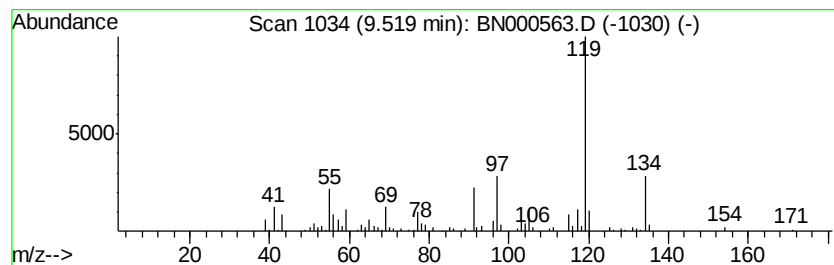
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 21 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 18

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 9.52 | 6.00 ng/ul | 302518 | 1,4-Dichlorobenzene-d4 | 8.41 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 4-ethyl-1,2-dimethyl-      | 134 | C10H14  | 000934-80-5 | 96   |
| 2    |    | Benzene, 1-methyl-2-(1-methyleth... | 134 | C10H14  | 000527-84-4 | 95   |
| 3    |    | Benzene, 1-methyl-2-(1-methyleth... | 134 | C10H14  | 000527-84-4 | 95   |
| 4    |    | Benzene, 1-methyl-4-(1-methyleth... | 134 | C10H14  | 000099-87-6 | 95   |
| 5    |    | Benzene, 1-methyl-2-(1-methyleth... | 134 | C10H14  | 000527-84-4 | 95   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

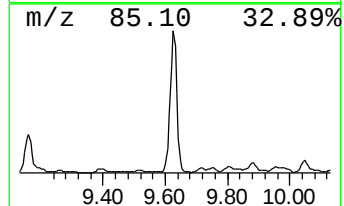
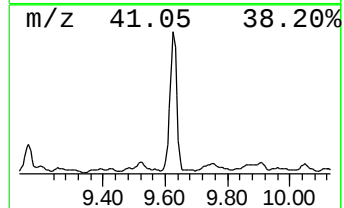
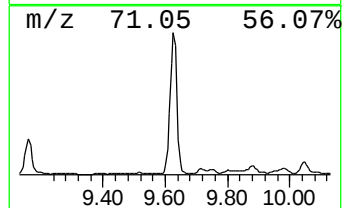
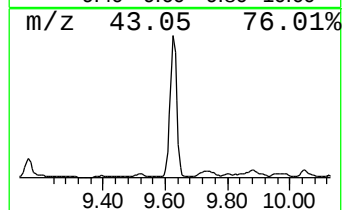
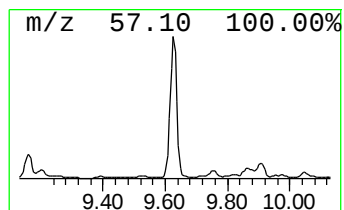
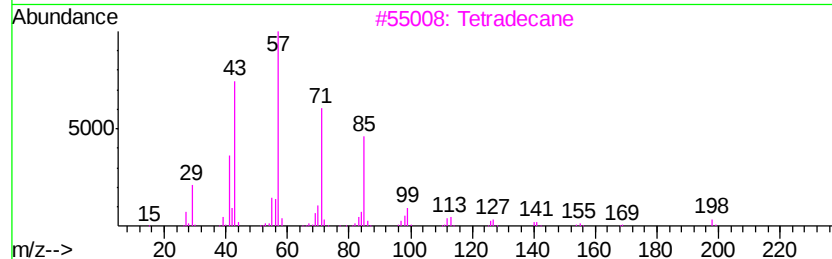
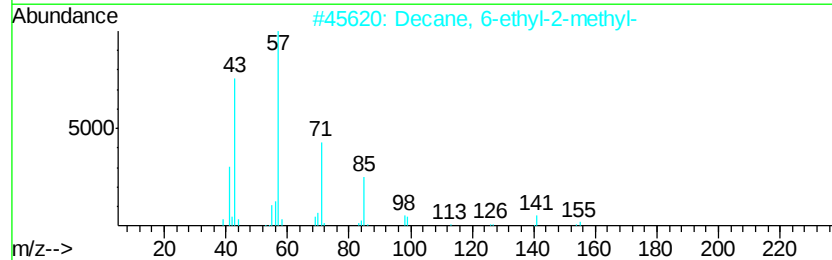
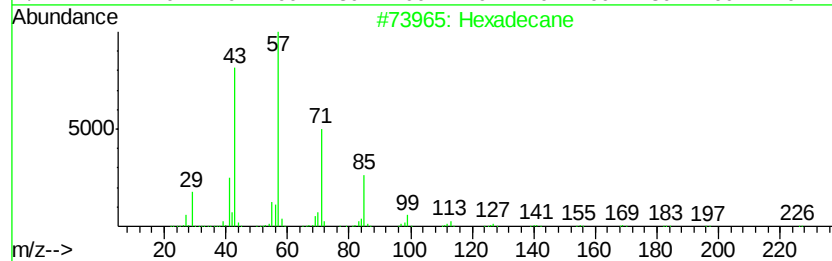
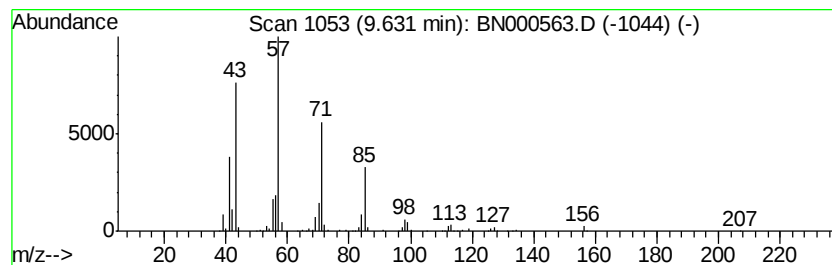
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 1

| R.T. | EstConc     | Area    | Relative to ISTD       | R.T. |
|------|-------------|---------|------------------------|------|
| 9.63 | 28.16 ng/ul | 1418830 | 1,4-Dichlorobenzene-d4 | 8.41 |

| Hit# | of | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---------------------------|-----|---------|-------------|------|
| 1    | 5  | Hexadecane                | 226 | C16H34  | 000544-76-3 | 86   |
| 2    |    | Decane, 6-ethyl-2-methyl- | 184 | C13H28  | 062108-21-8 | 83   |
| 3    |    | Tetradecane               | 198 | C14H30  | 000629-59-4 | 83   |
| 4    |    | Tridecane                 | 184 | C13H28  | 000629-50-5 | 78   |
| 5    |    | Hexadecane                | 226 | C16H34  | 000544-76-3 | 78   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

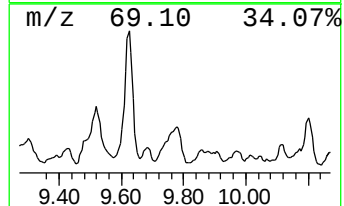
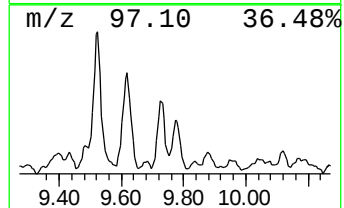
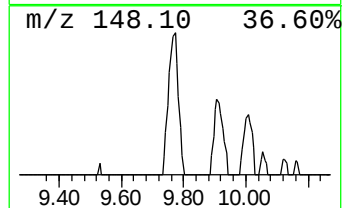
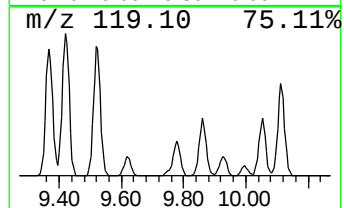
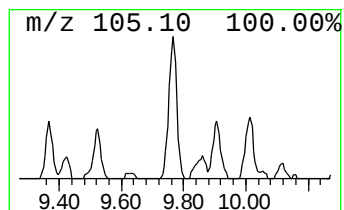
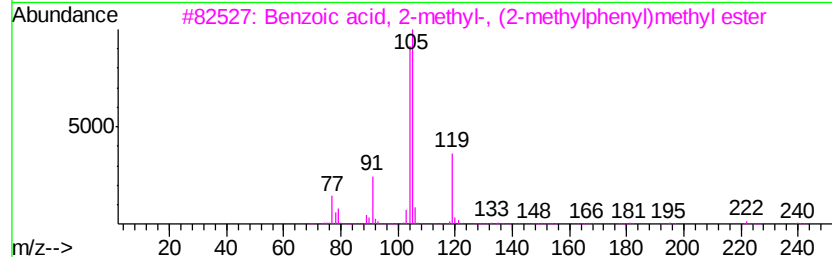
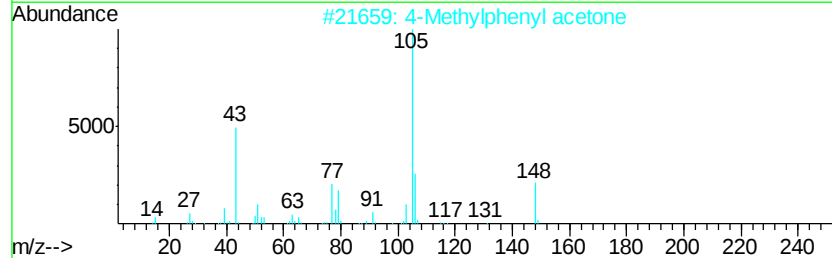
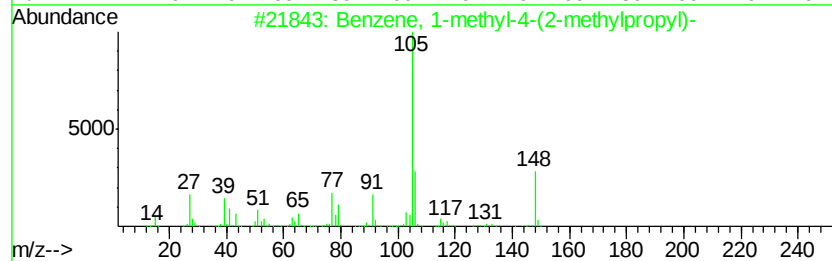
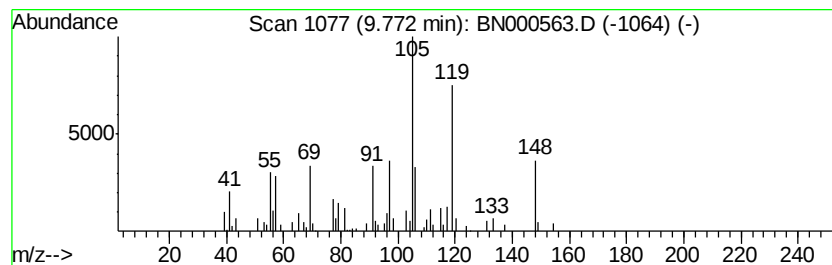
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 23 unknown-01 Concentration Rank 20

| R.T. | EstConc    | Area   | Relative to ISTD       | R.T. |
|------|------------|--------|------------------------|------|
| 9.77 | 5.66 ng/ul | 284965 | 1,4-Dichlorobenzene-d4 | 8.41 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Benzene, 1-methyl-4-(2-methylpro... | 148 | C11H16   | 005161-04-6 | 46   |
| 2    |    |   | 4-Methylphenyl acetone              | 148 | C10H12O  | 002096-86-8 | 38   |
| 3    |    |   | Benzoic acid, 2-methyl-, (2-meth... | 240 | C16H16O2 | 055133-99-8 | 35   |
| 4    |    |   | Benzenepropanal, .beta.-methyl-     | 148 | C10H12O  | 016251-77-7 | 35   |
| 5    |    |   | 2-Methylindan-2-ol                  | 148 | C10H12O  | 033223-84-6 | 35   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampled :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

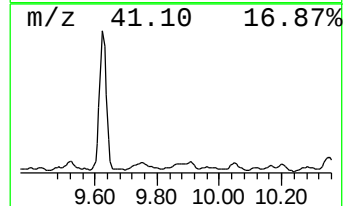
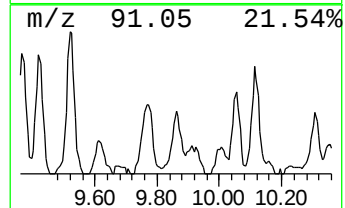
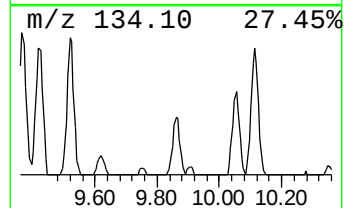
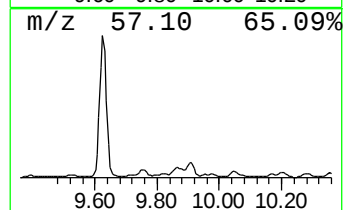
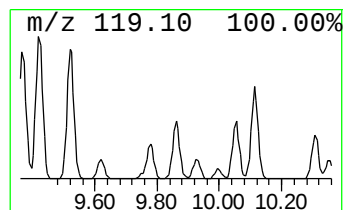
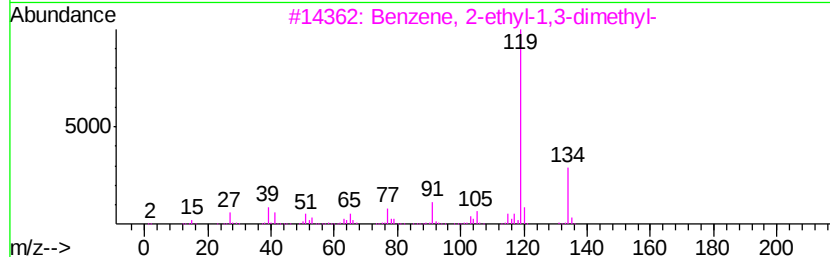
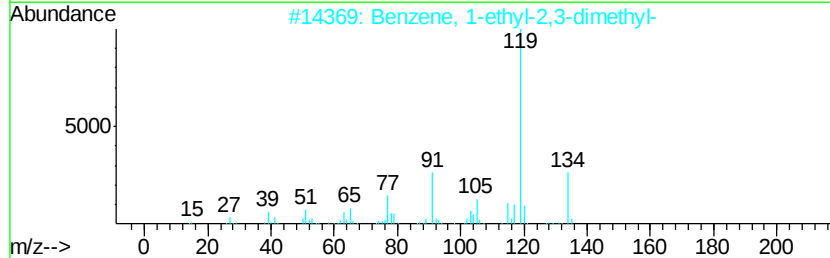
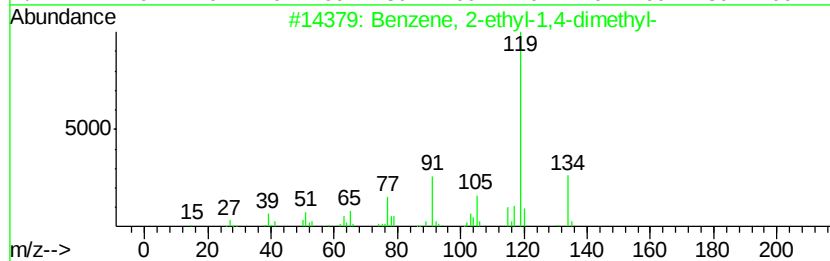
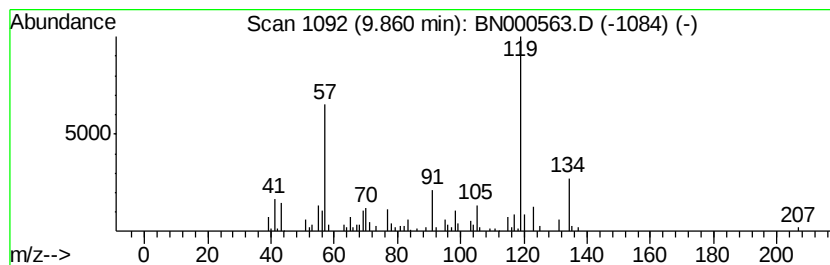
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 24 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 51

| R.T. | EstConc    | Area   | Relative to ISTD | R.T.  |
|------|------------|--------|------------------|-------|
| 9.86 | 2.69 ng/ul | 218205 | Napthalene-d8    | 11.25 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 2-ethyl-1,4-dimethyl-      | 134 | C10H14  | 001758-88-9 | 95   |
| 2    |    | Benzene, 1-ethyl-2,3-dimethyl-      | 134 | C10H14  | 000933-98-2 | 95   |
| 3    |    | Benzene, 2-ethyl-1,3-dimethyl-      | 134 | C10H14  | 002870-04-4 | 95   |
| 4    |    | Benzene, 1-methyl-2-(1-methyleth... | 134 | C10H14  | 000527-84-4 | 94   |
| 5    |    | Benzene, 1-methyl-2-(1-methyleth... | 134 | C10H14  | 000527-84-4 | 94   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

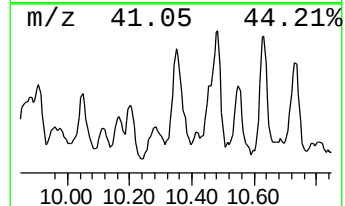
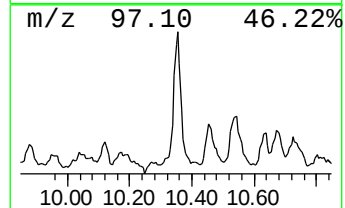
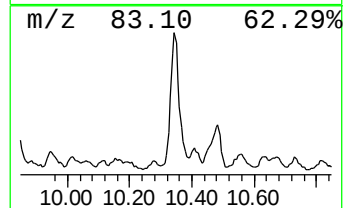
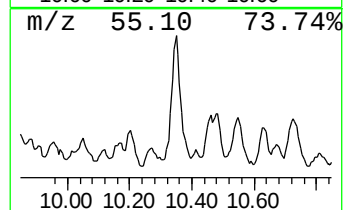
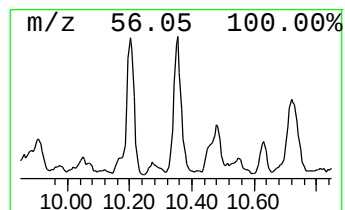
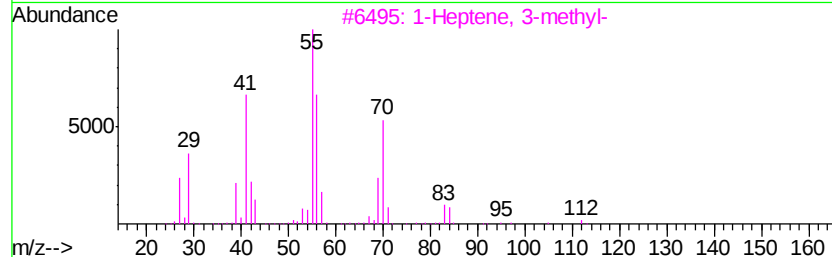
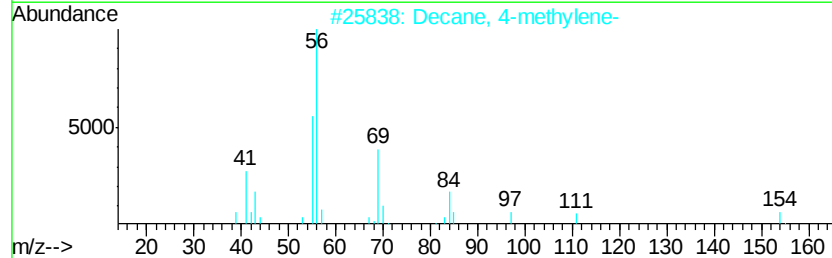
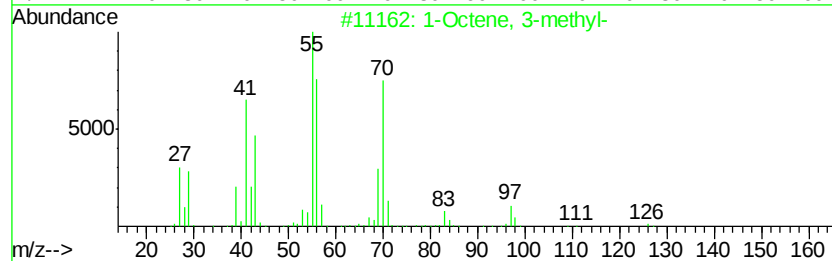
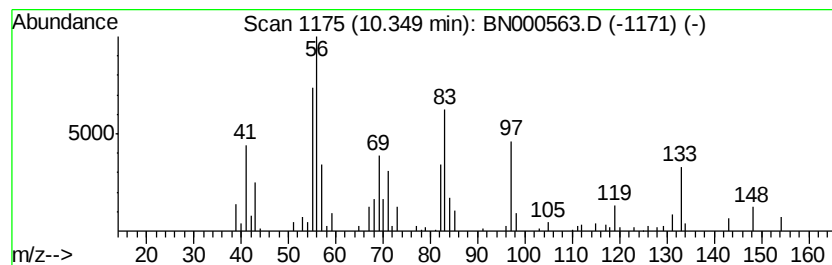
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 25 (DEL) Alkane: Cyclic10.35 Concentration Rank 42

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 10.35 | 3.11 ng/ul | 251834 | Napthalene-d8    | 11.25 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | 1-Octene, 3-methyl-            | 126 | C9H18   | 013151-08-1 | 35   |
| 2    |    |   | Decane, 4-methylene-           | 154 | C11H22  | 024949-41-5 | 30   |
| 3    |    |   | 1-Heptene, 3-methyl-           | 112 | C8H16   | 004810-09-7 | 27   |
| 4    |    |   | 1-Hexene, 2-methyl-            | 98  | C7H14   | 006094-02-6 | 25   |
| 5    |    |   | Cyclopentane, 1,1,3-trimethyl- | 112 | C8H16   | 004516-69-2 | 25   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

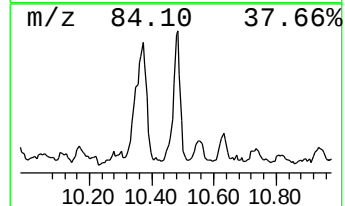
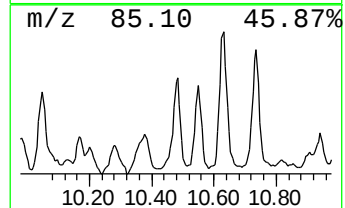
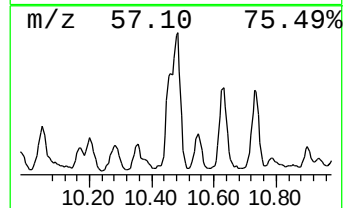
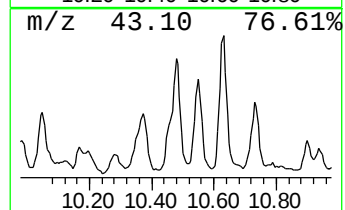
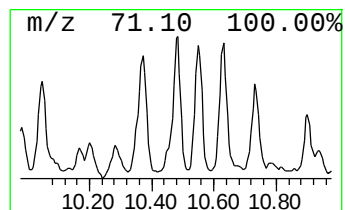
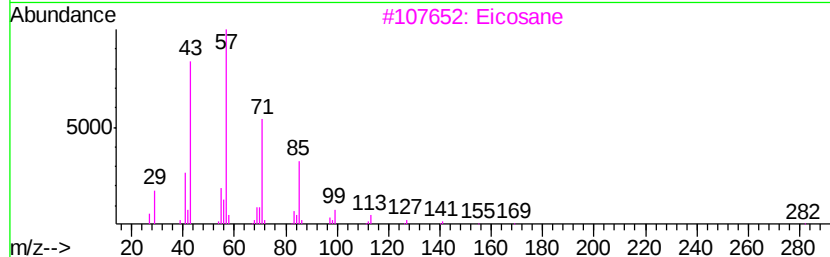
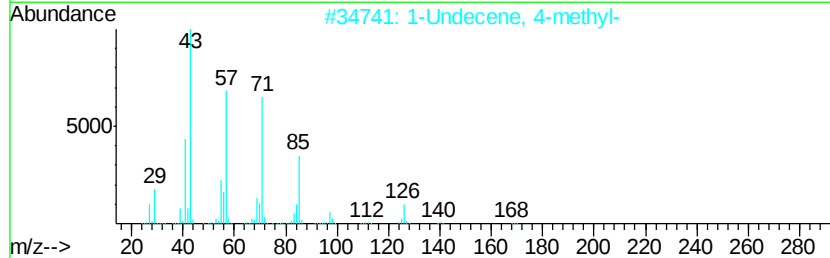
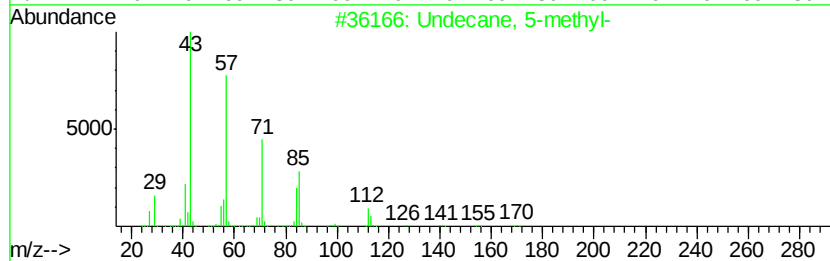
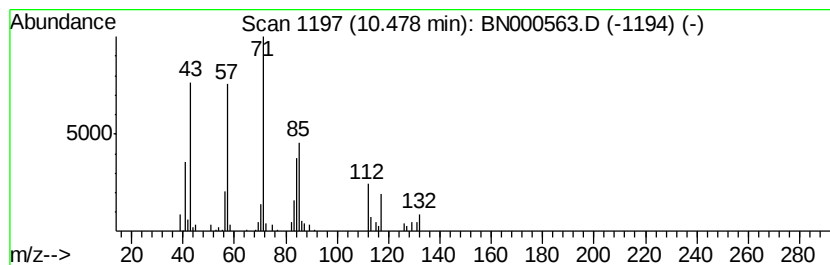
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 26 (DEL) Alkane: Straight-Chai... Concentration Rank 58

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 10.48 | 2.26 ng/ul | 182997 | Naphthalene-d8   | 11.25 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|-------------------------------------|-----|----------|--------------|------|
| 1    | 5  | Undecane, 5-methyl-                 | 170 | C12H26   | 001632-70-8  | 50   |
| 2    |    | 1-Undecene, 4-methyl-               | 168 | C12H24   | 074630-39-0  | 50   |
| 3    |    | Eicosane                            | 282 | C20H42   | 000112-95-8  | 47   |
| 4    |    | Methoxyacetic acid, 2-ethylhexyl... | 202 | C11H22O3 | 1000282-41-4 | 43   |
| 5    |    | Nonane, 5-butyl-                    | 184 | C13H28   | 017312-63-9  | 43   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

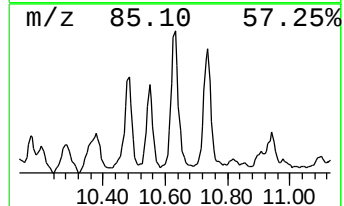
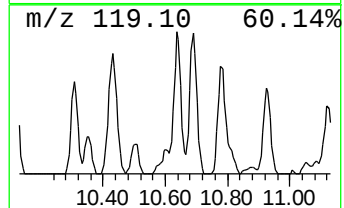
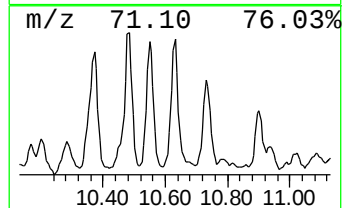
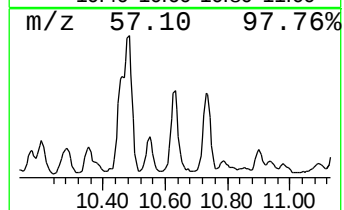
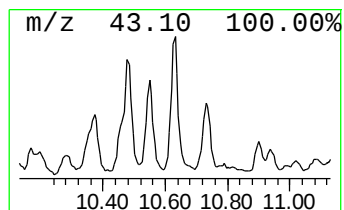
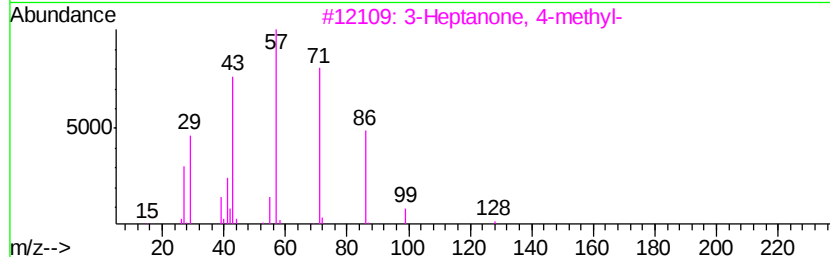
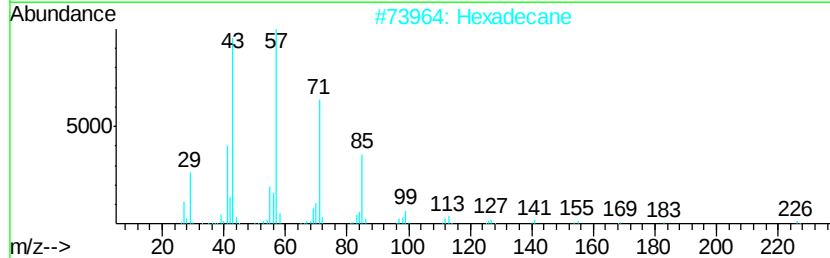
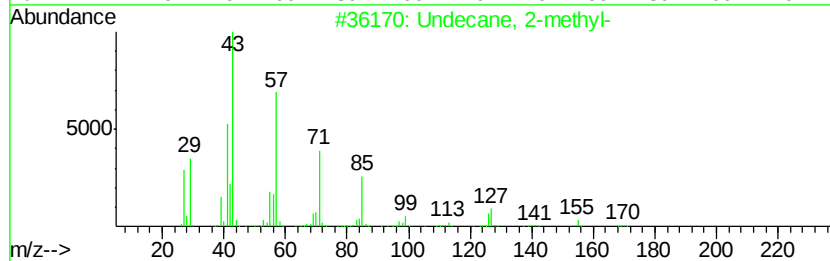
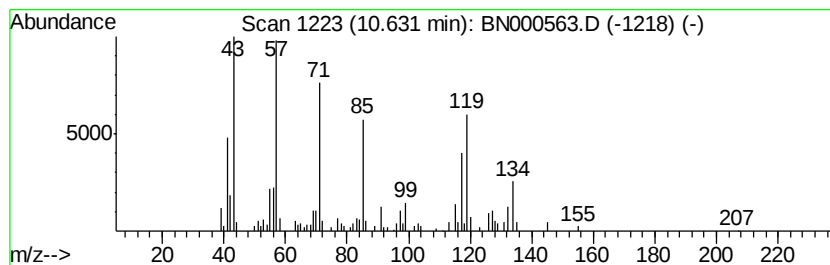
Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 27 (DEL) Alkane: Straight-Chai... Concentration Rank 47

| R.T.    | EstConc    | Area                                | Relative to ISTD | R.T.           |
|---------|------------|-------------------------------------|------------------|----------------|
| 10.63   | 2.79 ng/ul | 226338                              | Napthalene-d8    | 11.25          |
| Hit# of | 5          | Tentative ID                        | MW MolForm       | CAS# Qual      |
| 1       |            | Undecane, 2-methyl-                 | 170 C12H26       | 007045-71-8 46 |
| 2       |            | Hexadecane                          | 226 C16H34       | 000544-76-3 30 |
| 3       |            | 3-Heptanone, 4-methyl-              | 128 C8H16O       | 006137-11-7 18 |
| 4       |            | 1,3-Cyclopentadiene, 1,2,3,4-tet... | 134 C10H14       | 076089-59-3 15 |
| 5       |            | Benzene, 1-ethyl-3,5-dimethyl-      | 134 C10H14       | 000934-74-7 11 |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

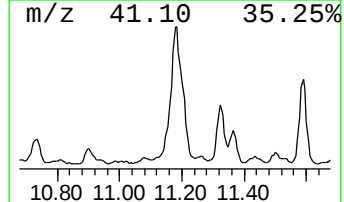
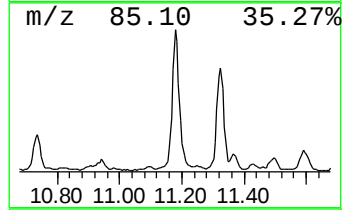
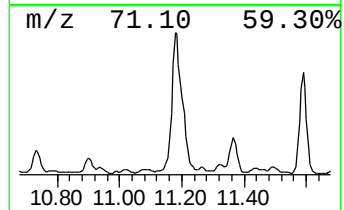
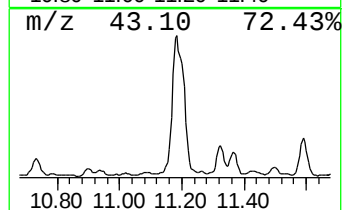
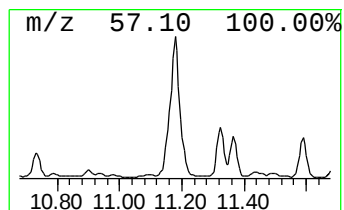
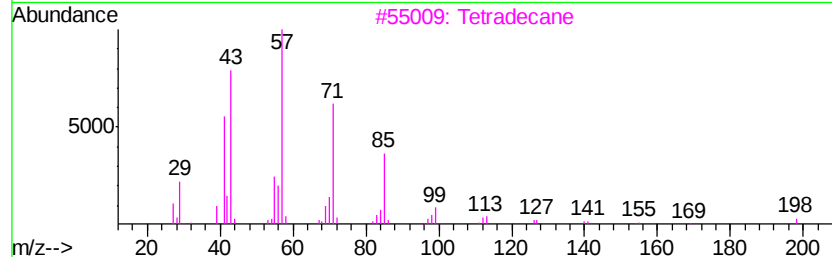
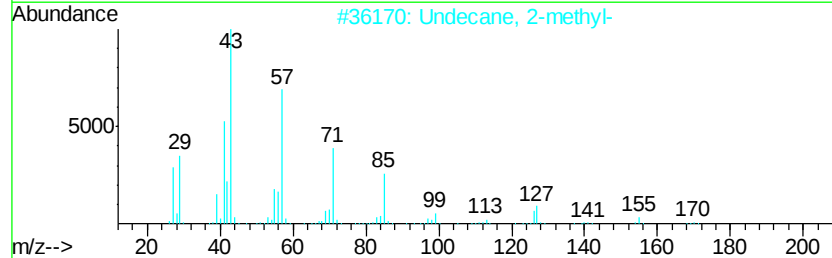
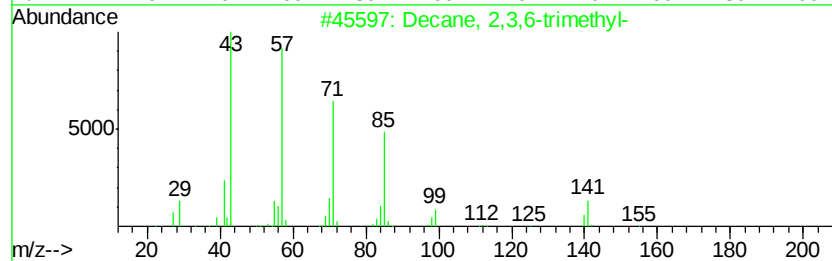
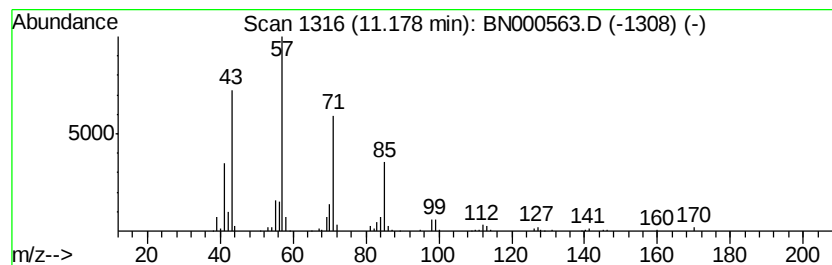
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 28 (DEL) Alkane: Straight-Chai... Concentration Rank 5

| R.T.  | EstConc     | Area   | Relative to ISTD | R.T.  |
|-------|-------------|--------|------------------|-------|
| 11.18 | 10.25 ng/ul | 831080 | Napthalene-d8    | 11.25 |

| Hit# | of | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------|-----|---------|-------------|------|
| 1    | 5  | Decane, 2,3,6-trimethyl- | 184 | C13H28  | 062238-12-4 | 80   |
| 2    |    | Undecane, 2-methyl-      | 170 | C12H26  | 007045-71-8 | 80   |
| 3    |    | Tetradecane              | 198 | C14H30  | 000629-59-4 | 80   |
| 4    |    | Hexadecane               | 226 | C16H34  | 000544-76-3 | 80   |
| 5    |    | Eicosane                 | 282 | C20H42  | 000112-95-8 | 80   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

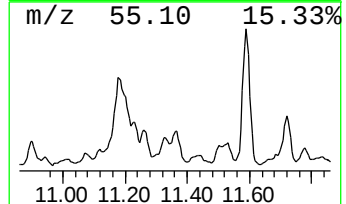
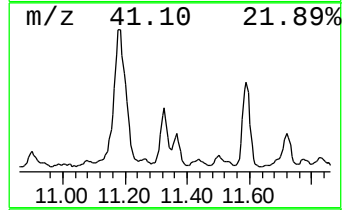
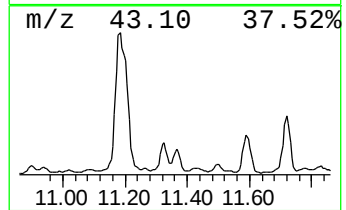
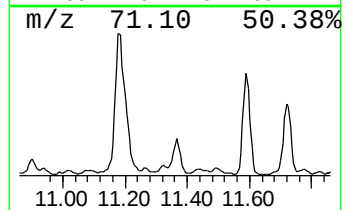
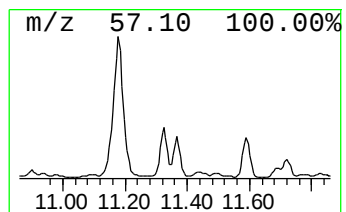
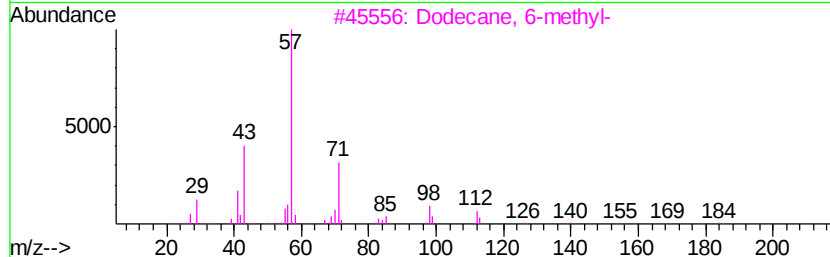
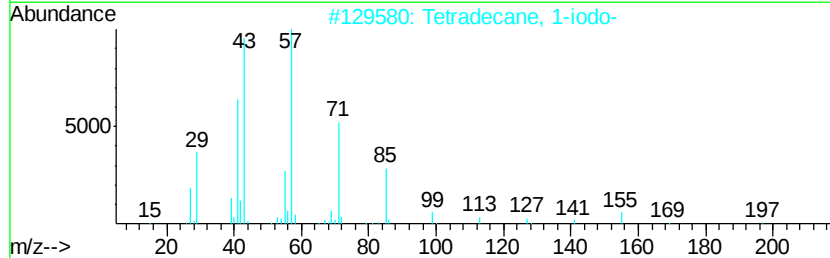
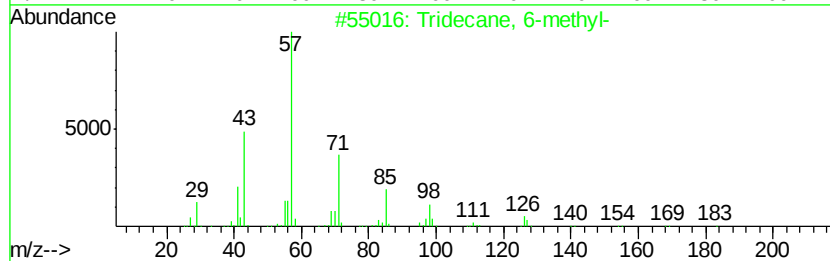
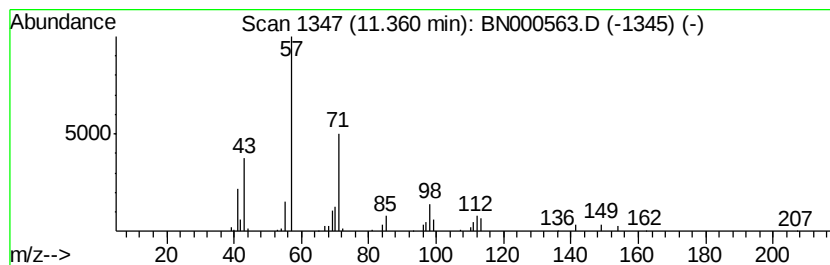
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 30 (DEL) Alkane: Straight-Chai... Concentration Rank 55

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 11.36 | 2.45 ng/ul | 198641 | Naphthalene-d8   | 11.25 |

| Hit# | of | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------|-----|---------|-------------|------|
| 1    | 5  | Tridecane, 6-methyl-    | 198 | C14H30  | 013287-21-3 | 59   |
| 2    |    | Tetradecane, 1-iodo-    | 324 | C14H29I | 019218-94-1 | 50   |
| 3    |    | Dodecane, 6-methyl-     | 184 | C13H28  | 006044-71-9 | 50   |
| 4    |    | Undecane, 2,6-dimethyl- | 184 | C13H28  | 017301-23-4 | 50   |
| 5    |    | Undecane, 2,6-dimethyl- | 184 | C13H28  | 017301-23-4 | 47   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

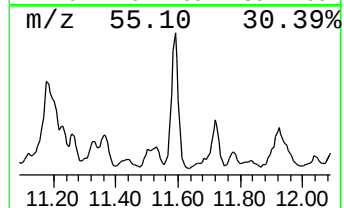
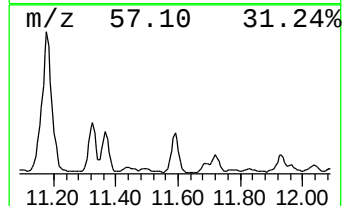
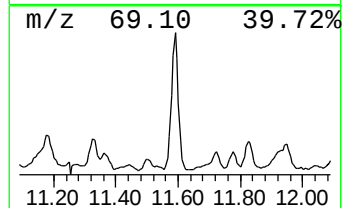
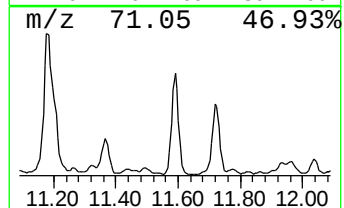
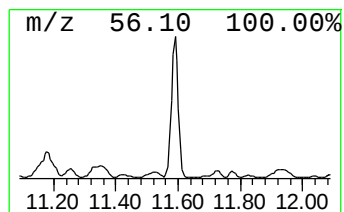
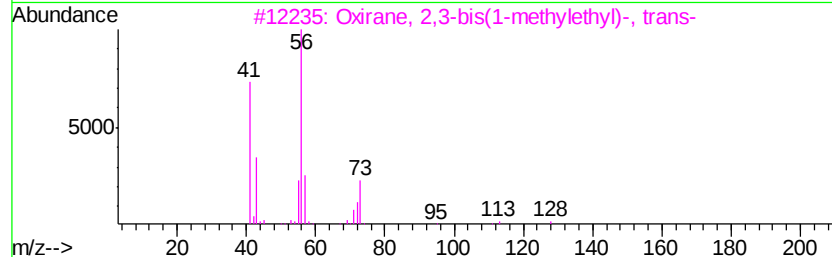
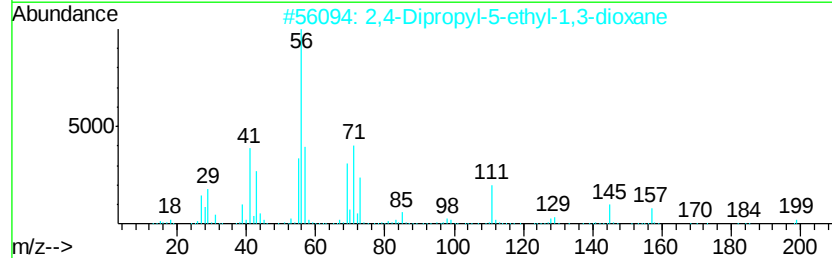
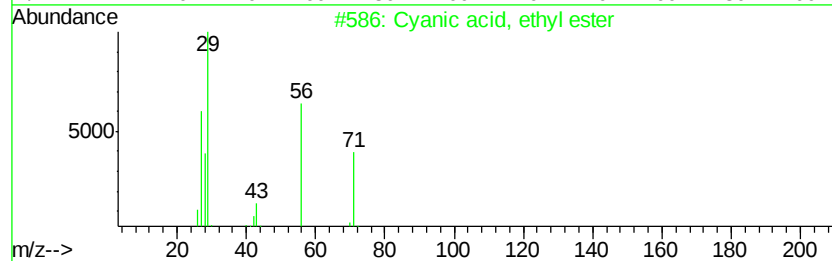
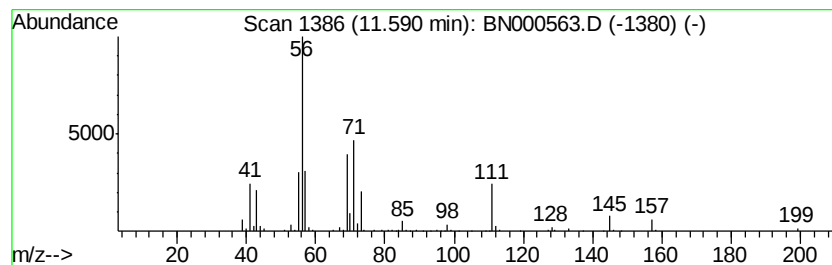
Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 31 Cyanic acid, ethyl ester Concentration Rank 11

| R.T.    | EstConc    | Area                                | Relative to ISTD | R.T.           |
|---------|------------|-------------------------------------|------------------|----------------|
| 11.59   | 7.36 ng/ul | 596699                              | Naphthalene-d8   | 11.25          |
| Hit# of | 5          | Tentative ID                        | MW MolForm       | CAS# Qual      |
| 1       |            | Cyanic acid, ethyl ester            | 71 C3H5NO        | 000627-48-5 58 |
| 2       |            | 2,4-Dipropyl-5-ethyl-1,3-dioxane    | 200 C12H24O2     | 006413-83-8 43 |
| 3       |            | Oxirane, 2,3-bis(1-methylethyl)-... | 128 C8H16O       | 054644-32-5 32 |
| 4       |            | 1,3-Pentanediol, 2,2,4-trimethyl-   | 146 C8H18O2      | 000144-19-4 25 |
| 5       |            | 1-Hexene, 2-methyl-                 | 98 C7H14         | 006094-02-6 17 |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

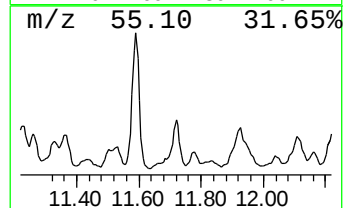
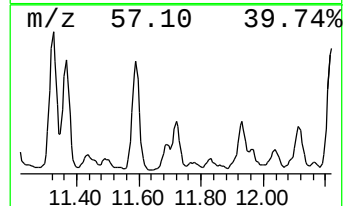
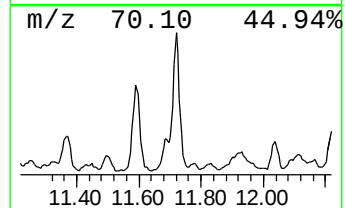
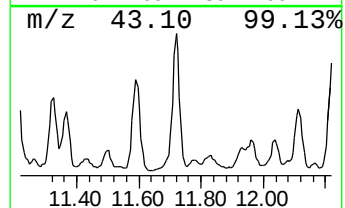
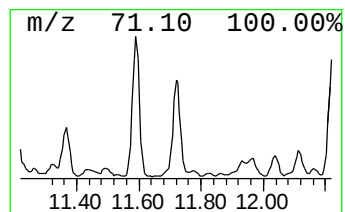
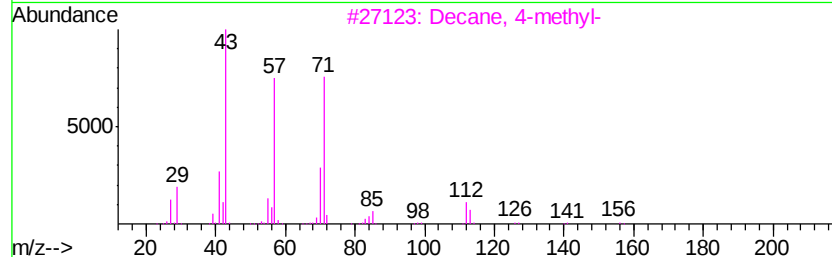
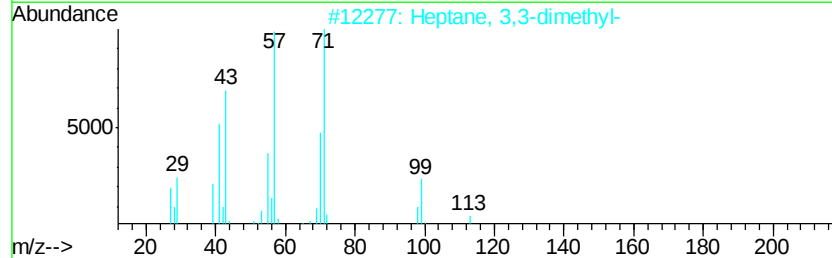
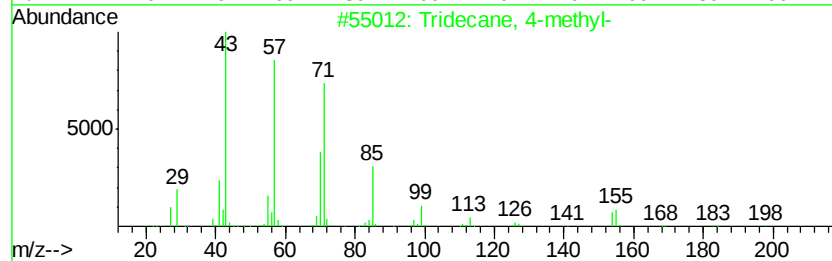
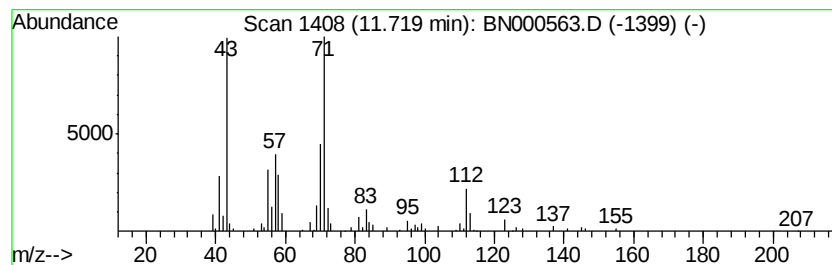
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 32 (DEL) Alkane: Straight-Chai... Concentration Rank 36

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 11.72 | 3.66 ng/ul | 296553 | Napthalene-d8    | 11.25 |

| Hit# | of | 5 | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------------|-----|---------|-------------|------|
| 1    |    |   | Tridecane, 4-methyl-      | 198 | C14H30  | 026730-12-1 | 50   |
| 2    |    |   | Heptane, 3,3-dimethyl-    | 128 | C9H20   | 004032-86-4 | 50   |
| 3    |    |   | Decane, 4-methyl-         | 156 | C11H24  | 002847-72-5 | 50   |
| 4    |    |   | Heptane, 3,3,5-trimethyl- | 142 | C10H22  | 007154-80-5 | 50   |
| 5    |    |   | Octane, 3-ethyl-          | 142 | C10H22  | 005881-17-4 | 50   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

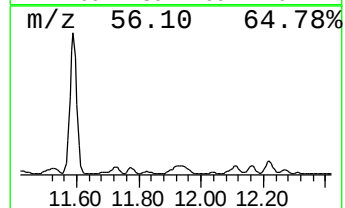
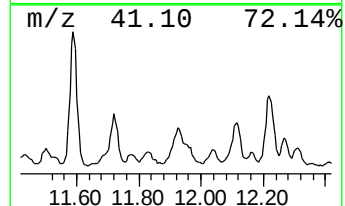
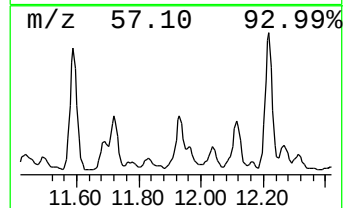
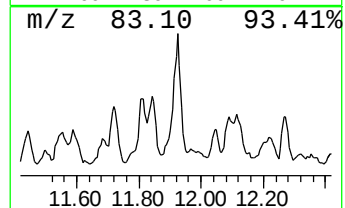
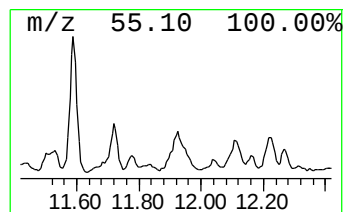
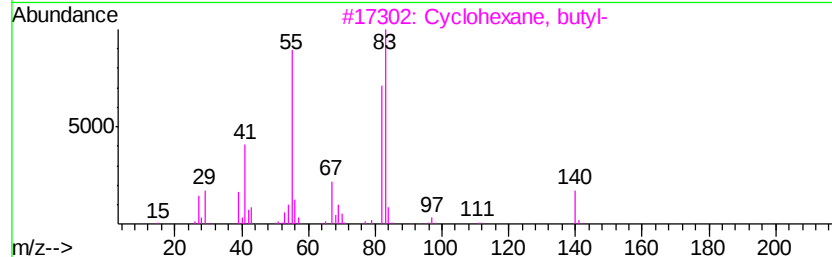
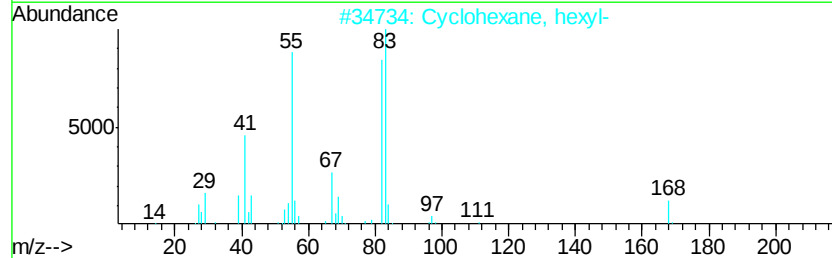
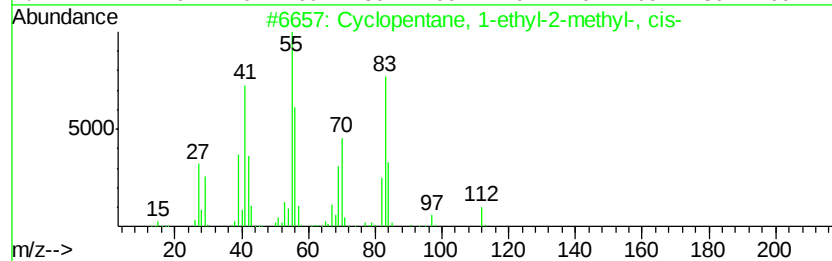
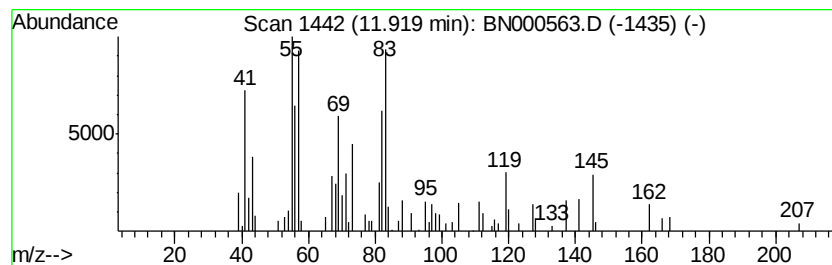
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 33 (DEL) Alkane: Cyclic11.92 Concentration Rank 37

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 11.92 | 3.57 ng/ul | 289593 | Napthalene-d8    | 11.25 |

| Hit# | of | Tentative ID                         | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclopentane, 1-ethyl-2-methyl-, ... | 112 | C8H16   | 000930-89-2 | 30   |
| 2    |    | Cyclohexane, hexyl-                  | 168 | C12H24  | 004292-75-5 | 30   |
| 3    |    | Cyclohexane, butyl-                  | 140 | C10H20  | 001678-93-9 | 27   |
| 4    |    | Cyclohexane, (1-methylethyl)-        | 126 | C9H18   | 000696-29-7 | 27   |
| 5    |    | Cyclohexane, (4-methylpentyl)-       | 168 | C12H24  | 061142-20-9 | 27   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

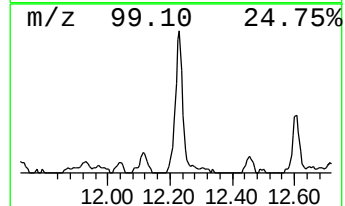
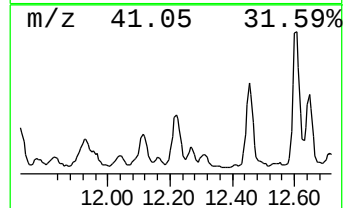
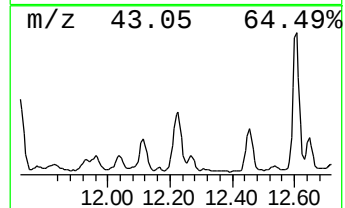
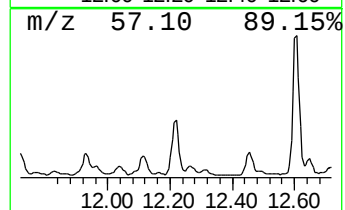
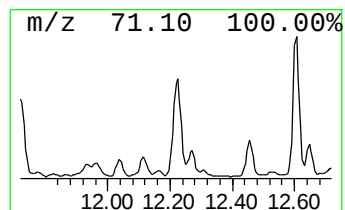
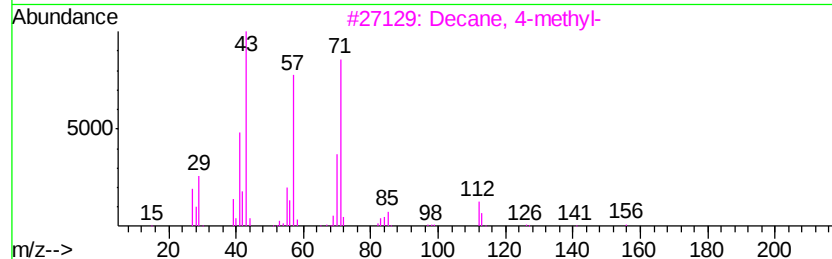
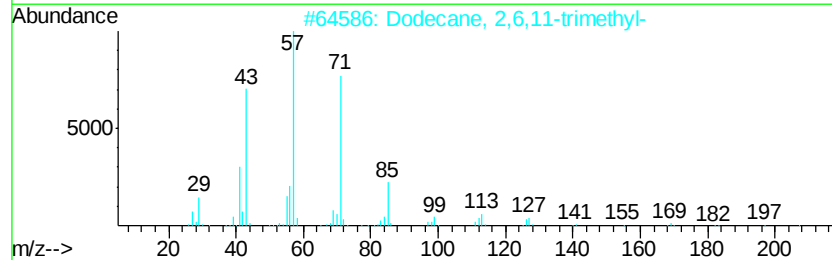
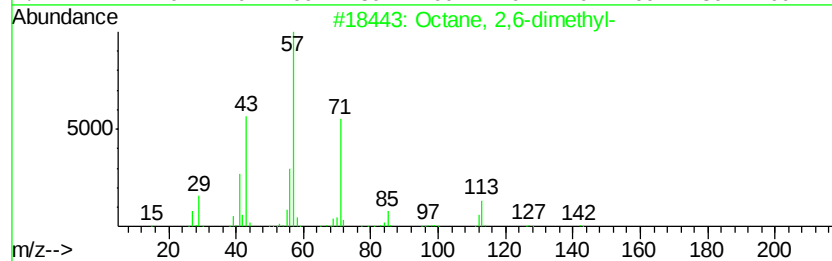
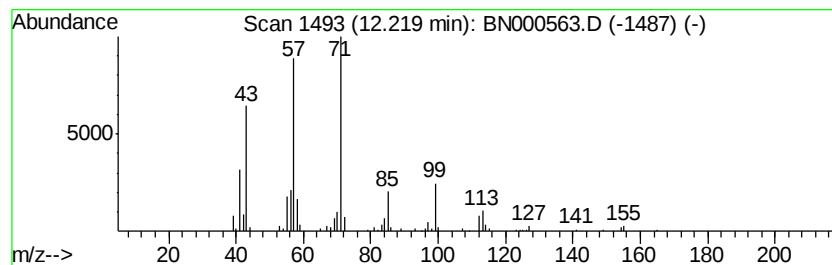
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 34 (DEL) Alkane: Straight-Chai... Concentration Rank 32

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 12.22 | 4.01 ng/ul | 325549 | Napthalene-d8    | 11.25 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | Octane, 2,6-dimethyl-       | 142 | C10H22  | 002051-30-1 | 80   |
| 2    |    |   | Dodecane, 2,6,11-trimethyl- | 212 | C15H32  | 031295-56-4 | 78   |
| 3    |    |   | Decane, 4-methyl-           | 156 | C11H24  | 002847-72-5 | 59   |
| 4    |    |   | Nonane, 3-methyl-           | 142 | C10H22  | 005911-04-6 | 43   |
| 5    |    |   | 1-Methoxycyclohexane        | 114 | C7H14O  | 000931-56-6 | 43   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

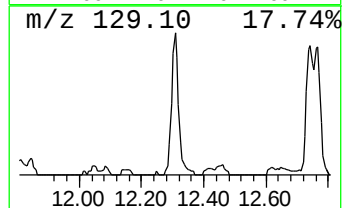
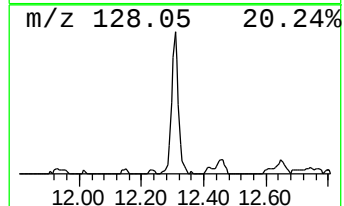
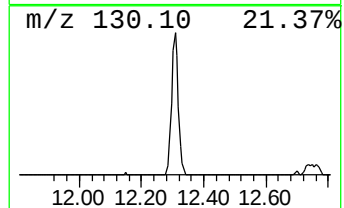
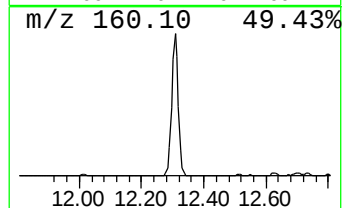
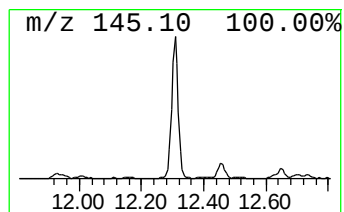
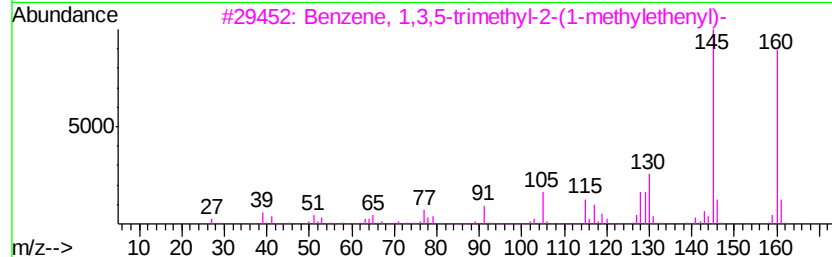
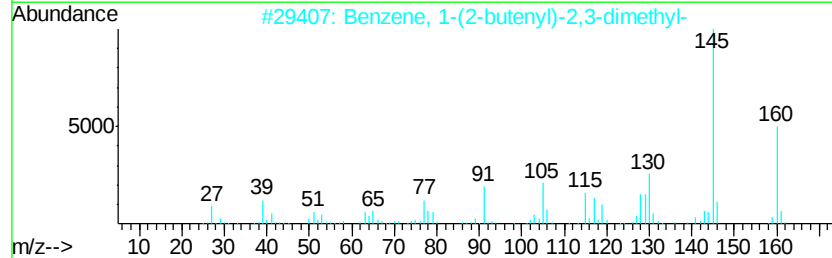
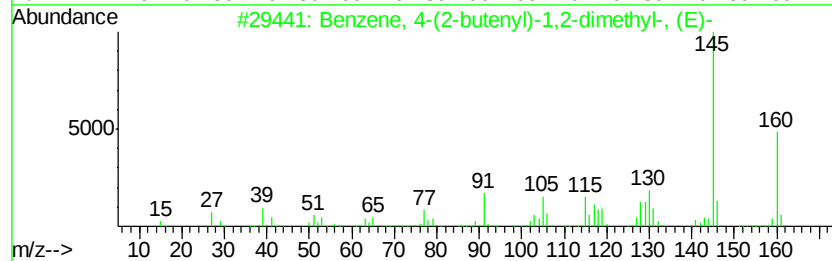
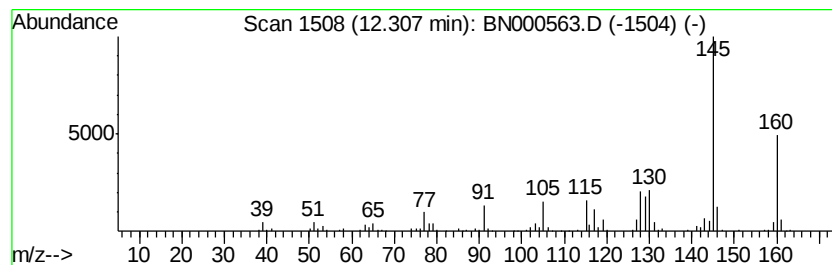
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 35 Benzene, 4-(2-butenyl)-1,2-... Concentration Rank 15

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 12.31 | 6.43 ng/ul | 521376 | Naphthalene-d8   | 11.25 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 4-(2-butenyl)-1,2-dimet... | 160 | C12H16  | 054340-86-2 | 94   |
| 2    |    | Benzene, 1-(2-butenyl)-2,3-dimet... | 160 | C12H16  | 054340-85-1 | 94   |
| 3    |    | Benzene, 1,3,5-trimethyl-2-(1-me... | 160 | C12H16  | 014679-13-1 | 93   |
| 4    |    | Benzene, 1-(2-butenyl)-2,3-dimet... | 160 | C12H16  | 054340-85-1 | 91   |
| 5    |    | Naphthalene, 1,2,3,4-tetrahydro-... | 160 | C12H16  | 004175-54-6 | 87   |



Data Path : Z:\HPCHEM1\BNA\_N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampled :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

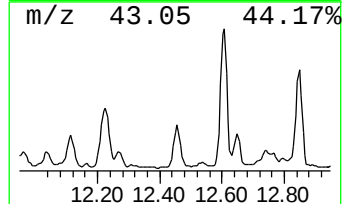
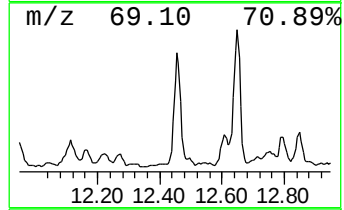
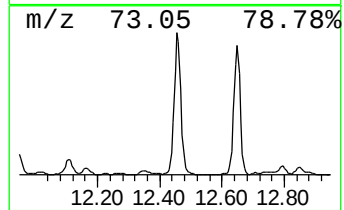
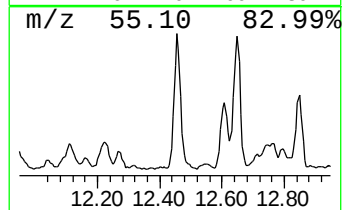
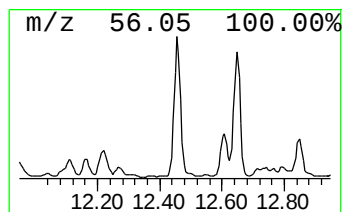
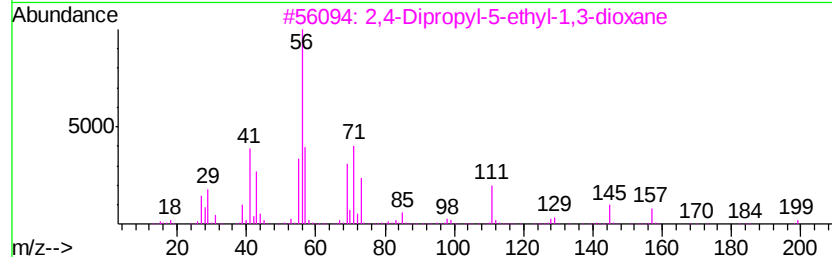
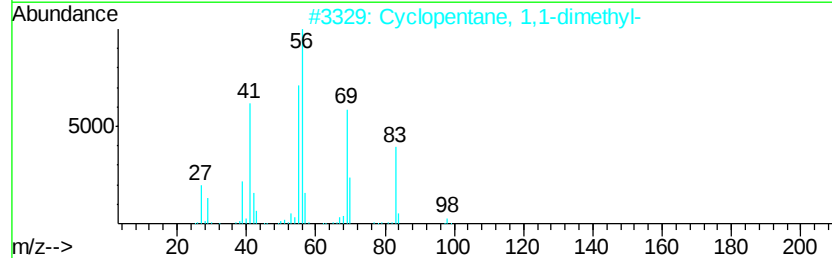
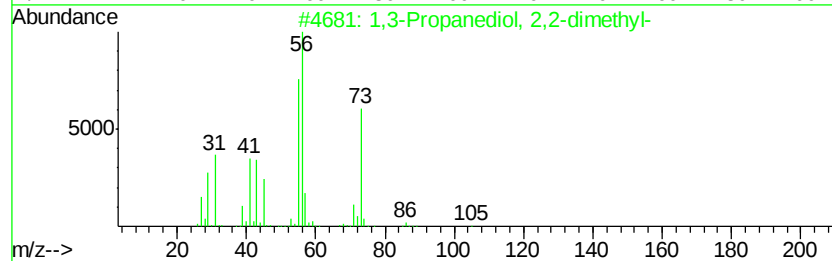
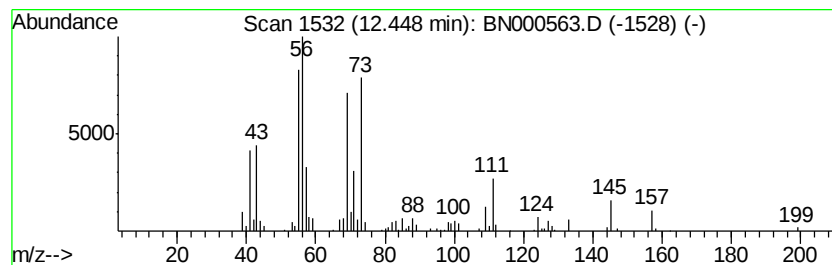
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 36 1,3-Propanediol, 2,2-dimethyl- Concentration Rank 22

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 12.45 | 5.14 ng/ul | 417139 | Napthalene-d8    | 11.25 |

| Hit# | of | Tentative ID                     | MW  | MolForm  | CAS#        | Qual |
|------|----|----------------------------------|-----|----------|-------------|------|
| 1    | 5  | 1,3-Propanediol, 2,2-dimethyl-   | 104 | C5H12O2  | 000126-30-7 | 50   |
| 2    |    | Cyclopentane, 1,1-dimethyl-      | 98  | C7H14    | 001638-26-2 | 43   |
| 3    |    | 2,4-Dipropyl-5-ethyl-1,3-dioxane | 200 | C12H24O2 | 006413-83-8 | 42   |
| 4    |    | 1,3-Propanediol, 2,2-dimethyl-   | 104 | C5H12O2  | 000126-30-7 | 40   |
| 5    |    | 1,3-Propanediol, 2,2-dimethyl-   | 104 | C5H12O2  | 000126-30-7 | 40   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

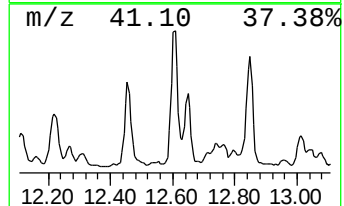
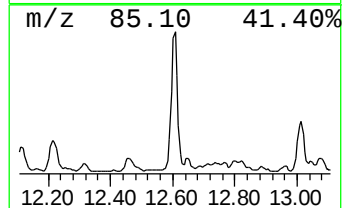
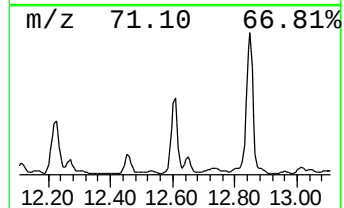
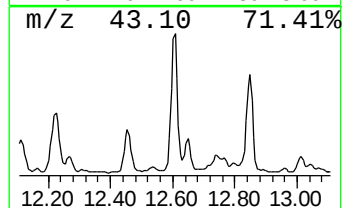
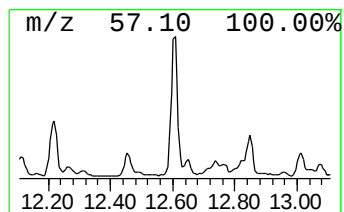
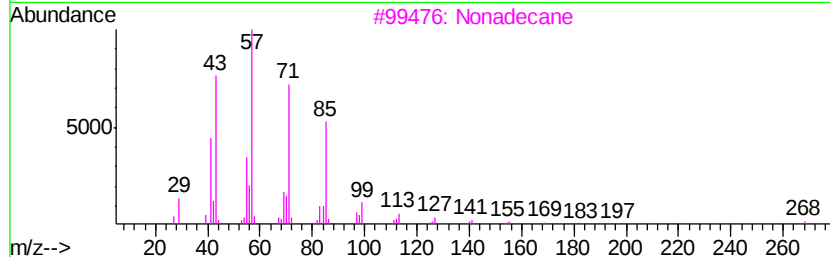
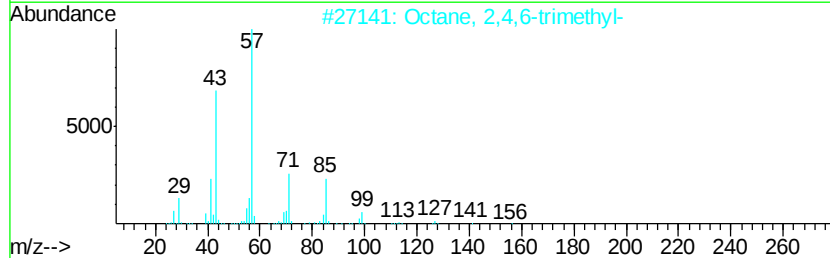
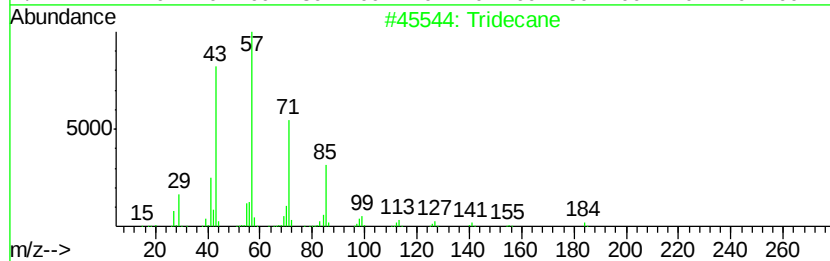
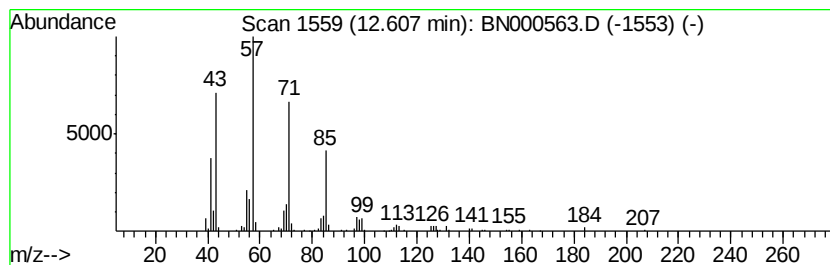
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 37 (DEL) Alkane: Straight-Chai... Concentration Rank 10

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 12.61 | 7.45 ng/ul | 604416 | Napthalene-d8    | 11.25 |

| Hit# | of | 5 | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------|-----|---------|-------------|------|
| 1    |    |   | Tridecane                | 184 | C13H28  | 000629-50-5 | 93   |
| 2    |    |   | Octane, 2,4,6-trimethyl- | 156 | C11H24  | 062016-37-9 | 86   |
| 3    |    |   | Nonadecane               | 268 | C19H40  | 000629-92-5 | 86   |
| 4    |    |   | Nonane, 5-butyl-         | 184 | C13H28  | 017312-63-9 | 80   |
| 5    |    |   | Undecane, 2,6-dimethyl-  | 184 | C13H28  | 017301-23-4 | 70   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

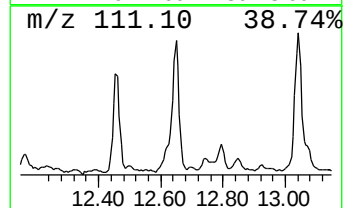
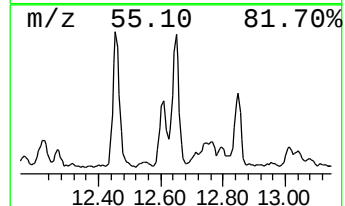
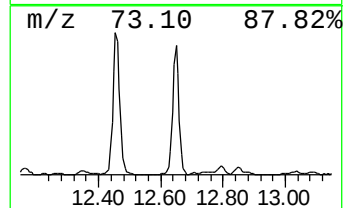
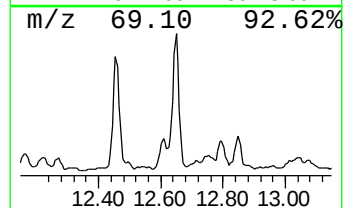
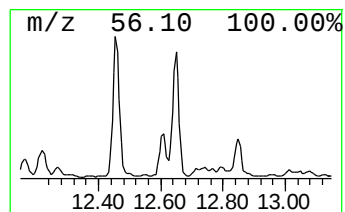
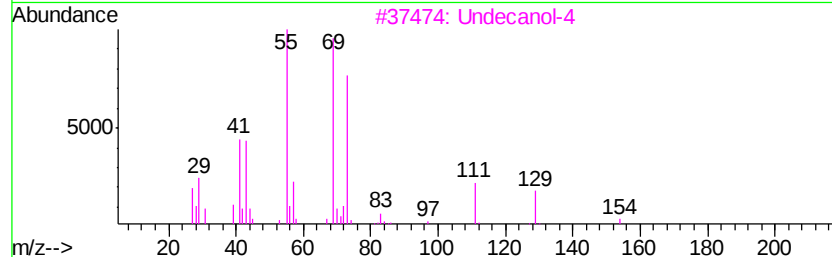
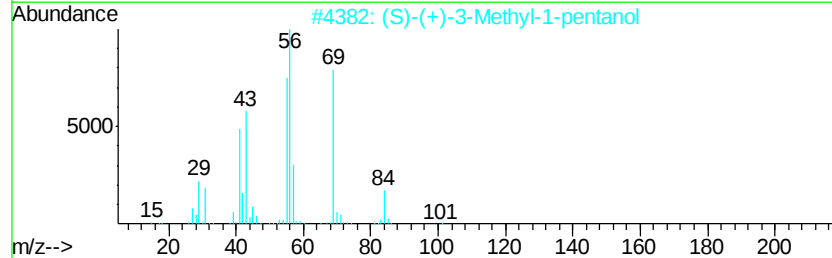
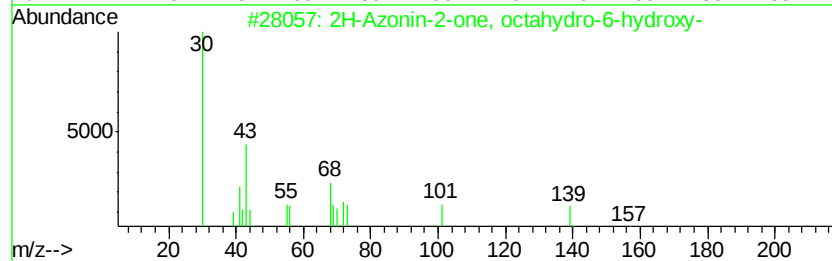
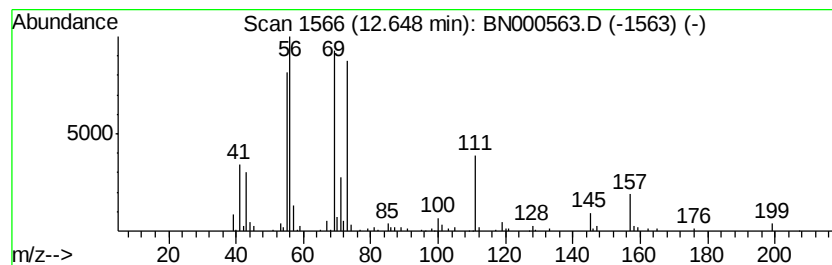
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TIC Integration Parameters: LSCINT.P

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Peak Number 38 unknown-02 Concentration Rank 28

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 12.65 | 4.21 ng/ul | 341589 | Naphthalene-d8   | 11.25 |

| Hit# | of | Tentative ID                          | MW  | MolForm  | CAS#        | Qual |
|------|----|---------------------------------------|-----|----------|-------------|------|
| 1    | 5  | 2H-Azonin-2-one, octahydro-6-hydroxy- | 157 | C8H15NO2 | 023435-07-6 | 42   |
| 2    |    | (S)-(+)-3-Methyl-1-pentanol           | 102 | C6H14O   | 042072-39-9 | 28   |
| 3    |    | Undecanol-4                           | 172 | C11H24O  | 004272-06-4 | 25   |
| 4    |    | Cyclooctane, (1-methylpropyl)-        | 168 | C12H24   | 016538-89-9 | 17   |
| 5    |    | 1,3-Propanediol, 2,2-dimethyl-        | 104 | C5H12O2  | 000126-30-7 | 16   |



Data Path : Z:\HPCHEM1\BNA\_N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

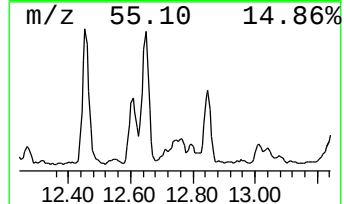
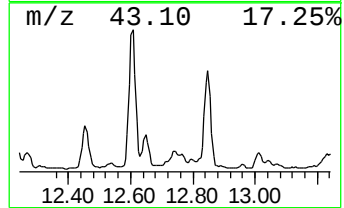
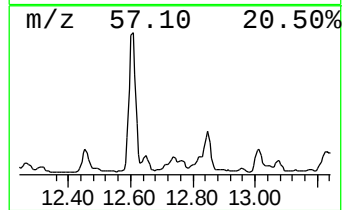
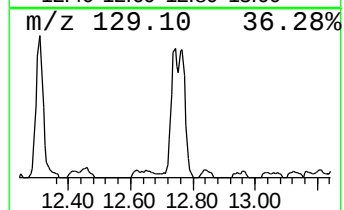
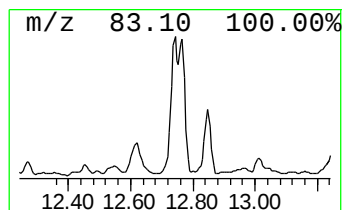
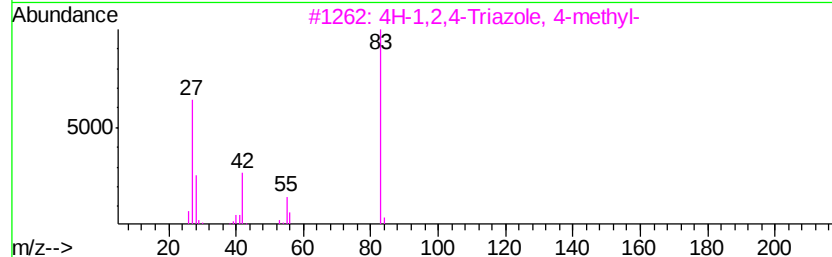
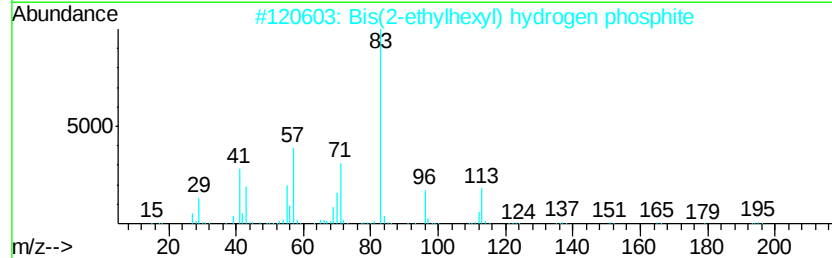
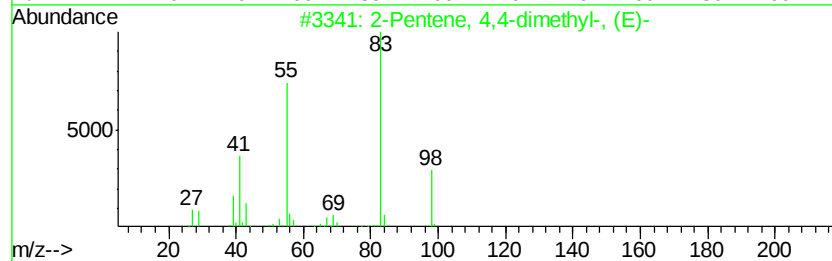
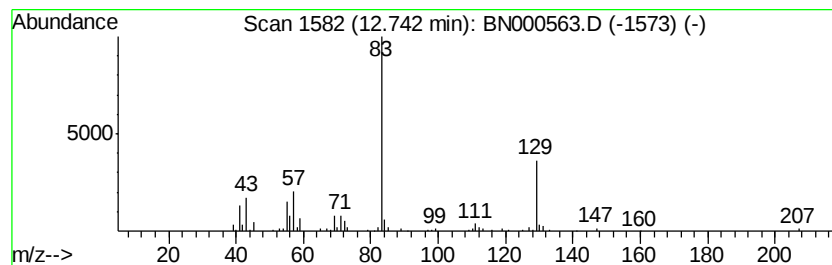
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 39 (DEL) Alkane: Cyclic12.74 Concentration Rank 56

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 12.74 | 2.41 ng/ul | 195827 | Naphthalene-d8   | 11.25 |

| Hit# | of | Tentative ID                         | MW  | MolForm   | CAS#        | Qual |
|------|----|--------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | 2-Pentene, 4,4-dimethyl-, (E)-       | 98  | C7H14     | 000690-08-4 | 40   |
| 2    |    | Bis(2-ethylhexyl) hydrogen phosph... | 306 | C16H35O3P | 003658-48-8 | 40   |
| 3    |    | 4H-1,2,4-Triazole, 4-methyl-         | 83  | C3H5N3    | 010570-40-8 | 40   |
| 4    |    | 1-Butene, 2,3,3-trimethyl-           | 98  | C7H14     | 000594-56-9 | 36   |
| 5    |    | 2-Hepten-4-one, 2-methyl-            | 126 | C8H14O    | 022319-24-0 | 36   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

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Quant Title : SVOA CALIBRATION

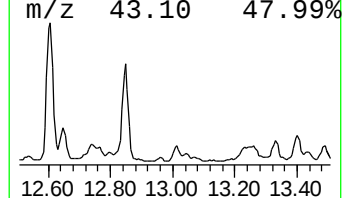
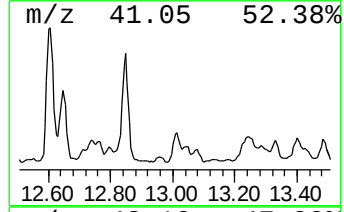
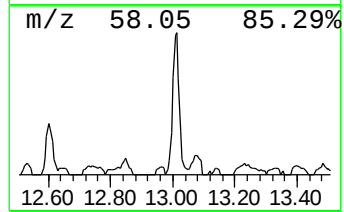
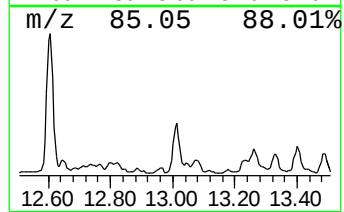
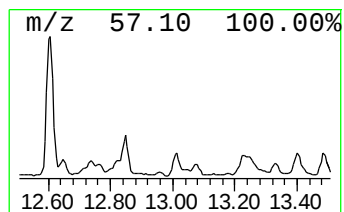
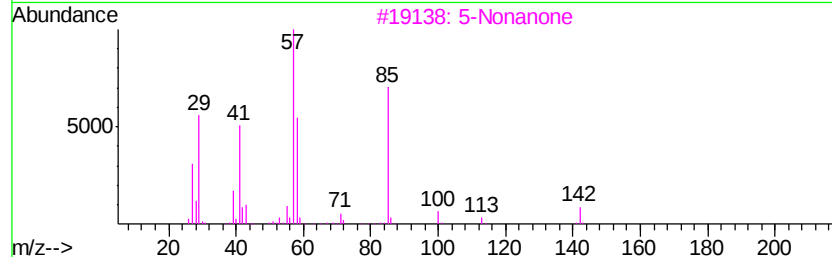
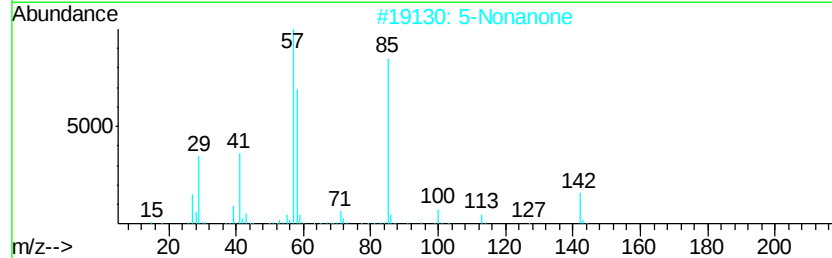
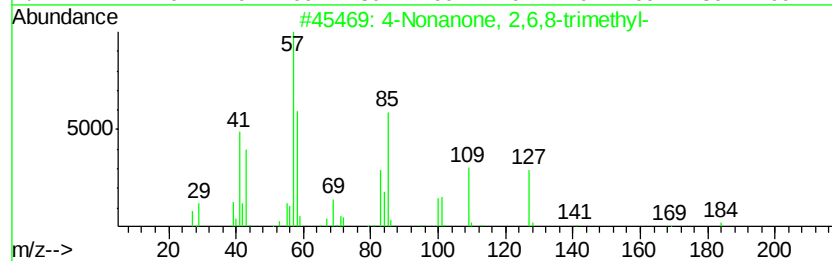
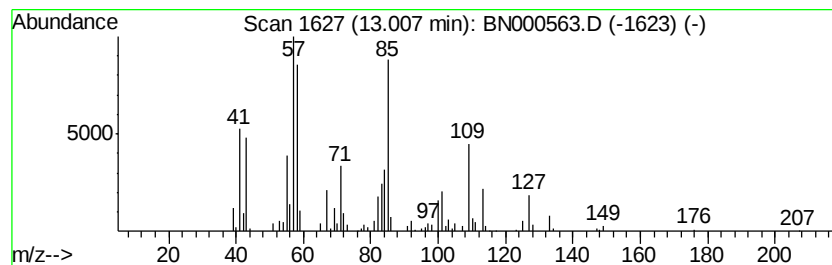
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 40 4-Nonanone, 2,6,8-trimethyl- Concentration Rank 59

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 13.01 | 2.23 ng/ul | 180607 | Napthalene-d8    | 11.25 |

| Hit# | of | Tentative ID                 | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------|-----|---------|-------------|------|
| 1    | 5  | 4-Nonanone, 2,6,8-trimethyl- | 184 | C12H24O | 000123-18-2 | 64   |
| 2    |    | 5-Nonanone                   | 142 | C9H18O  | 000502-56-7 | 47   |
| 3    |    | 5-Nonanone                   | 142 | C9H18O  | 000502-56-7 | 43   |
| 4    |    | N-Ethyl trimethylenediamine  | 102 | C5H14N2 | 038571-87-8 | 43   |
| 5    |    | 5-Nonanone                   | 142 | C9H18O  | 000502-56-7 | 43   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

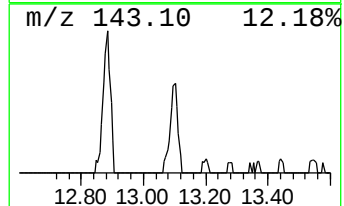
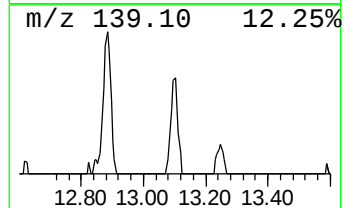
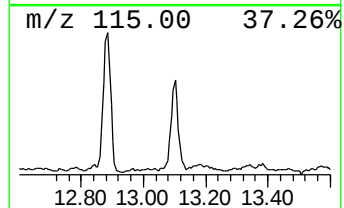
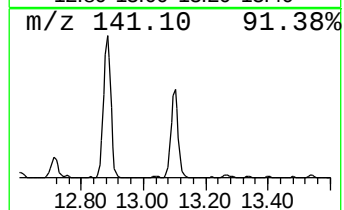
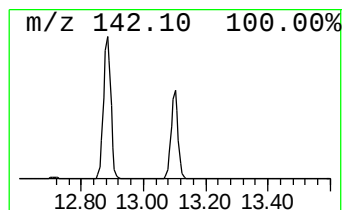
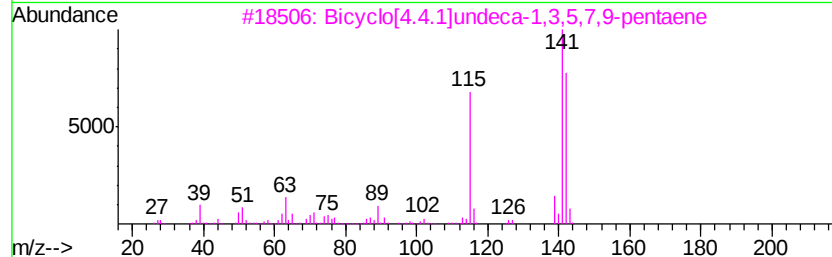
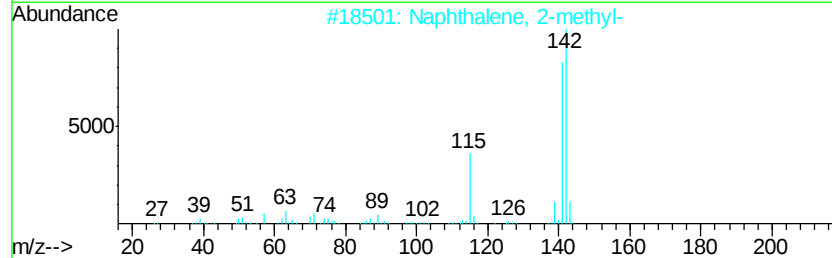
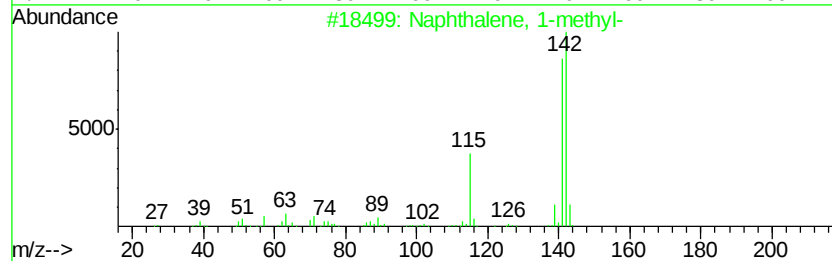
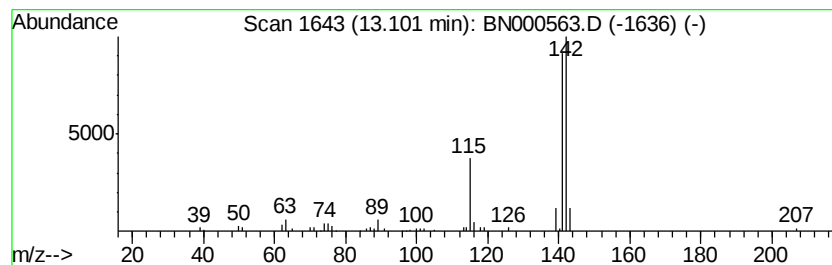
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 41 Naphthalene, 1-methyl- Concentration Rank 62

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 13.10 | 2.12 ng/ul | 172187 | Naphthalene-d8   | 11.25 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 1-methyl-              | 142 | C11H10  | 000090-12-0 | 95   |
| 2    |    | Naphthalene, 2-methyl-              | 142 | C11H10  | 000091-57-6 | 95   |
| 3    |    | Bicyclo[4.4.1]undeca-1,3,5,7,9-p... | 142 | C11H10  | 002443-46-1 | 95   |
| 4    |    | 1,4-Methanonaphthalene, 1,4-dihy... | 142 | C11H10  | 004453-90-1 | 93   |
| 5    |    | Benzocycloheptatriene               | 142 | C11H10  | 000264-09-5 | 91   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampled :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

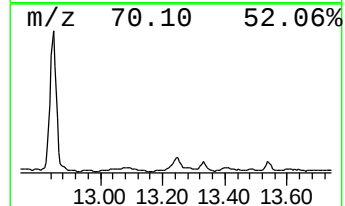
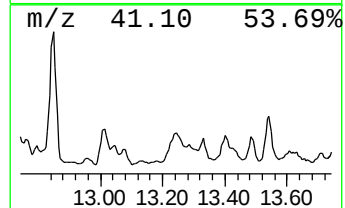
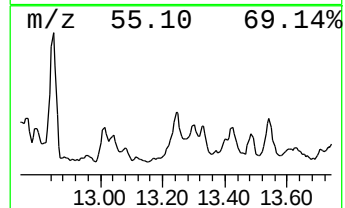
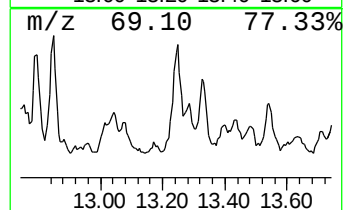
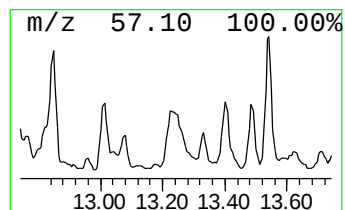
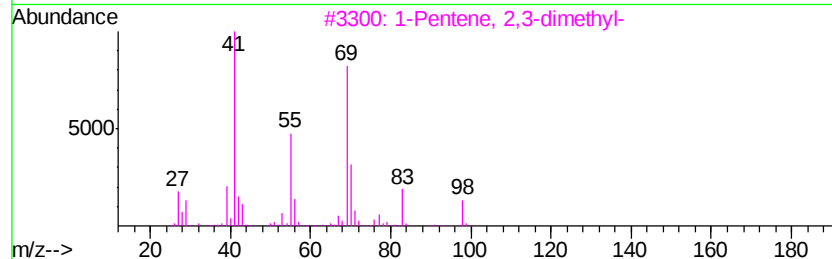
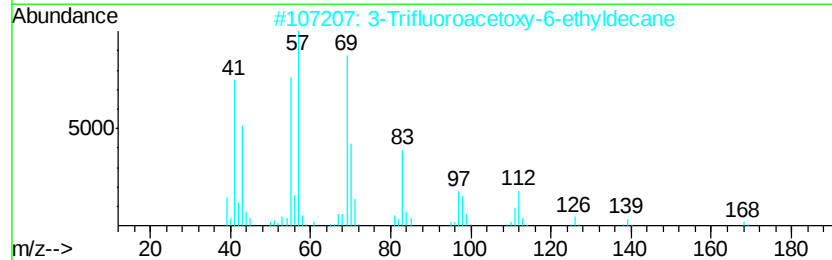
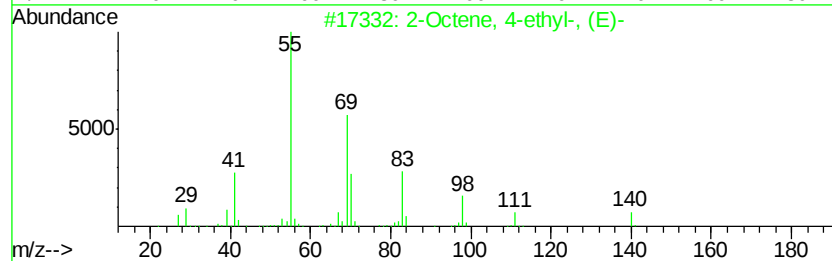
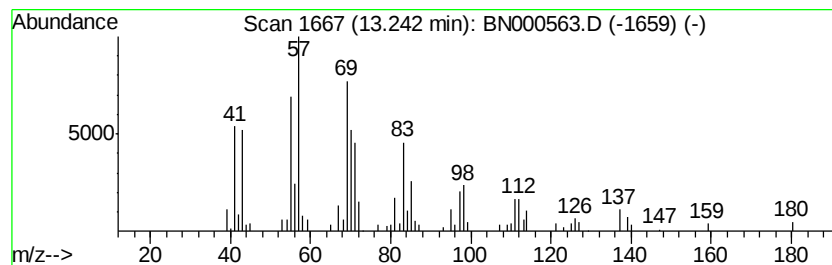
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 42 (DEL) Alkane: Cyclic13.24 Concentration Rank 61

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 13.24 | 2.18 ng/ul | 245385 | Acenaphthene-d10 | 15.03 |

| Hit# | of | Tentative ID                     | MW  | MolForm    | CAS#        | Qual |
|------|----|----------------------------------|-----|------------|-------------|------|
| 1    | 5  | 2-Octene, 4-ethyl-, (E)-         | 140 | C10H20     | 074630-09-4 | 46   |
| 2    |    | 3-Trifluoroacetoxy-6-ethyldecane | 282 | C14H25F3O2 | 116436-59-0 | 43   |
| 3    |    | 1-Pentene, 2,3-dimethyl-         | 98  | C7H14      | 003404-72-6 | 43   |
| 4    |    | Heptane, 4-methylene-            | 112 | C8H16      | 015918-08-8 | 30   |
| 5    |    | 3-Methyl-3-hexene                | 98  | C7H14      | 003404-65-7 | 30   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

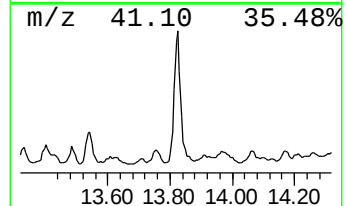
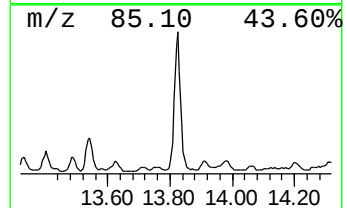
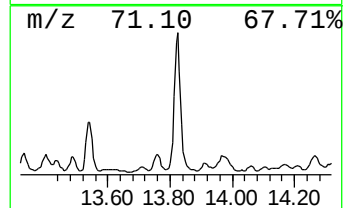
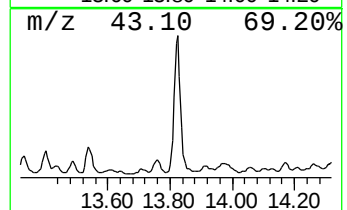
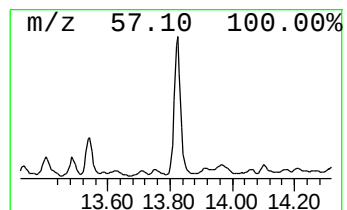
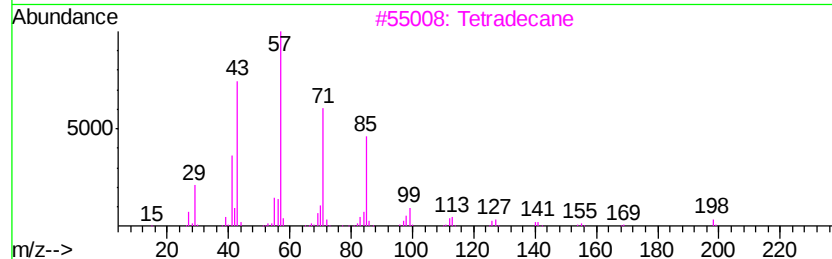
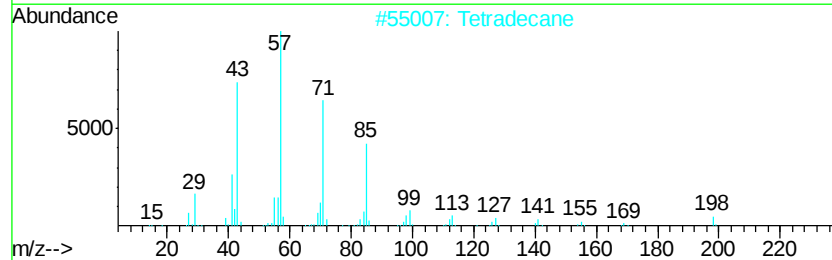
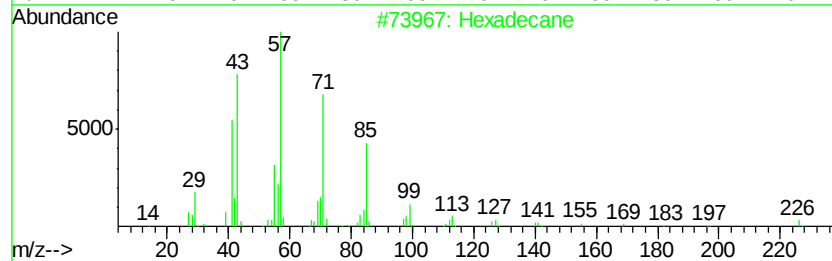
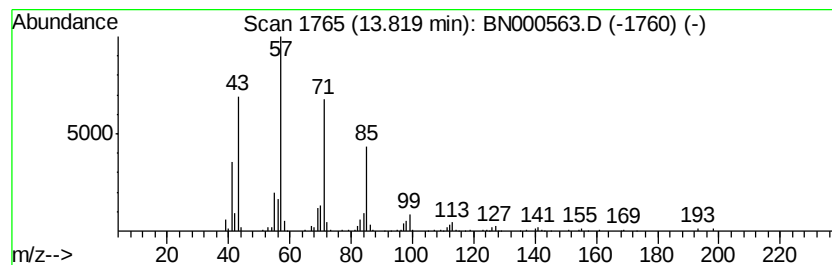
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 43 (DEL) Alkane: Straight-Chai... Concentration Rank 16

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 13.82 | 6.32 ng/ul | 711119 | Acenaphthene-d10 | 15.03 |

| Hit# | of | 5 | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------|-----|---------|-------------|------|
| 1    |    |   | Hexadecane               | 226 | C16H34  | 000544-76-3 | 91   |
| 2    |    |   | Tetradecane              | 198 | C14H30  | 000629-59-4 | 90   |
| 3    |    |   | Tetradecane              | 198 | C14H30  | 000629-59-4 | 87   |
| 4    |    |   | Tridecane                | 184 | C13H28  | 000629-50-5 | 86   |
| 5    |    |   | Decane, 2,3,5-trimethyl- | 184 | C13H28  | 062238-11-3 | 83   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

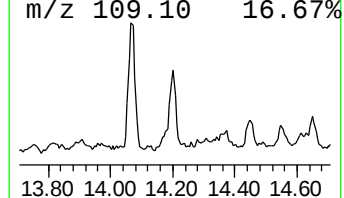
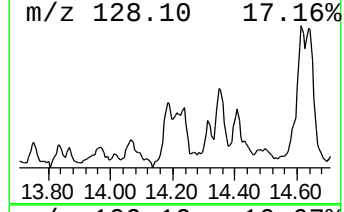
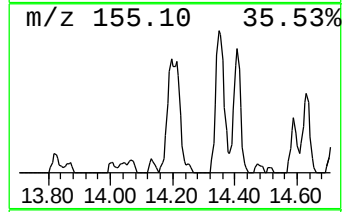
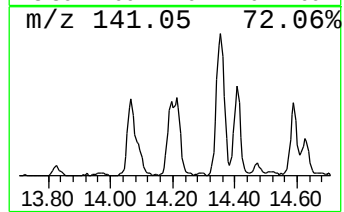
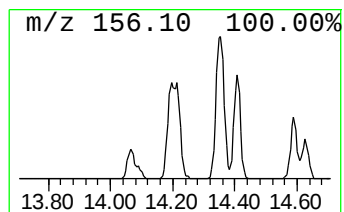
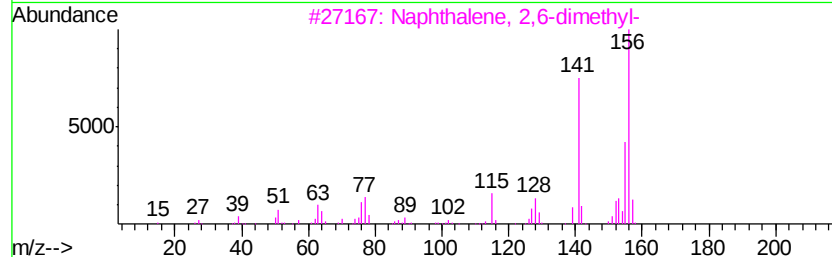
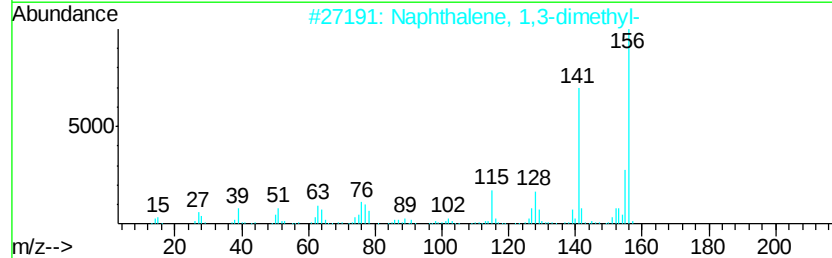
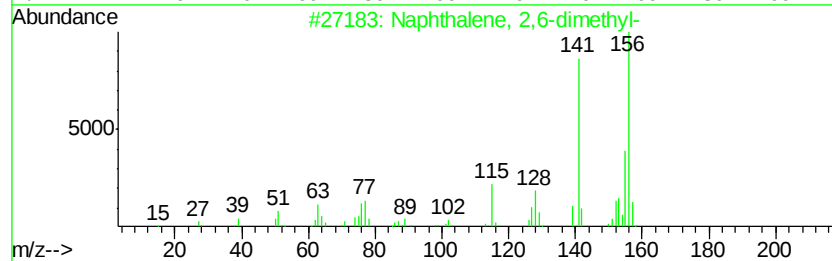
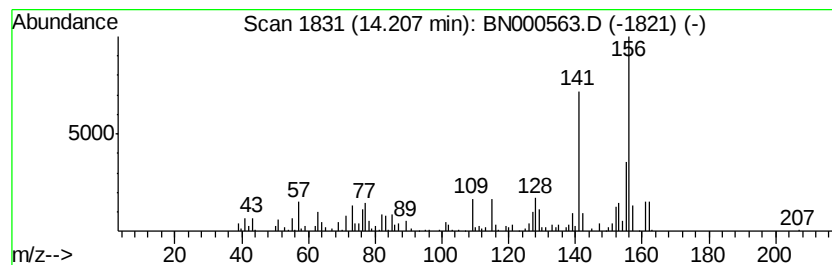
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 44 Naphthalene, 2,6-dimethyl- Concentration Rank 38

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 14.21 | 3.56 ng/ul | 400764 | Acenaphthene-d10 | 15.03 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 97   |
| 2    |    |   | Naphthalene, 1,3-dimethyl- | 156 | C12H12  | 000575-41-7 | 96   |
| 3    |    |   | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 96   |
| 4    |    |   | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 96   |
| 5    |    |   | Naphthalene, 2,7-dimethyl- | 156 | C12H12  | 000582-16-1 | 96   |



Data Path : Z:\HPCHEM1\BNA\_N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

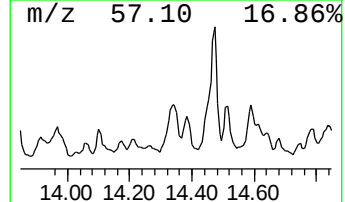
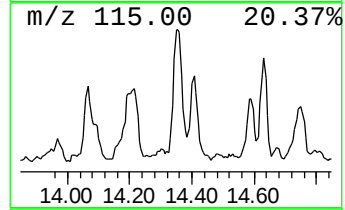
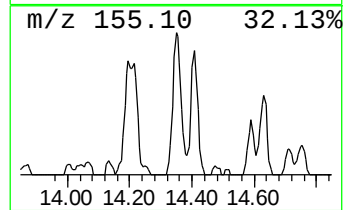
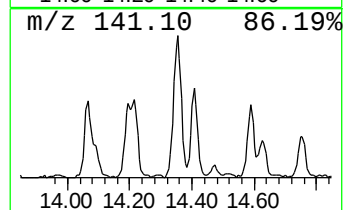
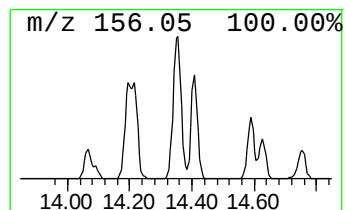
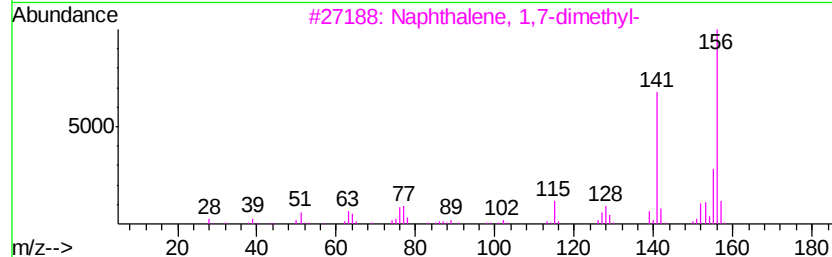
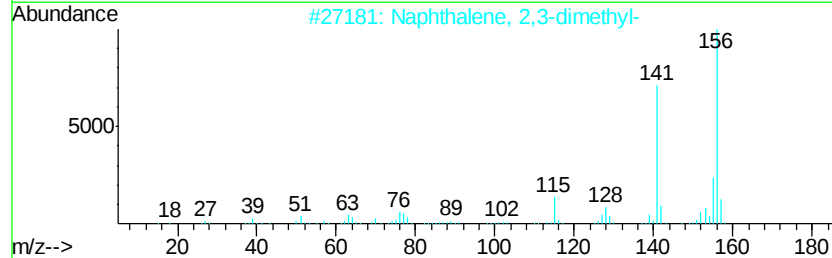
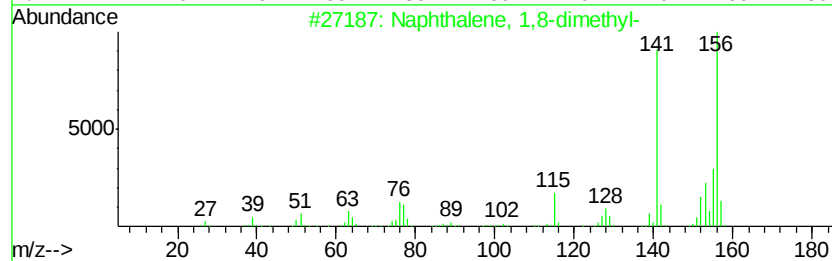
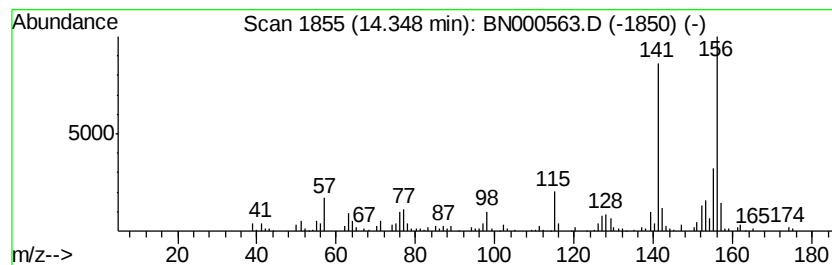
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 45 Naphthalene, 1,8-dimethyl- Concentration Rank 33

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 14.35 | 4.00 ng/ul | 450259 | Acenaphthene-d10 | 15.03 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 1,8-dimethyl- | 156 | C12H12  | 000569-41-5 | 98   |
| 2    |    |   | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 96   |
| 3    |    |   | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 96   |
| 4    |    |   | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 96   |
| 5    |    |   | Naphthalene, 1,2-dimethyl- | 156 | C12H12  | 000573-98-8 | 96   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

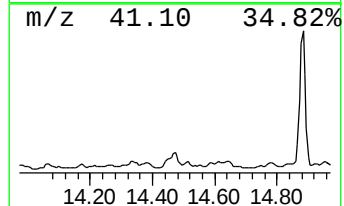
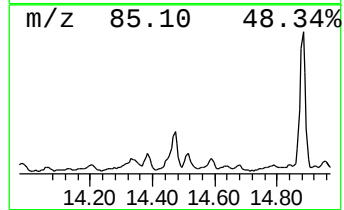
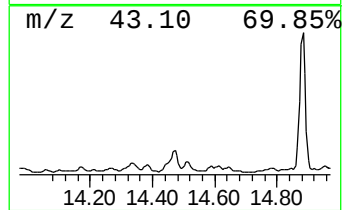
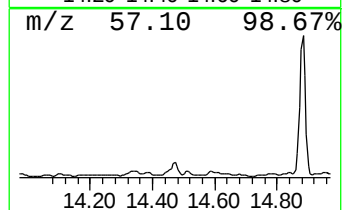
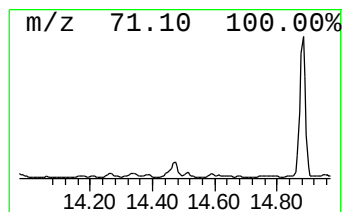
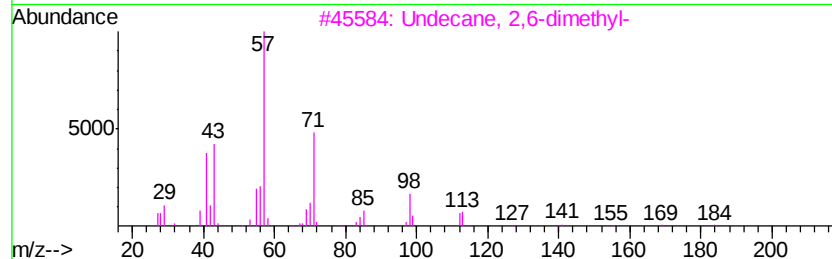
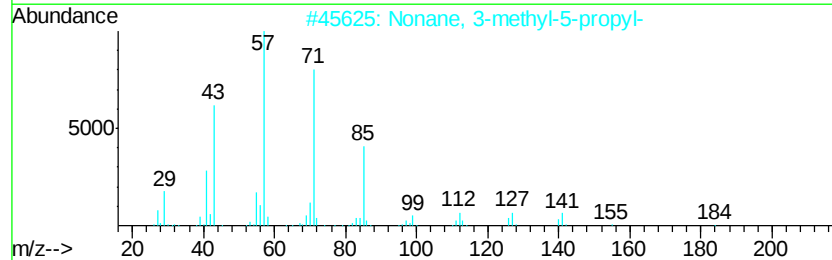
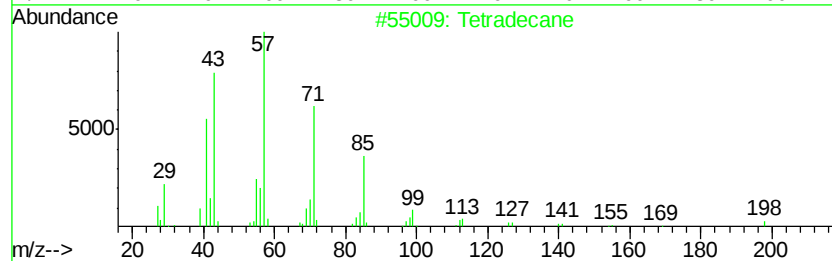
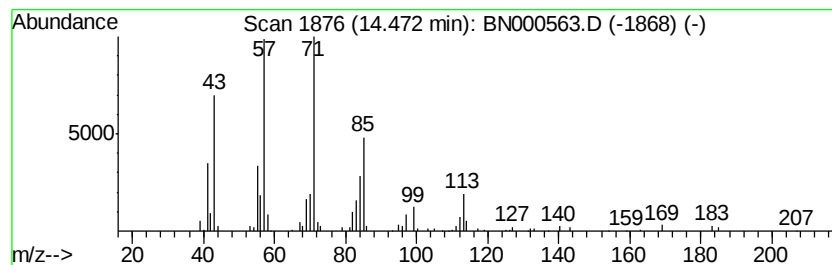
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 46 (DEL) Alkane: Straight-Chai... Concentration Rank 45

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 14.47 | 2.87 ng/ul | 322713 | Acenaphthene-d10 | 15.03 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#         | Qual |
|------|----|---|----------------------------|-----|---------|--------------|------|
| 1    |    |   | Tetradecane                | 198 | C14H30  | 000629-59-4  | 72   |
| 2    |    |   | Nonane, 3-methyl-5-propyl- | 184 | C13H28  | 031081-18-2  | 59   |
| 3    |    |   | Undecane, 2,6-dimethyl-    | 184 | C13H28  | 017301-23-4  | 53   |
| 4    |    |   | 1-Iodo-2-methylnonane      | 268 | C10H21I | 1000101-47-9 | 50   |
| 5    |    |   | Octane, 3,6-dimethyl-      | 142 | C10H22  | 015869-94-0  | 50   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

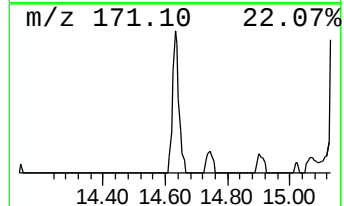
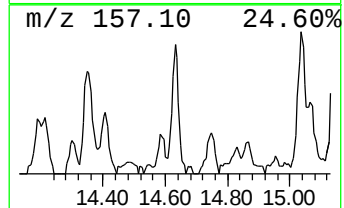
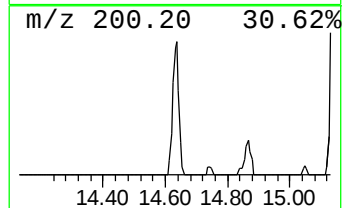
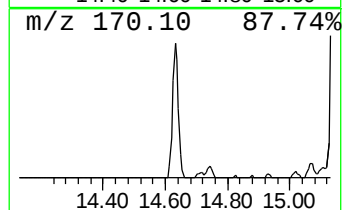
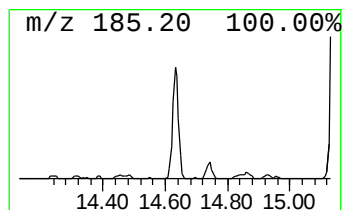
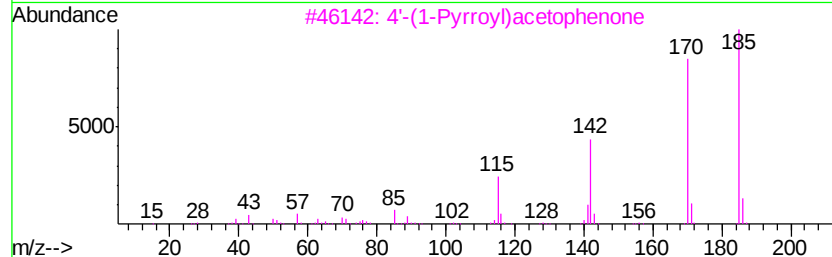
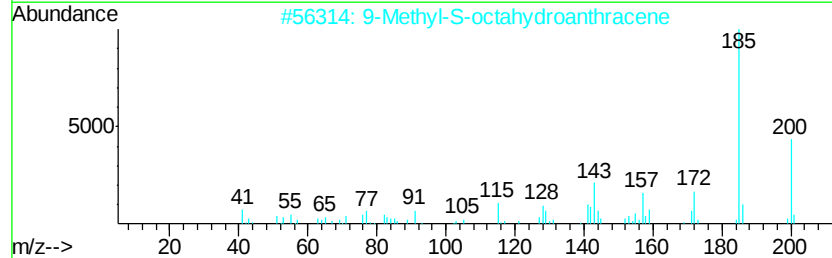
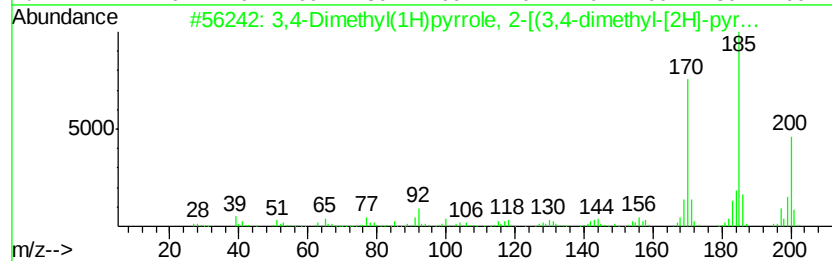
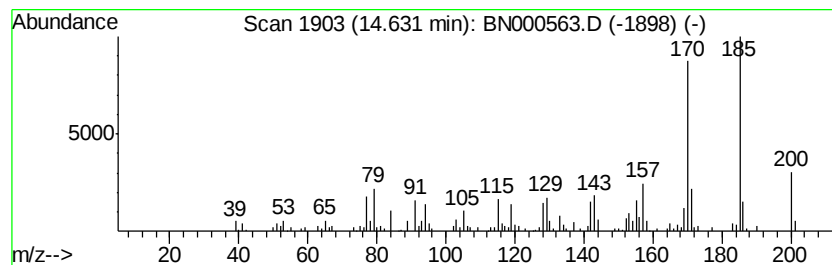
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 47 3,4-Dimethyl(1H)pyrrole, 2-... Concentration Rank 34

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 14.63 | 3.74 ng/ul | 421468 | Acenaphthene-d10 | 15.03 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | 3,4-Dimethyl(1H)pyrrole, 2-[(3,4... | 200 | C13H16N2 | 065539-79-9  | 74   |
| 2    |    |   | 9-Methyl-S-octahydroanthracene      | 200 | C15H20   | 004948-51-0  | 50   |
| 3    |    |   | 4'-(1-Pyrrovl)acetophenone          | 185 | C12H11NO | 022106-37-2  | 50   |
| 4    |    |   | 1,2,3,4,5,6,7,8-Octahydro-1-meth... | 200 | C15H20   | 1000080-19-4 | 46   |
| 5    |    |   | Naphthalene, 1,4,6-trimethyl-       | 170 | C13H14   | 002131-42-2  | 38   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

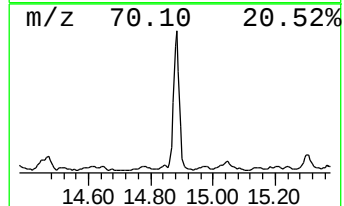
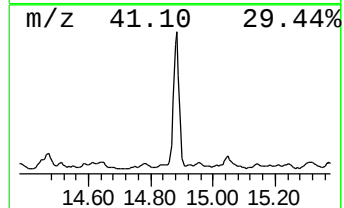
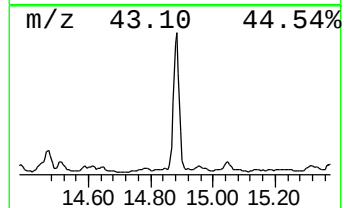
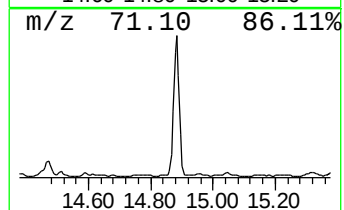
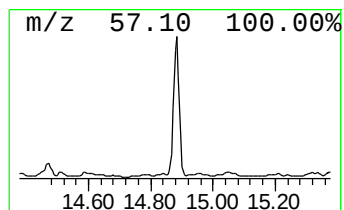
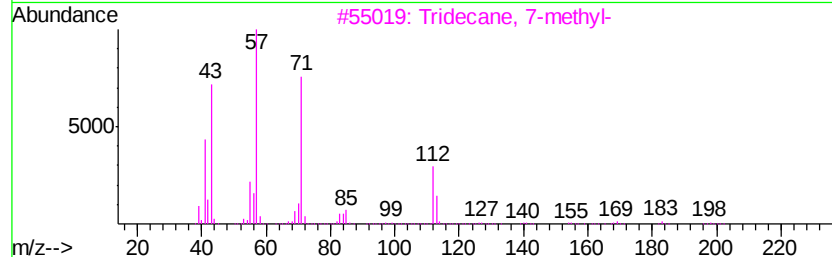
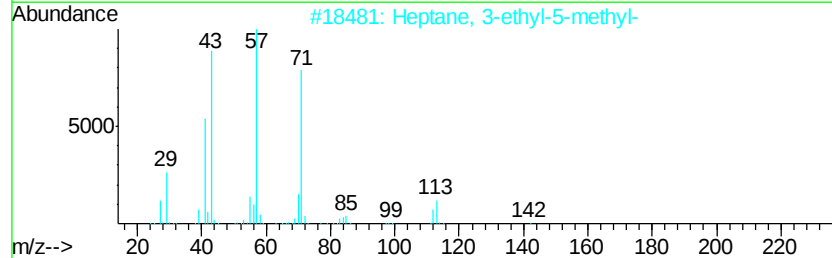
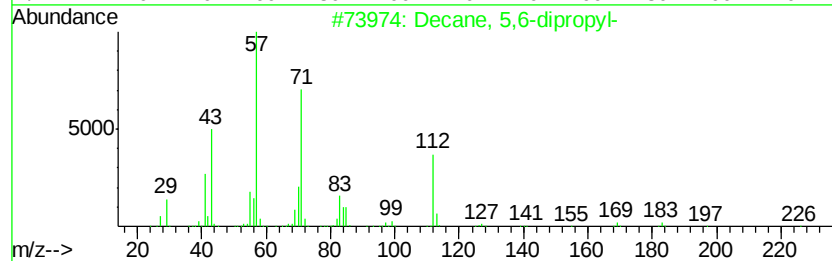
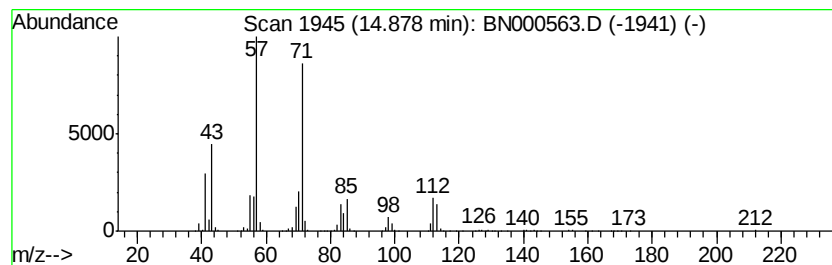
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 48 (DEL) Alkane: Straight-Chai... Concentration Rank 3

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.  |
|-------|-------------|---------|------------------|-------|
| 14.88 | 15.61 ng/ul | 1757570 | Acenaphthene-d10 | 15.03 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Decane, 5,6-dipropyl-      | 226 | C16H34  | 119209-20-0 | 78   |
| 2    |    |   | Heptane, 3-ethyl-5-methyl- | 142 | C10H22  | 052896-90-9 | 72   |
| 3    |    |   | Tridecane, 7-methyl-       | 198 | C14H30  | 026730-14-3 | 64   |
| 4    |    |   | Dodecane, 4,6-dimethyl-    | 198 | C14H30  | 061141-72-8 | 59   |
| 5    |    |   | Undecane, 6,6-dimethyl-    | 184 | C13H28  | 017312-76-4 | 53   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

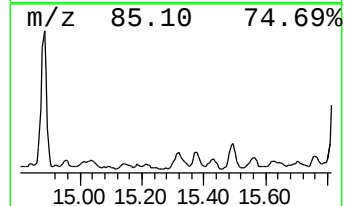
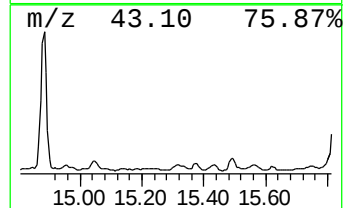
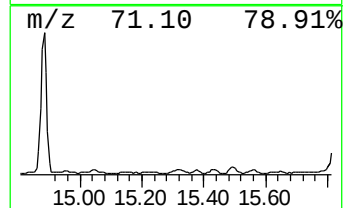
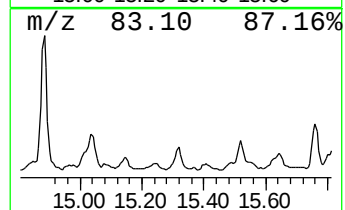
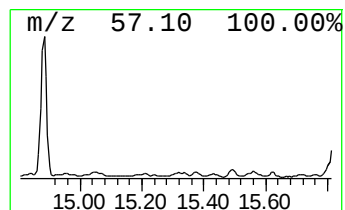
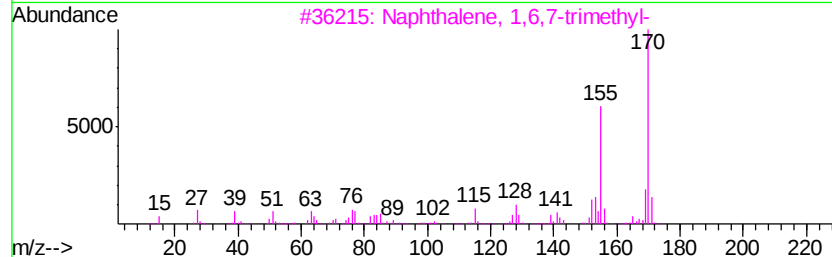
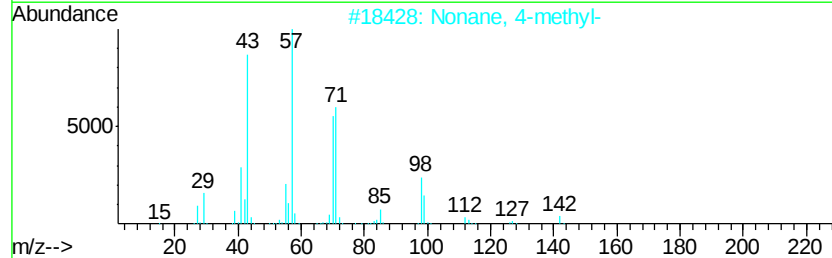
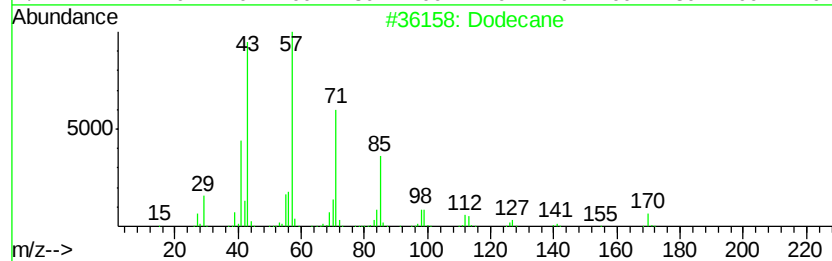
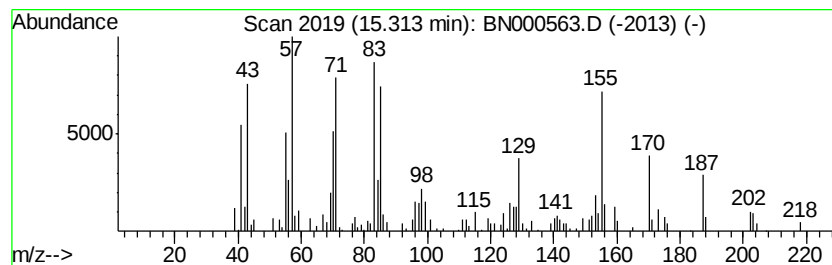
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 50 (DEL) Alkane: Straight-Chai... Concentration Rank 63

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 15.31 | 2.08 ng/ul | 233954 | Acenaphthene-d10 | 15.03 |

| Hit# | of | 5 | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Dodecane                           | 170 | C12H26  | 000112-40-3 | 46   |
| 2    |    |   | Nonane, 4-methyl-                  | 142 | C10H22  | 017301-94-9 | 43   |
| 3    |    |   | Naphthalene, 1,6,7-trimethyl-      | 170 | C13H14  | 002245-38-7 | 35   |
| 4    |    |   | 5-Methyl-1-phenylhexa-1,3,4-triene | 170 | C13H14  | 103240-79-5 | 35   |
| 5    |    |   | 4-Heptanone, 2-methyl-             | 128 | C8H16O  | 000626-33-5 | 25   |



Data Path : Z:\HPCHEM1\BNA\_N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

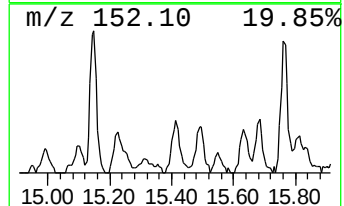
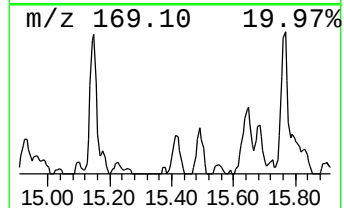
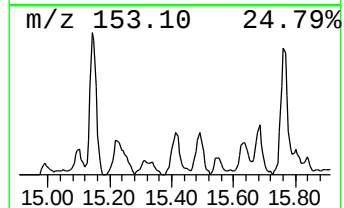
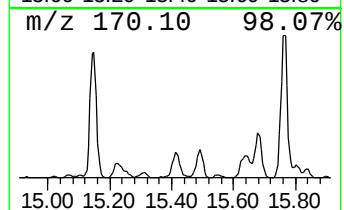
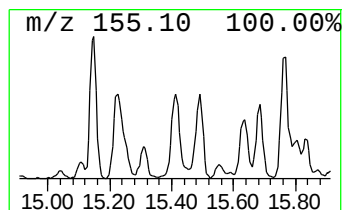
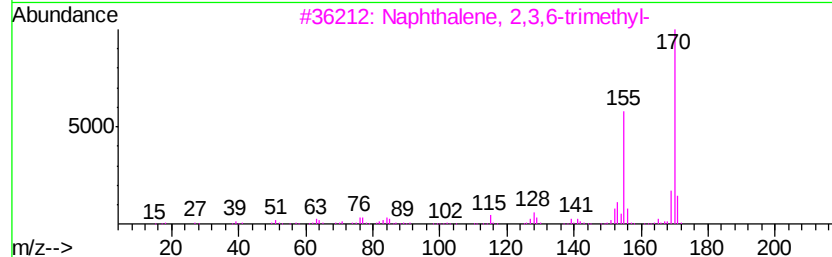
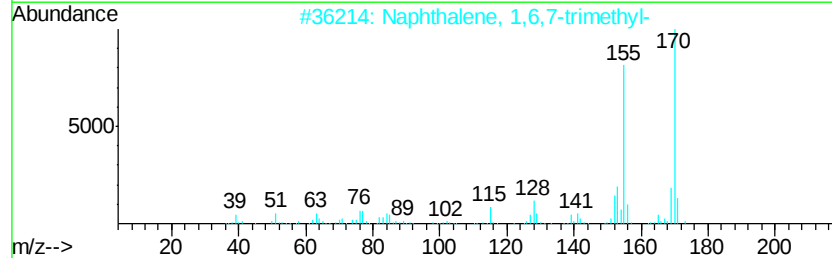
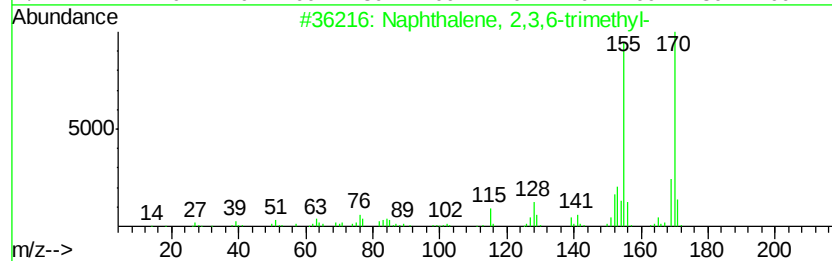
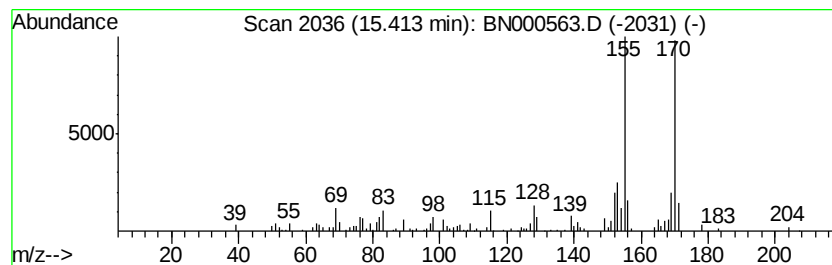
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 51 Naphthalene, 2,3,6-trimethyl- Concentration Rank 52

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 15.41 | 2.66 ng/ul | 299246 | Acenaphthene-d10 | 15.03 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 96   |
| 2    |    |   | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 96   |
| 3    |    |   | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 96   |
| 4    |    |   | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 95   |
| 5    |    |   | Naphthalene, 1,4,5-trimethyl- | 170 | C13H14  | 002131-41-1 | 95   |



Data Path : Z:\HPCHEM1\BNA\_N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

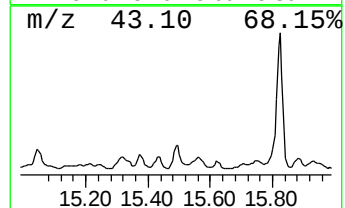
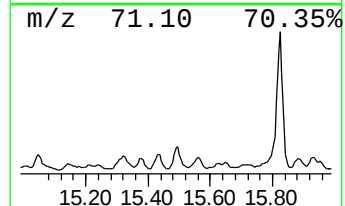
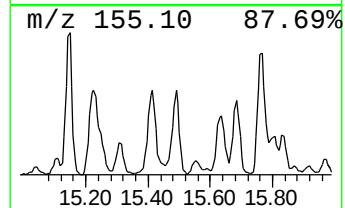
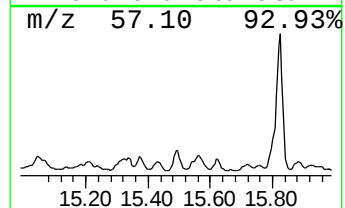
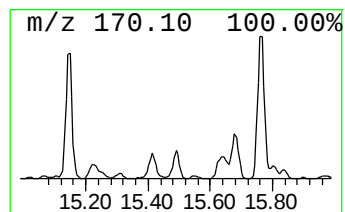
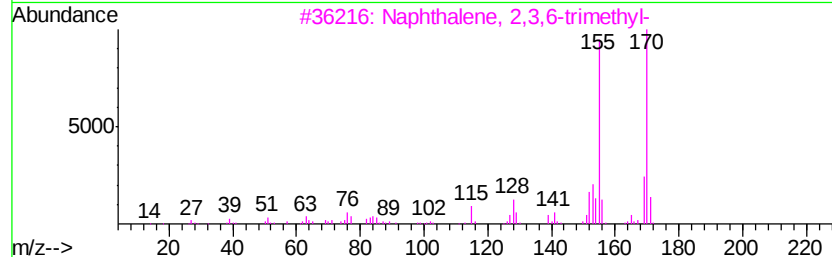
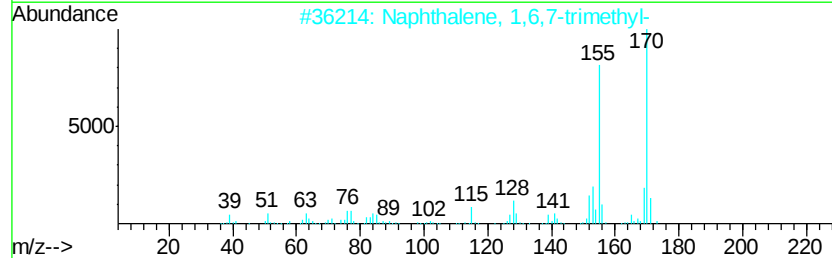
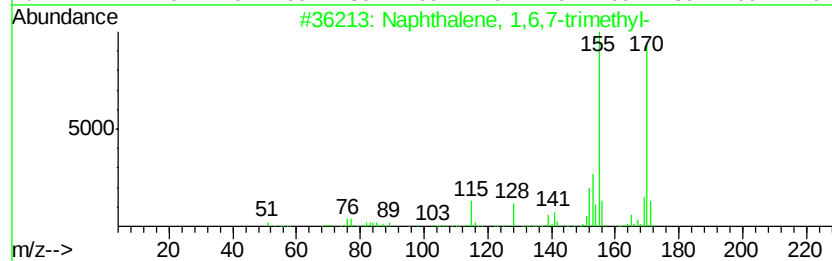
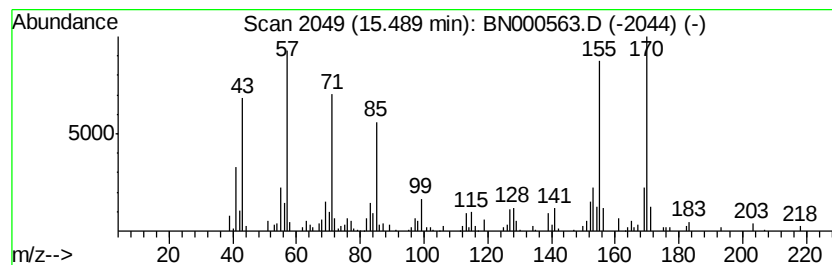
Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

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

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Peak Number 52 Naphthalene, 1,6,7-trimethyl- Concentration Rank 60

| R.T.    | EstConc    | Area                          | Relative to ISTD | R.T.           |
|---------|------------|-------------------------------|------------------|----------------|
| 15.49   | 2.20 ng/ul | 248282                        | Acenaphthene-d10 | 15.03          |
| Hit# of | 5          | Tentative ID                  | MW MolForm       | CAS# Qual      |
| 1       |            | Naphthalene, 1,6,7-trimethyl- | 170 C13H14       | 002245-38-7 96 |
| 2       |            | Naphthalene, 1,6,7-trimethyl- | 170 C13H14       | 002245-38-7 96 |
| 3       |            | Naphthalene, 2,3,6-trimethyl- | 170 C13H14       | 000829-26-5 96 |
| 4       |            | Naphthalene, 1,4,6-trimethyl- | 170 C13H14       | 002131-42-2 93 |
| 5       |            | Naphthalene, 2,3,6-trimethyl- | 170 C13H14       | 000829-26-5 93 |



Data Path : Z:\HPCHEM1\BNA\_N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

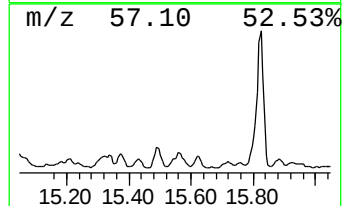
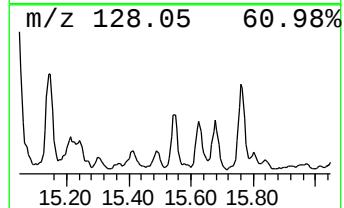
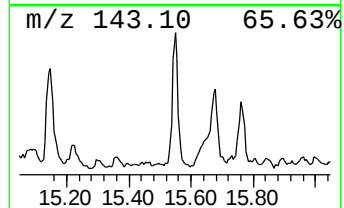
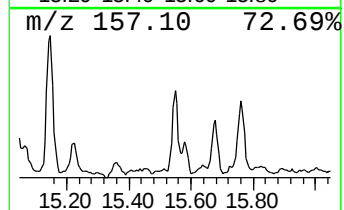
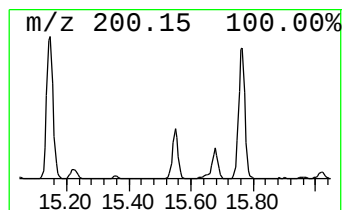
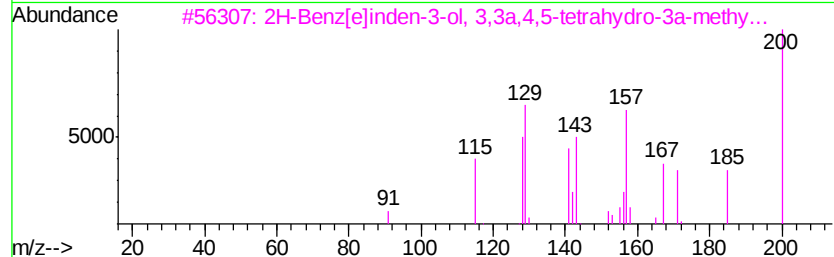
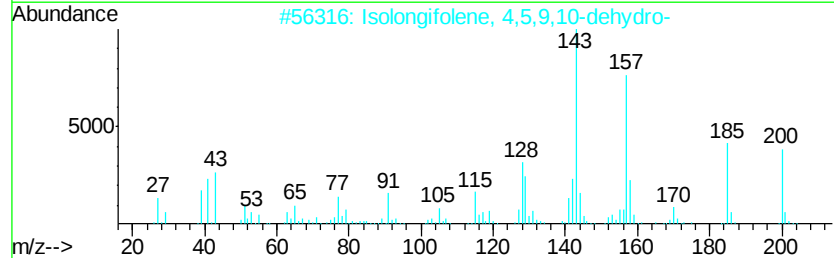
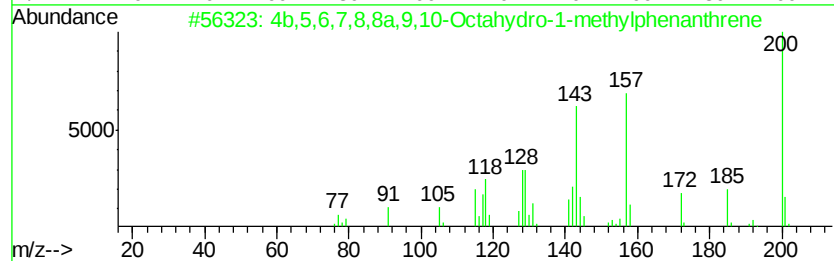
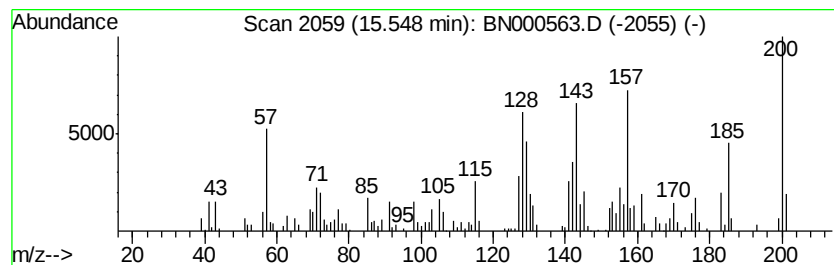
Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 53 4b,5,6,7,8,8a,9,10-Octahydr... Concentration Rank 44

| R.T.    | EstConc                              | Area          | Relative to ISTD | R.T.      |
|---------|--------------------------------------|---------------|------------------|-----------|
| 15.55   | 2.91 ng/ul                           | 328060        | Acenaphthene-d10 | 15.03     |
| Hit# of | 5                                    | Tentative ID  | MW MolForm       | CAS# Qual |
| 1       | 4b,5,6,7,8,8a,9,10-Octahydro-1-m...  | 200 C15H20    | 1000080-19-3     | 64        |
| 2       | Isolongifolene, 4,5,9,10-dehydro-    | 200 C15H20    | 156747-45-4      | 58        |
| 3       | 2H-Benz[e]indene-3-ol, 3,3a,4,5-t... | 200 C14H16O   | 071805-91-9      | 52        |
| 4       | 1H-Indene-1,3(2H)-dione, 2-(2-me...  | 200 C13H12O2  | 022123-54-2      | 47        |
| 5       | 3,1-Benzoxazepine-2-carbonitrile...  | 200 C11H8N2O2 | 019062-88-5      | 41        |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

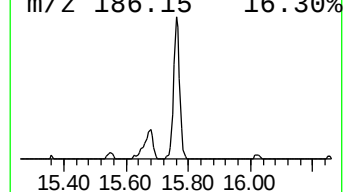
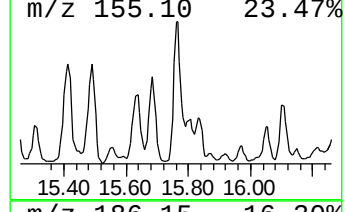
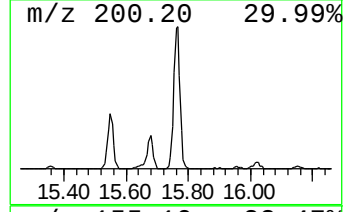
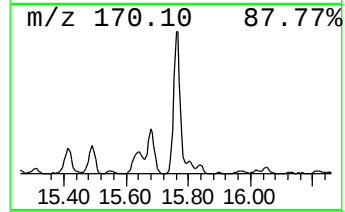
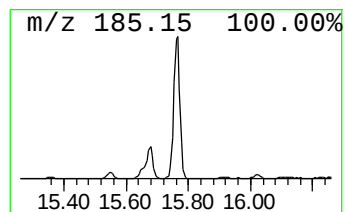
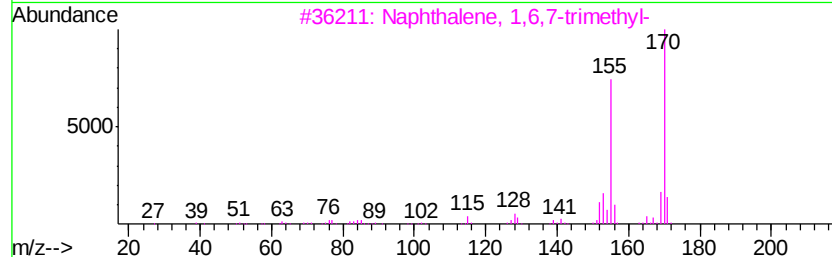
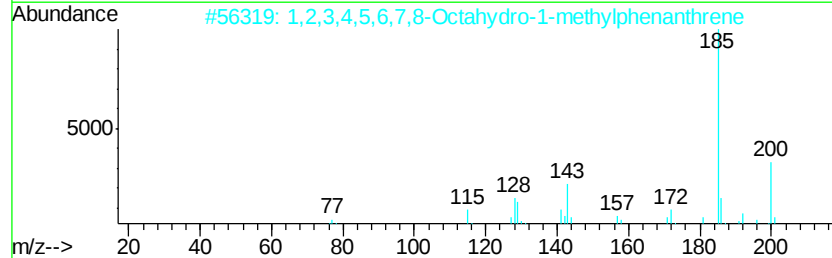
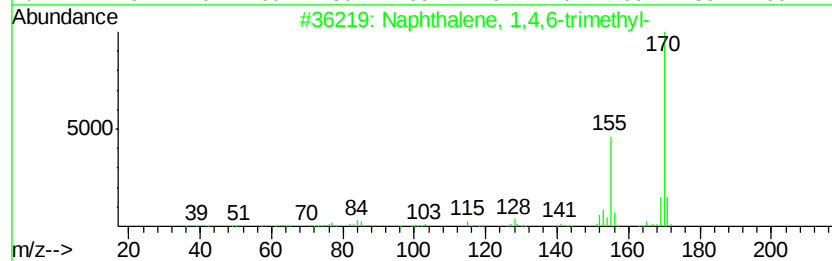
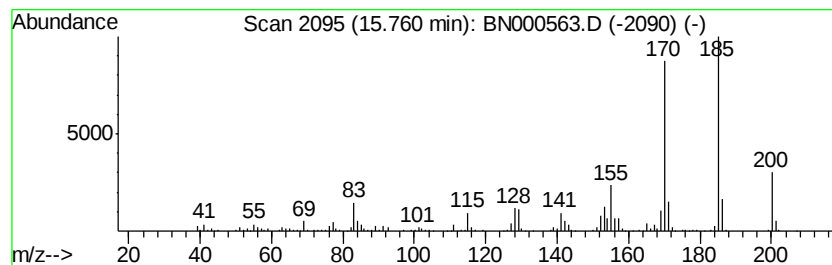
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 54 Naphthalene, 1,4,6-trimethyl- Concentration Rank 12

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 15.76 | 6.69 ng/ul | 753415 | Acenaphthene-d10 | 15.03 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | Naphthalene, 1,4,6-trimethyl-       | 170 | C13H14  | 002131-42-2  | 50   |
| 2    |    |   | 1,2,3,4,5,6,7,8-Octahydro-1-meth... | 200 | C15H20  | 1000080-19-4 | 50   |
| 3    |    |   | Naphthalene, 1,6,7-trimethyl-       | 170 | C13H14  | 002245-38-7  | 46   |
| 4    |    |   | Azulene, 4,6,8-trimethyl-           | 170 | C13H14  | 000941-81-1  | 46   |
| 5    |    |   | Naphthalene, 2,3,6-trimethyl-       | 170 | C13H14  | 000829-26-5  | 43   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

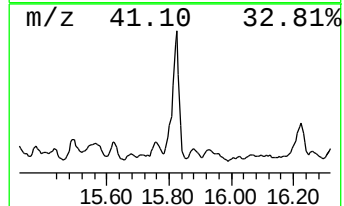
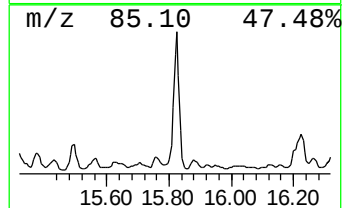
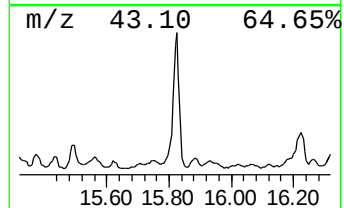
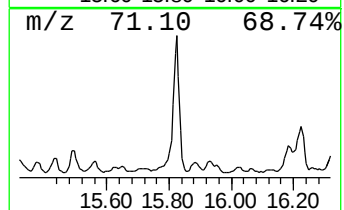
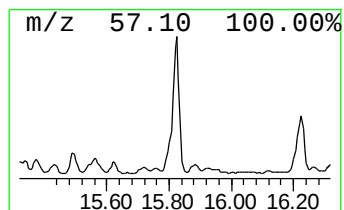
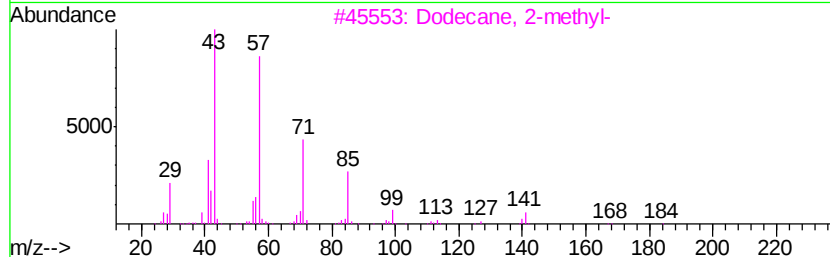
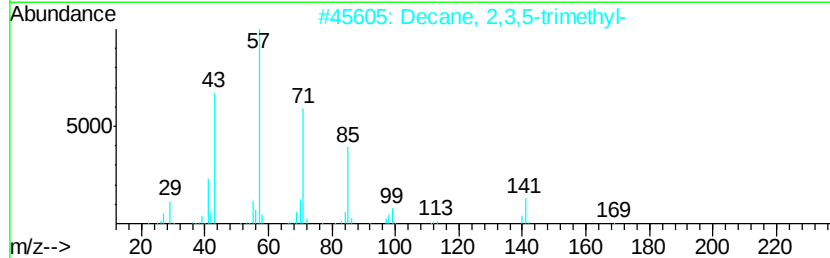
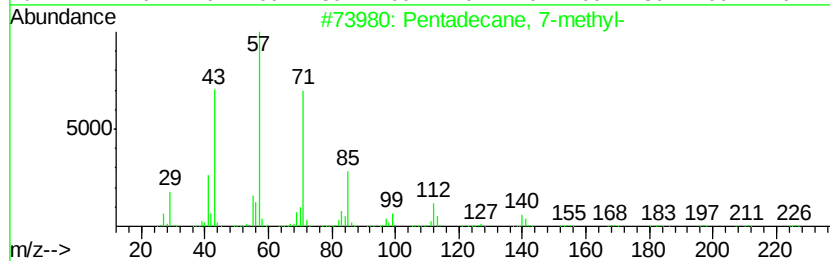
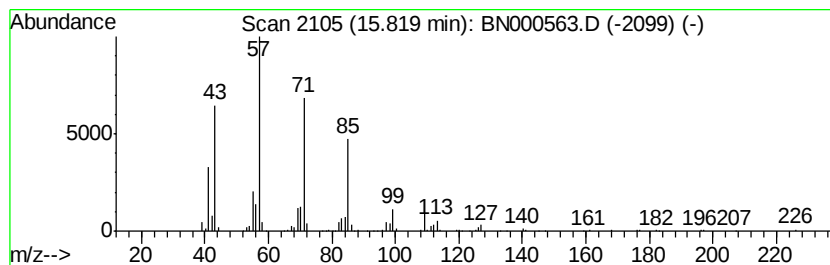
Instrument :  
BNA\_N  
ClientSampleID :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

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

\*\*\*\*\*  
Peak Number 55 (DEL) Alkane: Straight-Chai... Concentration Rank 6

| R.T.      | EstConc                  | Area   | Relative to ISTD | R.T.        |      |
|-----------|--------------------------|--------|------------------|-------------|------|
| 15.82     | 8.16 ng/ul               | 918483 | Acenaphthene-d10 | 15.03       |      |
| Hit# of 5 | Tentative ID             | MW     | MolForm          | CAS#        | Qual |
| 1         | Pentadecane, 7-methyl-   | 226    | C16H34           | 006165-40-8 | 90   |
| 2         | Decane, 2,3,5-trimethyl- | 184    | C13H28           | 062238-11-3 | 90   |
| 3         | Dodecane, 2-methyl-      | 184    | C13H28           | 001560-97-0 | 86   |
| 4         | Eicosane                 | 282    | C20H42           | 000112-95-8 | 72   |
| 5         | Nonane, 5-butyl-         | 184    | C13H28           | 017312-63-9 | 64   |





Data Path : Z:\HPCHEM1\BNA\_N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

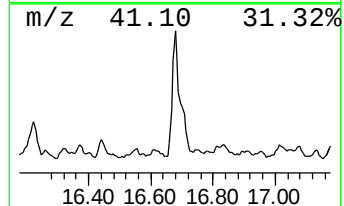
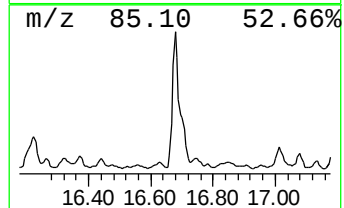
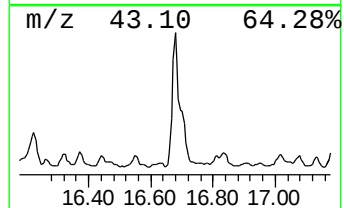
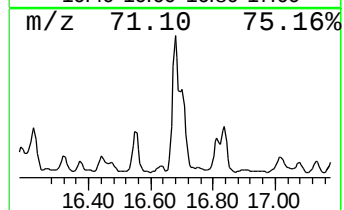
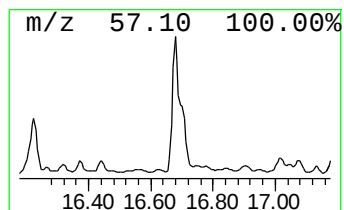
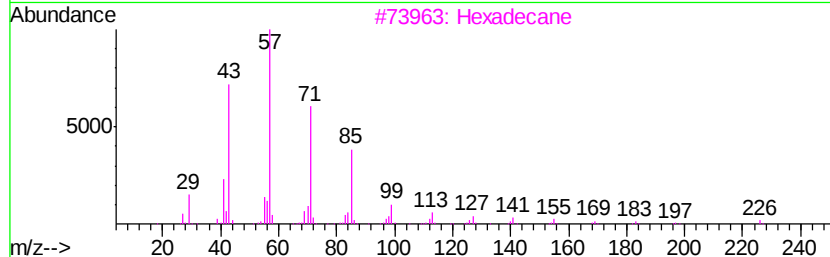
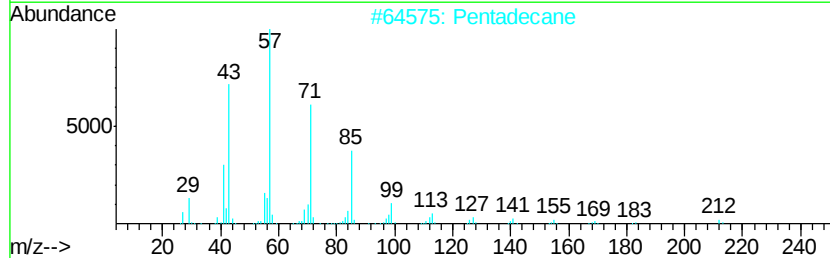
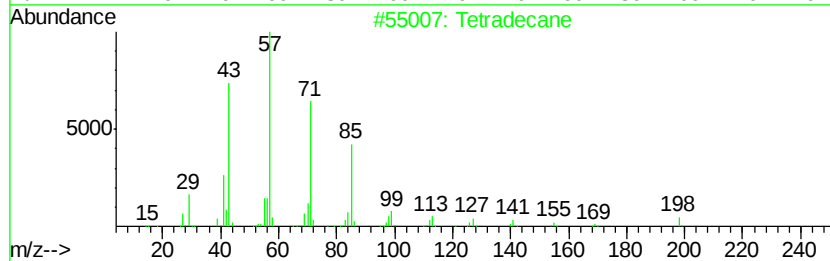
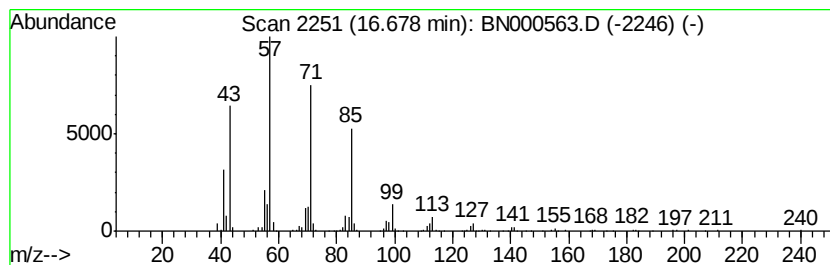
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 57 (DEL) Alkane: Straight-Chai... Concentration Rank 19

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 16.68 | 5.95 ng/ul | 736802 | Phenanthrene-d10 | 17.76 |

| Hit# | of | 5 | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------|-----|---------|-------------|------|
| 1    |    |   | Tetradecane              | 198 | C14H30  | 000629-59-4 | 90   |
| 2    |    |   | Pentadecane              | 212 | C15H32  | 000629-62-9 | 90   |
| 3    |    |   | Hexadecane               | 226 | C16H34  | 000544-76-3 | 90   |
| 4    |    |   | Dodecane, 1-iodo-        | 296 | C12H25I | 004292-19-7 | 86   |
| 5    |    |   | Decane, 2,3,5-trimethyl- | 184 | C13H28  | 062238-11-3 | 86   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

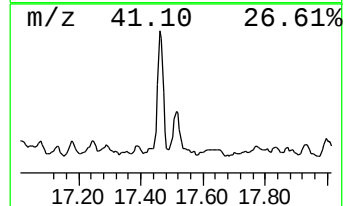
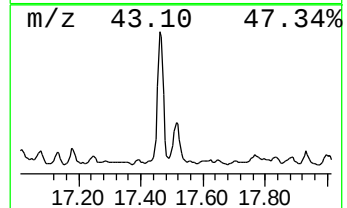
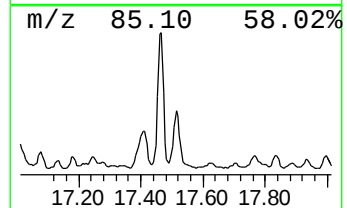
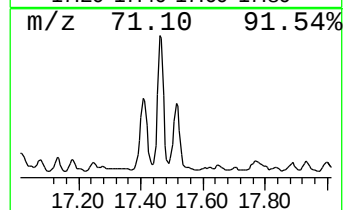
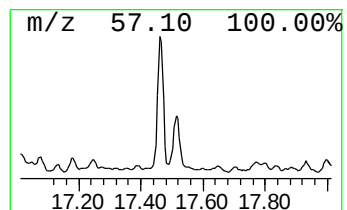
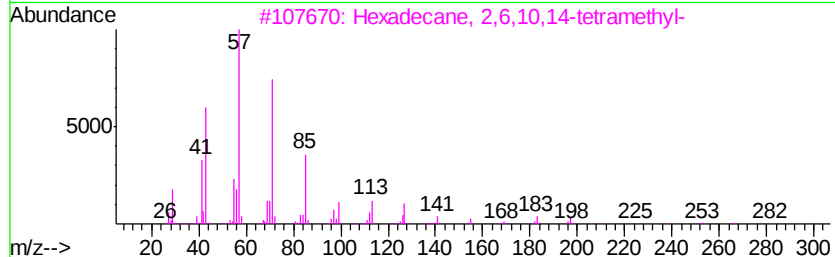
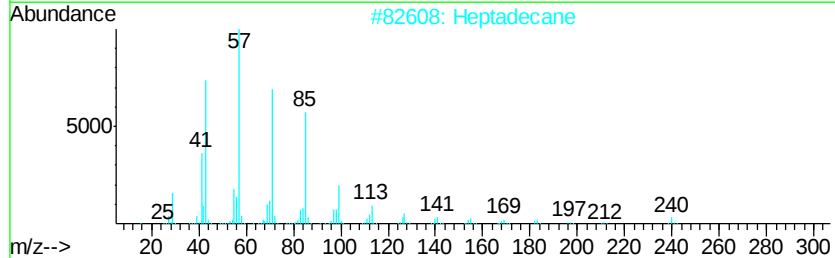
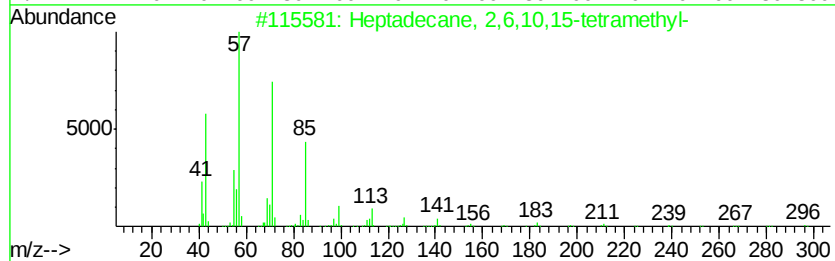
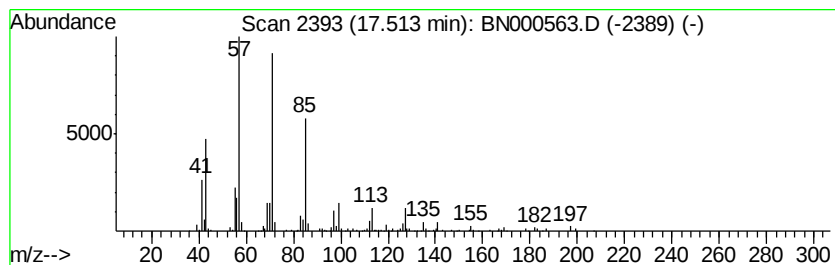
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 58 (DEL) Alkane: Straight-Chai... Concentration Rank 48

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 17.51 | 2.78 ng/ul | 344733 | Phenanthrene-d10 | 17.76 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44  | 054833-48-6 | 87   |
| 2    |    |   | Heptadecane                         | 240 | C17H36  | 000629-78-7 | 86   |
| 3    |    |   | Hexadecane, 2,6,10,14-tetramethyl-  | 282 | C20H42  | 000638-36-8 | 83   |
| 4    |    |   | Tetracosane                         | 338 | C24H50  | 000646-31-1 | 78   |
| 5    |    |   | Tetradecane, 1-iodo-                | 324 | C14H29I | 019218-94-1 | 78   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

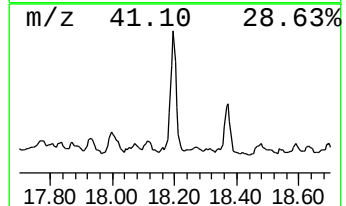
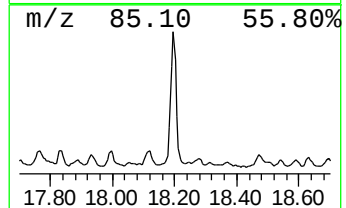
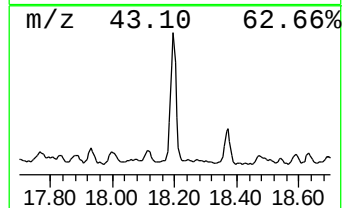
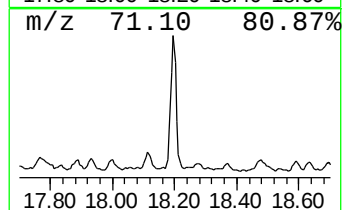
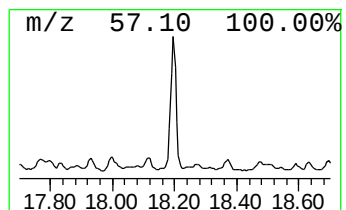
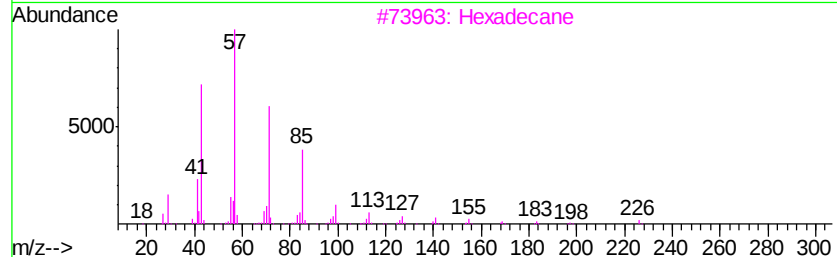
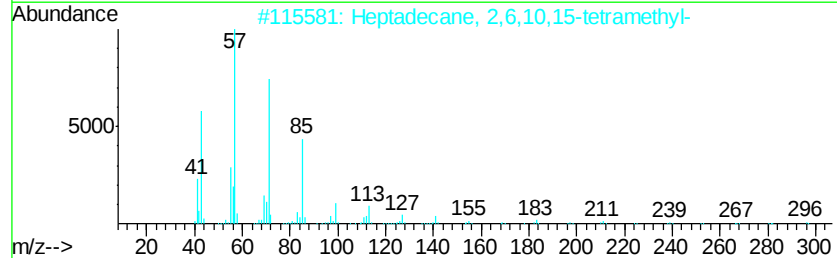
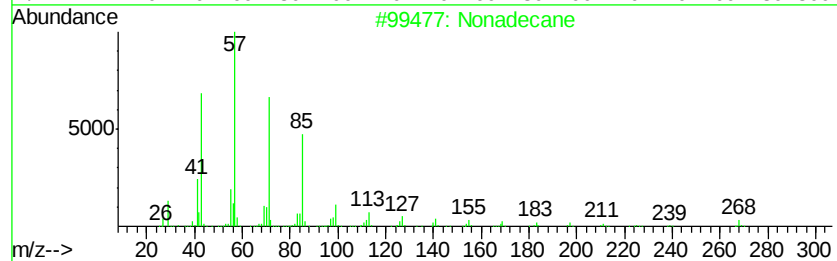
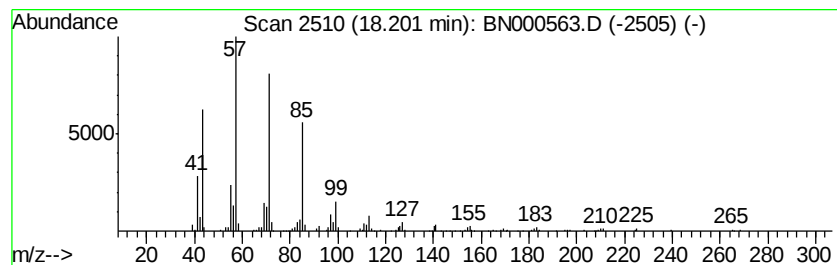
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 59 (DEL) Alkane: Straight-Chai... Concentration Rank 25

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.20 | 4.37 ng/ul | 541342 | Phenanthrene-d10 | 17.76 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Nonadecane                          | 268 | C19H40  | 000629-92-5 | 90   |
| 2    |    |   | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44  | 054833-48-6 | 90   |
| 3    |    |   | Hexadecane                          | 226 | C16H34  | 000544-76-3 | 86   |
| 4    |    |   | Tetracosane                         | 338 | C24H50  | 000646-31-1 | 86   |
| 5    |    |   | Nonane, 3,7-dimethyl-               | 156 | C11H24  | 017302-32-8 | 80   |



Data Path : Z:\HPCHEM1\BNA\_N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

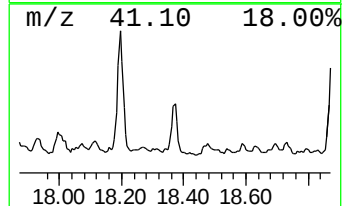
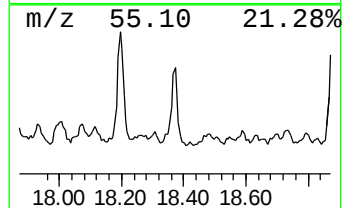
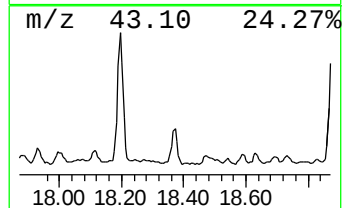
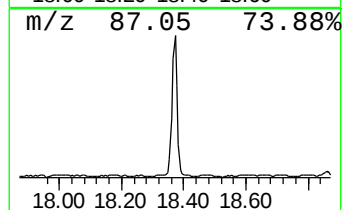
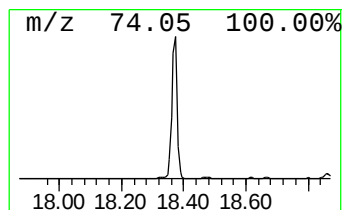
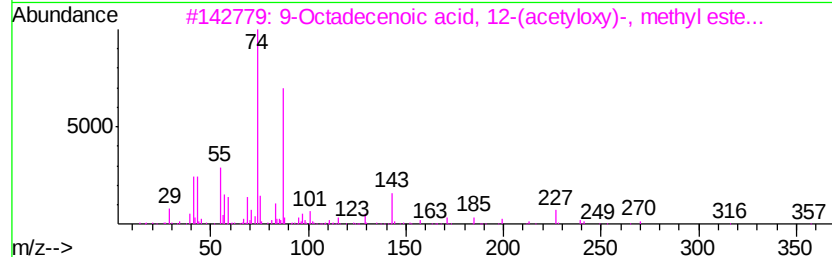
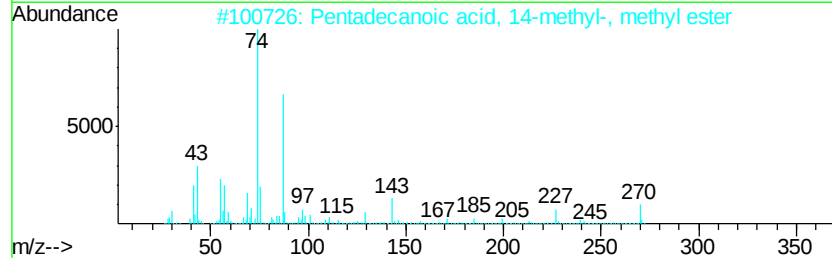
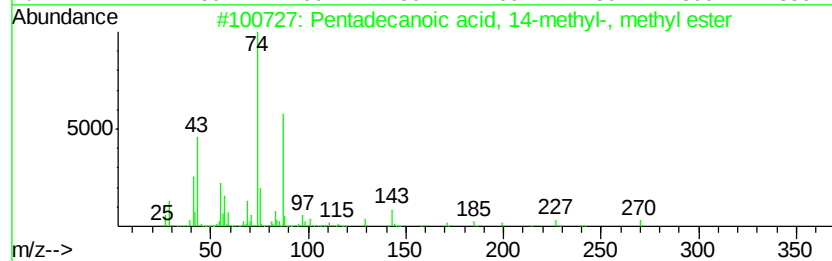
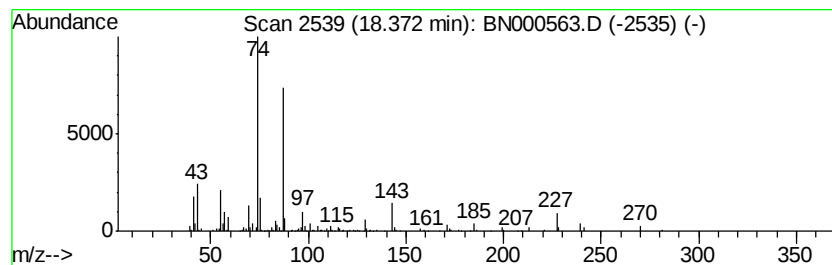
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 60 Pentadecanoic acid, 14-meth... Concentration Rank 57

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.37 | 2.41 ng/ul | 297920 | Phenanthrene-d10 | 17.76 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Pentadecanoic acid, 14-methyl-, ... | 270 | C17H34O2 | 005129-60-2 | 94   |
| 2    |    |   | Pentadecanoic acid, 14-methyl-, ... | 270 | C17H34O2 | 005129-60-2 | 91   |
| 3    |    |   | 9-Octadecenoic acid, 12-(acetylo... | 354 | C21H38O4 | 000140-03-4 | 90   |
| 4    |    |   | Hexadecanoic acid, methyl ester     | 270 | C17H34O2 | 000112-39-0 | 90   |
| 5    |    |   | Tridecanoic acid, methyl ester      | 228 | C14H28O2 | 001731-88-0 | 87   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

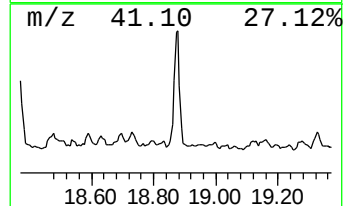
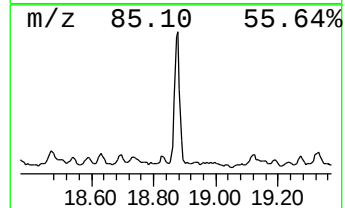
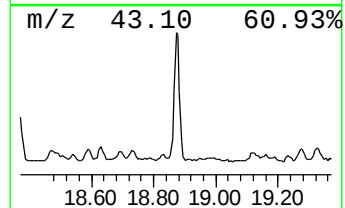
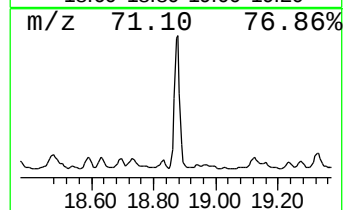
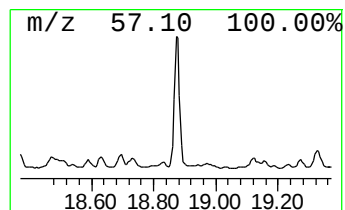
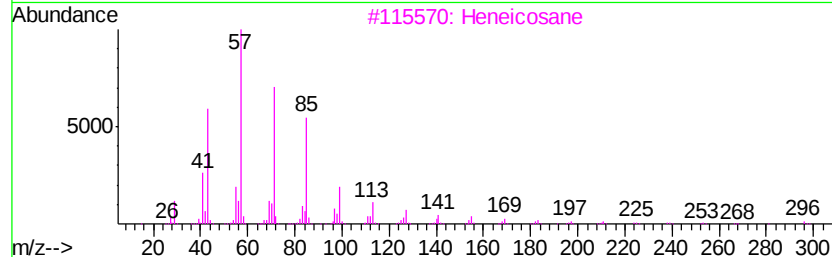
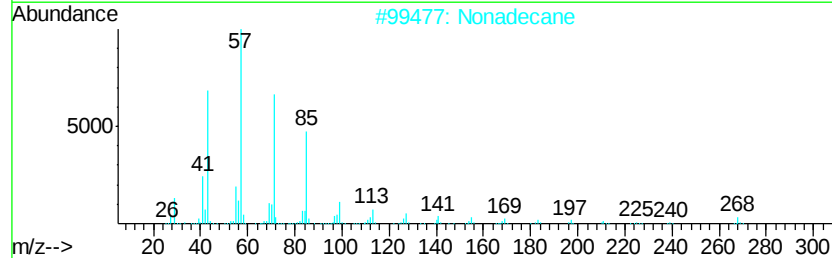
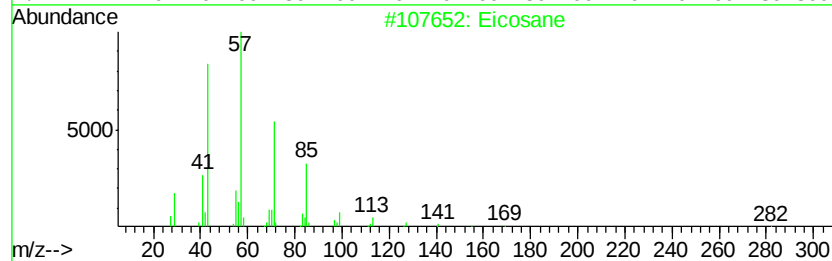
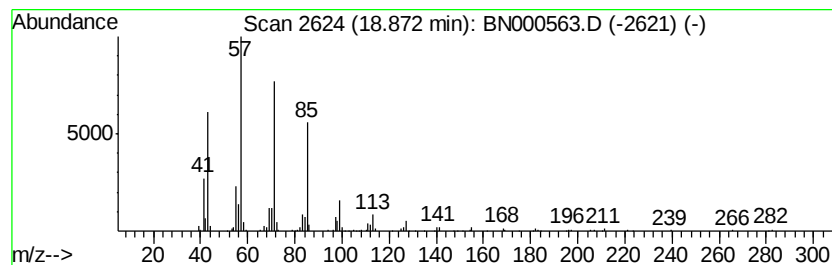
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 61 (DEL) Alkane: Straight-Chai... Concentration Rank 50

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 18.87 | 2.72 ng/ul | 336721 | Phenanthrene-d10 | 17.76 |

| Hit# | of | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------|-----|---------|-------------|------|
| 1    |    |   | Eicosane     | 282 | C20H42  | 000112-95-8 | 93   |
| 2    |    |   | Nonadecane   | 268 | C19H40  | 000629-92-5 | 91   |
| 3    |    |   | Heneicosane  | 296 | C21H44  | 000629-94-7 | 91   |
| 4    |    |   | Hexadecane   | 226 | C16H34  | 000544-76-3 | 90   |
| 5    |    |   | Tetracosane  | 338 | C24H50  | 000646-31-1 | 90   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

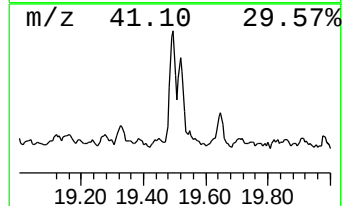
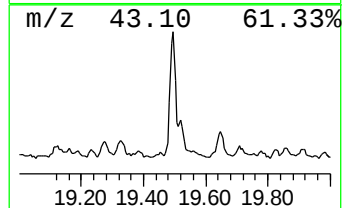
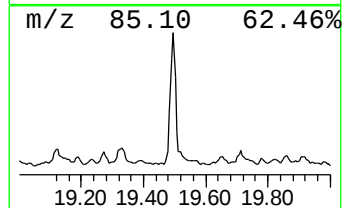
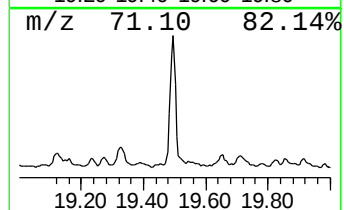
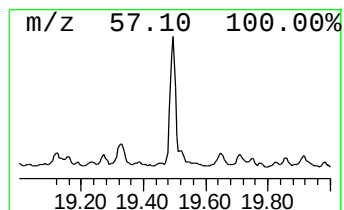
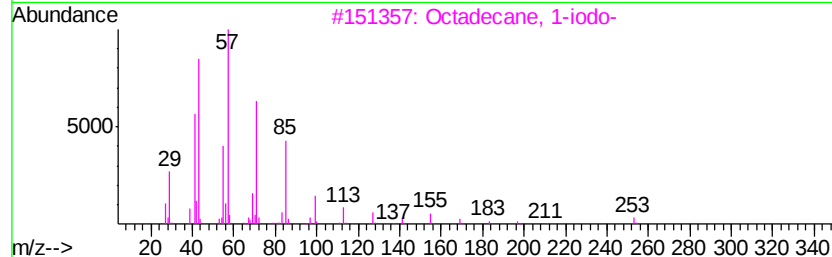
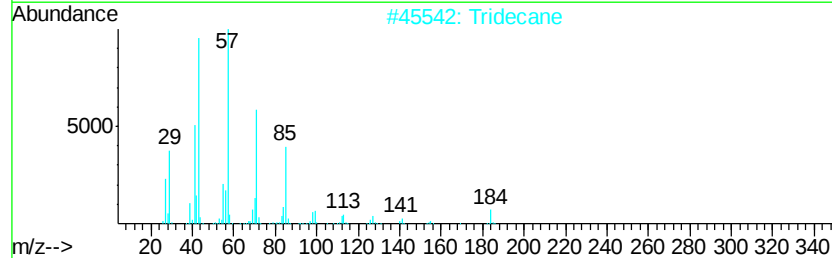
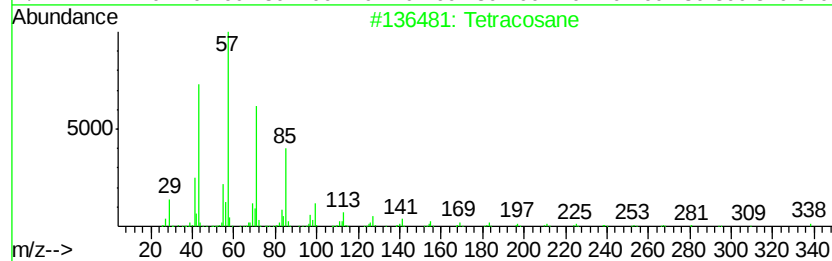
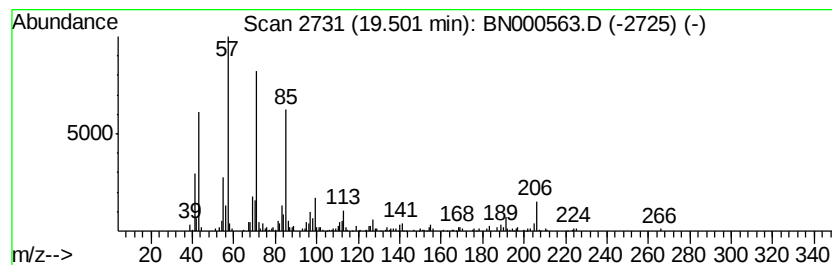
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 62 (DEL) Alkane: Straight-Chai... Concentration Rank 46

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 19.50 | 2.80 ng/ul | 346437 | Phenanthrene-d10 | 17.76 |

| Hit# | of | 5 | Tentative ID        | MW  | MolForm | CAS#        | Qual |
|------|----|---|---------------------|-----|---------|-------------|------|
| 1    |    |   | Tetracosane         | 338 | C24H50  | 000646-31-1 | 91   |
| 2    |    |   | Tridecane           | 184 | C13H28  | 000629-50-5 | 87   |
| 3    |    |   | Octadecane, 1-iodo- | 380 | C18H37I | 000629-93-6 | 86   |
| 4    |    |   | Dodecane, 1-iodo-   | 296 | C12H25I | 004292-19-7 | 86   |
| 5    |    |   | Octadecane          | 254 | C18H38  | 000593-45-3 | 80   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampled :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

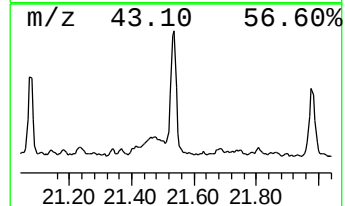
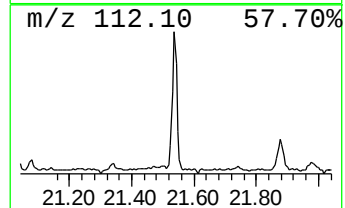
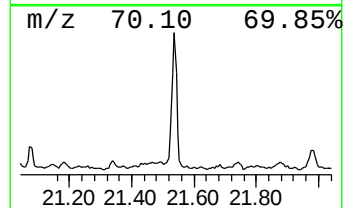
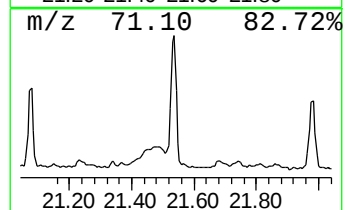
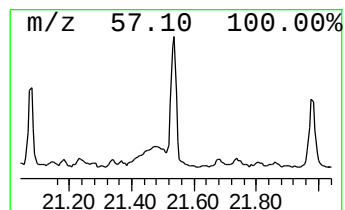
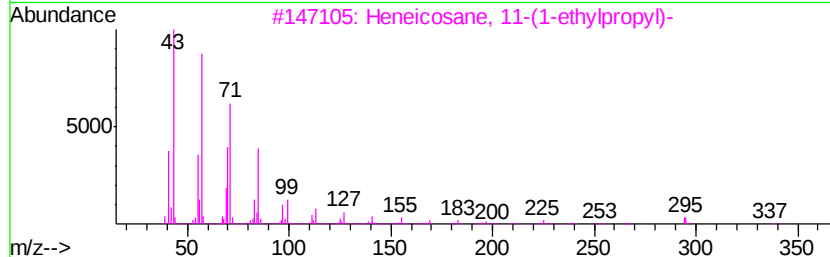
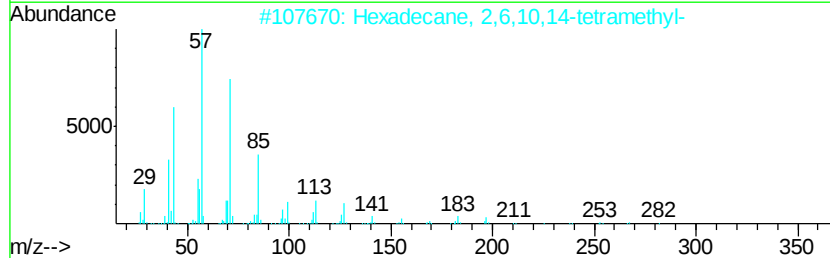
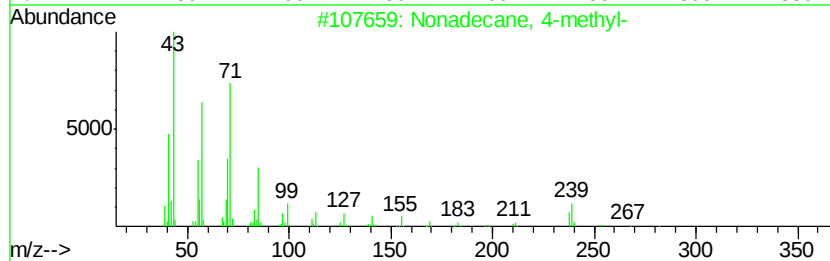
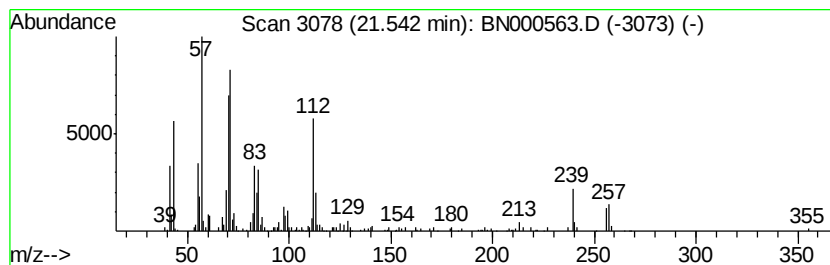
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 64 (DEL) Alkane: Straight-Chai... Concentration Rank 54

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 21.54 | 2.58 ng/ul | 343216 | Chrysene-d12     | 21.88 |

| Hit# | of | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Nonadecane, 4-methyl-              | 282 | C20H42  | 025117-27-5 | 89   |
| 2    |    | Hexadecane, 2,6,10,14-tetramethyl- | 282 | C20H42  | 000638-36-8 | 60   |
| 3    |    | Heneicosane, 11-(1-ethylpropyl)-   | 366 | C26H54  | 055282-11-6 | 59   |
| 4    |    | Decane, 2-methyl-                  | 156 | C11H24  | 006975-98-0 | 55   |
| 5    |    | Hexatriacontane                    | 507 | C36H74  | 000630-06-8 | 50   |



Data Path : Z:\HPCHEM1\BNA N\DATA\BN032218\  
Data File : BN000563.D  
Acq On : 23 Mar 2018 04:28  
Operator : JU/SJ  
Sample : J1905-03DL2 200X  
Misc :  
ALS Vial : 31 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2

Quant Method : Z:\HPCHEM1\BNA N\METHODS\SOM-EPA-BN031918.M  
Quant Title : SVOA CALIBRATION

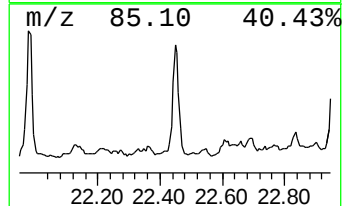
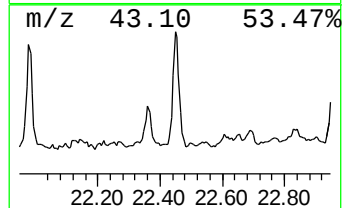
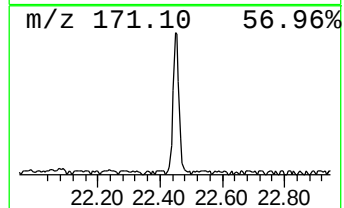
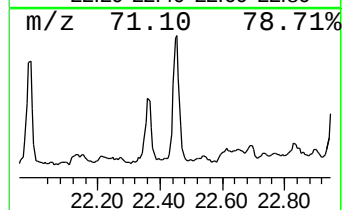
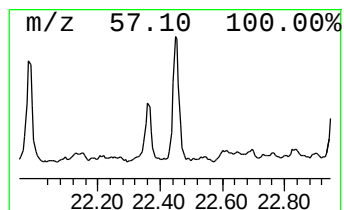
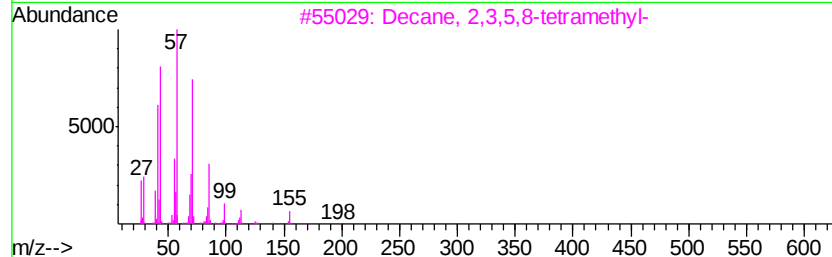
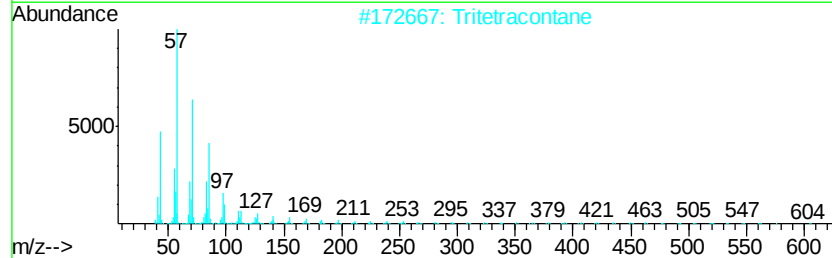
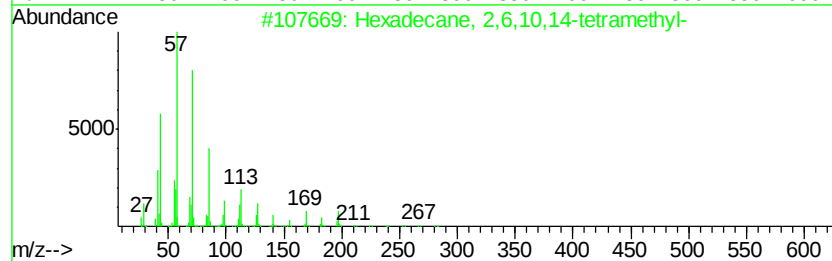
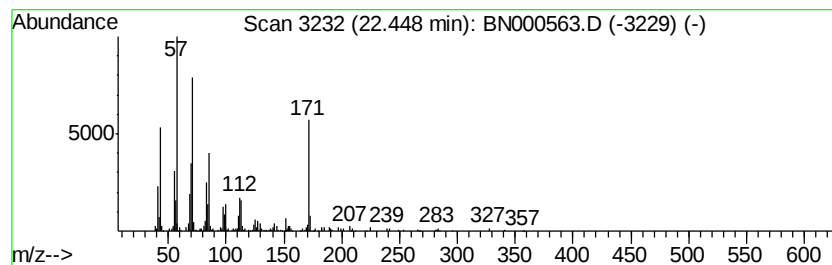
TIC Library : C:\DATABASE\NIST02.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 65 (DEL) Alkane: Straight-Chai... Concentration Rank 65

| R.T.  | EstConc    | Area   | Relative to ISTD | R.T.  |
|-------|------------|--------|------------------|-------|
| 22.45 | 2.04 ng/ul | 271473 | Chrysene-d12     | 21.88 |

| Hit# | of | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Hexadecane, 2,6,10,14-tetramethyl- | 282 | C20H42  | 000638-36-8 | 52   |
| 2    |    | Tritetracontane                    | 605 | C43H88  | 007098-21-7 | 49   |
| 3    |    | Decane, 2,3,5,8-tetramethyl-       | 198 | C14H30  | 192823-15-7 | 46   |
| 4    |    | Hexadecane                         | 226 | C16H34  | 000544-76-3 | 46   |
| 5    |    | Dodecane, 4-methyl-                | 184 | C13H28  | 006117-97-1 | 43   |



Data Path : Z:\HPCHEM1\BNA\_N\DATA\BN032218\  
 Data File : BN000563.D  
 Acq On : 23 Mar 2018 04:28  
 Operator : JU/SJ  
 Sample : J1905-03DL2 200X  
 Misc :  
 ALS Vial : 31 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 DAM40DL2

Quant Method : Z:\HPCHEM1\BNA\_N\METHODS\SOM-EPA-BN031918.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 |
| (DEL) Alkane: Str...  | 6.35  | 8.0     | ng/ul | 401921   | 1                     | 8.41  | 1007730 | 20.0 |
| (DEL) Alkane: Cyc...  | 7.00  | 4.2     | ng/ul | 211343   | 1                     | 8.41  | 1007730 | 20.0 |
| (DEL) Alkane: Str...  | 7.35  | 4.3     | ng/ul | 214458   | 1                     | 8.41  | 1007730 | 20.0 |
| (DEL) Alkane: Str...  | 7.41  | 4.2     | ng/ul | 212987   | 1                     | 8.41  | 1007730 | 20.0 |
| Benzene, 1-ethyl-...  | 7.47  | 4.2     | ng/ul | 209383   | 1                     | 8.41  | 1007730 | 20.0 |
| (DEL) Alkane: Str...  | 7.53  | 7.7     | ng/ul | 388101   | 1                     | 8.41  | 1007730 | 20.0 |
| Benzene, 1-ethyl-...  | 7.78  | 3.4     | ng/ul | 170953   | 1                     | 8.41  | 1007730 | 20.0 |
| (DEL) Alkane: Str...  | 8.00  | 22.7    | ng/ul | 1141830  | 1                     | 8.41  | 1007730 | 20.0 |
| (DEL) Alkane: Str...  | 8.36  | 4.2     | ng/ul | 211907   | 1                     | 8.41  | 1007730 | 20.0 |
| Benzene, 1,2,3-tr...  | 8.53  | 4.7     | ng/ul | 235286   | 1                     | 8.41  | 1007730 | 20.0 |
| (DEL) Alkane: Cyc...  | 8.69  | 2.6     | ng/ul | 131942   | 1                     | 8.41  | 1007730 | 20.0 |
| (DEL) Alkane: Str...  | 8.92  | 5.0     | ng/ul | 249193   | 1                     | 8.41  | 1007730 | 20.0 |
| (DEL) Alkane: Str...  | 8.98  | 7.8     | ng/ul | 391756   | 1                     | 8.41  | 1007730 | 20.0 |
| Benzene, 1,4-diet...  | 9.05  | 12.2    | ng/ul | 616280   | 1                     | 8.41  | 1007730 | 20.0 |
| (DEL) Alkane: Str...  | 9.16  | 5.5     | ng/ul | 275076   | 1                     | 8.41  | 1007730 | 20.0 |
| Benzene, 1-methyl...  | 9.21  | 3.4     | ng/ul | 168681   | 1                     | 8.41  | 1007730 | 20.0 |
| Benzene, 1-methyl...  | 9.42  | 3.7     | ng/ul | 184990   | 1                     | 8.41  | 1007730 | 20.0 |
| Benzene, 4-ethyl-...  | 9.52  | 6.0     | ng/ul | 302518   | 1                     | 8.41  | 1007730 | 20.0 |
| (DEL) Alkane: Str...  | 9.63  | 28.2    | ng/ul | 1418830  | 1                     | 8.41  | 1007730 | 20.0 |
| unknown-01            | 9.77  | 5.7     | ng/ul | 284965   | 1                     | 8.41  | 1007730 | 20.0 |
| Benzene, 2-ethyl-...  | 9.86  | 2.7     | ng/ul | 218205   | 2                     | 11.25 | 1621940 | 20.0 |
| (DEL) Alkane: Cyc...  | 10.35 | 3.1     | ng/ul | 251834   | 2                     | 11.25 | 1621940 | 20.0 |
| (DEL) Alkane: Str...  | 10.48 | 2.3     | ng/ul | 182997   | 2                     | 11.25 | 1621940 | 20.0 |
| (DEL) Alkane: Str...  | 10.63 | 2.8     | ng/ul | 226338   | 2                     | 11.25 | 1621940 | 20.0 |
| (DEL) Alkane: Str...  | 11.18 | 10.3    | ng/ul | 831080   | 2                     | 11.25 | 1621940 | 20.0 |
| (DEL) Alkane: Str...  | 11.36 | 2.5     | ng/ul | 198641   | 2                     | 11.25 | 1621940 | 20.0 |
| Cyanoic acid, ethy... | 11.59 | 7.4     | ng/ul | 596699   | 2                     | 11.25 | 1621940 | 20.0 |
| (DEL) Alkane: Str...  | 11.72 | 3.7     | ng/ul | 296553   | 2                     | 11.25 | 1621940 | 20.0 |
| (DEL) Alkane: Cyc...  | 11.92 | 3.6     | ng/ul | 289593   | 2                     | 11.25 | 1621940 | 20.0 |
| (DEL) Alkane: Str...  | 12.22 | 4.0     | ng/ul | 325549   | 2                     | 11.25 | 1621940 | 20.0 |
| Benzene, 4-(2-but...  | 12.31 | 6.4     | ng/ul | 521376   | 2                     | 11.25 | 1621940 | 20.0 |
| 1,3-Propanediol, ...  | 12.45 | 5.1     | ng/ul | 417139   | 2                     | 11.25 | 1621940 | 20.0 |
| (DEL) Alkane: Str...  | 12.61 | 7.5     | ng/ul | 604416   | 2                     | 11.25 | 1621940 | 20.0 |
| unknown-02            | 12.65 | 4.2     | ng/ul | 341589   | 2                     | 11.25 | 1621940 | 20.0 |
| (DEL) Alkane: Cyc...  | 12.74 | 2.4     | ng/ul | 195827   | 2                     | 11.25 | 1621940 | 20.0 |
| 4-Nonanone, 2,6,8...  | 13.01 | 2.2     | ng/ul | 180607   | 2                     | 11.25 | 1621940 | 20.0 |
| Naphthalene, 1-me...  | 13.10 | 2.1     | ng/ul | 172187   | 2                     | 11.25 | 1621940 | 20.0 |
| (DEL) Alkane: Cyc...  | 13.24 | 2.2     | ng/ul | 245385   | 3                     | 15.03 | 2252140 | 20.0 |
| (DEL) Alkane: Str...  | 13.82 | 6.3     | ng/ul | 711119   | 3                     | 15.03 | 2252140 | 20.0 |
| Naphthalene, 2,6-...  | 14.21 | 3.6     | ng/ul | 400764   | 3                     | 15.03 | 2252140 | 20.0 |
| Naphthalene, 1,8-...  | 14.35 | 4.0     | ng/ul | 450259   | 3                     | 15.03 | 2252140 | 20.0 |
| (DEL) Alkane: Str...  | 14.47 | 2.9     | ng/ul | 322713   | 3                     | 15.03 | 2252140 | 20.0 |
| 3,4-Dimethyl(1H)p...  | 14.63 | 3.7     | ng/ul | 421468   | 3                     | 15.03 | 2252140 | 20.0 |
| (DEL) Alkane: Str...  | 14.88 | 15.6    | ng/ul | 1757570  | 3                     | 15.03 | 2252140 | 20.0 |
| (DEL) Alkane: Str...  | 15.31 | 2.1     | ng/ul | 233954   | 3                     | 15.03 | 2252140 | 20.0 |
| Naphthalene, 2,3,...  | 15.41 | 2.7     | ng/ul | 299246   | 3                     | 15.03 | 2252140 | 20.0 |
| Naphthalene, 1,6,...  | 15.49 | 2.2     | ng/ul | 248282   | 3                     | 15.03 | 2252140 | 20.0 |
| 4b,5,6,7,8,8a,9,1...  | 15.55 | 2.9     | ng/ul | 328060   | 3                     | 15.03 | 2252140 | 20.0 |
| Naphthalene, 1,4,...  | 15.76 | 6.7     | ng/ul | 753415   | 3                     | 15.03 | 2252140 | 20.0 |

|                      |       |           |        |   |       |         |      |
|----------------------|-------|-----------|--------|---|-------|---------|------|
| (DEL) Alkane: Str... | 15.82 | 8.2 ng/ul | 918483 | 3 | 15.03 | 2252140 | 20.0 |
| unknown-03           | 16.61 | 2.7 ng/ul | 339208 | 4 | 17.76 | 2477090 | 20.0 |
| (DEL) Alkane: Str... | 16.68 | 6.0 ng/ul | 736802 | 4 | 17.76 | 2477090 | 20.0 |
| (DEL) Alkane: Str... | 17.51 | 2.8 ng/ul | 344733 | 4 | 17.76 | 2477090 | 20.0 |
| (DEL) Alkane: Str... | 18.20 | 4.4 ng/ul | 541342 | 4 | 17.76 | 2477090 | 20.0 |
| Pentadecanoic aci... | 18.37 | 2.4 ng/ul | 297920 | 4 | 17.76 | 2477090 | 20.0 |
| (DEL) Alkane: Str... | 18.87 | 2.7 ng/ul | 336721 | 4 | 17.76 | 2477090 | 20.0 |
| (DEL) Alkane: Str... | 19.50 | 2.8 ng/ul | 346437 | 4 | 17.76 | 2477090 | 20.0 |
| (DEL) Alkane: Str... | 21.54 | 2.6 ng/ul | 343216 | 5 | 21.88 | 2664600 | 20.0 |
| (DEL) Alkane: Str... | 22.45 | 2.0 ng/ul | 271473 | 5 | 21.88 | 2664600 | 20.0 |

Instrument :  
BNA\_N  
ClientSampleId :  
DAM40DL2