

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
 Data File : BG051787.D  
 Acq On : 23 Dec 2021 16:23  
 Operator : CG/JU  
 Sample : M5215-02  
 Misc :  
 ALS Vial : 38 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BGLD4

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
 Title : SVOA CALIBRATION

Signal : TIC: BG051787.D\data.ms

| 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         | 5.680       | 368           | 373         | 376          | rBV2     | 125964         | 222854        | 3.72%           | 0.213%        |
| 2         | 5.745       | 379           | 384         | 392          | rVB      | 2461475        | 4107866       | 68.51%          | 3.927%        |
| 3         | 5.821       | 392           | 397         | 402          | rVB      | 160010         | 256580        | 4.28%           | 0.245%        |
| 4         | 5.903       | 402           | 411         | 425          | rVB      | 1960219        | 3546739       | 59.16%          | 3.390%        |
| 5         | 6.127       | 442           | 449         | 455          | rBV3     | 75793          | 168703        | 2.81%           | 0.161%        |
| 6         | 6.209       | 457           | 463         | 467          | rBV      | 150421         | 259035        | 4.32%           | 0.248%        |
| 7         | 6.268       | 467           | 473         | 482          | rVB5     | 112577         | 270179        | 4.51%           | 0.258%        |
| 8         | 6.526       | 510           | 517         | 525          | rBV4     | 398030         | 1001610       | 16.71%          | 0.957%        |
| 9         | 6.644       | 531           | 537         | 544          | rVB      | 369203         | 742963        | 12.39%          | 0.710%        |
| 10        | 6.738       | 544           | 553         | 561          | rBV6     | 441657         | 1004655       | 16.76%          | 0.960%        |
| 11        | 6.814       | 561           | 566         | 571          | rVB      | 770075         | 1095272       | 18.27%          | 1.047%        |
| 12        | 6.867       | 571           | 575         | 576          | rBV      | 201088         | 259485        | 4.33%           | 0.248%        |
| 13        | 6.902       | 576           | 581         | 595          | rVB4     | 637405         | 1747159       | 29.14%          | 1.670%        |
| 14        | 7.020       | 597           | 601         | 607          | rVB2     | 156586         | 261601        | 4.36%           | 0.250%        |
| 15        | 7.090       | 607           | 613         | 617          | rBV2     | 187412         | 339342        | 5.66%           | 0.324%        |
| 16        | 7.161       | 617           | 625         | 630          | rVB3     | 475179         | 1140399       | 19.02%          | 1.090%        |
| 17        | 7.219       | 630           | 635         | 636          | rBV      | 428272         | 552764        | 9.22%           | 0.528%        |
| 18        | 7.255       | 636           | 641         | 646          | rVB2     | 1220196        | 2159965       | 36.03%          | 2.065%        |
| 19        | 7.313       | 646           | 651         | 656          | rVB      | 1098653        | 1572267       | 26.22%          | 1.503%        |
| 20        | 7.360       | 656           | 659         | 663          | rBV2     | 221819         | 322913        | 5.39%           | 0.309%        |
| 21        | 7.419       | 663           | 669         | 676          | rBV3     | 2014397        | 3717976       | 62.01%          | 3.554%        |
| 22        | 7.478       | 676           | 679         | 689          | rVB5     | 150700         | 340688        | 5.68%           | 0.326%        |
| 23        | 7.601       | 689           | 700         | 706          | rBV5     | 342840         | 1012159       | 16.88%          | 0.968%        |
| 24        | 7.666       | 706           | 711         | 715          | rBV2     | 341118         | 582265        | 9.71%           | 0.557%        |
| 25        | 7.730       | 715           | 722         | 724          | rBV5     | 169676         | 309469        | 5.16%           | 0.296%        |
| 26        | 7.766       | 724           | 728         | 732          | rBV      | 364613         | 569487        | 9.50%           | 0.544%        |
| 27        | 7.842       | 737           | 741         | 749          | rVB5     | 345687         | 672348        | 11.21%          | 0.643%        |
| 28        | 7.930       | 749           | 756         | 762          | rBV      | 3586600        | 5995628       | 100.00%         | 5.731%        |
| 29        | 8.007       | 765           | 769         | 774          | rVB3     | 234746         | 355689        | 5.93%           | 0.340%        |
| 30        | 8.053       | 774           | 777         | 779          | rBV      | 121068         | 170091        | 2.84%           | 0.163%        |
| 31        | 8.106       | 779           | 786         | 790          | rBV3     | 334634         | 618190        | 10.31%          | 0.591%        |
| 32        | 8.194       | 793           | 801         | 807          | rBV4     | 521517         | 1312072       | 21.88%          | 1.254%        |
| 33        | 8.259       | 807           | 812         | 820          | rVB      | 2272693        | 4038741       | 67.36%          | 3.861%        |
| 34        | 8.353       | 820           | 828         | 832          | rBV2     | 721102         | 1568820       | 26.17%          | 1.500%        |
| 35        | 8.406       | 832           | 837         | 844          | rVB2     | 1419272        | 2589570       | 43.19%          | 2.475%        |
| 36        | 8.471       | 844           | 848         | 853          | rBV2     | 800827         | 1258792       | 21.00%          | 1.203%        |

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Instrument :  
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 ClientSampleId :  
 BGLD4

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
 Title : SVOA CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 37 | 8.535  | 853  | 859  | 863  | rBV2 | 628192  | 1261729 | 21.04% | 1.206% |
| 38 | 8.582  | 863  | 867  | 873  | rVB2 | 999920  | 1647531 | 27.48% | 1.575% |
| 39 | 8.664  | 874  | 881  | 885  | rVB3 | 447119  | 958504  | 15.99% | 0.916% |
| 40 | 8.717  | 885  | 890  | 896  | rBV3 | 266459  | 577022  | 9.62%  | 0.552% |
| 41 | 8.782  | 896  | 901  | 902  | rBV  | 388428  | 519100  | 8.66%  | 0.496% |
| 42 | 8.817  | 903  | 907  | 913  | rVV  | 1087085 | 1970696 | 32.87% | 1.884% |
| 43 | 8.876  | 913  | 917  | 922  | rVB  | 1207303 | 1935948 | 32.29% | 1.851% |
| 44 | 8.946  | 922  | 929  | 941  | rVB3 | 1720238 | 4630971 | 77.24% | 4.427% |
| 45 | 9.058  | 941  | 948  | 952  | rBV  | 1377694 | 2362513 | 39.40% | 2.258% |
| 46 | 9.099  | 952  | 955  | 958  | rVB  | 425583  | 538685  | 8.98%  | 0.515% |
| 47 | 9.140  | 958  | 962  | 967  | rBV2 | 412576  | 676811  | 11.29% | 0.647% |
| 48 | 9.246  | 976  | 980  | 985  | rBV  | 477910  | 742508  | 12.38% | 0.710% |
| 49 | 9.299  | 985  | 989  | 997  | rVB  | 781707  | 1514267 | 25.26% | 1.448% |
| 50 | 9.405  | 997  | 1007 | 1013 | rBV2 | 1565088 | 3631530 | 60.57% | 3.471% |
| 51 | 9.499  | 1018 | 1023 | 1030 | rVB5 | 664139  | 1296517 | 21.62% | 1.239% |
| 52 | 9.563  | 1031 | 1034 | 1038 | rVB4 | 206083  | 315171  | 5.26%  | 0.301% |
| 53 | 9.640  | 1038 | 1047 | 1048 | rBV5 | 513545  | 1278566 | 21.32% | 1.222% |
| 54 | 9.781  | 1059 | 1071 | 1079 | rVB5 | 652272  | 2446831 | 40.81% | 2.339% |
| 55 | 9.880  | 1079 | 1088 | 1092 | rVV6 | 430410  | 1030260 | 17.18% | 0.985% |
| 56 | 9.939  | 1092 | 1098 | 1102 | rVV2 | 914756  | 1801784 | 30.05% | 1.722% |
| 57 | 9.992  | 1102 | 1107 | 1112 | rVV  | 921052  | 1711775 | 28.55% | 1.636% |
| 58 | 10.039 | 1112 | 1115 | 1125 | rVB2 | 369094  | 905674  | 15.11% | 0.866% |
| 59 | 10.186 | 1132 | 1140 | 1144 | rBV6 | 517778  | 1305502 | 21.77% | 1.248% |
| 60 | 10.233 | 1144 | 1148 | 1153 | rVV2 | 592441  | 1006810 | 16.79% | 0.962% |
| 61 | 10.298 | 1153 | 1159 | 1164 | rVV3 | 492777  | 1023814 | 17.08% | 0.979% |
| 62 | 10.374 | 1164 | 1172 | 1178 | rVB3 | 573342  | 1572772 | 26.23% | 1.503% |
| 63 | 10.444 | 1178 | 1184 | 1189 | rBV2 | 484108  | 818893  | 13.66% | 0.783% |
| 64 | 10.521 | 1190 | 1197 | 1202 | rBV3 | 978531  | 1925390 | 32.11% | 1.840% |
| 65 | 10.574 | 1202 | 1206 | 1211 | rVB2 | 327062  | 558973  | 9.32%  | 0.534% |
| 66 | 10.632 | 1211 | 1216 | 1219 | rBV  | 493160  | 743657  | 12.40% | 0.711% |
| 67 | 10.691 | 1223 | 1226 | 1229 | rVB2 | 128185  | 170133  | 2.84%  | 0.163% |
| 68 | 10.809 | 1241 | 1246 | 1256 | rVB2 | 221278  | 505187  | 8.43%  | 0.483% |
| 69 | 10.973 | 1270 | 1274 | 1276 | rVV2 | 128331  | 211378  | 3.53%  | 0.202% |
| 70 | 11.008 | 1276 | 1280 | 1287 | rVV4 | 241159  | 591465  | 9.86%  | 0.565% |
| 71 | 11.096 | 1287 | 1295 | 1301 | rVV5 | 268926  | 1001176 | 16.70% | 0.957% |
| 72 | 11.173 | 1302 | 1308 | 1314 | rVV2 | 656272  | 1508133 | 25.15% | 1.442% |
| 73 | 11.267 | 1314 | 1324 | 1330 | rVV2 | 680795  | 1594202 | 26.59% | 1.524% |
| 74 | 11.331 | 1330 | 1335 | 1342 | rVV2 | 216545  | 410757  | 6.85%  | 0.393% |
| 75 | 11.390 | 1342 | 1345 | 1355 | rVB3 | 115264  | 273038  | 4.55%  | 0.261% |
| 76 | 11.490 | 1356 | 1362 | 1367 | rBV6 | 101643  | 219344  | 3.66%  | 0.210% |

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 Sample : M5215-02  
 Misc :  
 ALS Vial : 38 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BGLD4

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
 Title : SVOA CALIBRATION

|     |        |      |      |      |      |        |        |        |        |
|-----|--------|------|------|------|------|--------|--------|--------|--------|
| 77  | 11.819 | 1413 | 1418 | 1423 | rBV4 | 203648 | 420658 | 7.02%  | 0.402% |
| 78  | 11.866 | 1423 | 1426 | 1433 | rVB3 | 121158 | 219057 | 3.65%  | 0.209% |
| 79  | 11.942 | 1433 | 1439 | 1446 | rVB3 | 192022 | 322564 | 5.38%  | 0.308% |
| 80  | 12.013 | 1446 | 1451 | 1461 | rBV2 | 246350 | 475949 | 7.94%  | 0.455% |
| 81  | 12.119 | 1463 | 1469 | 1474 | rBV  | 492138 | 852593 | 14.22% | 0.815% |
| 82  | 12.213 | 1481 | 1485 | 1490 | rVB5 | 98540  | 172931 | 2.88%  | 0.165% |
| 83  | 12.724 | 1566 | 1572 | 1574 | rVV2 | 130281 | 246908 | 4.12%  | 0.236% |
| 84  | 12.759 | 1574 | 1578 | 1587 | rVB  | 296243 | 486009 | 8.11%  | 0.465% |
| 85  | 12.976 | 1610 | 1615 | 1621 | rVB  | 187320 | 318293 | 5.31%  | 0.304% |
| 86  | 13.129 | 1636 | 1641 | 1645 | rBV2 | 111443 | 194050 | 3.24%  | 0.185% |
| 87  | 13.199 | 1650 | 1653 | 1656 | rVV2 | 109865 | 177370 | 2.96%  | 0.170% |
| 88  | 13.305 | 1667 | 1671 | 1677 | rVV  | 135347 | 211732 | 3.53%  | 0.202% |
| 89  | 13.446 | 1690 | 1695 | 1701 | rVB  | 224181 | 329582 | 5.50%  | 0.315% |
| 90  | 14.075 | 1797 | 1802 | 1813 | rVB2 | 62406  | 175588 | 2.93%  | 0.168% |
| 91  | 14.298 | 1834 | 1840 | 1845 | rVB  | 317889 | 504438 | 8.41%  | 0.482% |
| 92  | 14.603 | 1886 | 1892 | 1901 | rVB  | 316385 | 525826 | 8.77%  | 0.503% |
| 93  | 14.909 | 1934 | 1944 | 1951 | rBV  | 243374 | 414529 | 6.91%  | 0.396% |
| 94  | 15.079 | 1967 | 1973 | 1978 | rBV2 | 119314 | 200017 | 3.34%  | 0.191% |
| 95  | 15.902 | 2107 | 2113 | 2119 | rVB  | 401832 | 621638 | 10.37% | 0.594% |
| 96  | 17.658 | 2406 | 2412 | 2418 | rBV  | 278983 | 442535 | 7.38%  | 0.423% |
| 97  | 17.758 | 2423 | 2429 | 2435 | rVB2 | 368881 | 595178 | 9.93%  | 0.569% |
| 98  | 20.031 | 2812 | 2816 | 2822 | rVB  | 431069 | 622336 | 10.38% | 0.595% |
| 99  | 21.841 | 3120 | 3124 | 3130 | rBV  | 165568 | 301247 | 5.02%  | 0.288% |
| 100 | 21.976 | 3144 | 3147 | 3155 | rVB2 | 263375 | 442143 | 7.37%  | 0.423% |

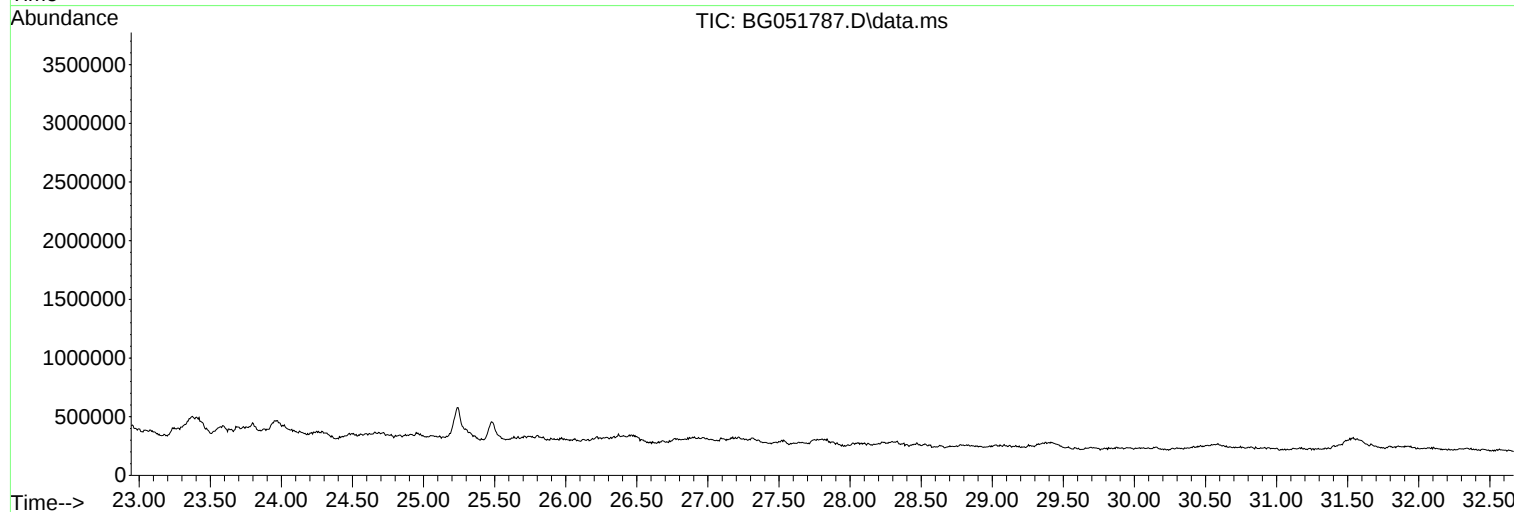
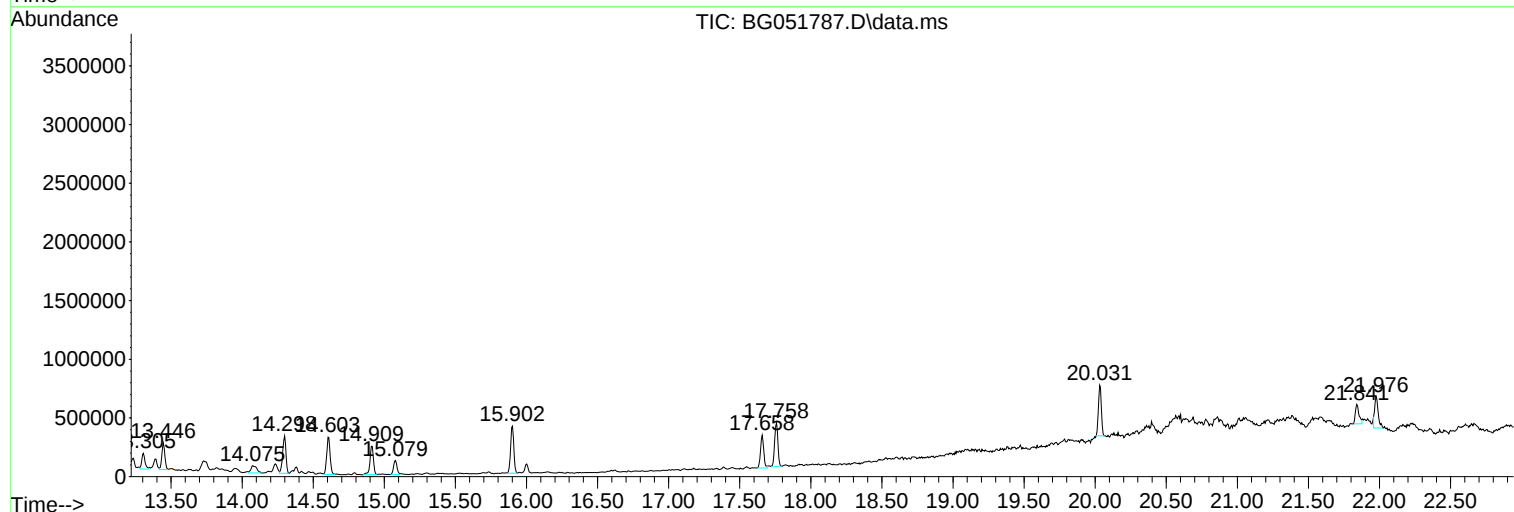
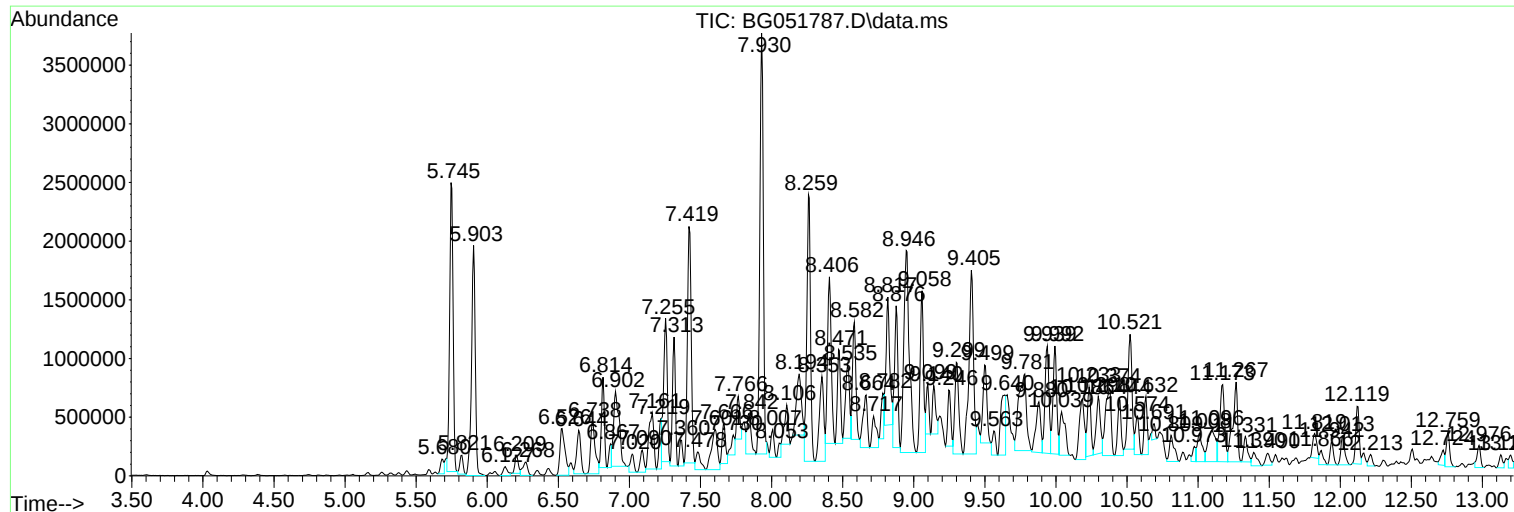
Sum of corrected areas: 104612524

Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
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Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
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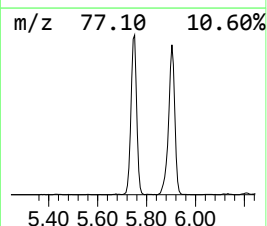
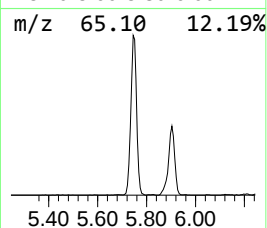
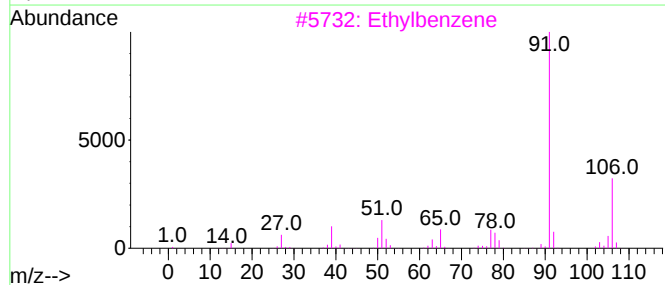
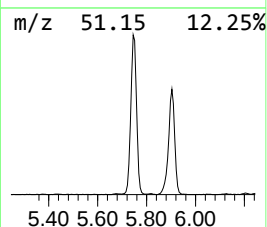
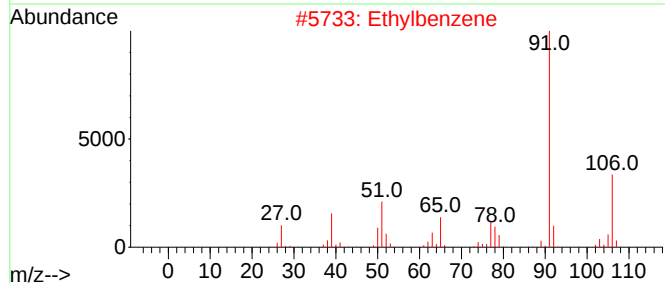
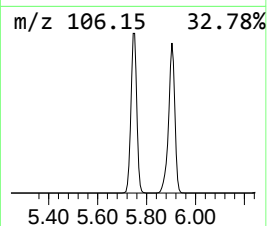
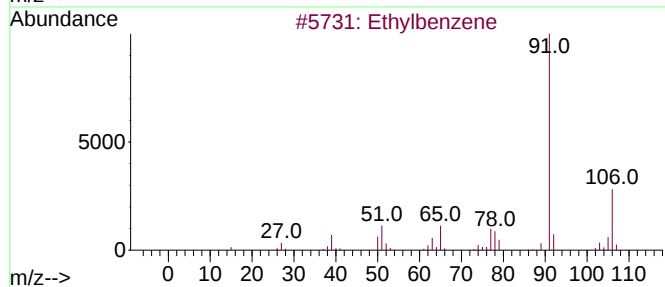
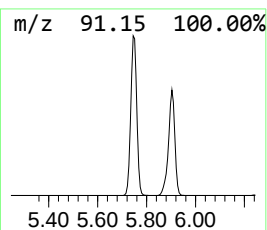
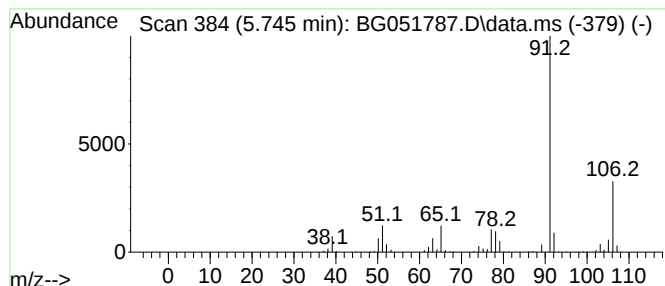
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 Ethylbenzene Concentration Rank 7

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 5.745 | 20.34 ng/ul | 4107870 | 1,4-Dichlorobenzene-d4 | 8.277 |

| Hit# | of           | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|--------------|---|--------------|-----|---------|-------------|------|
| 1    | Ethylbenzene |   |              | 106 | C8H10   | 000100-41-4 | 95   |
| 2    | Ethylbenzene |   |              | 106 | C8H10   | 000100-41-4 | 91   |
| 3    | Ethylbenzene |   |              | 106 | C8H10   | 000100-41-4 | 91   |
| 4    | Ethylbenzene |   |              | 106 | C8H10   | 000100-41-4 | 91   |
| 5    | Ethylbenzene |   |              | 106 | C8H10   | 000100-41-4 | 91   |



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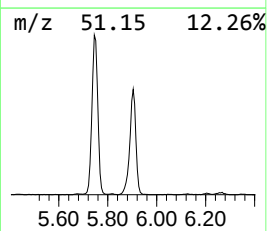
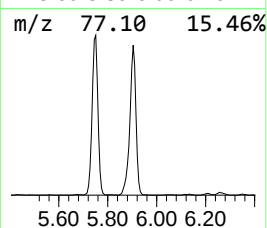
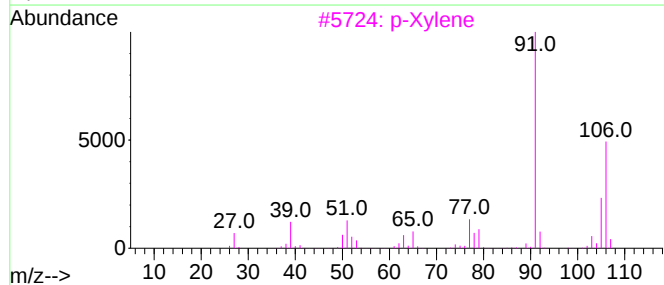
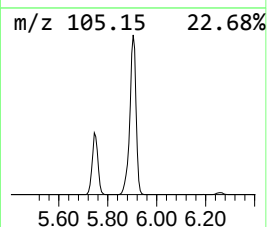
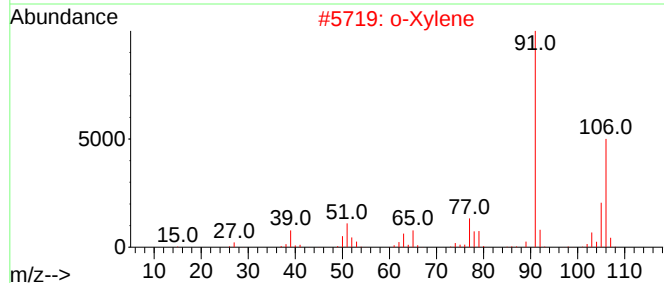
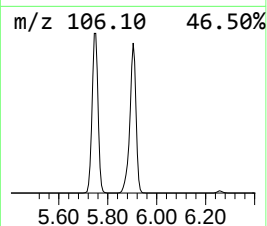
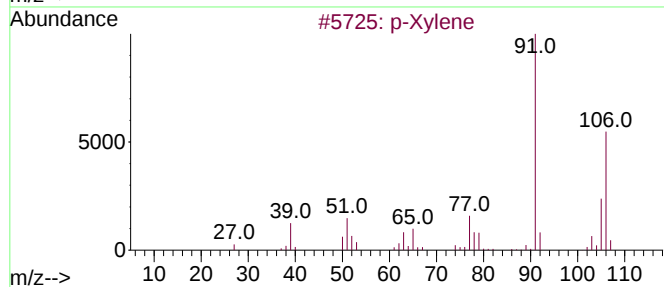
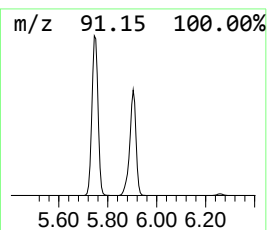
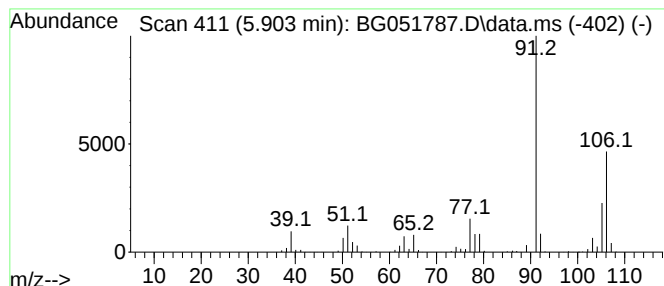
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 p-Xylene Concentration Rank 10

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 5.903 | 17.56 ng/ul | 3546740 | 1,4-Dichlorobenzene-d4 | 8.277 |

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



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

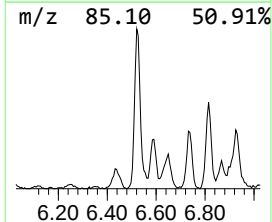
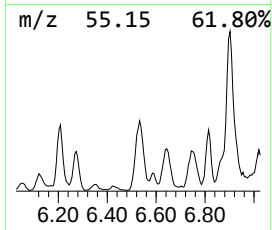
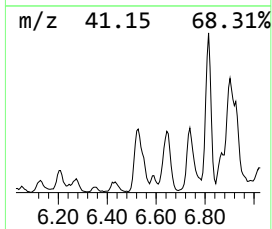
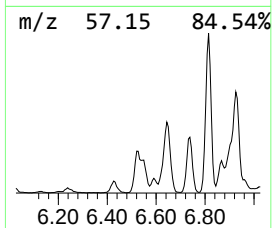
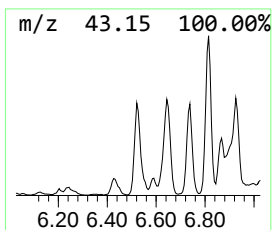
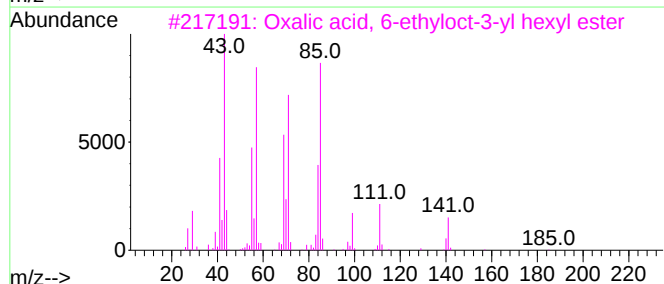
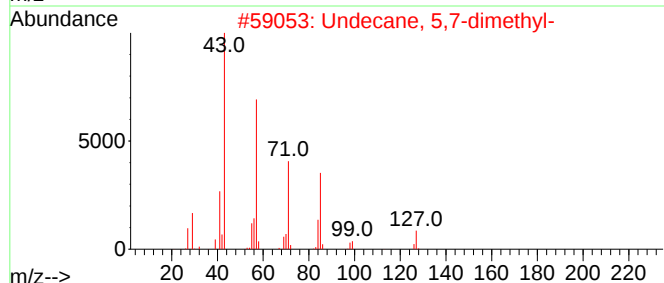
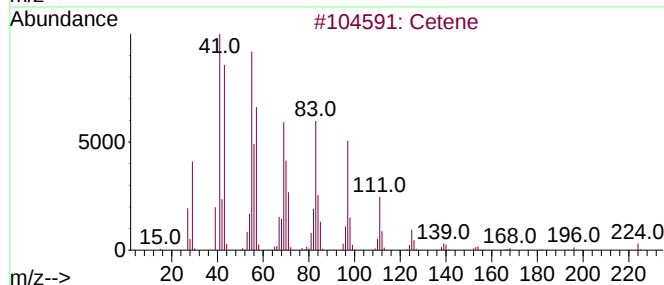
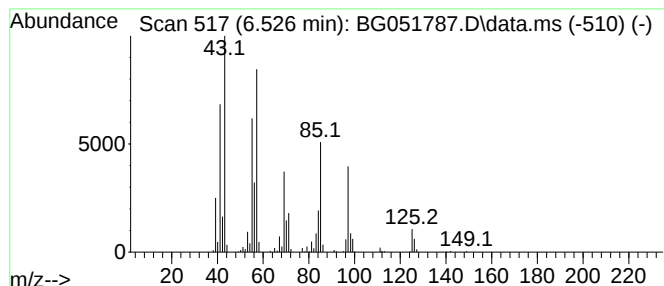
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 44

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 6.526 | 4.96 ng/ul | 1001610 | 1,4-Dichlorobenzene-d4 | 8.277 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Cetene                              | 224 | C16H32   | 000629-73-2  | 47   |
| 2    |    |   | Undecane, 5,7-dimethyl-             | 184 | C13H28   | 017312-83-3  | 43   |
| 3    |    |   | Oxalic acid, 6-ethyloct-3-yl hex... | 314 | C18H34O4 | 1000309-34-4 | 43   |
| 4    |    |   | Hexane, 2,3,3-trimethyl-            | 128 | C9H20    | 016747-28-7  | 38   |
| 5    |    |   | Cyclohexane, 1,3-dimethyl-, trans-  | 112 | C8H16    | 002207-03-6  | 35   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

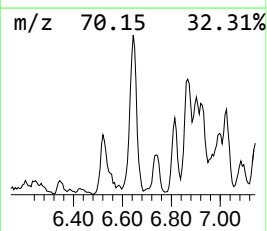
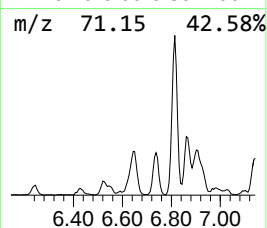
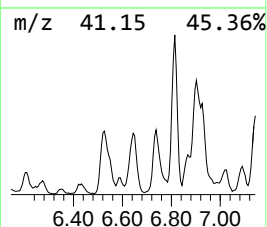
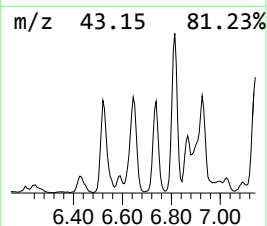
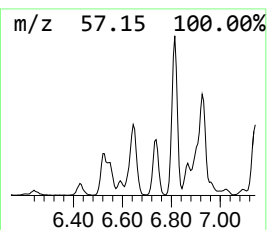
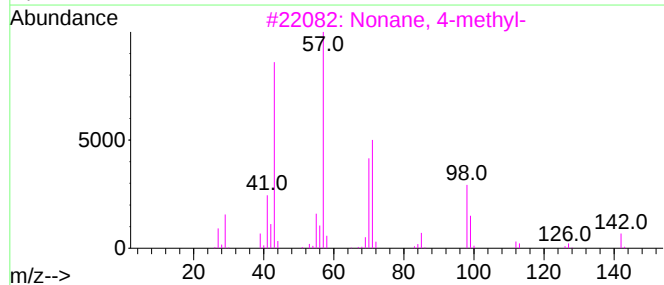
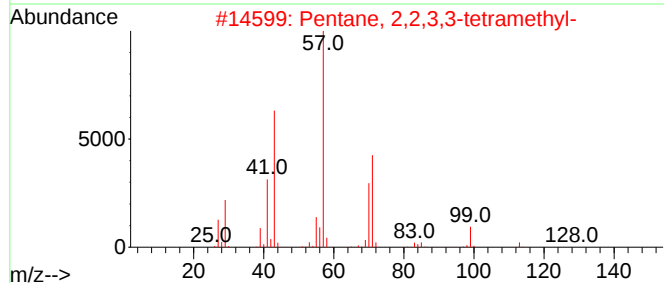
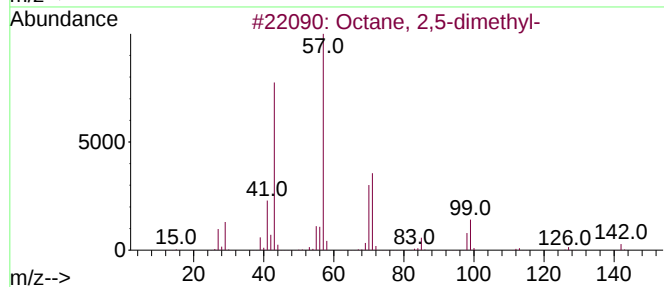
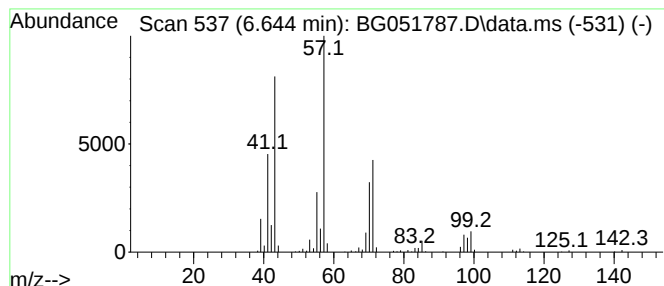
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 49

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 6.644 | 3.68 ng/ul | 742963 | 1,4-Dichlorobenzene-d4 | 8.277 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | Octane, 2,5-dimethyl-         | 142 | C10H22  | 015869-89-3 | 78   |
| 2    |    |   | Pentane, 2,2,3,3-tetramethyl- | 128 | C9H20   | 007154-79-2 | 72   |
| 3    |    |   | Nonane, 4-methyl-             | 142 | C10H22  | 017301-94-9 | 64   |
| 4    |    |   | Undecane, 4,5-dimethyl-       | 184 | C13H28  | 017312-79-7 | 64   |
| 5    |    |   | Hexane, 3,3-dimethyl-         | 114 | C8H18   | 000563-16-6 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

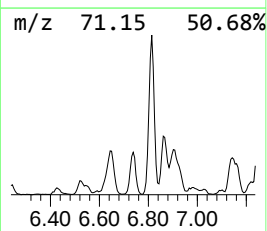
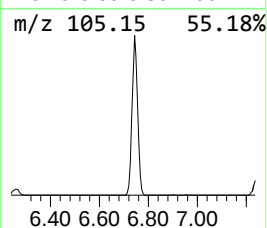
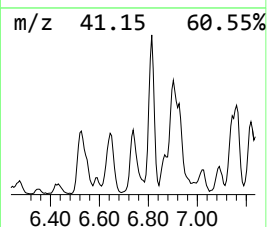
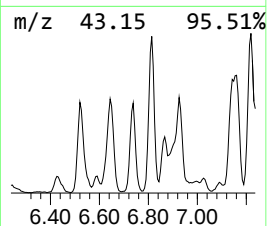
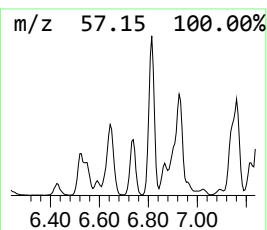
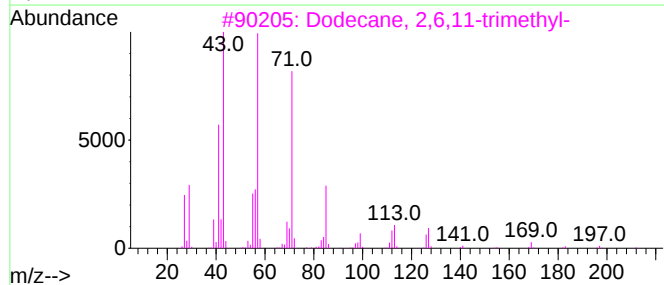
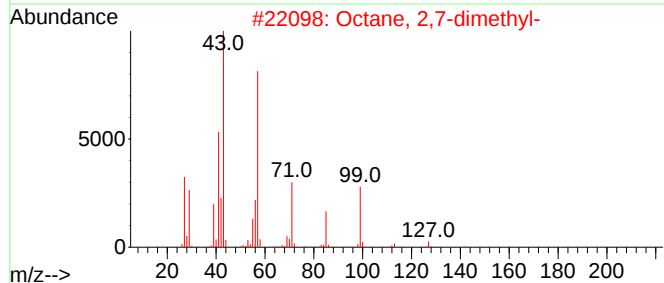
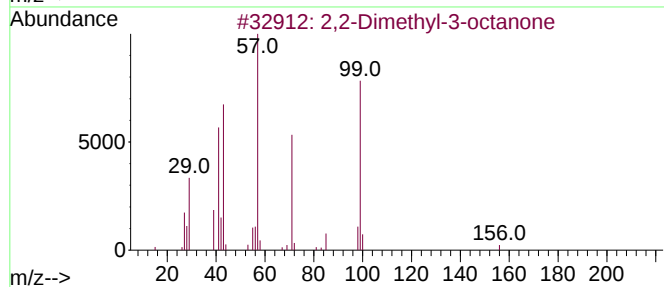
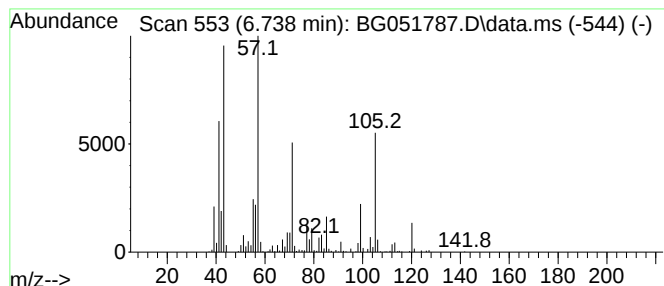
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 5 2,2-Dimethyl-3-octanone Concentration Rank 43

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 6.738 | 4.98 ng/ul | 1004660 | 1,4-Dichlorobenzene-d4 | 8.277 |

| Hit# | of | 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------------|-----|---------|-------------|------|
| 1    |    |   | 2,2-Dimethyl-3-octanone     | 156 | C10H20O | 005340-64-7 | 59   |
| 2    |    |   | Octane, 2,7-dimethyl-       | 142 | C10H22  | 001072-16-8 | 52   |
| 3    |    |   | Dodecane, 2,6,11-trimethyl- | 212 | C15H32  | 031295-56-4 | 50   |
| 4    |    |   | Octane, 2,7-dimethyl-       | 142 | C10H22  | 001072-16-8 | 50   |
| 5    |    |   | Dodecane, 2,7,10-trimethyl- | 212 | C15H32  | 074645-98-0 | 50   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

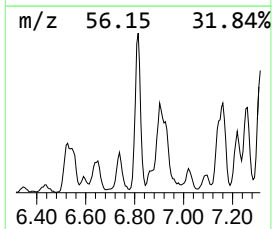
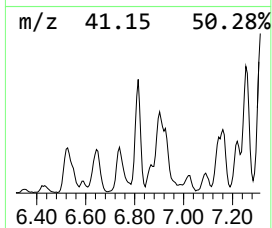
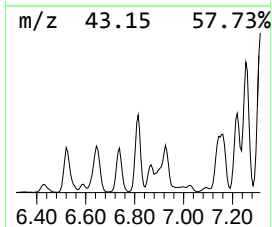
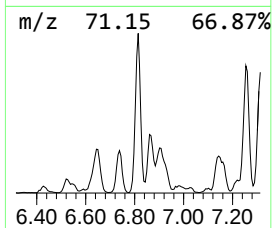
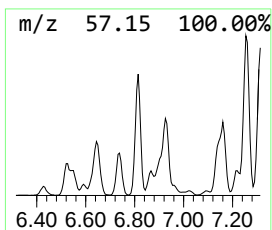
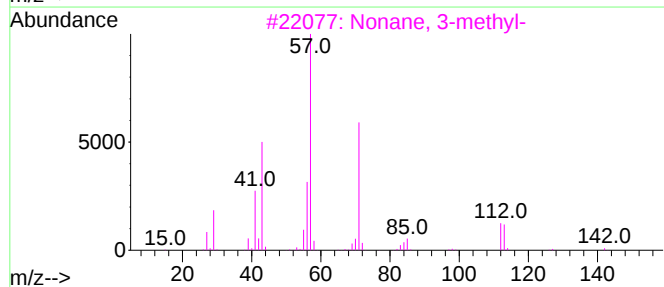
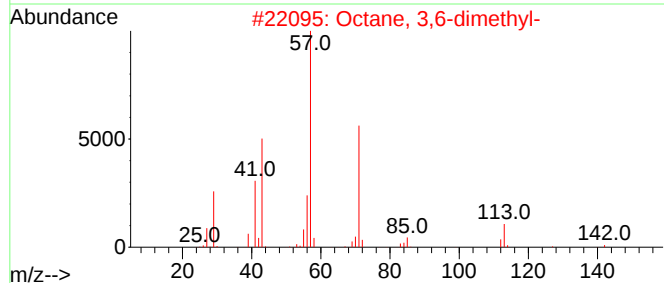
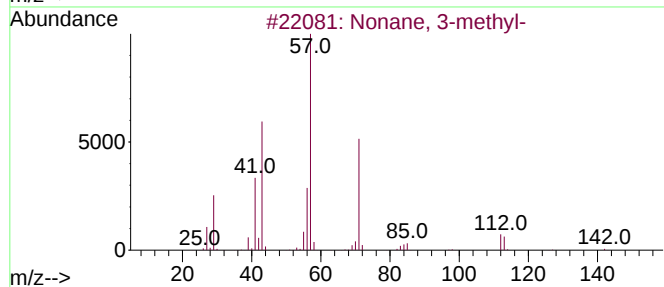
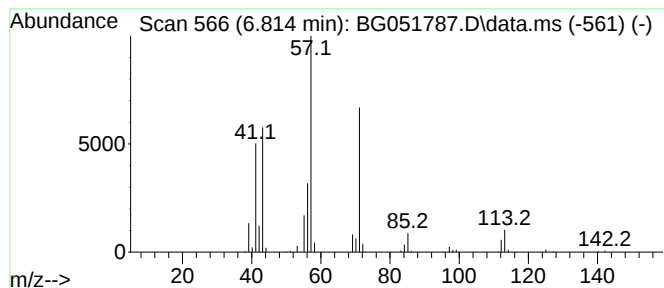
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TIC Integration Parameters: LSCINT.P

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Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 42

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 6.814 | 5.42 ng/ul | 1095270 | 1,4-Dichlorobenzene-d4 | 8.277 |

| Hit# | of | 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|---|-----------------------|-----|---------|-------------|------|
| 1    |    |   | Nonane, 3-methyl-     | 142 | C10H22  | 005911-04-6 | 87   |
| 2    |    |   | Octane, 3,6-dimethyl- | 142 | C10H22  | 015869-94-0 | 86   |
| 3    |    |   | Nonane, 3-methyl-     | 142 | C10H22  | 005911-04-6 | 86   |
| 4    |    |   | Octane, 3,6-dimethyl- | 142 | C10H22  | 015869-94-0 | 83   |
| 5    |    |   | Octane, 2,6-dimethyl- | 142 | C10H22  | 002051-30-1 | 81   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

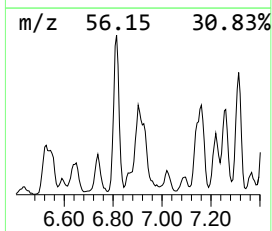
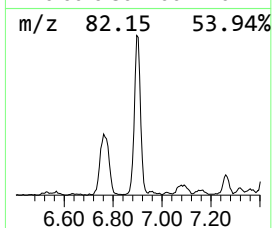
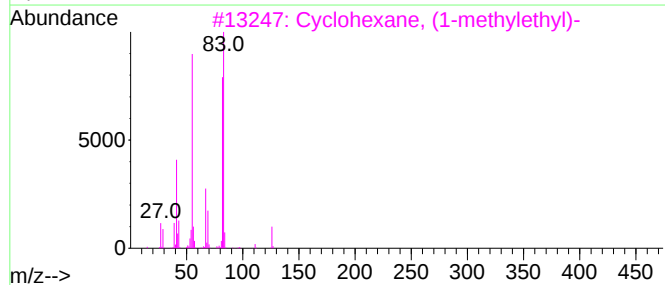
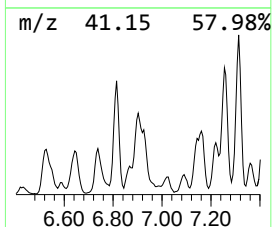
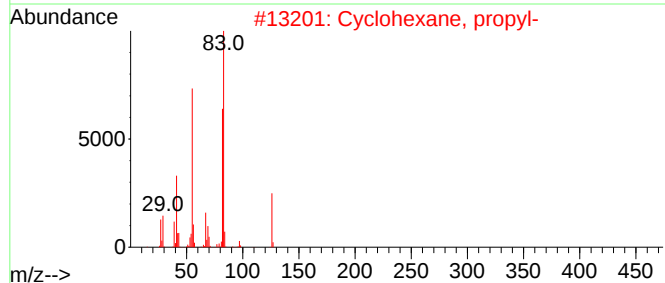
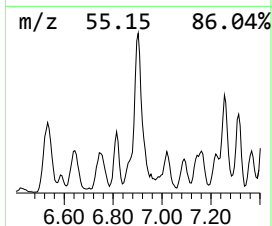
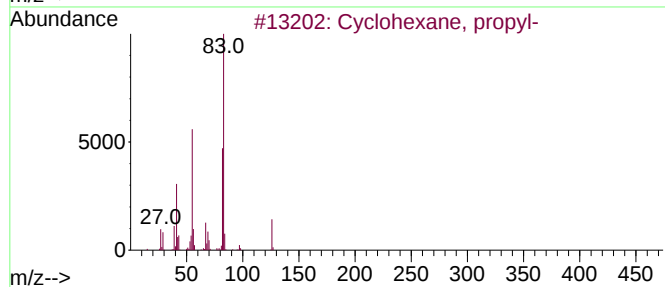
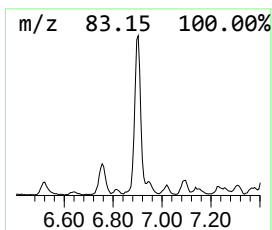
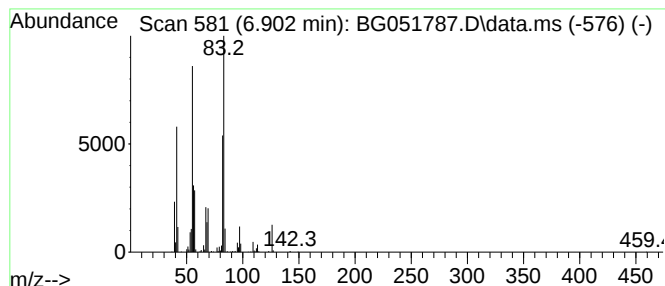
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 (DEL) Alkane: Cyclic6.902 Concentration Rank 26

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 6.902 | 8.65 ng/ul | 1747160 | 1,4-Dichlorobenzene-d4 | 8.277 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, propyl-          | 126 | C9H18   | 001678-92-8 | 83   |
| 2    |    |   | Cyclohexane, propyl-          | 126 | C9H18   | 001678-92-8 | 80   |
| 3    |    |   | Cyclohexane, (1-methylethyl)- | 126 | C9H18   | 000696-29-7 | 72   |
| 4    |    |   | Cyclohexane, propyl-          | 126 | C9H18   | 001678-92-8 | 72   |
| 5    |    |   | Cyclohexane, propyl-          | 126 | C9H18   | 001678-92-8 | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

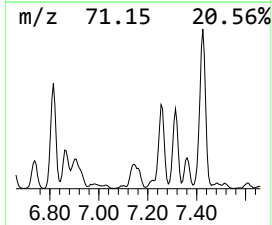
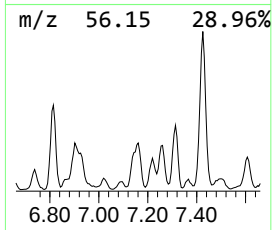
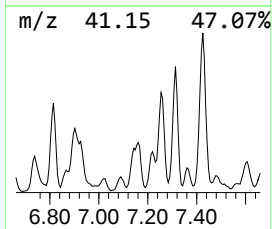
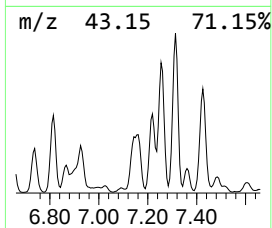
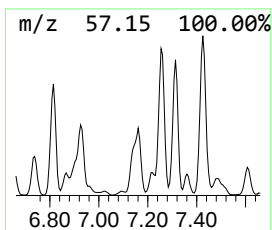
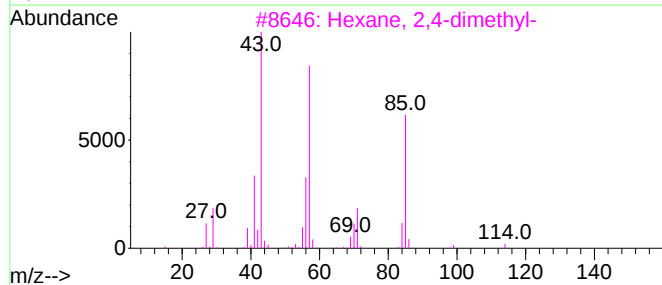
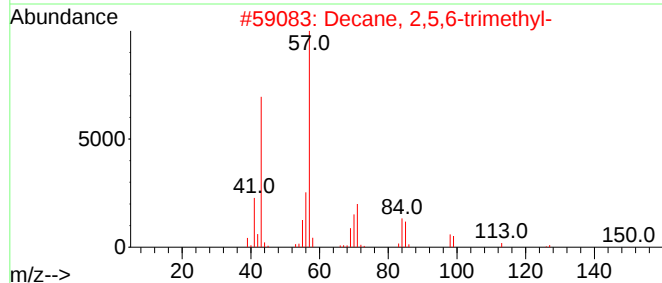
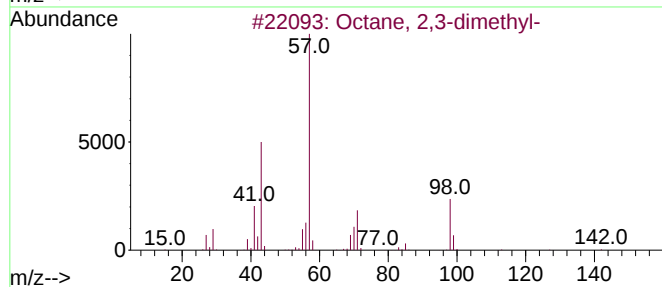
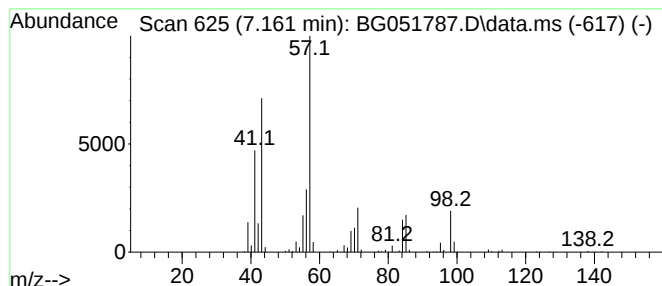
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 40

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 7.161 | 5.65 ng/ul | 1140400 | 1,4-Dichlorobenzene-d4 | 8.277 |

| Hit# | of                         | 5 | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|----------------------------|---|--------------|-----|---------|-------------|------|
| 1    | Octane, 2,3-dimethyl-      |   |              | 142 | C10H22  | 007146-60-3 | 59   |
| 2    | Decane, 2,5,6-trimethyl-   |   |              | 184 | C13H28  | 062108-23-0 | 59   |
| 3    | Hexane, 2,4-dimethyl-      |   |              | 114 | C8H18   | 000589-43-5 | 53   |
| 4    | Heptane, 3-ethyl-2-methyl- |   |              | 142 | C10H22  | 014676-29-0 | 53   |
| 5    | Decane, 5-methyl-          |   |              | 156 | C11H24  | 013151-35-4 | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

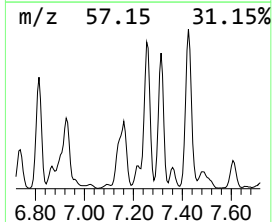
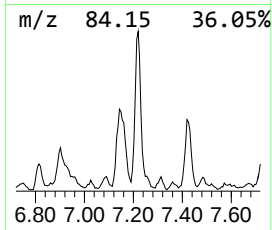
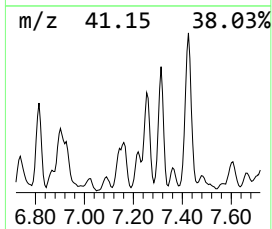
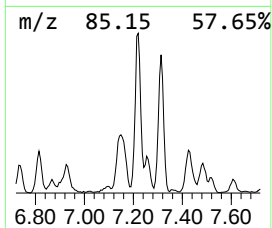
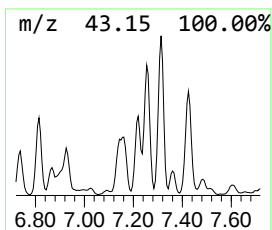
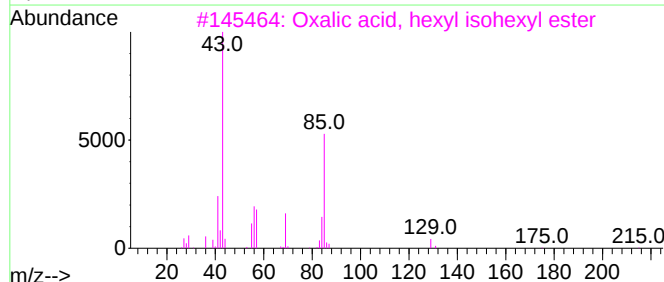
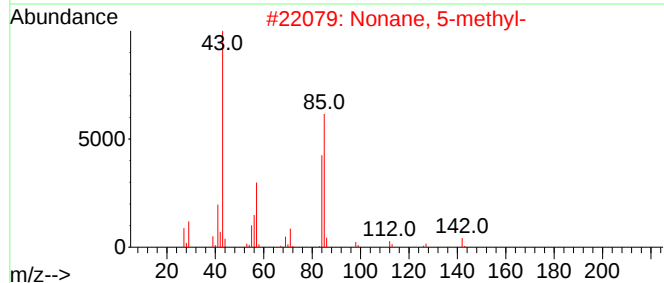
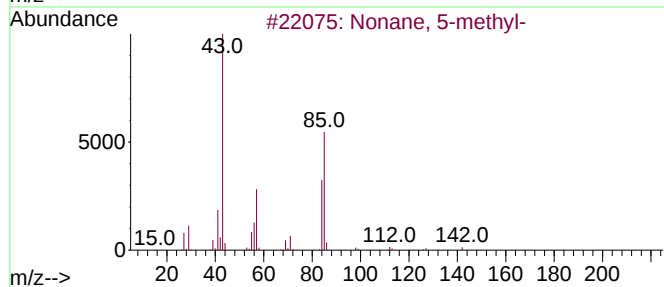
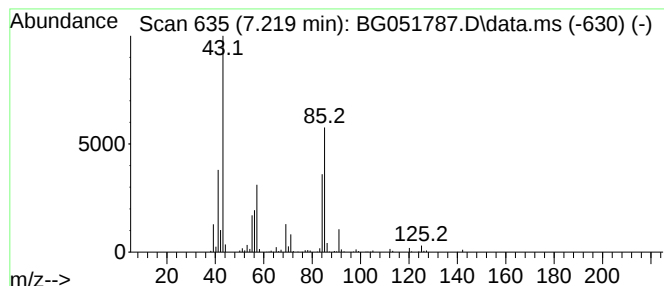
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TIC Integration Parameters: LSCINT.P

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Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 56

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 7.219 | 2.74 ng/ul | 552764 | 1,4-Dichlorobenzene-d4 | 8.277 |

| Hit# | of | 5 | Tentative ID                      | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-----------------------------------|-----|----------|--------------|------|
| 1    |    |   | Nonane, 5-methyl-                 | 142 | C10H22   | 015869-85-9  | 81   |
| 2    |    |   | Nonane, 5-methyl-                 | 142 | C10H22   | 015869-85-9  | 72   |
| 3    |    |   | Oxalic acid, hexyl isohexyl ester | 258 | C14H26O4 | 1010309-33-0 | 64   |
| 4    |    |   | Heptane, 3,4,5-trimethyl-         | 142 | C10H22   | 020278-89-1  | 64   |
| 5    |    |   | Hexane, 2,3,3-trimethyl-          | 128 | C9H20    | 016747-28-7  | 59   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

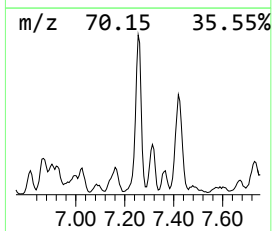
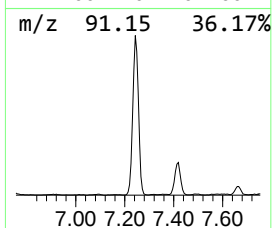
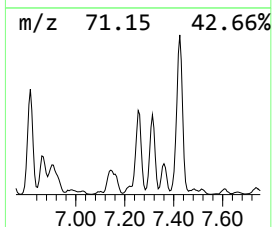
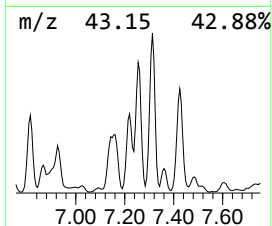
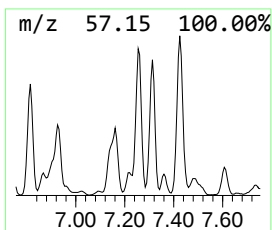
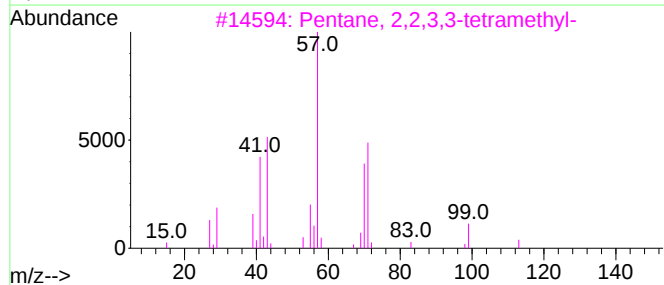
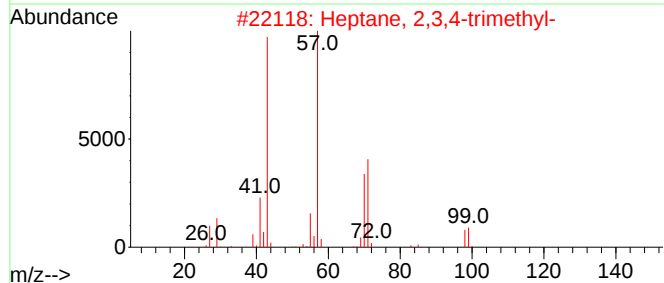
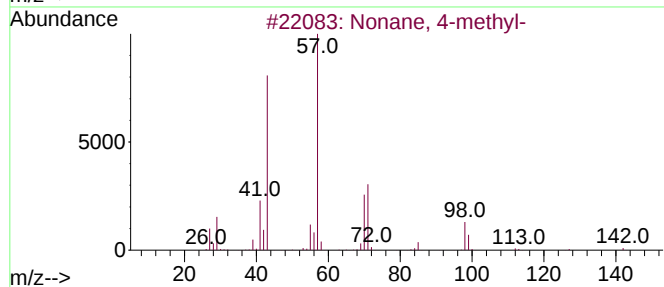
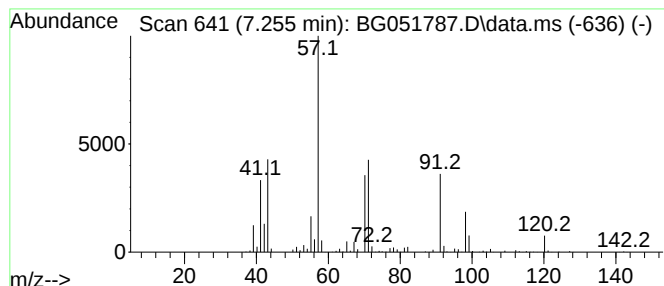
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 19

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 7.255 | 10.70 ng/ul | 2159970 | 1,4-Dichlorobenzene-d4 | 8.277 |

| Hit# | of 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------|-----|---------|-------------|------|
| 1    |      | Nonane, 4-methyl-             | 142 | C10H22  | 017301-94-9 | 80   |
| 2    |      | Heptane, 2,3,4-trimethyl-     | 142 | C10H22  | 052896-95-4 | 59   |
| 3    |      | Pentane, 2,2,3,3-tetramethyl- | 128 | C9H20   | 007154-79-2 | 45   |
| 4    |      | Undecane, 2,6-dimethyl-       | 184 | C13H28  | 017301-23-4 | 45   |
| 5    |      | Nonane, 4-methyl-             | 142 | C10H22  | 017301-94-9 | 45   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

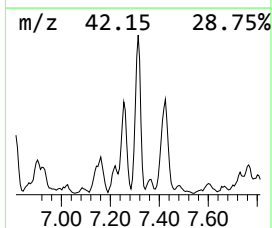
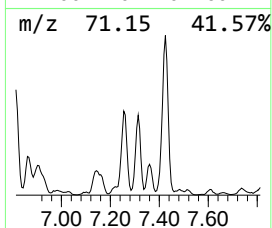
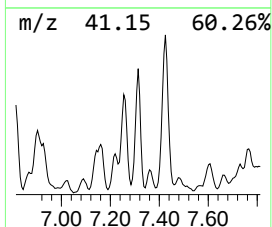
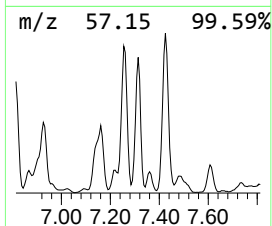
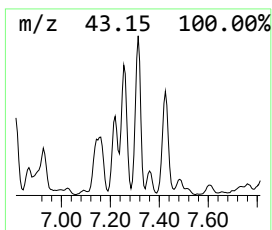
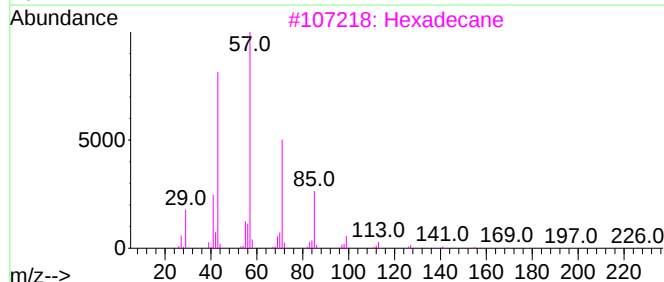
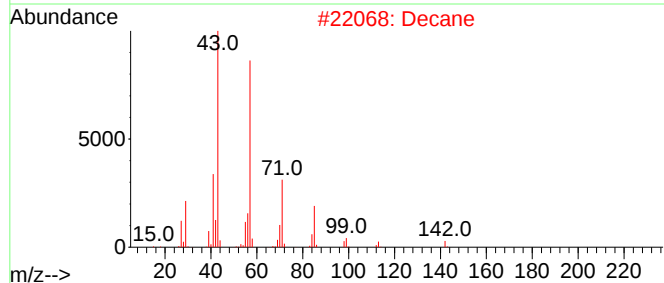
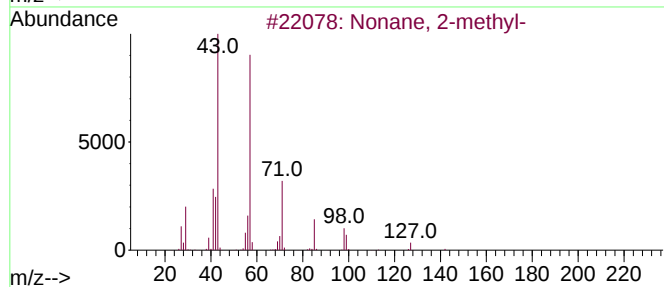
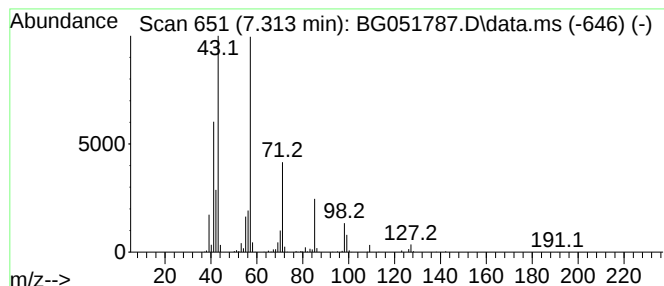
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 32

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 7.313 | 7.79 ng/ul | 1572270 | 1,4-Dichlorobenzene-d4 | 8.277 |

| Hit# | of 5 | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|------|-----------------------|-----|---------|-------------|------|
| 1    |      | Nonane, 2-methyl-     | 142 | C10H22  | 000871-83-0 | 90   |
| 2    |      | Decane                | 142 | C10H22  | 000124-18-5 | 72   |
| 3    |      | Hexadecane            | 226 | C16H34  | 000544-76-3 | 64   |
| 4    |      | Decane, 2,4-dimethyl- | 170 | C12H26  | 002801-84-5 | 53   |
| 5    |      | Undecane, 5-methyl-   | 170 | C12H26  | 001632-70-8 | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

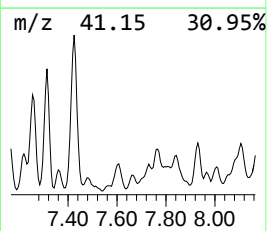
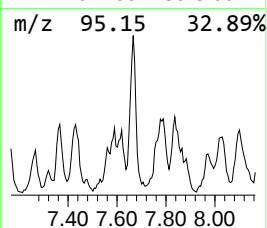
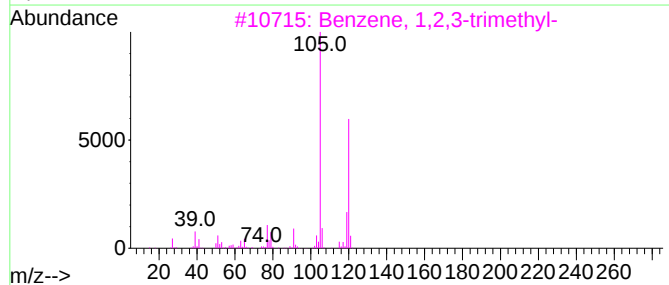
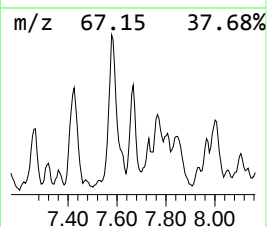
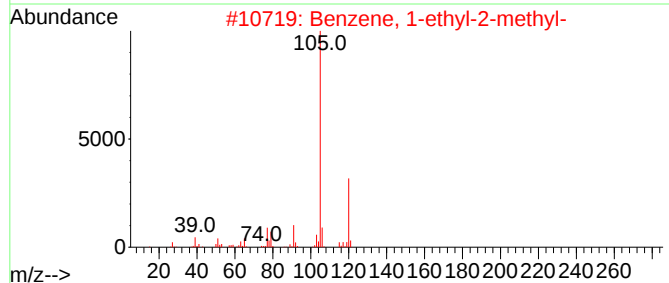
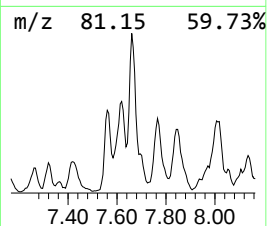
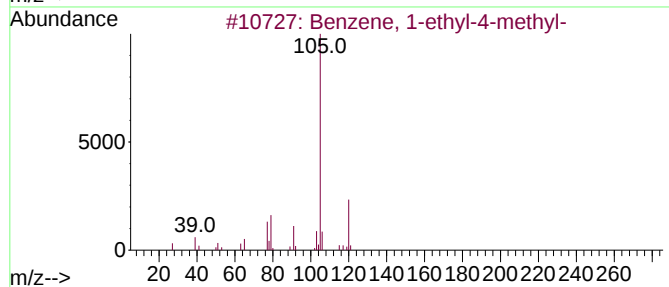
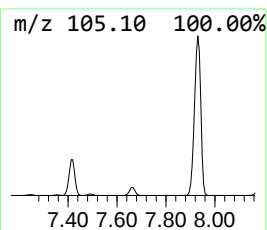
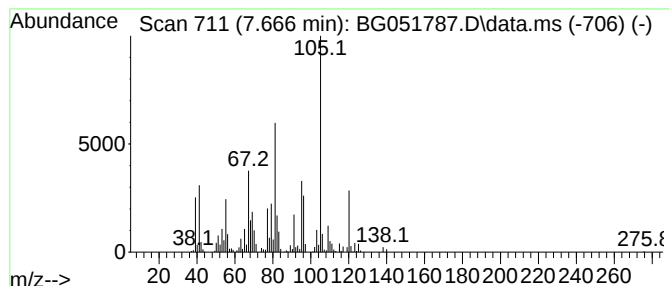
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 12 Benzene, 1-ethyl-4-methyl- Concentration Rank 54

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 7.666 | 2.88 ng/ul | 582265 | 1,4-Dichlorobenzene-d4 | 8.277 |

| Hit# | of | 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 70   |
| 2    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 55   |
| 3    |    |   | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 46   |
| 4    |    |   | Benzene, 1-ethyl-2-methyl- | 120 | C9H12   | 000611-14-3 | 46   |
| 5    |    |   | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 41   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

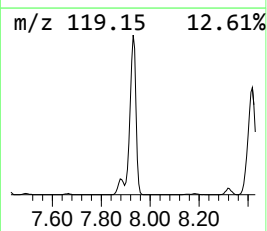
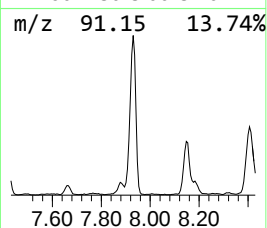
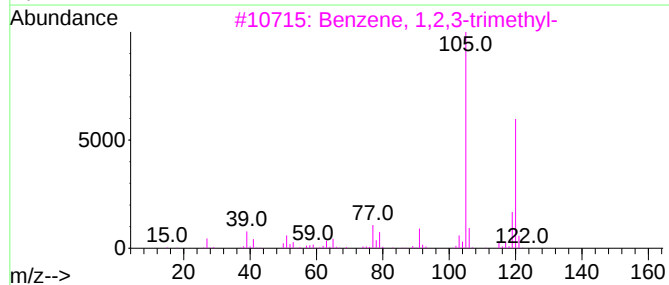
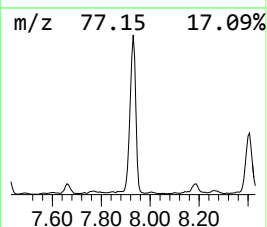
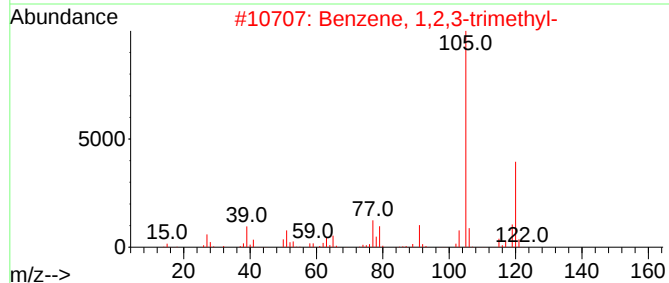
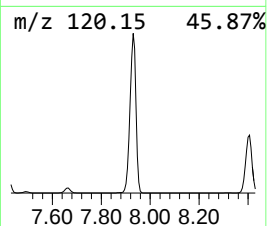
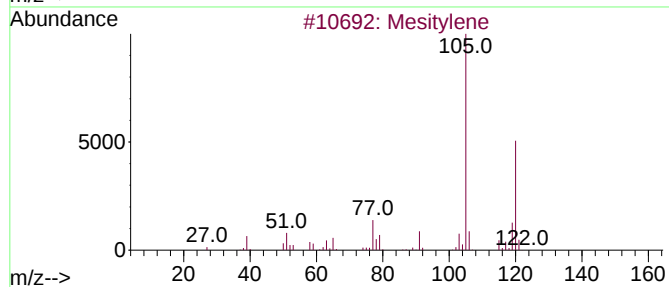
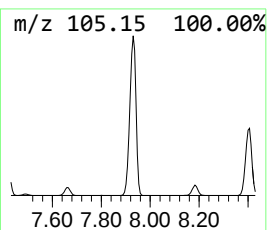
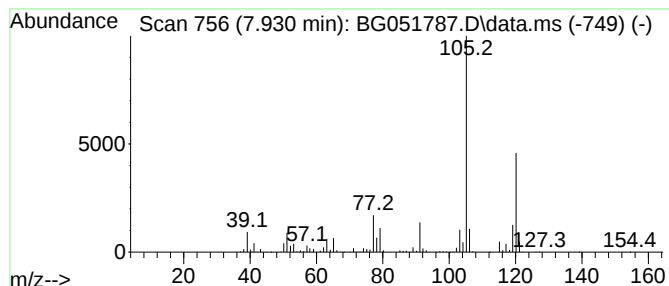
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 13 Mesitylene Concentration Rank 5

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 7.930 | 29.69 ng/ul | 5995630 | 1,4-Dichlorobenzene-d4 | 8.277 |

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



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

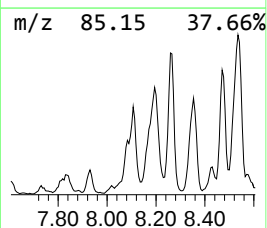
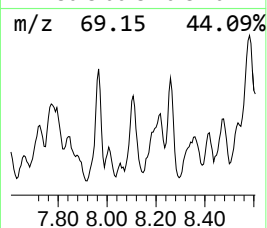
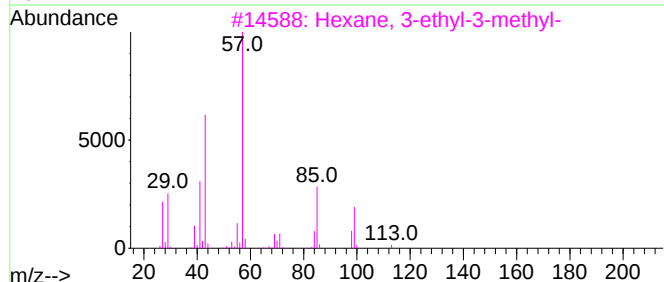
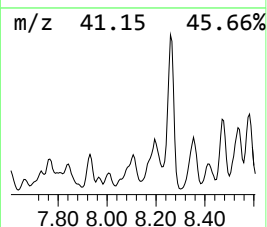
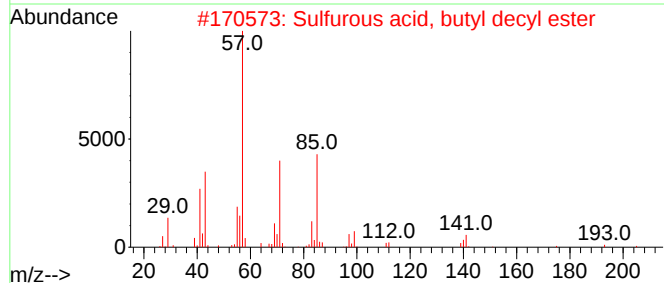
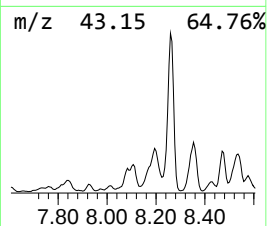
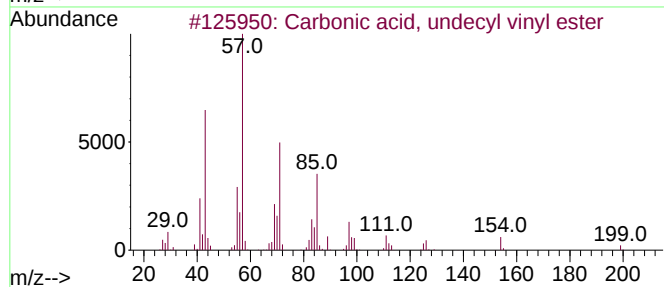
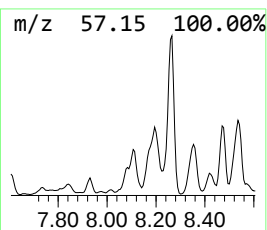
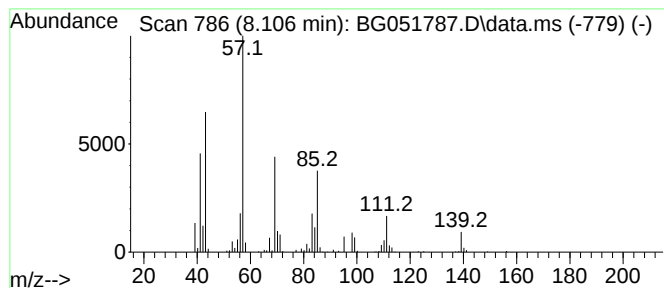
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 14 Carbonic acid, undecyl vinyl... Concentration Rank 53

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 8.106 | 3.06 ng/ul | 618190 | 1,4-Dichlorobenzene-d4 | 8.277 |

| Hit# | of | 5 | Tentative ID                       | MW  | MolForm   | CAS#         | Qual |
|------|----|---|------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Carbonic acid, undecyl vinyl ester | 242 | C14H26O3  | 1000382-54-9 | 59   |
| 2    |    |   | Sulfurous acid, butyl decyl ester  | 278 | C14H30O3S | 1000309-17-7 | 47   |
| 3    |    |   | Hexane, 3-ethyl-3-methyl-          | 128 | C9H20     | 003074-76-8  | 43   |
| 4    |    |   | Undecane, 4-ethyl-                 | 184 | C13H28    | 017312-59-3  | 43   |
| 5    |    |   | Tetradecane                        | 198 | C14H30    | 000629-59-4  | 43   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

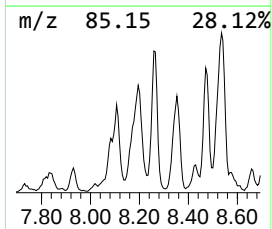
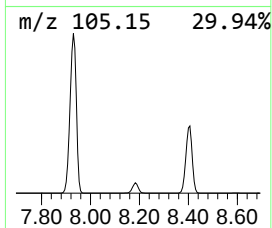
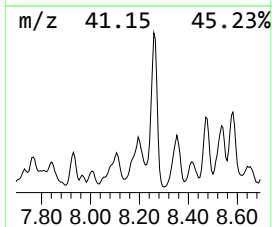
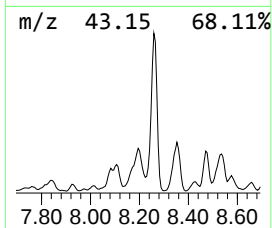
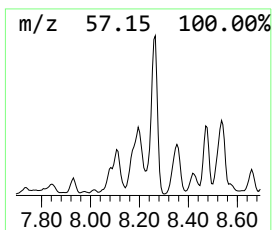
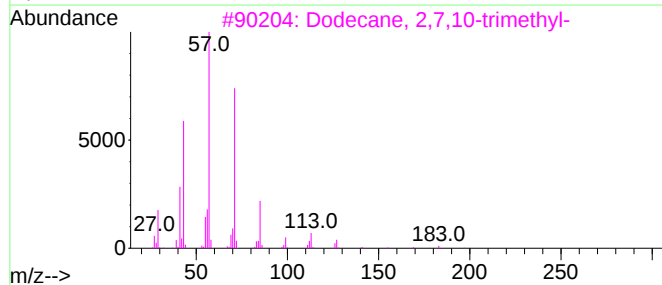
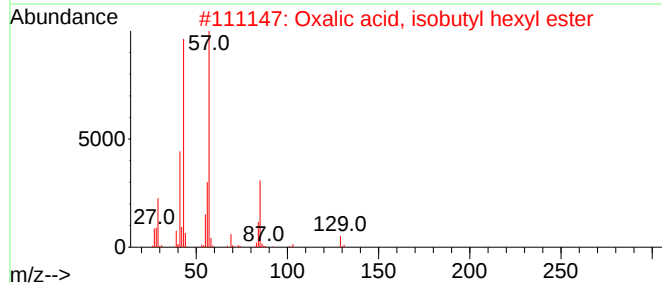
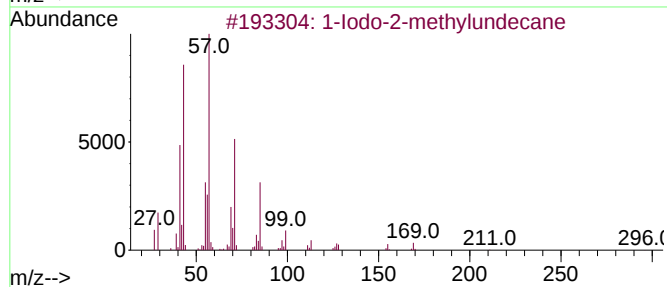
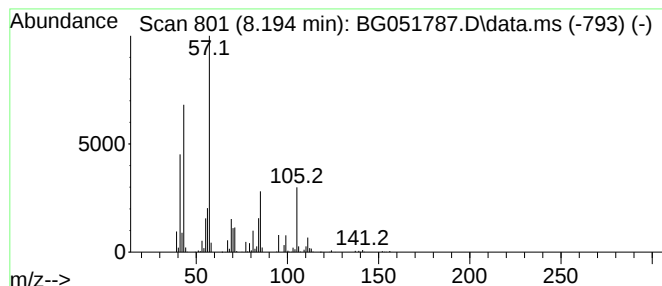
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 15 unknown-01 Concentration Rank 35

| R.T.      | EstConc                           | Area    | Relative to ISTD       | R.T.         |      |
|-----------|-----------------------------------|---------|------------------------|--------------|------|
| 8.194     | 6.50 ng/ul                        | 1312070 | 1,4-Dichlorobenzene-d4 | 8.277        |      |
| Hit# of 5 | Tentative ID                      | MW      | MolForm                | CAS#         | Qual |
| 1         | 1-Iodo-2-methylundecane           | 296     | C12H25I                | 073105-67-6  | 43   |
| 2         | Oxalic acid, isobutyl hexyl ester | 230     | C12H22O4               | 1000309-37-1 | 43   |
| 3         | Dodecane, 2,7,10-trimethyl-       | 212     | C15H32                 | 074645-98-0  | 43   |
| 4         | Hexane, 2,2,3,3-tetramethyl-      | 142     | C10H22                 | 013475-81-5  | 38   |
| 5         | Hexane, 2,4-dimethyl-             | 114     | C8H18                  | 000589-43-5  | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

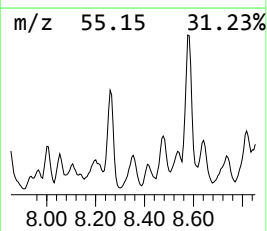
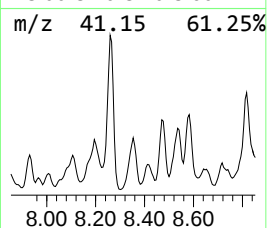
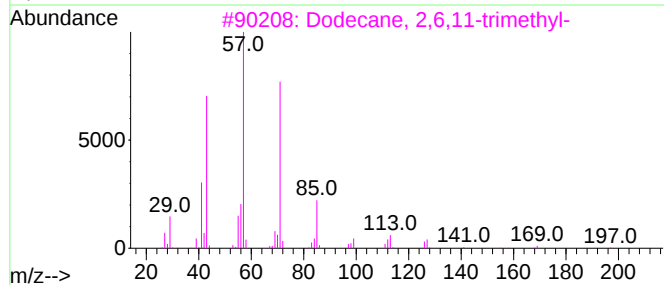
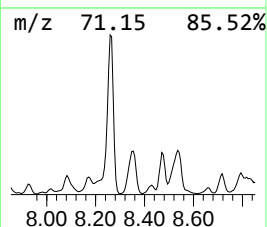
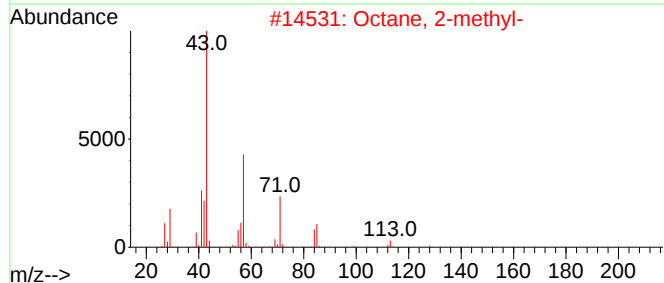
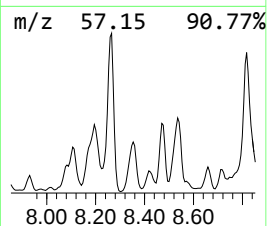
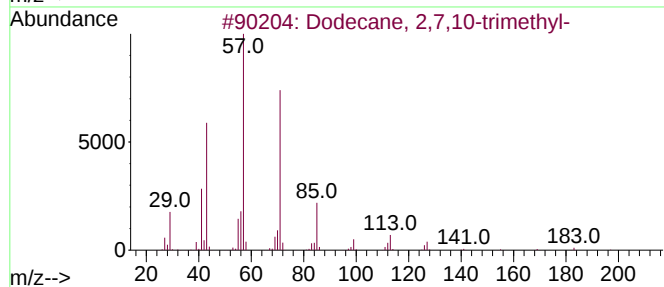
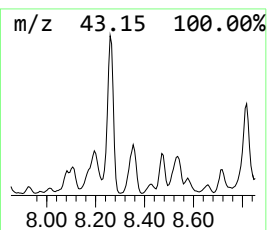
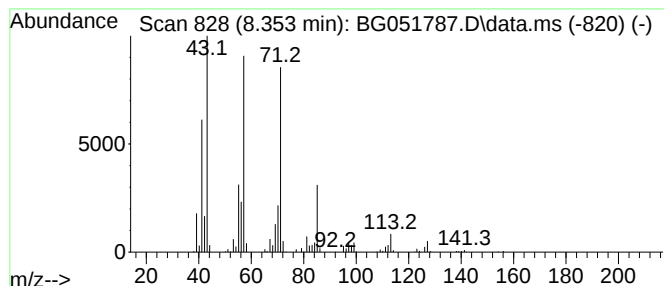
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TIC Integration Parameters: LSCINT.P

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Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 33

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 8.353 | 7.77 ng/ul | 1568820 | 1,4-Dichlorobenzene-d4 | 8.277 |

| Hit# | of 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|------|-----------------------------|-----|---------|-------------|------|
| 1    |      | Dodecane, 2,7,10-trimethyl- | 212 | C15H32  | 074645-98-0 | 86   |
| 2    |      | Octane, 2-methyl-           | 128 | C9H20   | 003221-61-2 | 81   |
| 3    |      | Dodecane, 2,6,11-trimethyl- | 212 | C15H32  | 031295-56-4 | 80   |
| 4    |      | Dodecane, 2,6,10-trimethyl- | 212 | C15H32  | 003891-98-3 | 72   |
| 5    |      | Nonane, 5-(2-methylpropyl)- | 184 | C13H28  | 062185-53-9 | 59   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

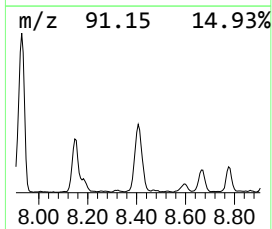
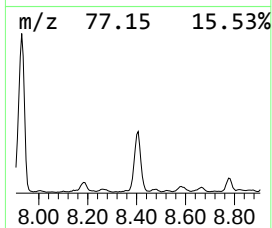
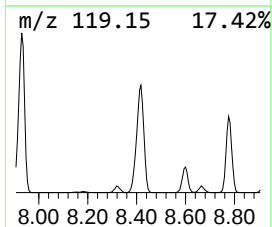
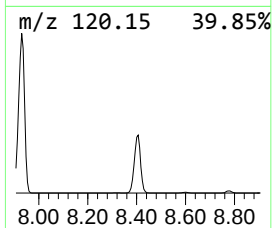
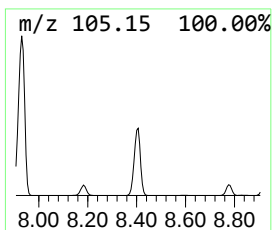
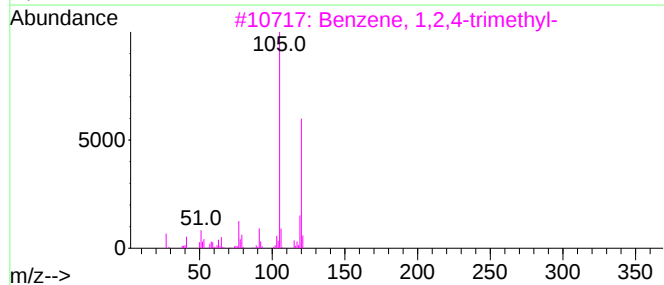
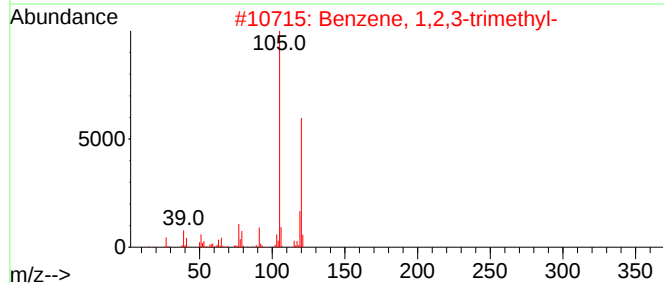
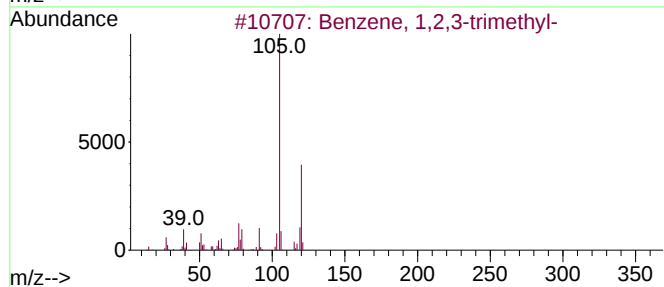
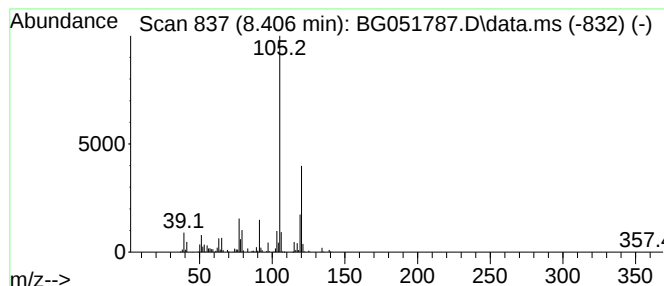
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 17 Benzene, 1,2,3-trimethyl- Concentration Rank 15

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 8.406 | 12.82 ng/ul | 2589570 | 1,4-Dichlorobenzene-d4 | 8.277 |

| Hit# | of 5 | Tentative ID              | MW  | MolForm | CAS#        | Qual |
|------|------|---------------------------|-----|---------|-------------|------|
| 1    |      | Benzene, 1,2,3-trimethyl- | 120 | C9H12   | 000526-73-8 | 95   |
| 2    |      | Benzene, 1,2,3-trimethyl- | 120 | C9H12   | 000526-73-8 | 94   |
| 3    |      | Benzene, 1,2,4-trimethyl- | 120 | C9H12   | 000095-63-6 | 91   |
| 4    |      | Mesitylene                | 120 | C9H12   | 000108-67-8 | 91   |
| 5    |      | Mesitylene                | 120 | C9H12   | 000108-67-8 | 91   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

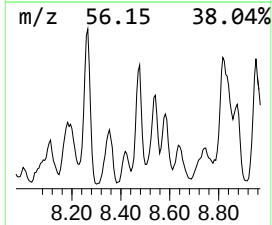
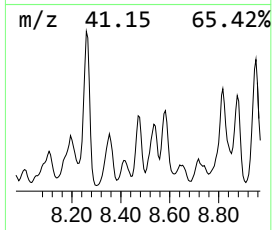
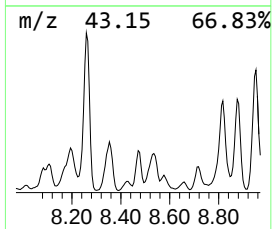
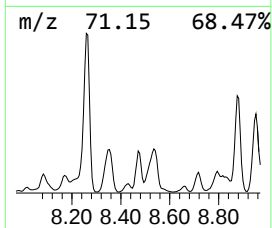
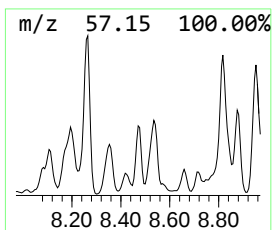
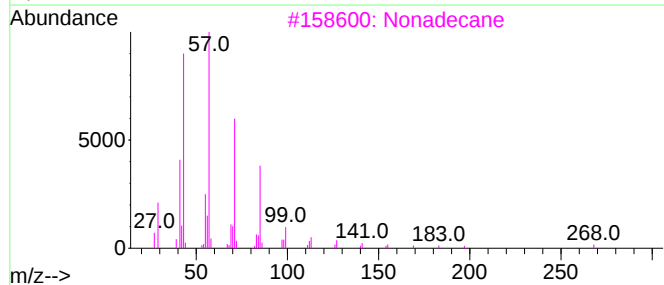
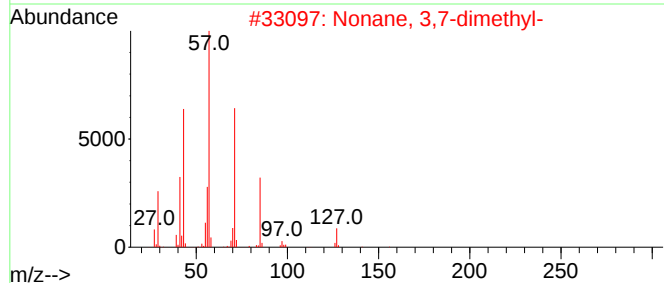
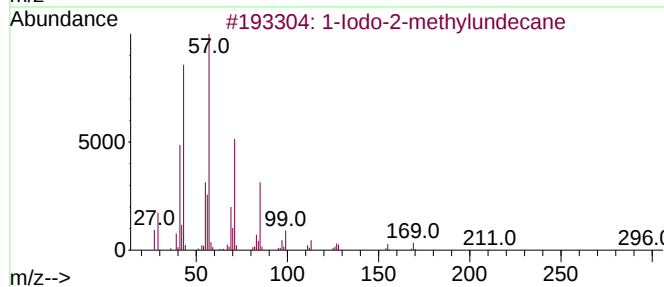
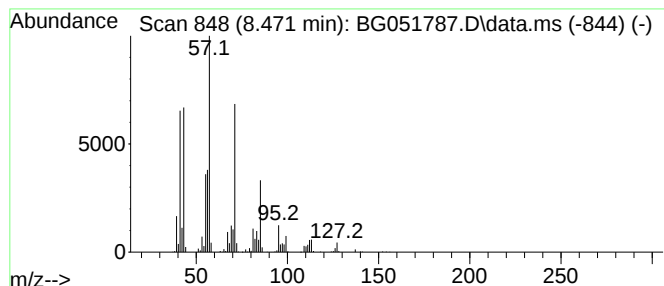
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 1-Iodo-2-methylundecane Concentration Rank 39

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 8.471 | 6.23 ng/ul | 1258790 | 1,4-Dichlorobenzene-d4 | 8.277 |

| Hit# | of | 5 | Tentative ID            | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------|-----|---------|-------------|------|
| 1    |    |   | 1-Iodo-2-methylundecane | 296 | C12H25I | 073105-67-6 | 80   |
| 2    |    |   | Nonane, 3,7-dimethyl-   | 156 | C11H24  | 017302-32-8 | 76   |
| 3    |    |   | Nonadecane              | 268 | C19H40  | 000629-92-5 | 72   |
| 4    |    |   | Pentadecane             | 212 | C15H32  | 000629-62-9 | 72   |
| 5    |    |   | Nonadecane              | 268 | C19H40  | 000629-92-5 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

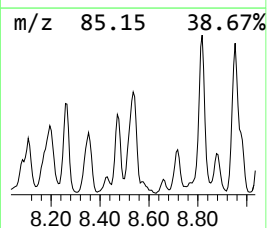
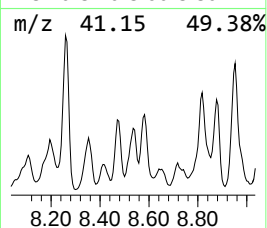
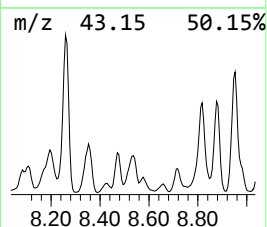
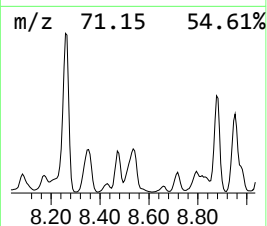
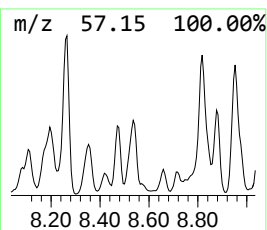
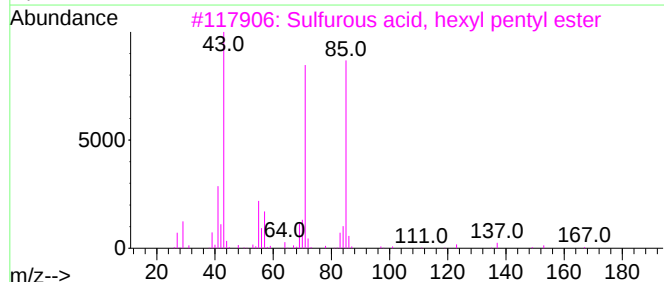
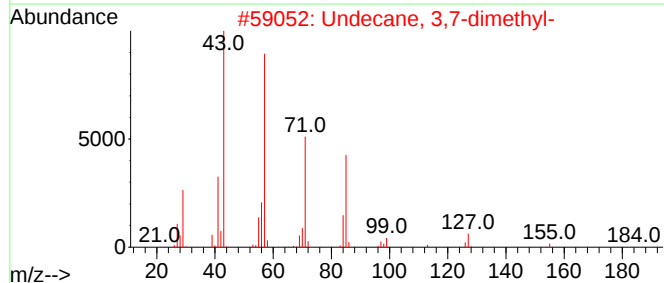
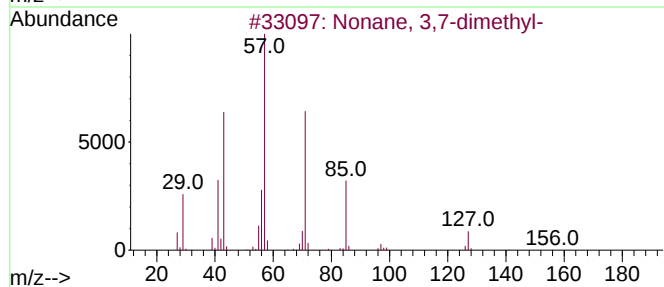
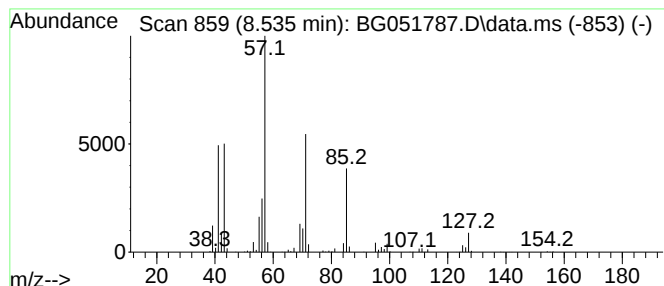
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 38

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 8.535 | 6.25 ng/ul | 1261730 | 1,4-Dichlorobenzene-d4 | 8.277 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|------|-------------------------------------|-----|-----------|--------------|------|
| 1    |      | Nonane, 3,7-dimethyl-               | 156 | C11H24    | 017302-32-8  | 76   |
| 2    |      | Undecane, 3,7-dimethyl-             | 184 | C13H28    | 017301-29-0  | 64   |
| 3    |      | Sulfurous acid, hexyl pentyl ester  | 236 | C11H24O3S | 1000309-14-1 | 64   |
| 4    |      | Sulfurous acid, 2-ethylhexyl hex... | 278 | C14H30O3S | 1000309-20-2 | 64   |
| 5    |      | Sulfurous acid, nonyl 2-propyl e... | 250 | C12H26O3S | 1000309-12-0 | 58   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

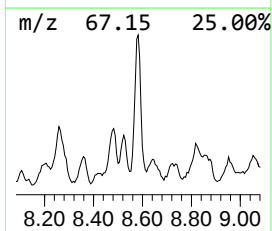
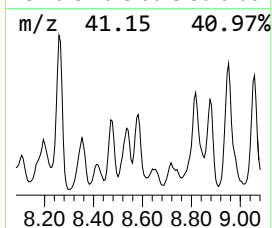
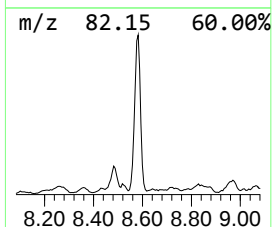
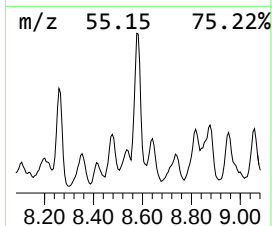
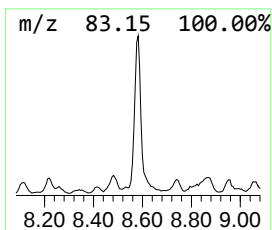
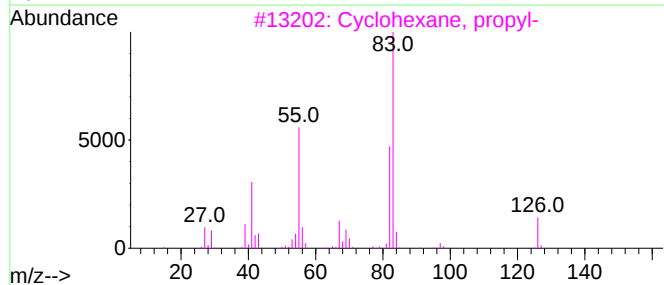
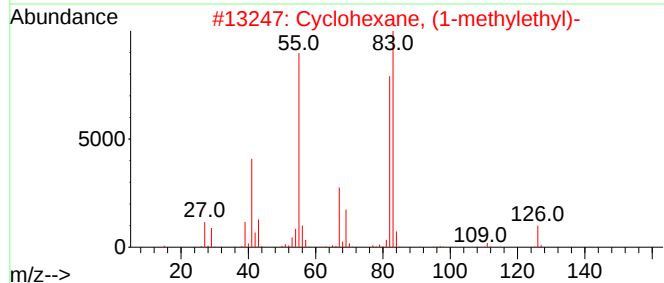
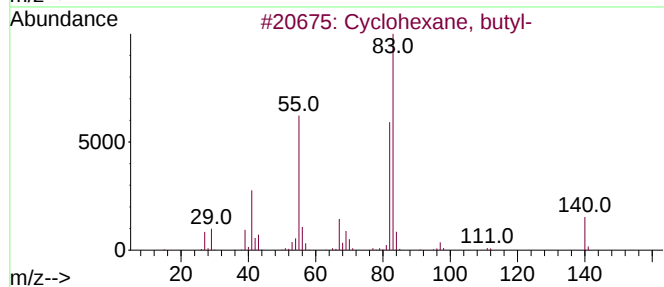
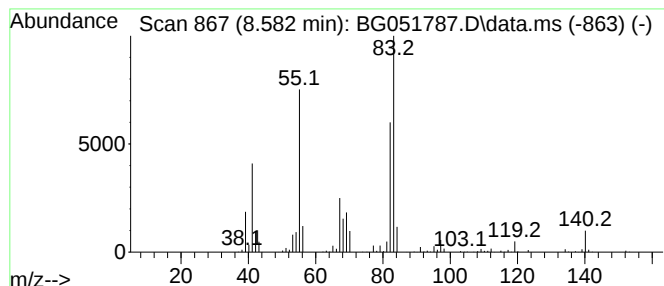
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 (DEL) Alkane: Cyclic8.582 Concentration Rank 31

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 8.582 | 8.16 ng/ul | 1647530 | 1,4-Dichlorobenzene-d4 | 8.277 |

| Hit# | of 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------|-----|---------|-------------|------|
| 1    |      | Cyclohexane, butyl-           | 140 | C10H20  | 001678-93-9 | 90   |
| 2    |      | Cyclohexane, (1-methylethyl)- | 126 | C9H18   | 000696-29-7 | 80   |
| 3    |      | Cyclohexane, propyl-          | 126 | C9H18   | 001678-92-8 | 72   |
| 4    |      | Cyclohexane, octyl-           | 196 | C14H28  | 001795-15-9 | 72   |
| 5    |      | Cyclohexane, octyl-           | 196 | C14H28  | 001795-15-9 | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
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Quant Title : SVOA CALIBRATION

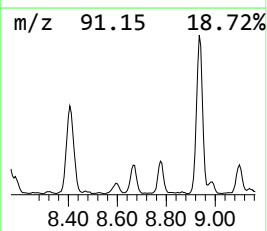
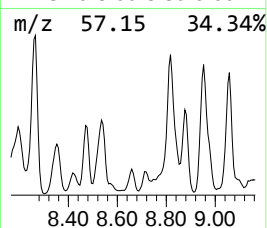
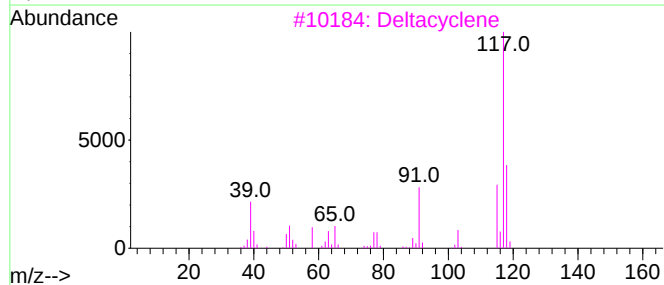
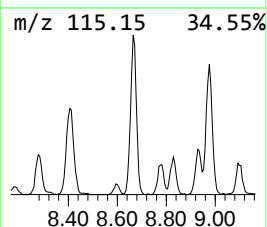
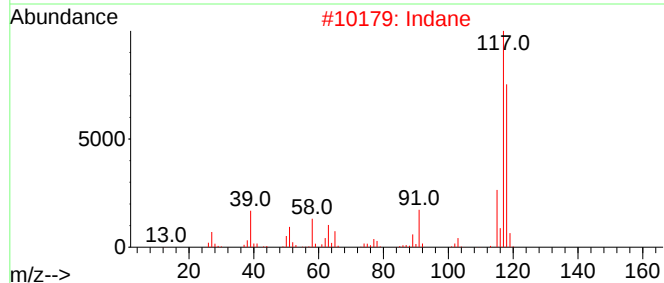
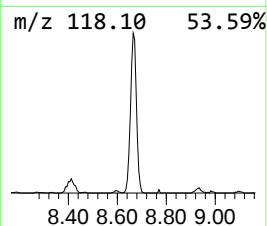
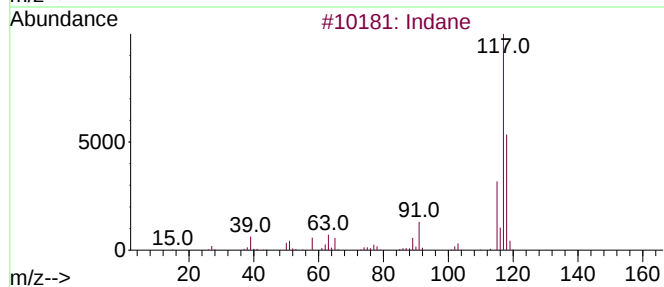
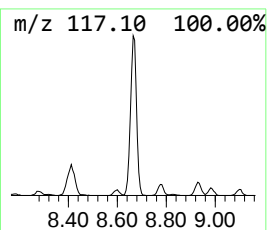
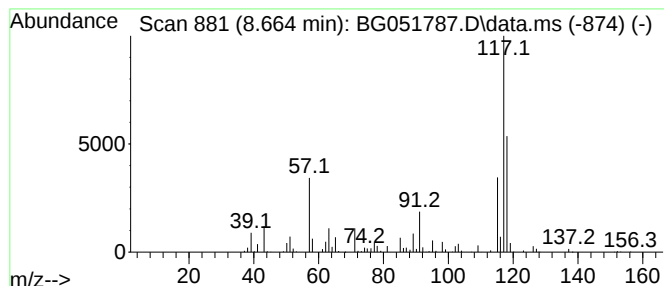
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 21 Indane Concentration Rank 45

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 8.664 | 4.75 ng/ul | 958504 | 1,4-Dichlorobenzene-d4 | 8.277 |

| Hit# | of 5                         | Tentative ID | MW  | MolForm | CAS#        | Qual |
|------|------------------------------|--------------|-----|---------|-------------|------|
| 1    | Indane                       |              | 118 | C9H10   | 000496-11-7 | 81   |
| 2    | Indane                       |              | 118 | C9H10   | 000496-11-7 | 81   |
| 3    | Deltacyclene                 |              | 118 | C9H10   | 007785-10-6 | 74   |
| 4    | Benzene, 1-ethenyl-2-methyl- |              | 118 | C9H10   | 000611-15-4 | 68   |
| 5    | Benzene, cyclopropyl-        |              | 118 | C9H10   | 000873-49-4 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

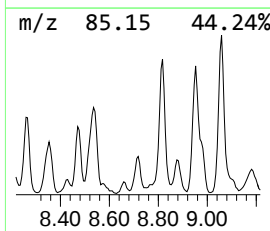
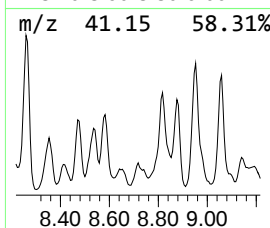
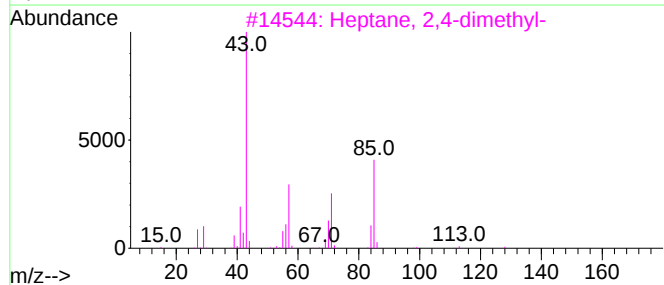
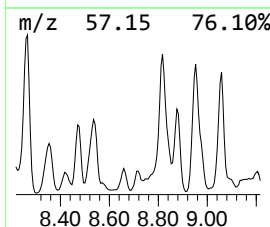
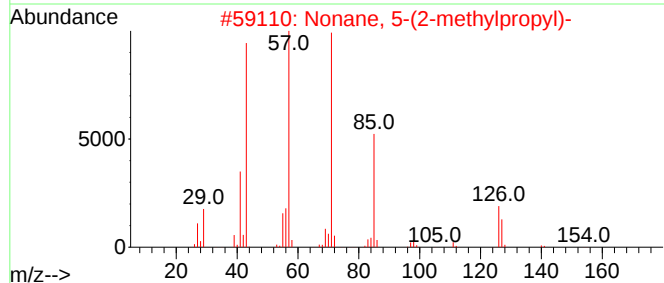
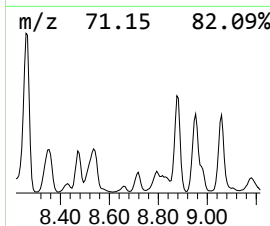
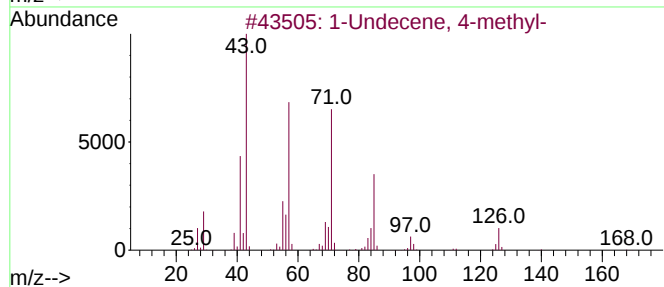
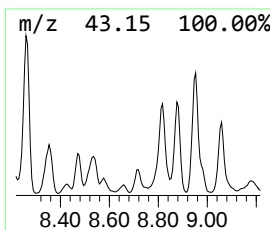
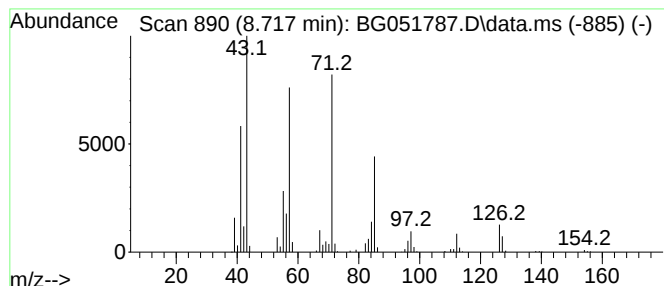
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Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 55

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 8.717 | 2.86 ng/ul | 577022 | 1,4-Dichlorobenzene-d4 | 8.277 |

| Hit# | of 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|------|-----------------------------|-----|---------|-------------|------|
| 1    |      | 1-Undecene, 4-methyl-       | 168 | C12H24  | 074630-39-0 | 78   |
| 2    |      | Nonane, 5-(2-methylpropyl)- | 184 | C13H28  | 062185-53-9 | 72   |
| 3    |      | Heptane, 2,4-dimethyl-      | 128 | C9H20   | 002213-23-2 | 64   |
| 4    |      | Dodecane, 4-methyl-         | 184 | C13H28  | 006117-97-1 | 64   |
| 5    |      | Decane, 5-propyl-           | 184 | C13H28  | 017312-62-8 | 59   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

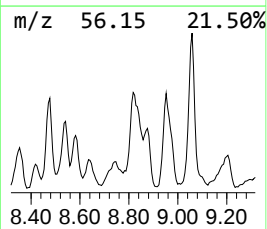
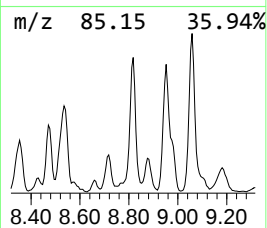
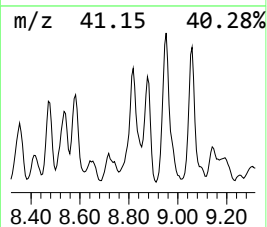
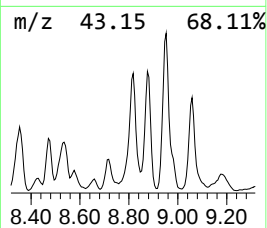
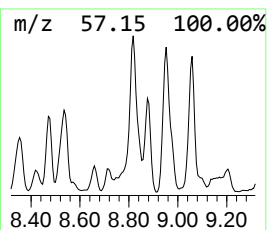
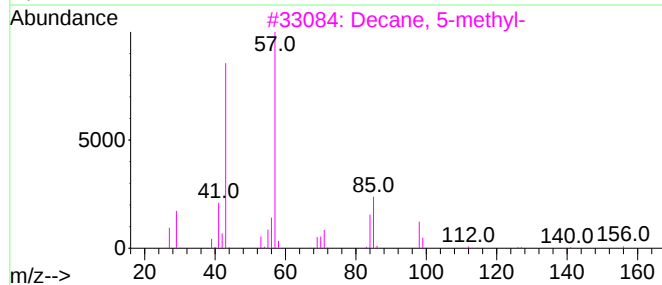
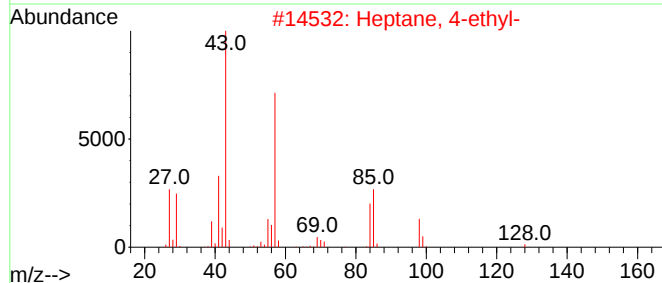
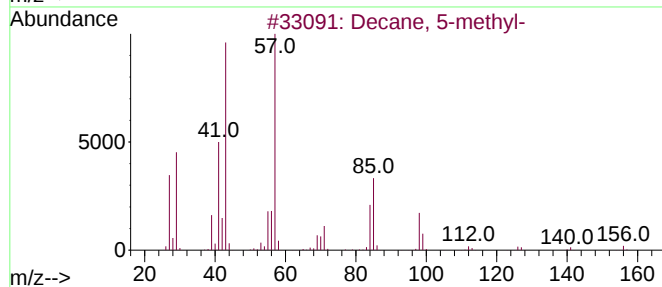
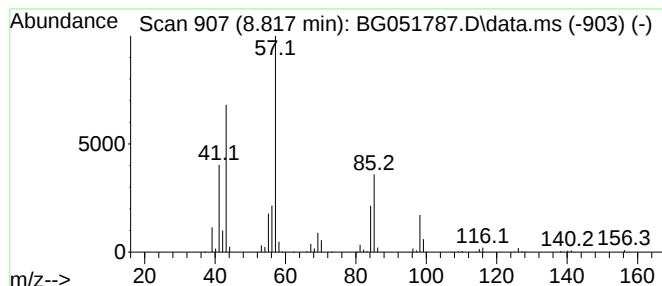
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 22

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 8.817 | 9.76 ng/ul | 1970700 | 1,4-Dichlorobenzene-d4 | 8.277 |

| Hit# | of 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|------|--------------------------------|-----|---------|-------------|------|
| 1    |      | Decane, 5-methyl-              | 156 | C11H24  | 013151-35-4 | 78   |
| 2    |      | Heptane, 4-ethyl-              | 128 | C9H20   | 002216-32-2 | 72   |
| 3    |      | Decane, 5-methyl-              | 156 | C11H24  | 013151-35-4 | 64   |
| 4    |      | Heptane, 4-ethyl-              | 128 | C9H20   | 002216-32-2 | 59   |
| 5    |      | Pentane, 3-ethyl-2,3-dimethyl- | 128 | C9H20   | 016747-33-4 | 50   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

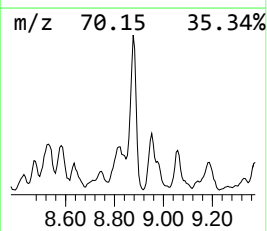
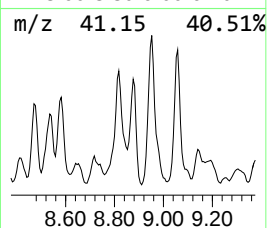
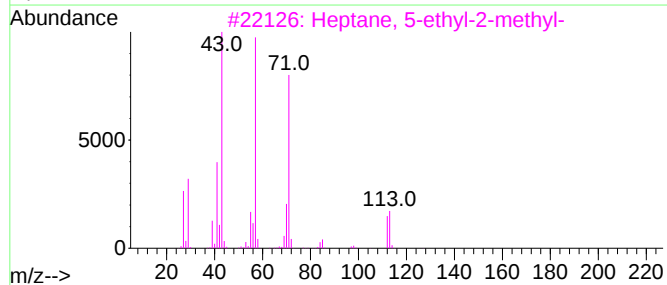
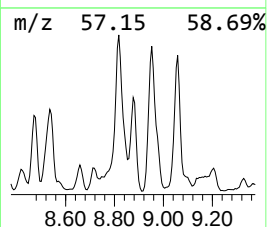
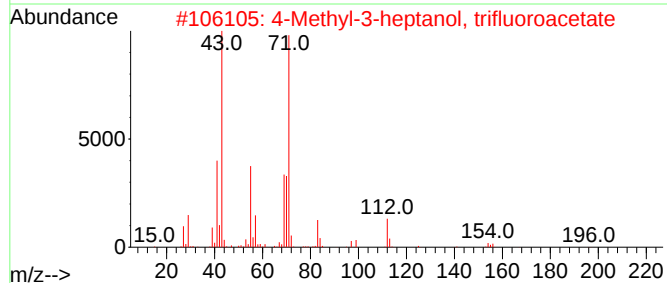
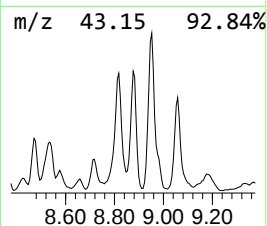
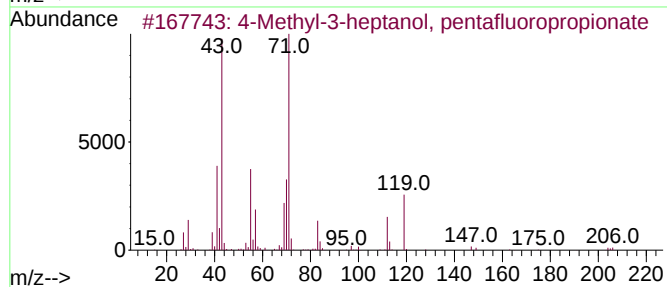
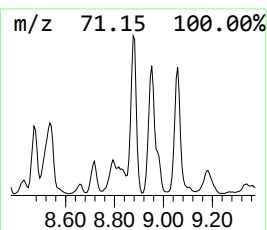
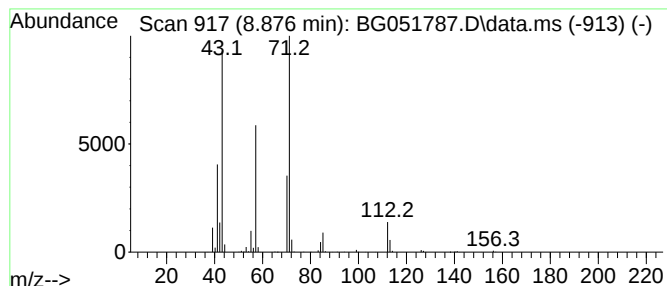
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 25 4-Methyl-3-heptanol, pentafl... Concentration Rank 23

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 8.876 | 9.59 ng/ul | 1935950 | 1,4-Dichlorobenzene-d4 | 8.277 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | 4-Methyl-3-heptanol, pentafluoro... | 276 | C11H17F5O2 | 1000365-51-7 | 83   |
| 2    |    |   | 4-Methyl-3-heptanol, trifluoroac... | 226 | C10H17F3O2 | 1000365-51-8 | 83   |
| 3    |    |   | Heptane, 5-ethyl-2-methyl-          | 142 | C10H22     | 013475-78-0  | 78   |
| 4    |    |   | Decane, 4-methyl-                   | 156 | C11H24     | 002847-72-5  | 74   |
| 5    |    |   | Decane, 4-methyl-                   | 156 | C11H24     | 002847-72-5  | 68   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

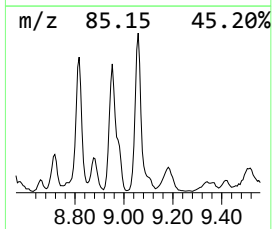
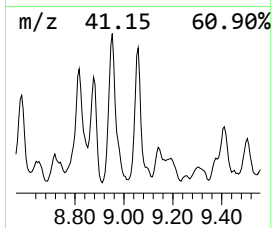
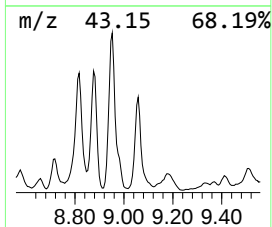
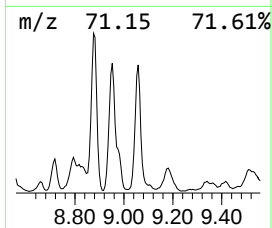
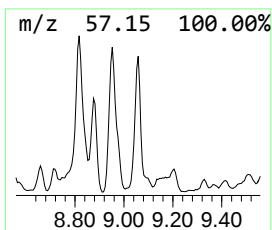
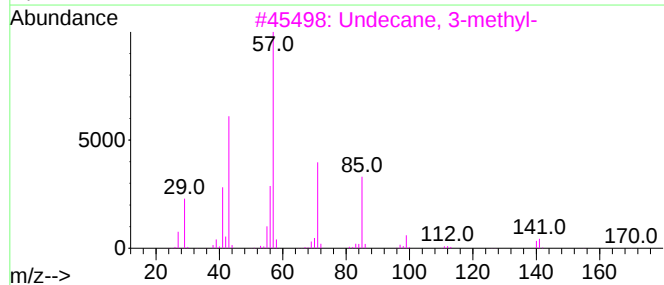
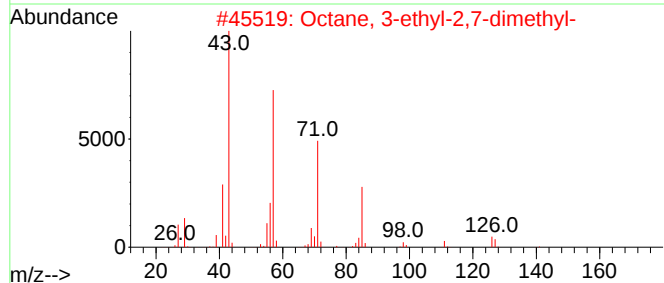
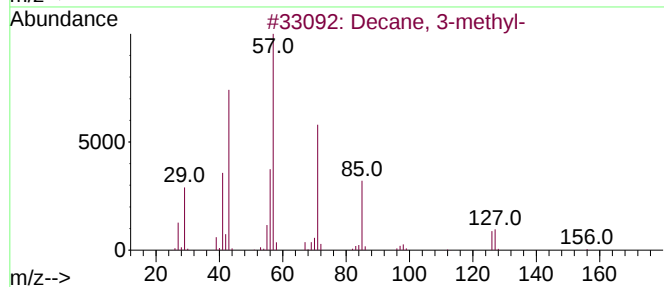
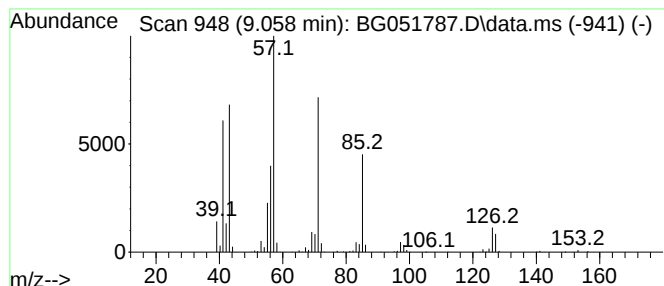
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Peak Number 26 (DEL) Alkane: Straight-Chai... Concentration Rank 17

| R.T.  | EstConc     | Area    | Relative to ISTD       | R.T.  |
|-------|-------------|---------|------------------------|-------|
| 9.058 | 11.70 ng/ul | 2362510 | 1,4-Dichlorobenzene-d4 | 8.277 |

| Hit# | of 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------|-----|---------|-------------|------|
| 1    |      | Decane, 3-methyl-             | 156 | C11H24  | 013151-34-3 | 91   |
| 2    |      | Octane, 3-ethyl-2,7-dimethyl- | 170 | C12H26  | 062183-55-5 | 83   |
| 3    |      | Undecane, 3-methyl-           | 170 | C12H26  | 001002-43-3 | 78   |
| 4    |      | Undecane, 3-methyl-           | 170 | C12H26  | 001002-43-3 | 72   |
| 5    |      | Nonane, 3,7-dimethyl-         | 156 | C11H24  | 017302-32-8 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

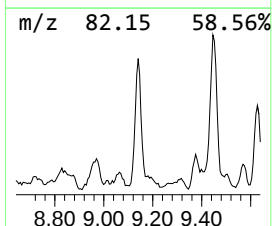
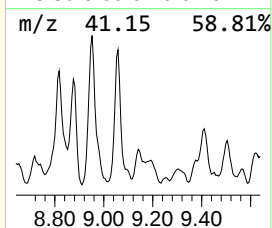
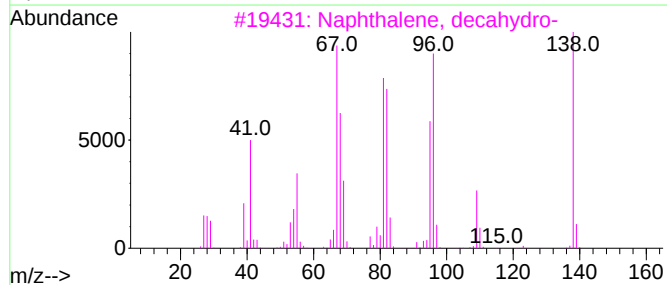
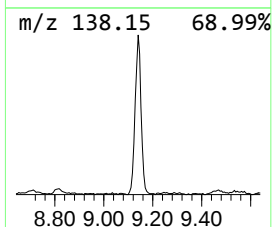
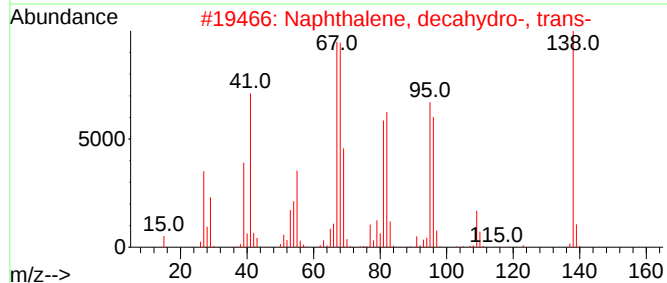
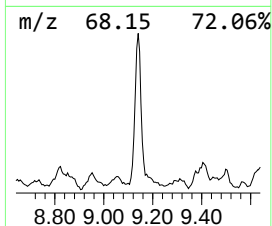
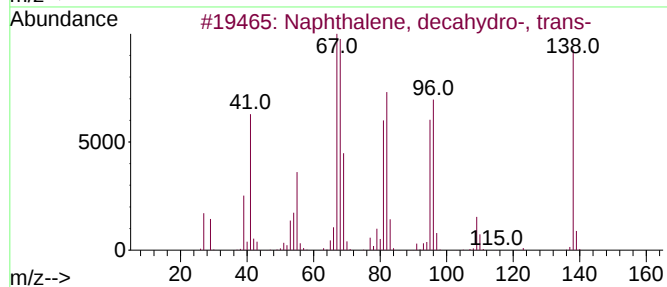
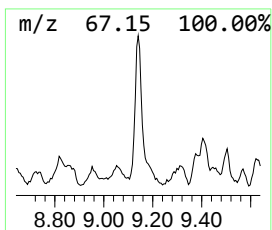
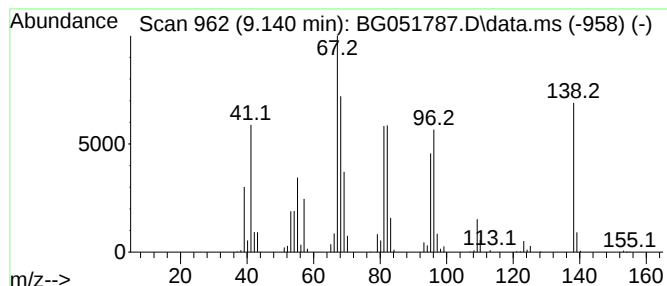
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 28 Naphthalene, decahydro-, tr... Concentration Rank 52

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 9.140 | 3.35 ng/ul | 676811 | 1,4-Dichlorobenzene-d4 | 8.277 |

| Hit# | of 5 | Tentative ID                    | MW  | MolForm | CAS#        | Qual |
|------|------|---------------------------------|-----|---------|-------------|------|
| 1    |      | Naphthalene, decahydro-, trans- | 138 | C10H18  | 000493-02-7 | 98   |
| 2    |      | Naphthalene, decahydro-, trans- | 138 | C10H18  | 000493-02-7 | 97   |
| 3    |      | Naphthalene, decahydro-         | 138 | C10H18  | 000091-17-8 | 97   |
| 4    |      | Naphthalene, decahydro-         | 138 | C10H18  | 000091-17-8 | 96   |
| 5    |      | Naphthalene, decahydro-         | 138 | C10H18  | 000091-17-8 | 96   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

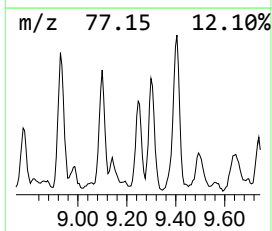
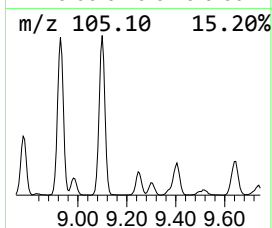
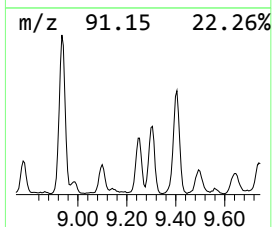
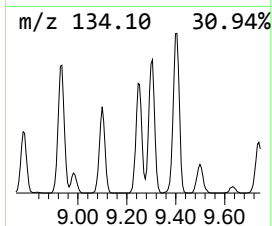
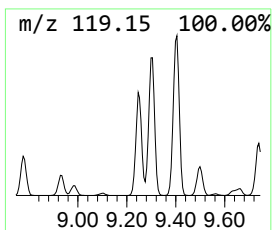
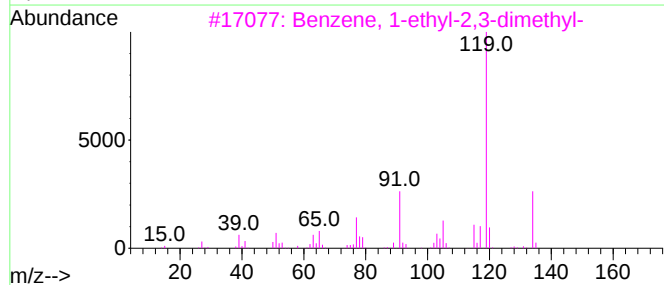
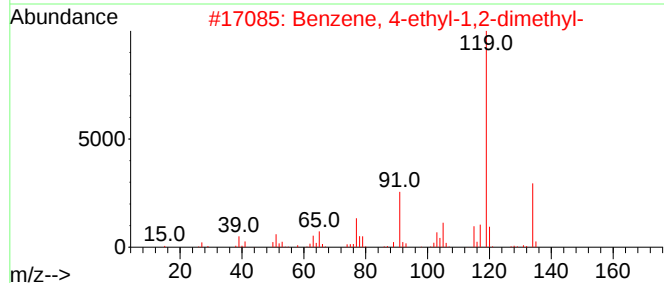
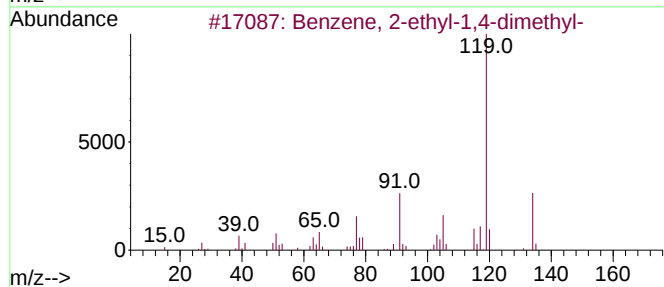
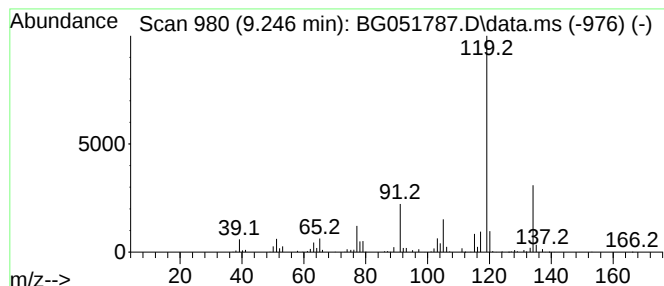
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TIC Integration Parameters: LSCINT.P

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Peak Number 29 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 50

| R.T.  | EstConc    | Area   | Relative to ISTD       | R.T.  |
|-------|------------|--------|------------------------|-------|
| 9.246 | 3.68 ng/ul | 742508 | 1,4-Dichlorobenzene-d4 | 8.277 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 96   |
| 2    |    |   | Benzene, 4-ethyl-1,2-dimethyl- | 134 | C10H14  | 000934-80-5 | 95   |
| 3    |    |   | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14  | 000933-98-2 | 95   |
| 4    |    |   | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 94   |
| 5    |    |   | o-Cymene                       | 134 | C10H14  | 000527-84-4 | 94   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

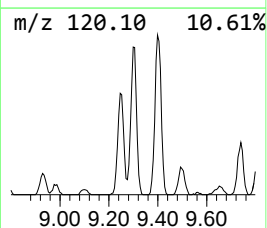
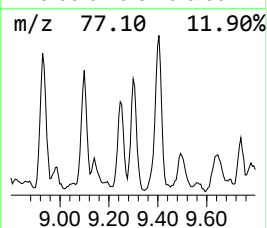
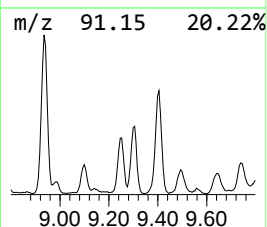
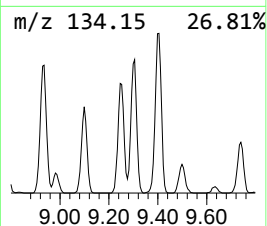
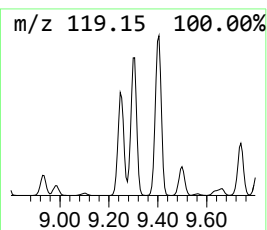
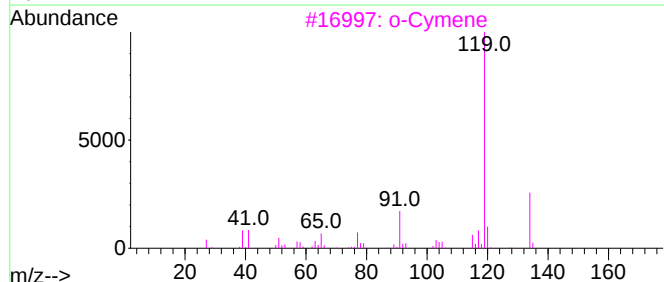
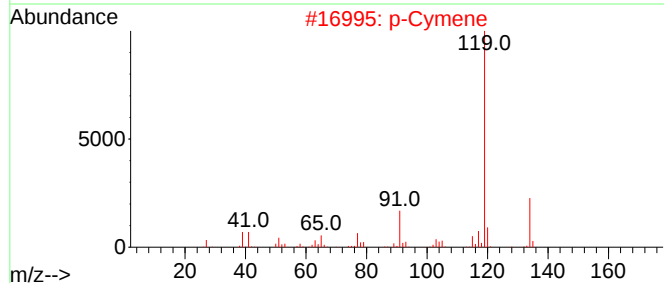
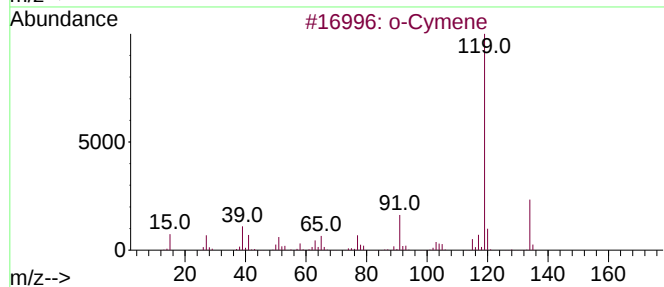
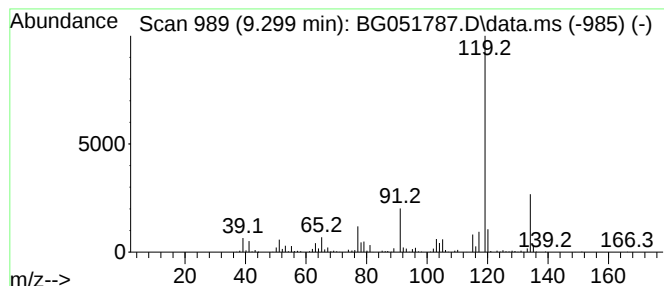
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 30 o-Cymene Concentration Rank 34

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 9.299 | 7.50 ng/ul | 1514270 | 1,4-Dichlorobenzene-d4 | 8.277 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | o-Cymene                            | 134 | C10H14  | 000527-84-4 | 95   |
| 2    |    |   | p-Cymene                            | 134 | C10H14  | 000099-87-6 | 95   |
| 3    |    |   | o-Cymene                            | 134 | C10H14  | 000527-84-4 | 95   |
| 4    |    |   | Benzene, 1-ethyl-2,4-dimethyl-      | 134 | C10H14  | 000874-41-9 | 95   |
| 5    |    |   | Benzene, 1-methyl-3-(1-methyleth... | 134 | C10H14  | 000535-77-3 | 94   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

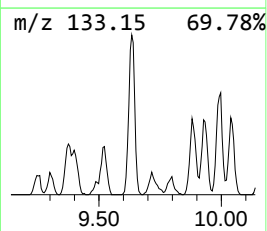
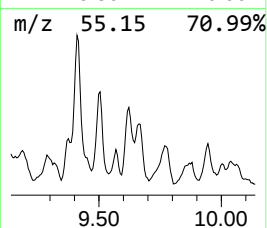
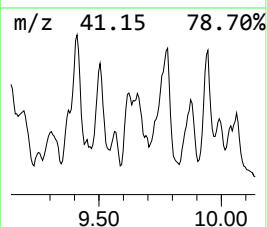
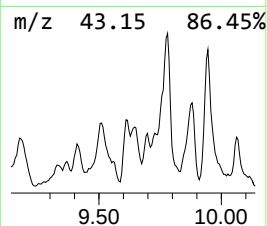
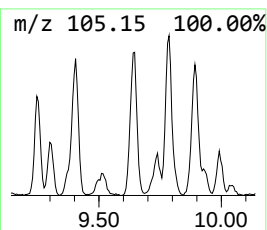
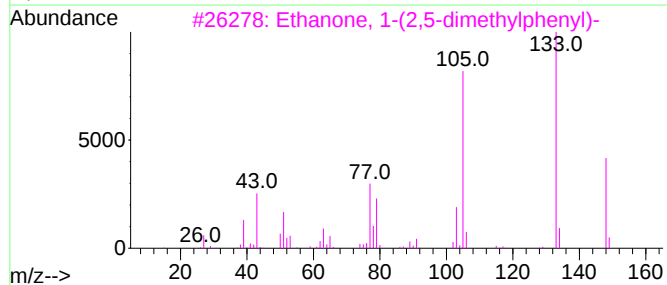
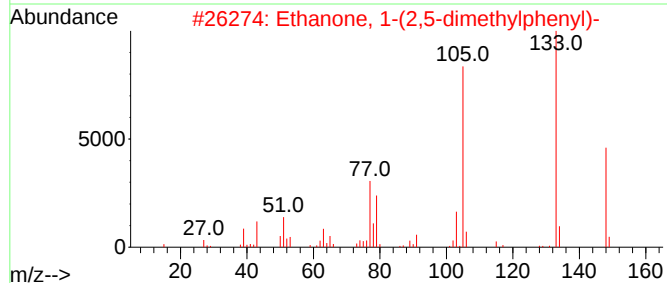
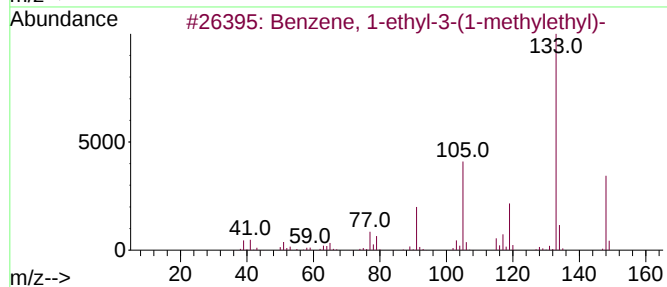
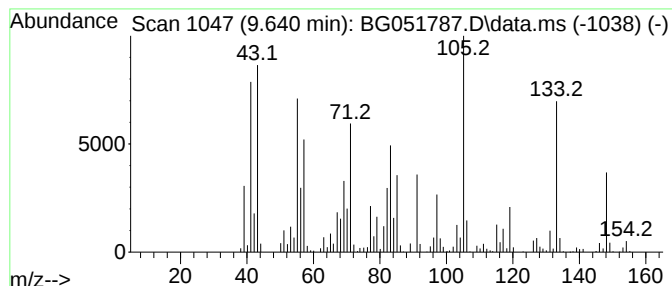
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 31 Benzene, 1-ethyl-3-(1-methy... Concentration Rank 37

| R.T.  | EstConc    | Area    | Relative to ISTD       | R.T.  |
|-------|------------|---------|------------------------|-------|
| 9.640 | 6.33 ng/ul | 1278570 | 1,4-Dichlorobenzene-d4 | 8.277 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|-------------|-------------|------|
| 1    |    |   | Benzene, 1-ethyl-3-(1-methylethyl)- | 148 | C11H16      | 004920-99-4 | 84   |
| 2    |    |   | Ethanone, 1-(2,5-dimethylphenyl)-   | 148 | C10H12O     | 002142-73-6 | 53   |
| 3    |    |   | Ethanone, 1-(2,5-dimethylphenyl)-   | 148 | C10H12O     | 002142-73-6 | 53   |
| 4    |    |   | Dichloroacetic acid, nonyl ester    | 254 | C11H20Cl2O2 | 083004-99-3 | 43   |
| 5    |    |   | Benzene, 1-methoxy-2-(1-methylet... | 148 | C10H12O     | 010278-02-1 | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

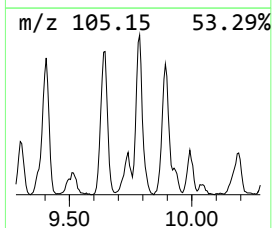
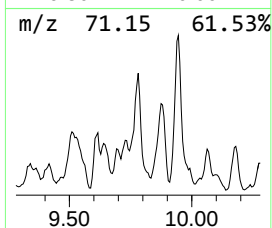
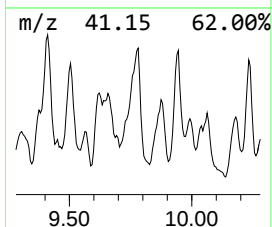
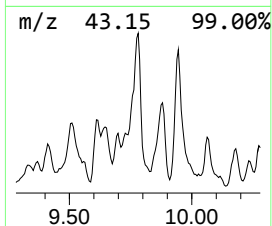
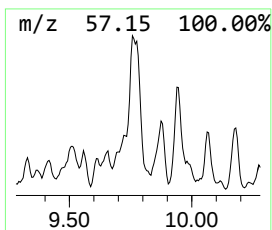
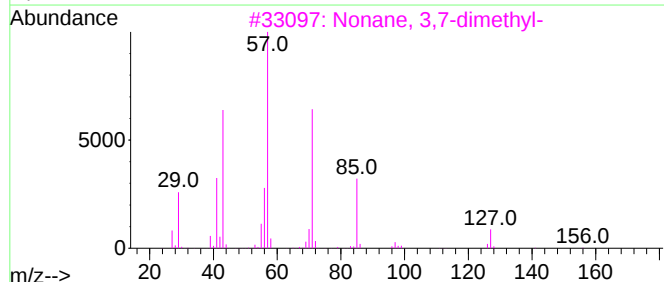
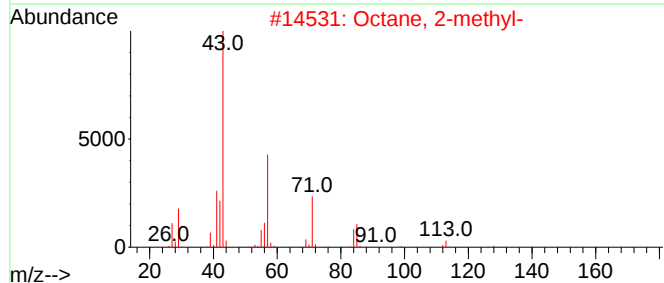
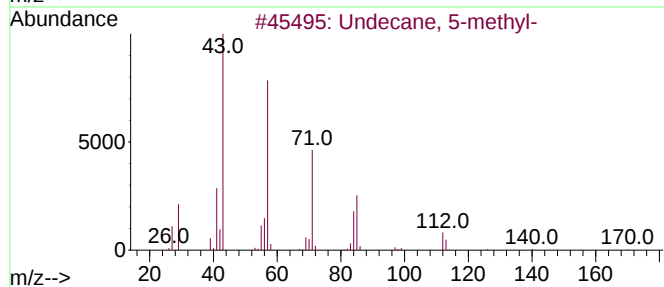
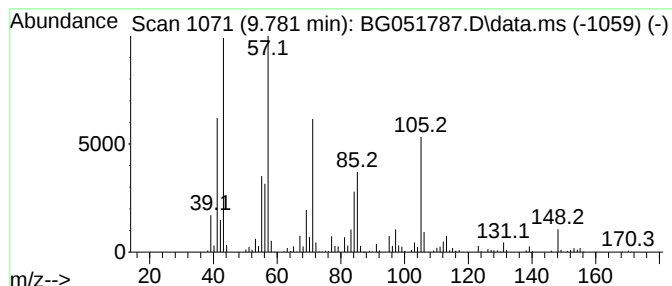
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 32 (DEL) Alkane: Straight-Chai... Concentration Rank 1

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.   |
|-------|-------------|---------|------------------|--------|
| 9.781 | 48.88 ng/ul | 2446830 | Naphthalene-d8   | 11.120 |

| Hit# | of 5 | Tentative ID             | MW  | MolForm | CAS#        | Qual |
|------|------|--------------------------|-----|---------|-------------|------|
| 1    |      | Undecane, 5-methyl-      | 170 | C12H26  | 001632-70-8 | 62   |
| 2    |      | Octane, 2-methyl-        | 128 | C9H20   | 003221-61-2 | 62   |
| 3    |      | Nonane, 3,7-dimethyl-    | 156 | C11H24  | 017302-32-8 | 52   |
| 4    |      | Undecane, 3-methyl-      | 170 | C12H26  | 001002-43-3 | 52   |
| 5    |      | Hexane, 2,3,4-trimethyl- | 128 | C9H20   | 000921-47-1 | 52   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

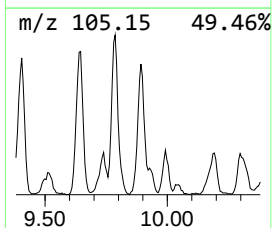
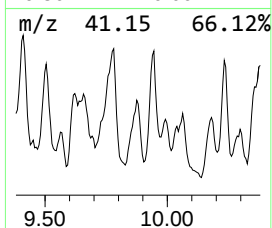
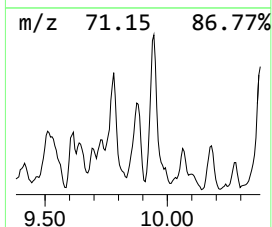
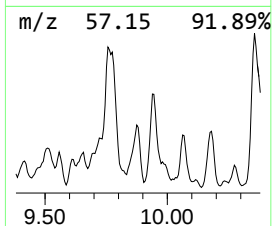
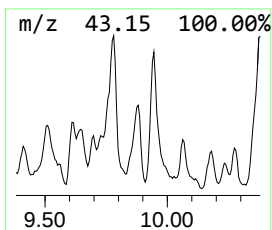
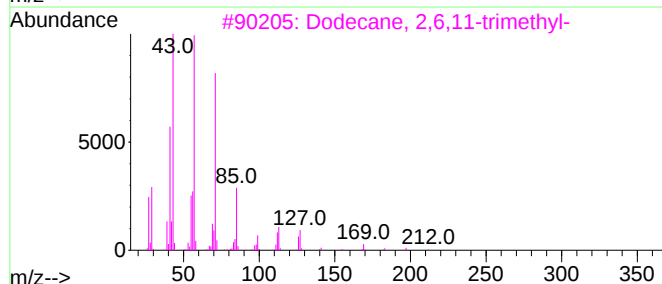
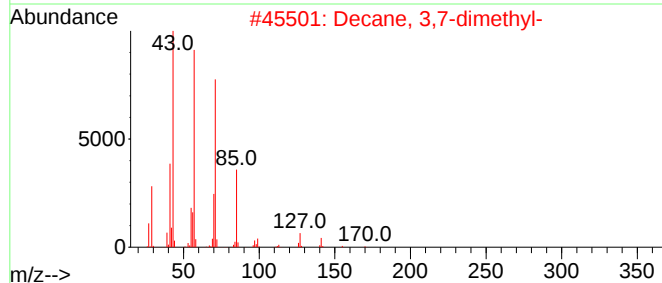
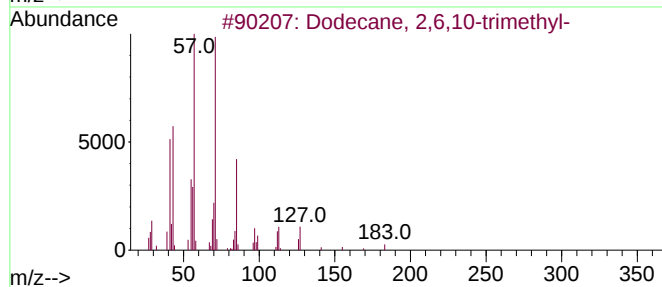
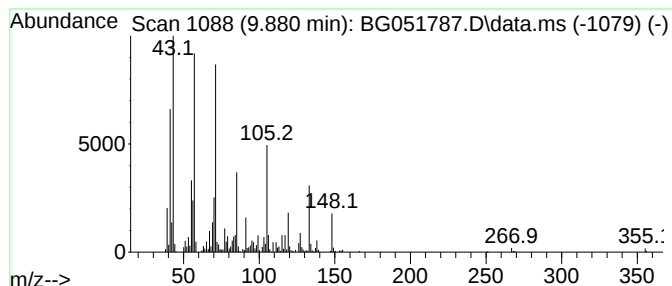
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Peak Number 33 (DEL) Alkane: Straight-Chai... Concentration Rank 6

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.   |
|-------|-------------|---------|------------------|--------|
| 9.880 | 20.58 ng/ul | 1030260 | Naphthalene-d8   | 11.120 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------------|-----|---------|-------------|------|
| 1    |      | Dodecane, 2,6,10-trimethyl-         | 212 | C15H32  | 003891-98-3 | 64   |
| 2    |      | Decane, 3,7-dimethyl-               | 170 | C12H26  | 017312-54-8 | 59   |
| 3    |      | Dodecane, 2,6,11-trimethyl-         | 212 | C15H32  | 031295-56-4 | 58   |
| 4    |      | Pentadecane, 2,6,10,14-tetramethyl- | 268 | C19H40  | 001921-70-6 | 58   |
| 5    |      | Heptadecane, 2,6-dimethyl-          | 268 | C19H40  | 054105-67-8 | 55   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

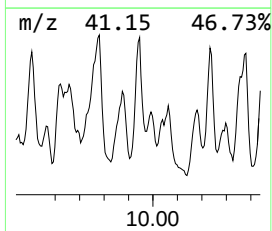
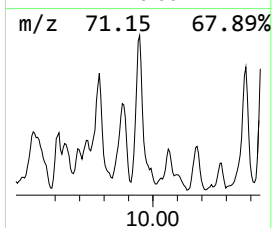
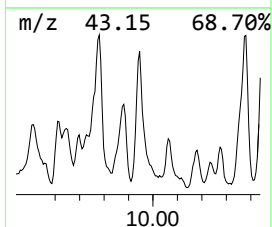
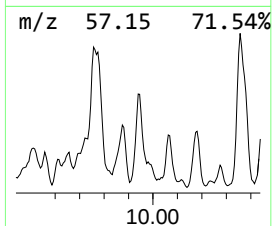
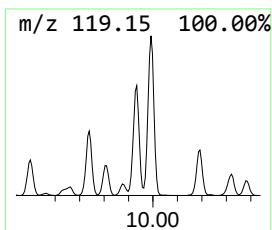
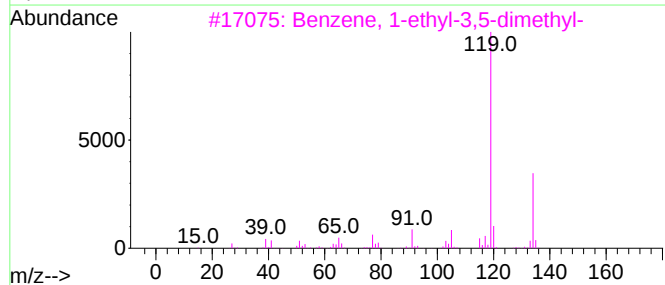
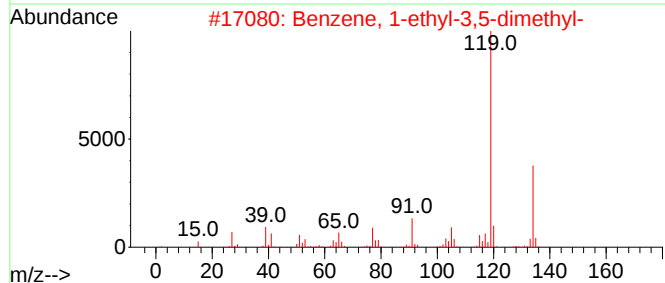
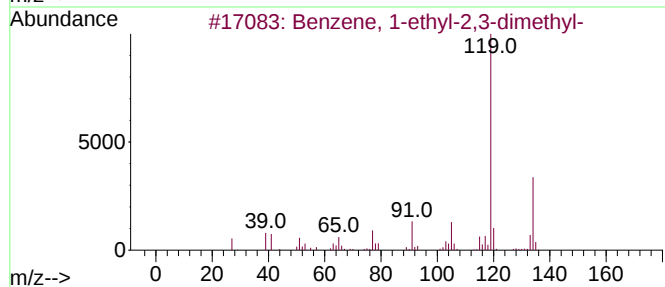
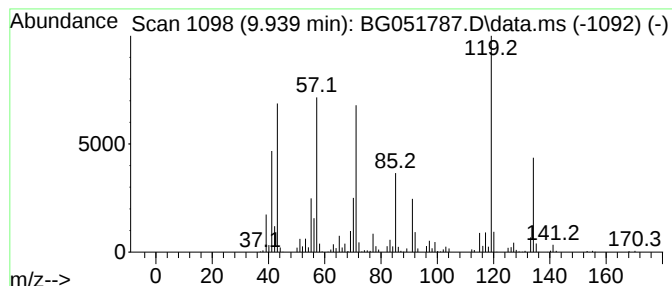
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Peak Number 34 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 3

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.   |
|-------|-------------|---------|------------------|--------|
| 9.939 | 35.99 ng/ul | 1801780 | Naphthalene-d8   | 11.120 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14  | 000933-98-2 | 60   |
| 2    |    |   | Benzene, 1-ethyl-3,5-dimethyl- | 134 | C10H14  | 000934-74-7 | 60   |
| 3    |    |   | Benzene, 1-ethyl-3,5-dimethyl- | 134 | C10H14  | 000934-74-7 | 60   |
| 4    |    |   | Benzene, 1,2,3,5-tetramethyl-  | 134 | C10H14  | 000527-53-7 | 60   |
| 5    |    |   | Benzene, 1-ethyl-3,5-dimethyl- | 134 | C10H14  | 000934-74-7 | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

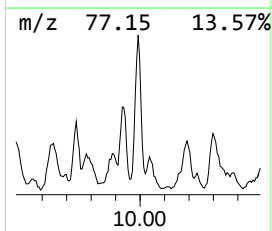
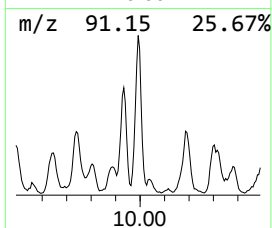
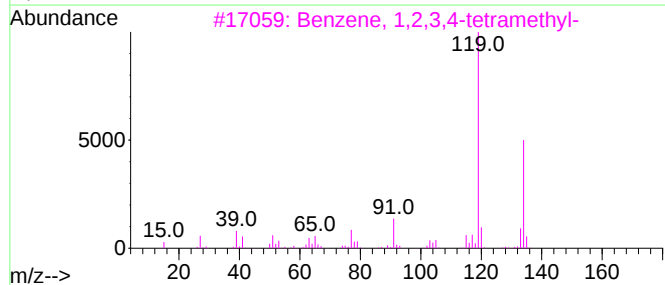
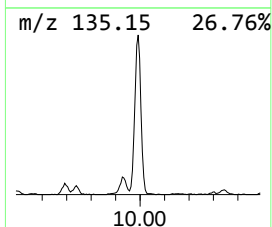
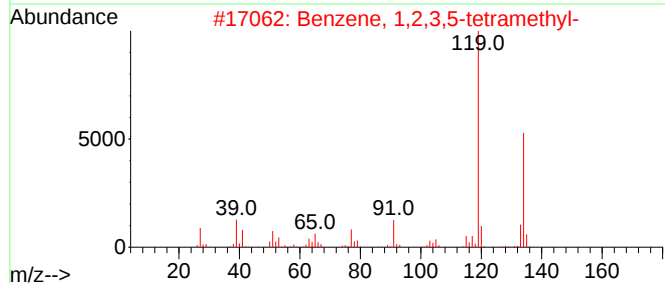
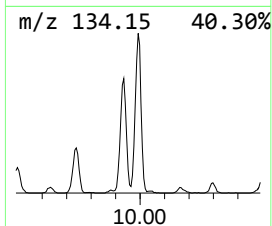
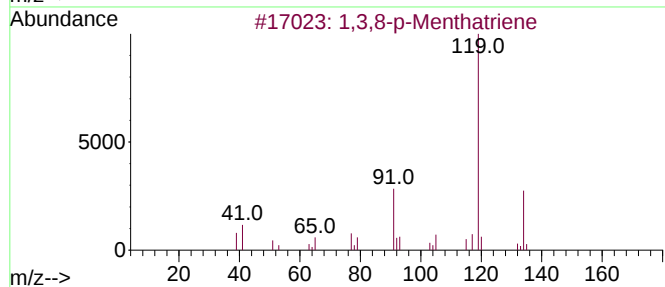
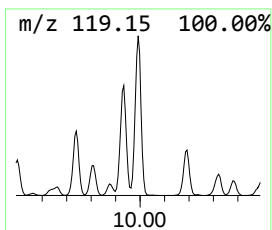
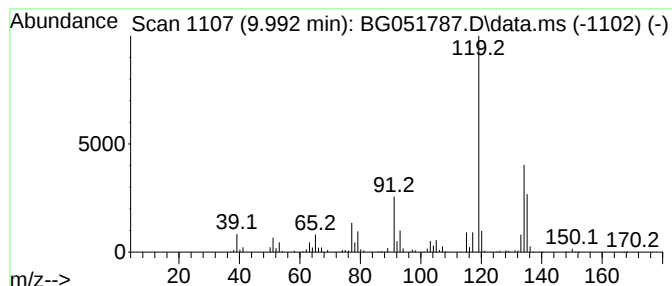
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 35 1,3,8-p-Menthatriene Concentration Rank 4

| R.T.  | EstConc     | Area    | Relative to ISTD | R.T.   |
|-------|-------------|---------|------------------|--------|
| 9.992 | 34.20 ng/ul | 1711780 | Naphthalene-d8   | 11.120 |

| Hit# | of 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|------|--------------------------------|-----|---------|-------------|------|
| 1    |      | 1,3,8-p-Menthatriene           | 134 | C10H14  | 018368-95-1 | 90   |
| 2    |      | Benzene, 1,2,3,5-tetramethyl-  | 134 | C10H14  | 000527-53-7 | 87   |
| 3    |      | Benzene, 1,2,3,4-tetramethyl-  | 134 | C10H14  | 000488-23-3 | 87   |
| 4    |      | o-Cymene                       | 134 | C10H14  | 000527-84-4 | 87   |
| 5    |      | Benzene, 1-ethyl-3,5-dimethyl- | 134 | C10H14  | 000934-74-7 | 87   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

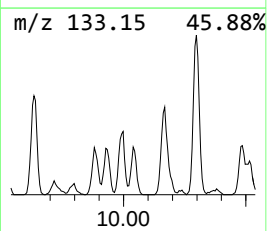
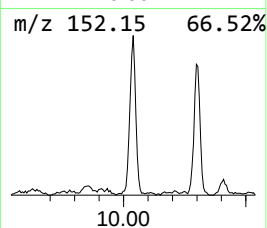
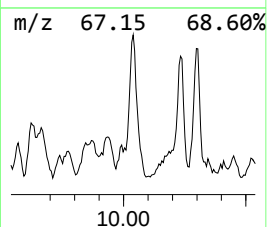
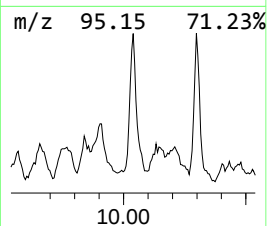
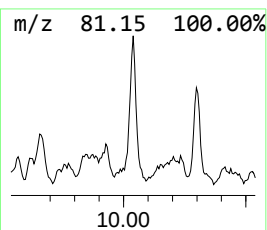
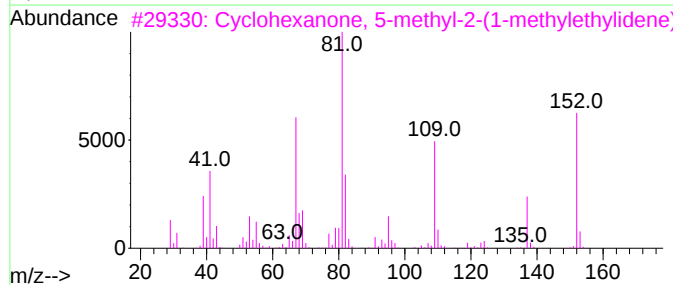
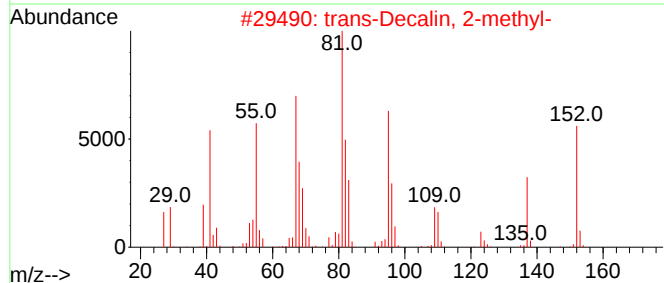
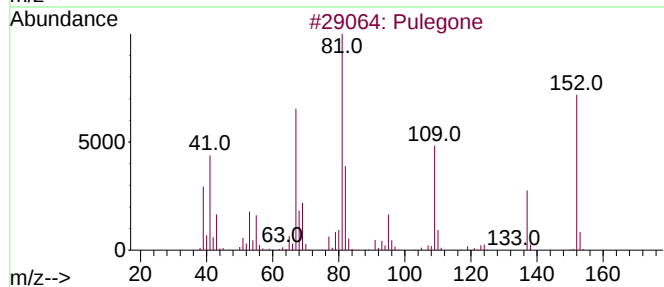
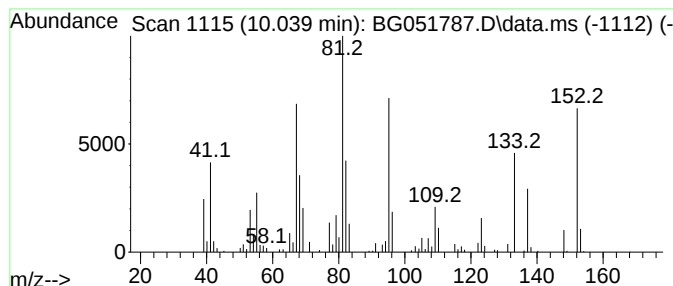
Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 36 Pulegone Concentration Rank 9

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 10.039    | 18.09 ng/ul                         | 905674 | Naphthalene-d8   | 11.120       |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | Pulegone                            | 152    | C10H16O          | 000089-82-7  | 86   |
| 2         | trans-Decalin, 2-methyl-            | 152    | C11H20           | 1000152-47-3 | 72   |
| 3         | Cyclohexanone, 5-methyl-2-(1-met... | 152    | C10H16O          | 015932-80-6  | 70   |
| 4         | cis-Decalin, 2-syn-methyl-          | 152    | C11H20           | 1000155-85-6 | 68   |
| 5         | Cyclohexanone, 5-methyl-2-(1-met... | 152    | C10H16O          | 015932-80-6  | 62   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

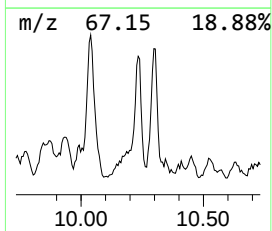
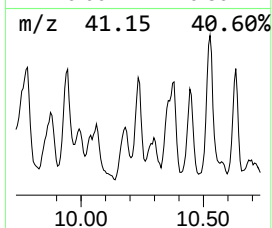
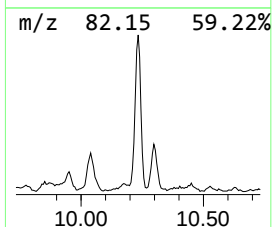
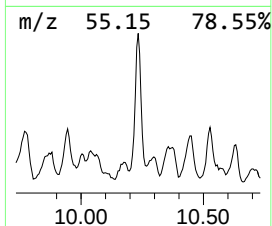
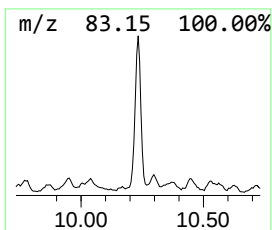
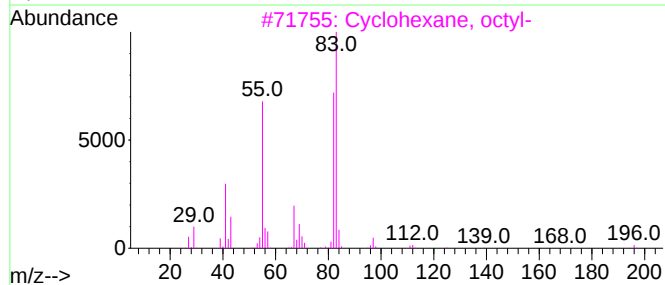
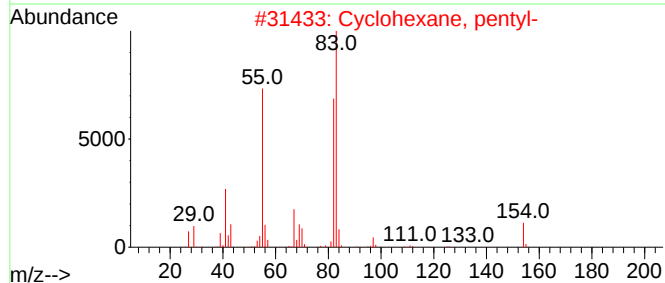
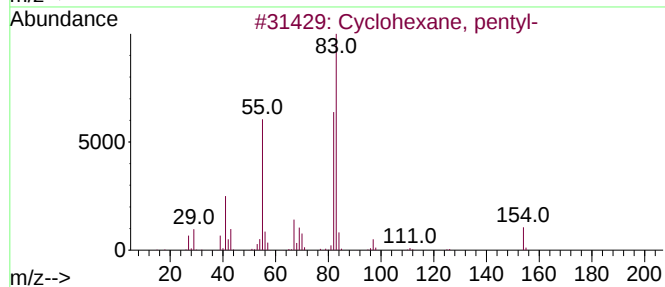
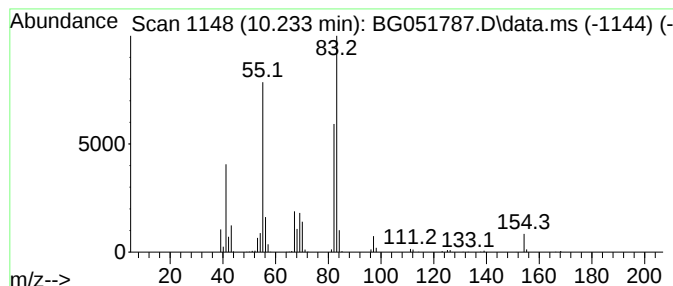
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 37 (DEL) Alkane: Cyclic10.233 Concentration Rank 8

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 10.233 | 20.11 ng/ul | 1006810 | Naphthalene-d8   | 11.120 |

| Hit# | of | 5 | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, pentyl- | 154 | C11H22  | 004292-92-6 | 91   |
| 2    |    |   | Cyclohexane, pentyl- | 154 | C11H22  | 004292-92-6 | 91   |
| 3    |    |   | Cyclohexane, octyl-  | 196 | C14H28  | 001795-15-9 | 78   |
| 4    |    |   | Cyclohexane, butyl-  | 140 | C10H20  | 001678-93-9 | 78   |
| 5    |    |   | Cyclohexane, propyl- | 126 | C9H18   | 001678-92-8 | 78   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

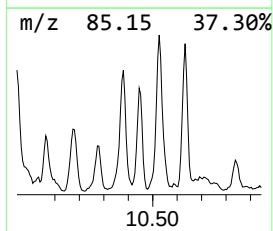
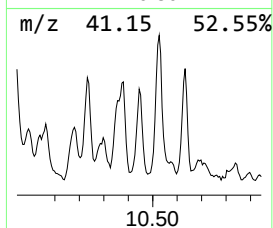
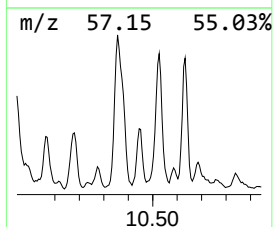
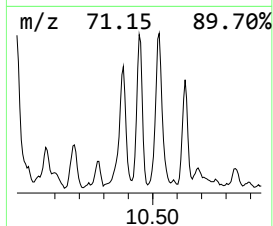
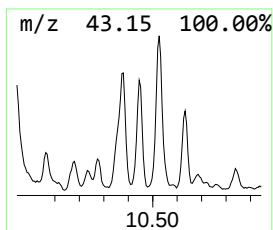
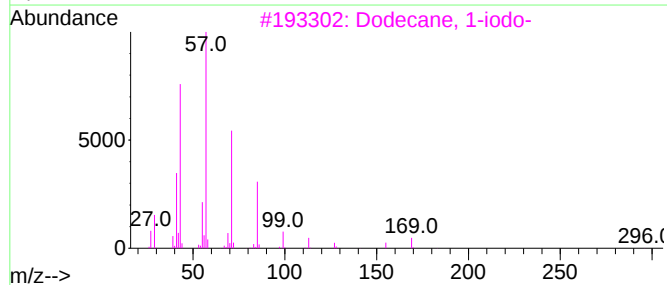
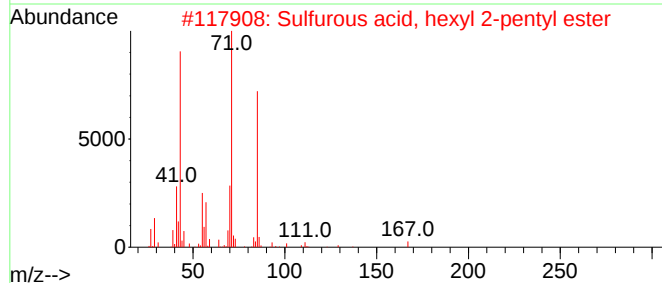
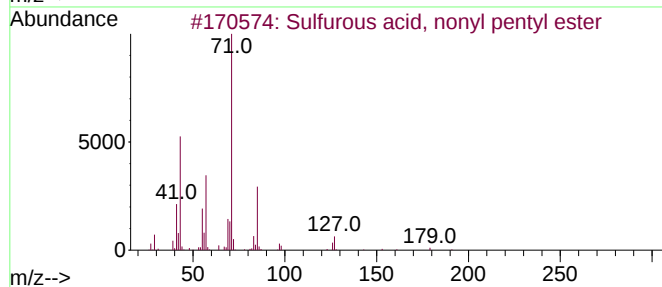
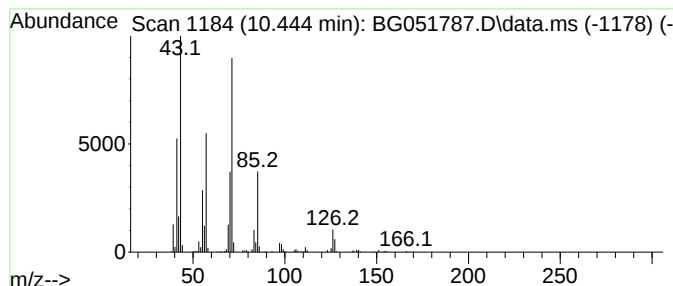
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 38 Sulfurous acid, nonyl penty... Concentration Rank 12

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 10.444 | 16.36 ng/ul | 818893 | Naphthalene-d8   | 11.120 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Sulfurous acid, nonyl pentyl ester  | 278 | C14H30O3S | 1000309-14-2 | 78   |
| 2    |    |   | Sulfurous acid, hexyl 2-pentyl e... | 236 | C11H24O3S | 1000309-15-6 | 59   |
| 3    |    |   | Dodecane, 1-iodo-                   | 296 | C12H25I   | 004292-19-7  | 53   |
| 4    |    |   | Decane, 4-methyl-                   | 156 | C11H24    | 002847-72-5  | 53   |
| 5    |    |   | Heptane, 3,3,5-trimethyl-           | 142 | C10H22    | 007154-80-5  | 50   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

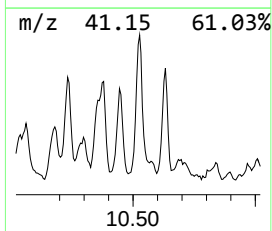
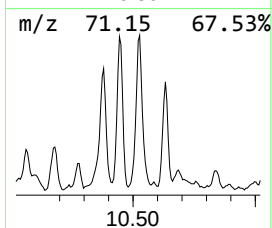
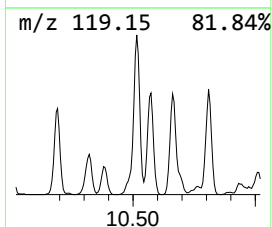
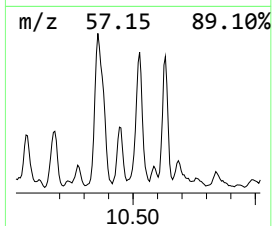
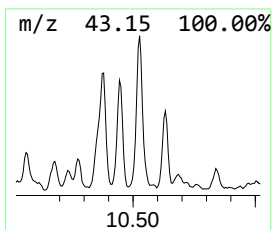
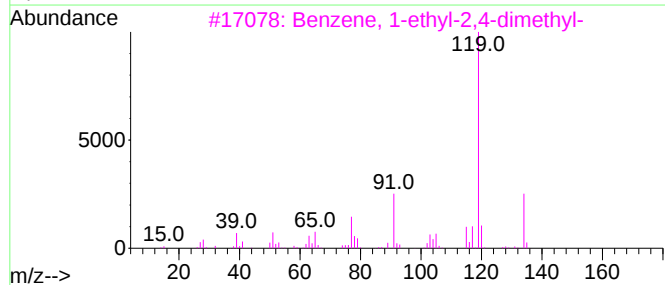
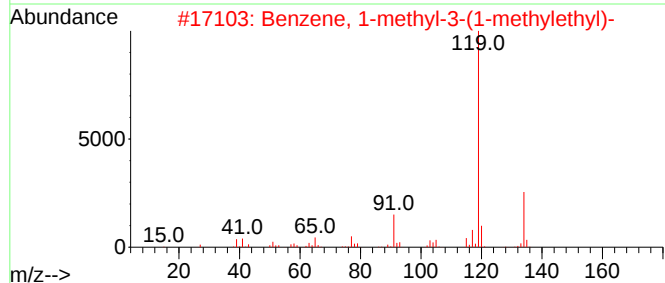
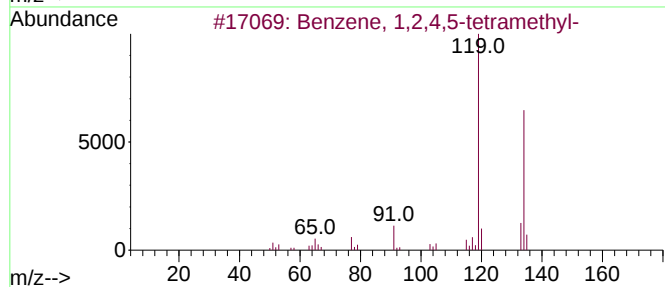
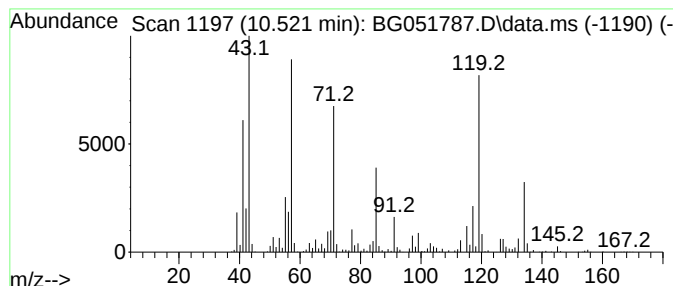
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TIC Integration Parameters: LSCINT.P

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Peak Number 39 unknown-02 Concentration Rank 2

| R.T.   | EstConc     | Area    | Relative to ISTD | R.T.   |
|--------|-------------|---------|------------------|--------|
| 10.521 | 38.46 ng/ul | 1925390 | Naphthalene-d8   | 11.120 |

| Hit# | of | 5 | Tentative ID                         | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1,2,4,5-tetramethyl-        | 134 | C10H14  | 000095-93-2 | 46   |
| 2    |    |   | Benzene, 1-methyl-3-(1-methylethyl)- | 134 | C10H14  | 000535-77-3 | 42   |
| 3    |    |   | Benzene, 1-ethyl-2,4-dimethyl-       | 134 | C10H14  | 000874-41-9 | 42   |
| 4    |    |   | Benzene, 1,2,3,4-tetramethyl-        | 134 | C10H14  | 000488-23-3 | 42   |
| 5    |    |   | o-Cymene                             | 134 | C10H14  | 000527-84-4 | 42   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

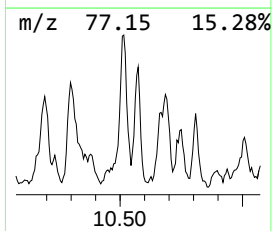
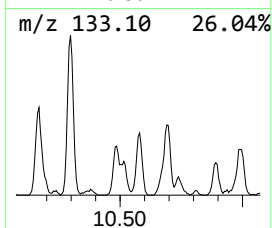
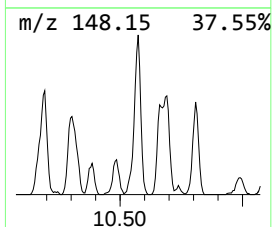
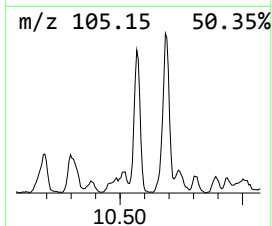
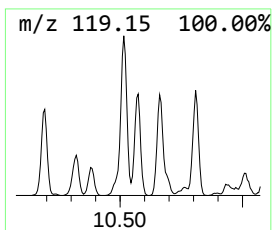
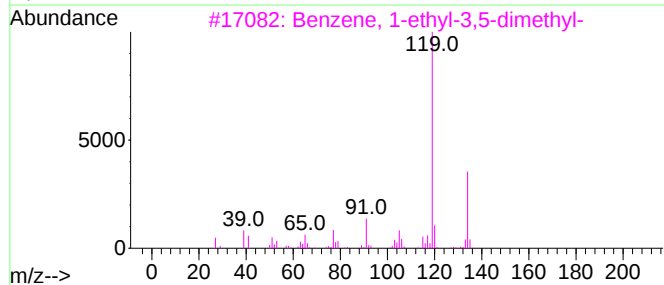
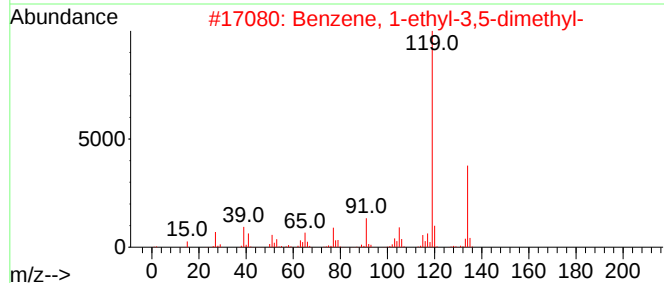
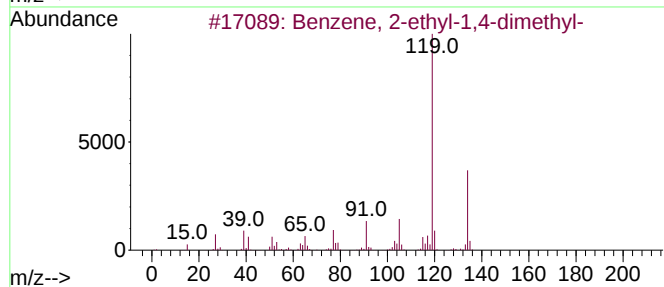
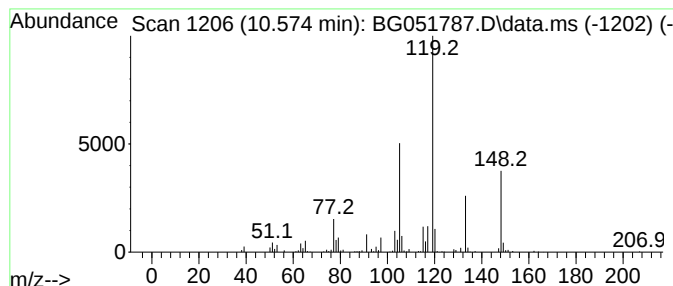
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 40 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 18

| R.T.      | EstConc                        | Area   | Relative to ISTD | R.T.        |      |
|-----------|--------------------------------|--------|------------------|-------------|------|
| 10.574    | 11.17 ng/ul                    | 558973 | Naphthalene-d8   | 11.120      |      |
| Hit# of 5 | Tentative ID                   | MW     | MolForm          | CAS#        | Qual |
| 1         | Benzene, 2-ethyl-1,4-dimethyl- | 134    | C10H14           | 001758-88-9 | 64   |
| 2         | Benzene, 1-ethyl-3,5-dimethyl- | 134    | C10H14           | 000934-74-7 | 59   |
| 3         | Benzene, 1-ethyl-3,5-dimethyl- | 134    | C10H14           | 000934-74-7 | 59   |
| 4         | 4'-Methylpropiofenone          | 148    | C10H12O          | 005337-93-9 | 58   |
| 5         | Benzene, 4-ethyl-1,2-dimethyl- | 134    | C10H14           | 000934-80-5 | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

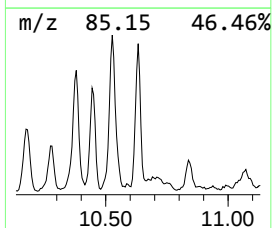
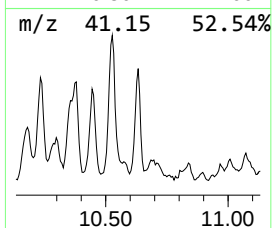
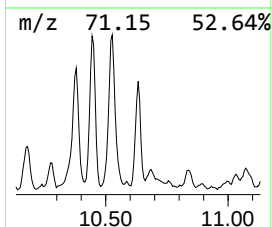
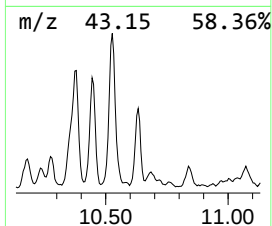
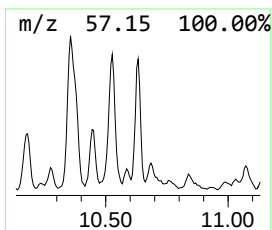
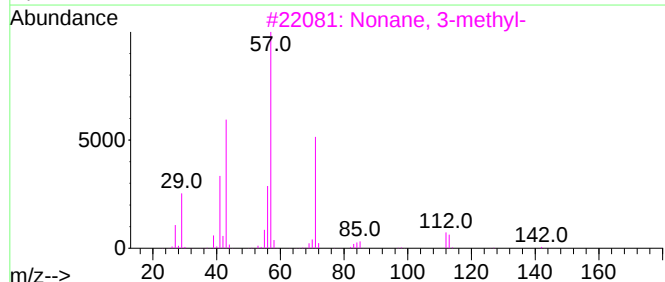
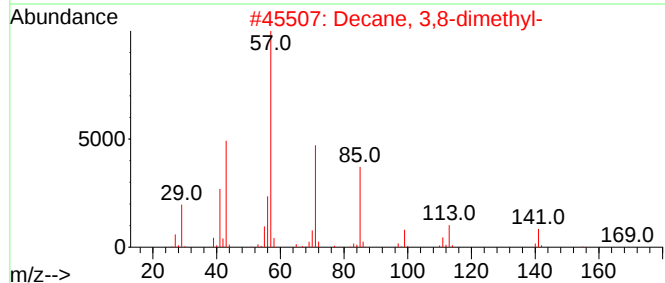
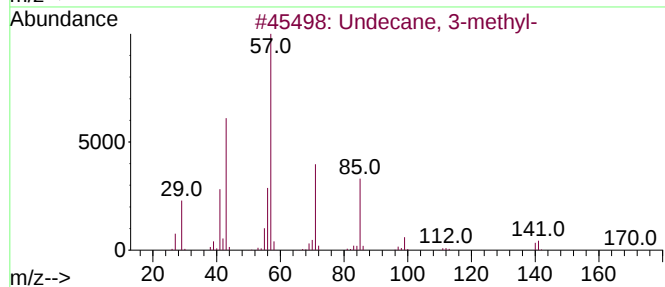
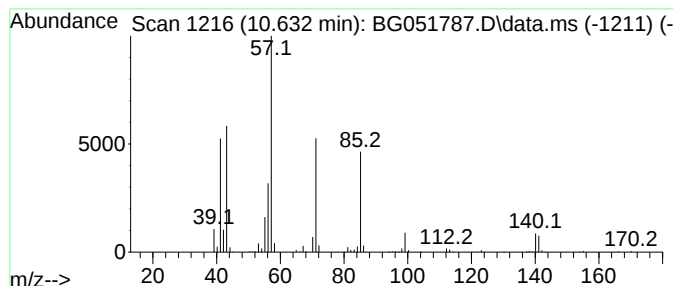
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 41 (DEL) Alkane: Straight-Chai... Concentration Rank 14

| R.T.      | EstConc                   | Area   | Relative to ISTD | R.T.        |      |
|-----------|---------------------------|--------|------------------|-------------|------|
| 10.632    | 14.86 ng/ul               | 743657 | Naphthalene-d8   | 11.120      |      |
| Hit# of 5 | Tentative ID              | MW     | MolForm          | CAS#        | Qual |
| 1         | Undecane, 3-methyl-       | 170    | C12H26           | 001002-43-3 | 80   |
| 2         | Decane, 3,8-dimethyl-     | 170    | C12H26           | 017312-55-9 | 72   |
| 3         | Nonane, 3-methyl-         | 142    | C10H22           | 005911-04-6 | 53   |
| 4         | Octane, 2-methyl-         | 128    | C9H20            | 003221-61-2 | 50   |
| 5         | Heptane, 3,4,5-trimethyl- | 142    | C10H22           | 020278-89-1 | 50   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

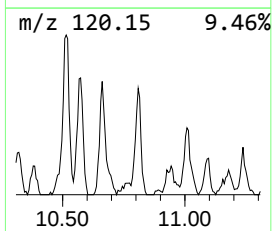
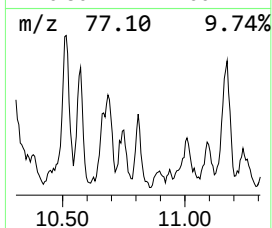
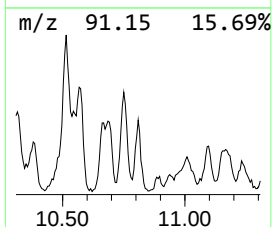
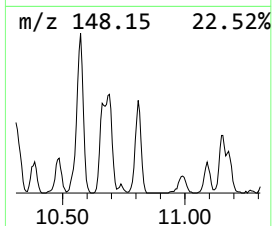
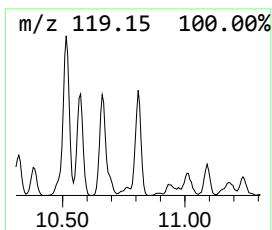
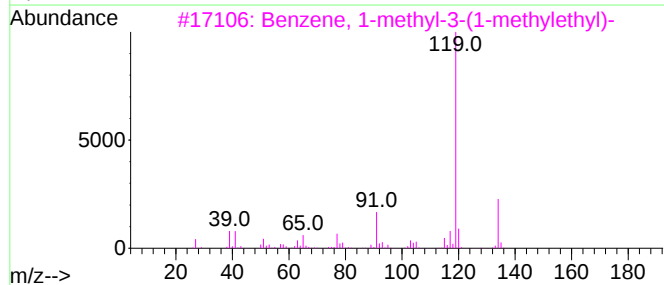
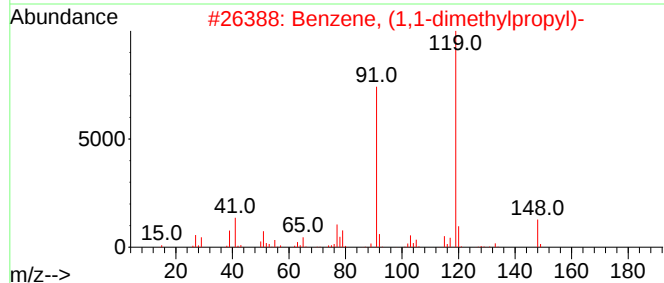
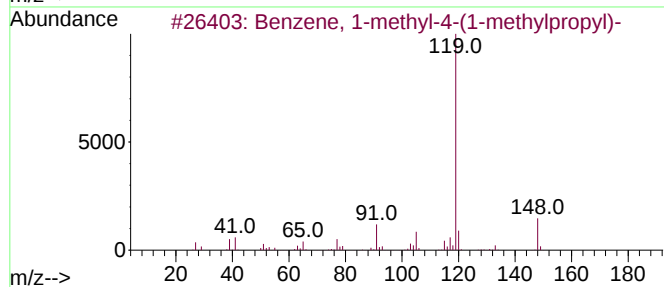
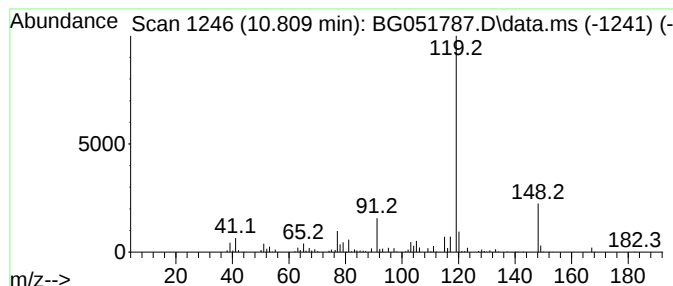
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 42 Benzene, 1-methyl-4-(1-meth... Concentration Rank 21

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|-------------|------|
| 10.809    | 10.09 ng/ul                         | 505187 | Naphthalene-d8   | 11.120      |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual |
| 1         | Benzene, 1-methyl-4-(1-methylpro... | 148    | C11H16           | 001595-16-0 | 90   |
| 2         | Benzene, (1,1-dimethylpropyl)-      | 148    | C11H16           | 002049-95-8 | 86   |
| 3         | Benzene, 1-methyl-3-(1-methyleth... | 134    | C10H14           | 000535-77-3 | 68   |
| 4         | o-Cymene                            | 134    | C10H14           | 000527-84-4 | 64   |
| 5         | Benzene, 1,1'-(1,1,2,2-tetrameth... | 238    | C18H22           | 001889-67-4 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

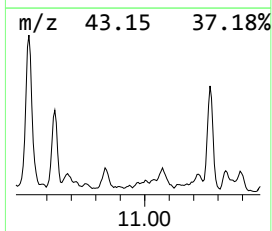
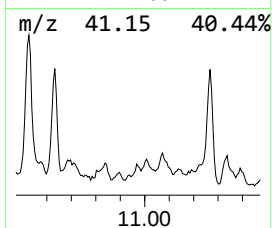
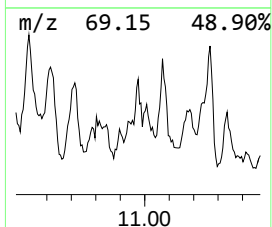
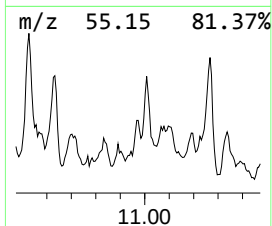
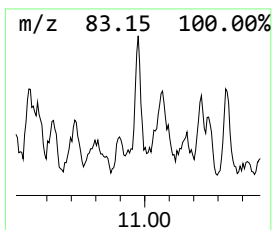
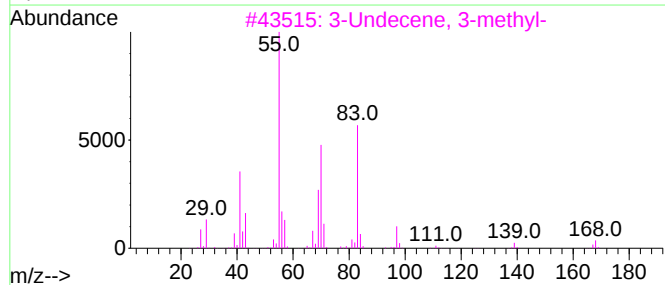
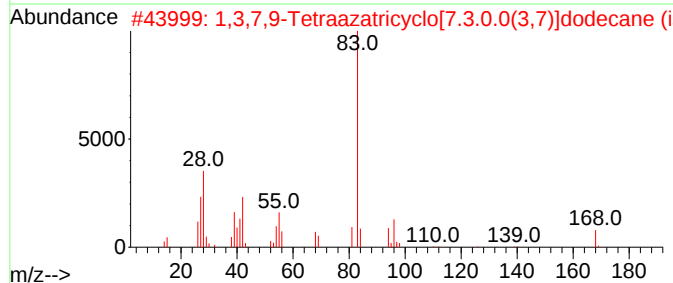
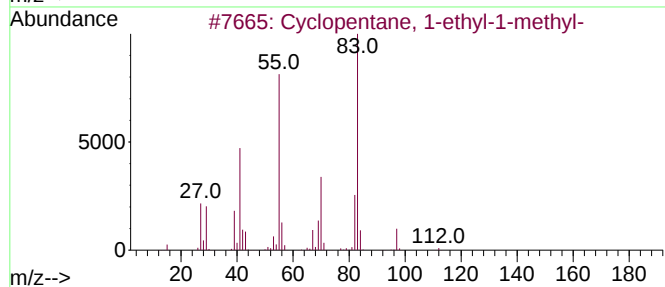
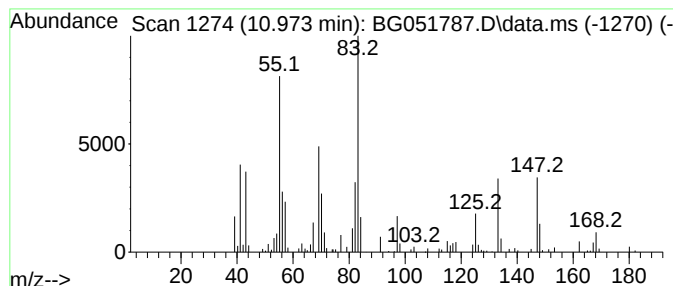
Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 43 (DEL) Alkane: Cyclic10.973 Concentration Rank 48

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 10.973    | 4.22 ng/ul                          | 211378 | Naphthalene-d8   | 11.120       |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | Cyclopentane, 1-ethyl-1-methyl-     | 112    | C8H16            | 016747-50-5  | 52   |
| 2         | 1,3,7,9-Tetraazatricyclo[7.3.0.0... | 168    | C8H16N4          | 1000444-63-7 | 38   |
| 3         | 3-Undecene, 3-methyl-               | 168    | C12H24           | 023381-94-4  | 38   |
| 4         | 2-Hexene, 4,4,5-trimethyl-          | 126    | C9H18            | 055702-61-9  | 35   |
| 5         | trans-4,4-Dimethyl-2-hexene         | 112    | C8H16            | 019550-83-5  | 30   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

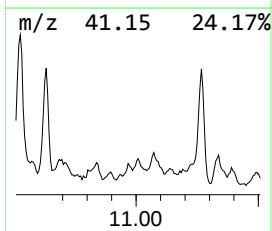
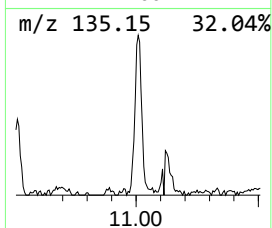
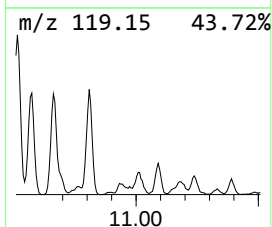
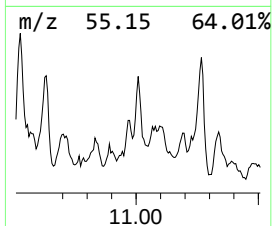
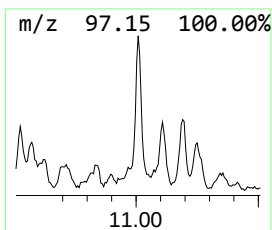
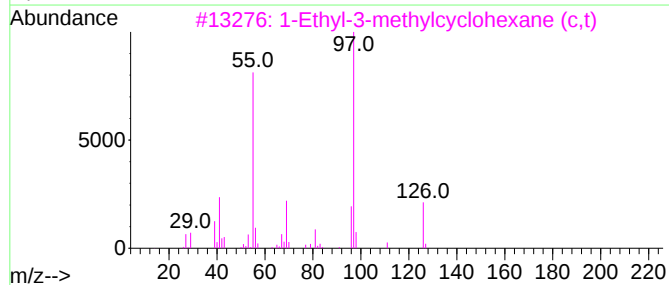
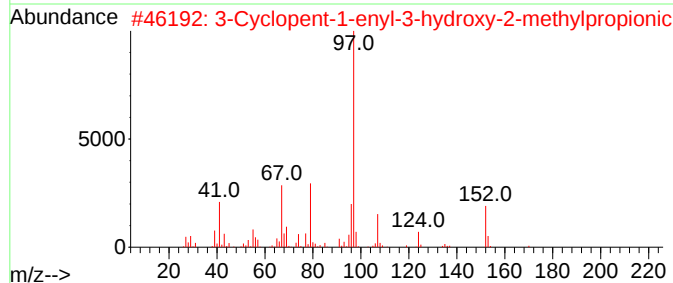
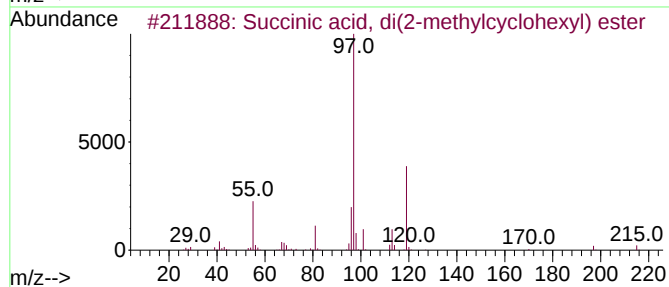
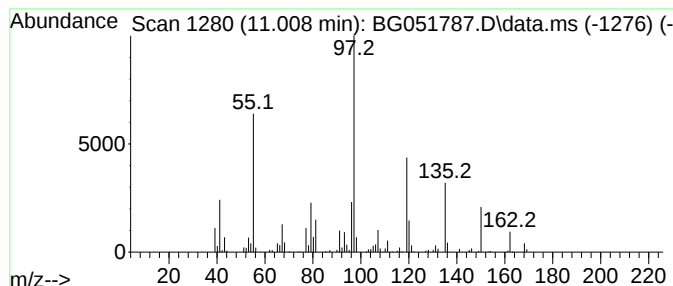
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 44 unknown-03 Concentration Rank 16

| R.T.   | EstConc     | Area   | Relative to ISTD | R.T.   |
|--------|-------------|--------|------------------|--------|
| 11.008 | 11.82 ng/ul | 591465 | Naphthalene-d8   | 11.120 |

| Hit# | of | 5 | Tentative ID                                        | MW  | MolForm  | CAS#         | Qual |
|------|----|---|-----------------------------------------------------|-----|----------|--------------|------|
| 1    |    |   | Succinic acid, di(2-methylcyclohexyl) ester         | 310 | C18H30O4 | 1000329-83-0 | 38   |
| 2    |    |   | 3-Cyclopent-1-enyl-3-hydroxy-2-methylpropionic acid | 170 | C9H14O3  | 1000191-19-7 | 38   |
| 3    |    |   | 1-Ethyl-3-methylcyclohexane (c,t)                   | 126 | C9H18    | 003728-55-0  | 27   |
| 4    |    |   | Oxalic acid, cyclohexylmethyl isomer                | 270 | C15H26O4 | 1000309-68-3 | 27   |
| 5    |    |   | Cyclopentane, 1,1'-ethylidenebis-                   | 166 | C12H22   | 004413-21-2  | 27   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

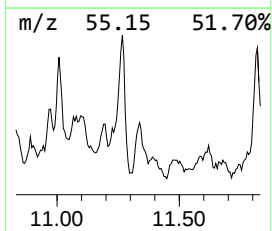
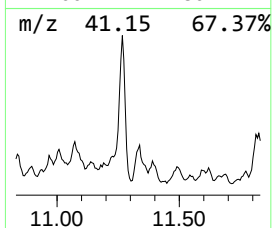
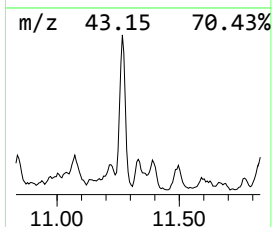
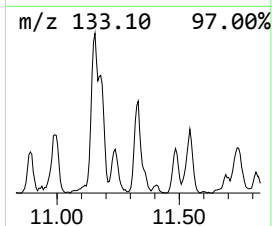
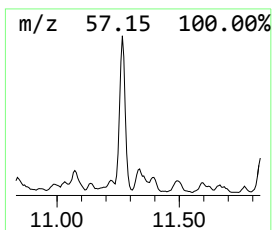
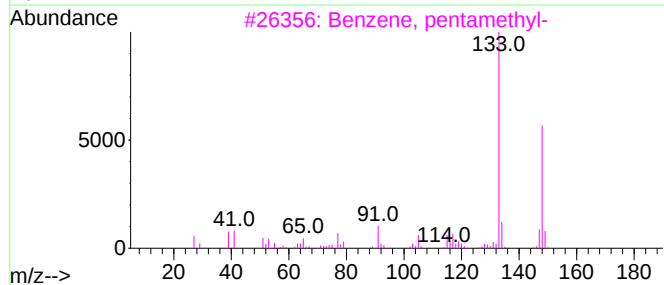
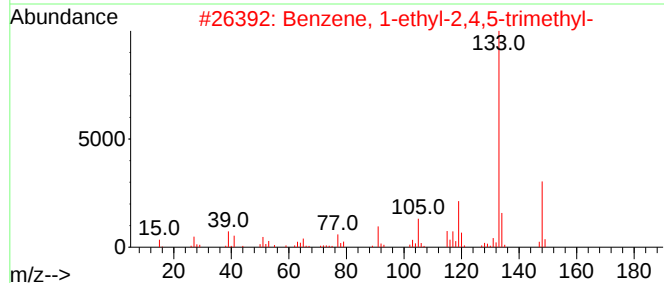
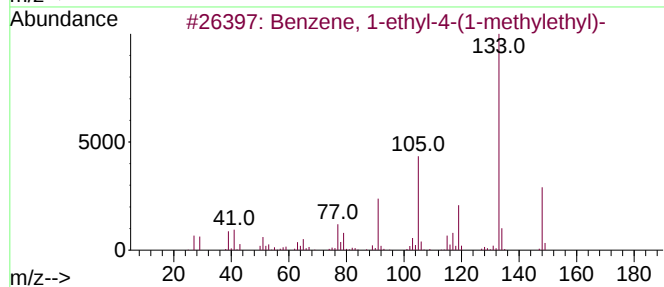
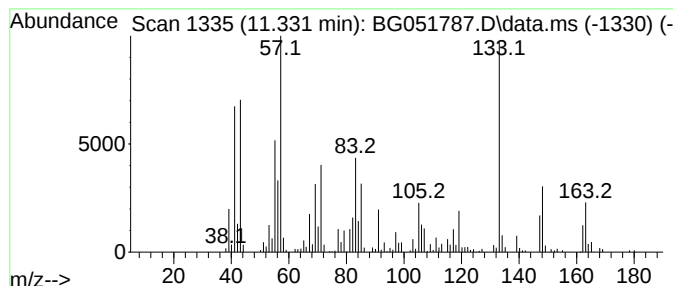
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 45 unknown-04 Concentration Rank 30

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 11.331 | 8.21 ng/ul | 410757 | Naphthalene-d8   | 11.120 |

| Hit# | of | 5 | Tentative ID                         | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Benzene, 1-ethyl-4-(1-methylethyl)-  | 148 | C11H16  | 004218-48-8 | 46   |
| 2    |    |   | Benzene, 1-ethyl-2,4,5-trimethyl-    | 148 | C11H16  | 017851-27-3 | 45   |
| 3    |    |   | Benzene, pentamethyl-                | 148 | C11H16  | 000700-12-9 | 42   |
| 4    |    |   | 1,5,6,7-Tetramethylbicyclo[3.2.0]... | 148 | C11H16  | 134329-46-7 | 38   |
| 5    |    |   | Benzene, (1-ethyl-1-methylpropyl)-   | 162 | C12H18  | 001985-97-3 | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

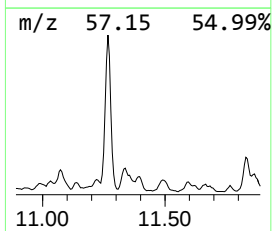
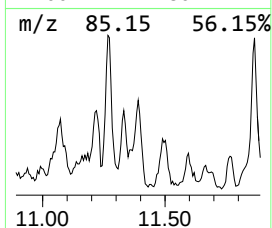
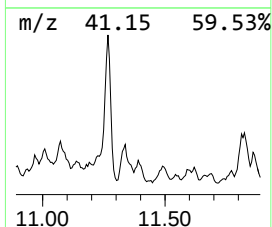
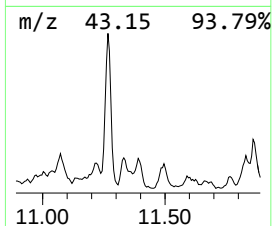
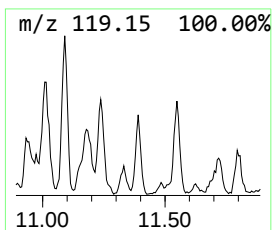
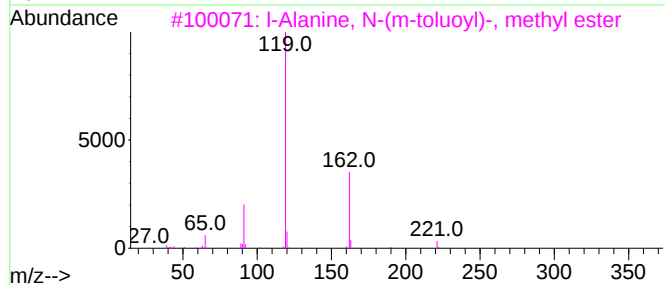
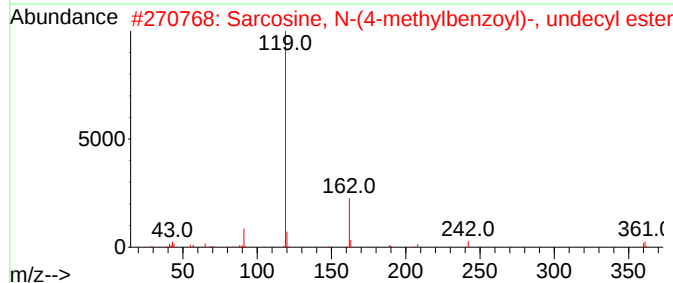
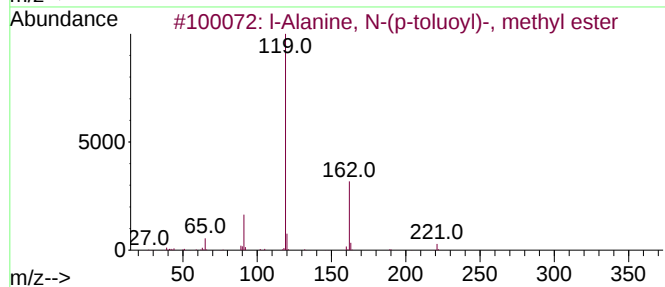
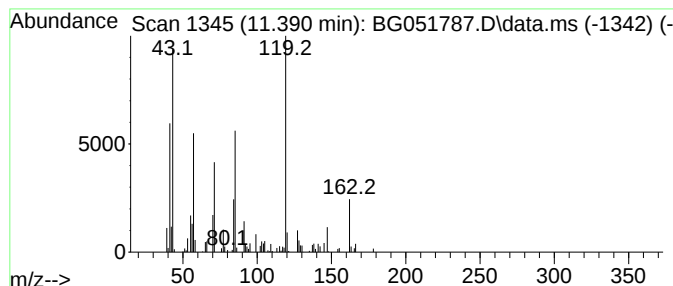
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 46 unknown-05 Concentration Rank 41

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 11.390 | 5.45 ng/ul | 273038 | Naphthalene-d8   | 11.120 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | l-Alanine, N-(p-toluoyl)-, methy... | 221 | C12H15NO3 | 1000299-64-6 | 35   |
| 2    |    |   | Sarcosine, N-(4-methylbenzoyl)-,... | 361 | C22H35NO3 | 1000321-22-3 | 35   |
| 3    |    |   | l-Alanine, N-(m-toluoyl)-, methy... | 221 | C12H15NO3 | 1010299-63-7 | 35   |
| 4    |    |   | Benzene, 1,4-dimethyl-2-(2-methy... | 162 | C12H18    | 055669-88-0  | 27   |
| 5    |    |   | Dodecane, 5-methyl-                 | 184 | C13H28    | 017453-93-9  | 27   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

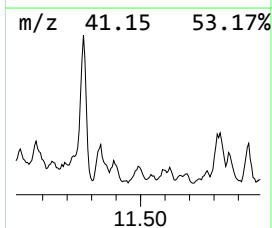
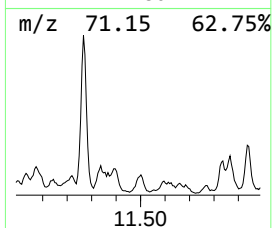
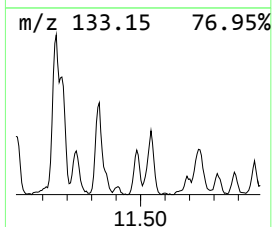
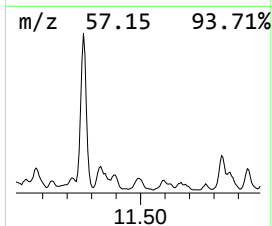
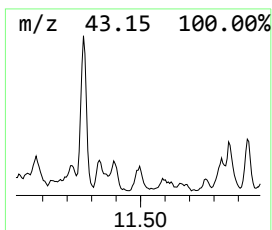
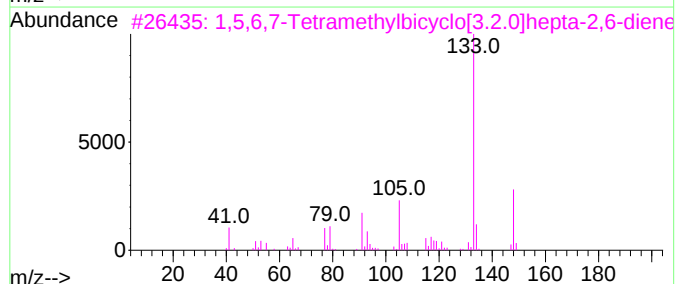
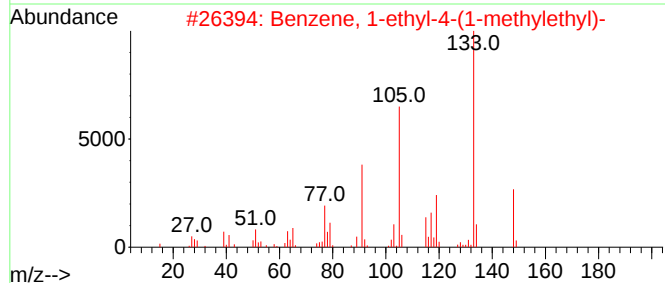
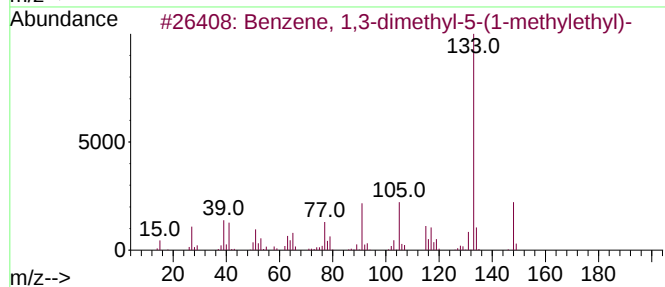
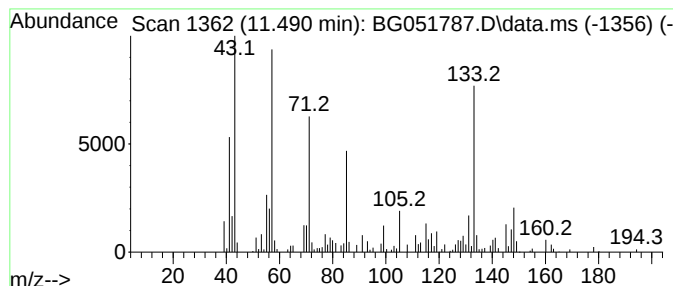
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 47 Benzene, 1,3-dimethyl-5-(1-... Concentration Rank 46

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|-------------|------|
| 11.490    | 4.38 ng/ul                          | 219344 | Naphthalene-d8   | 11.120      |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual |
| 1         | Benzene, 1,3-dimethyl-5-(1-methy... | 148    | C11H16           | 004706-90-5 | 86   |
| 2         | Benzene, 1-ethyl-4-(1-methylethyl)- | 148    | C11H16           | 004218-48-8 | 45   |
| 3         | 1,5,6,7-Tetramethylbicyclo[3.2.0... | 148    | C11H16           | 134329-46-7 | 43   |
| 4         | Ethanone, 1-(4-ethylphenyl)-        | 148    | C10H12O          | 000937-30-4 | 41   |
| 5         | Benzene, 1,4-diethyl-2-methyl-      | 148    | C11H16           | 013632-94-5 | 41   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

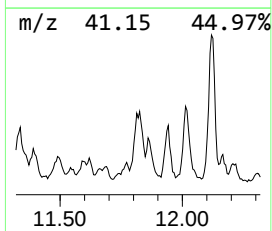
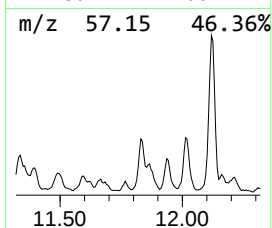
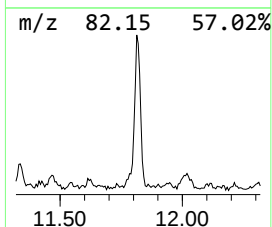
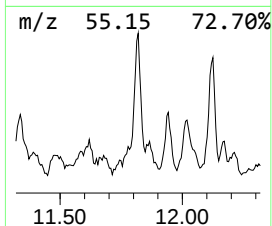
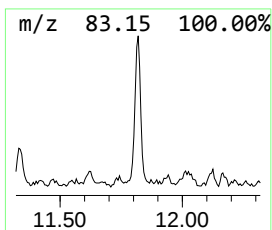
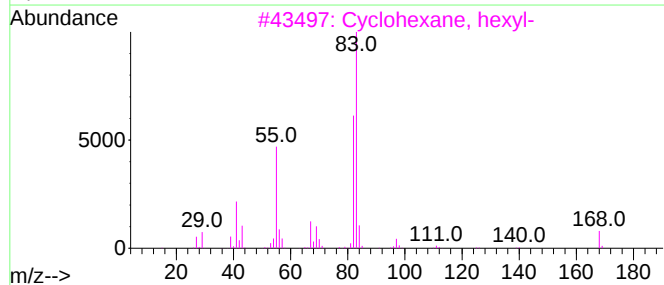
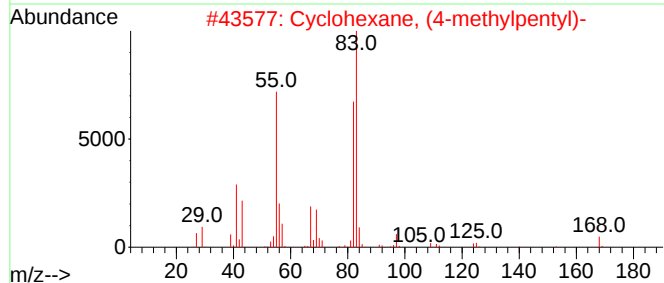
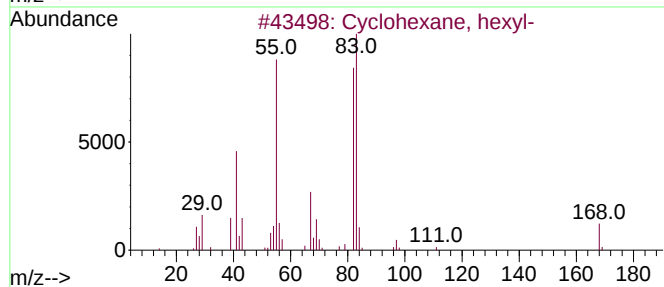
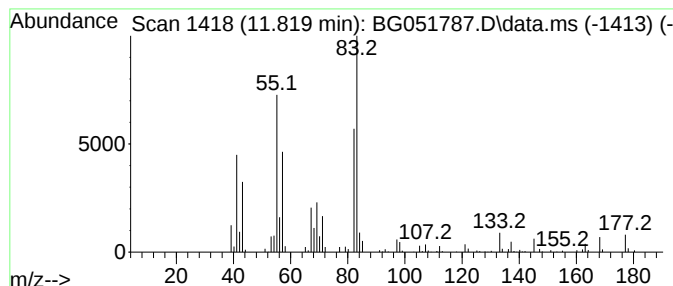
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TIC Integration Parameters: LSCINT.P

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Peak Number 48 (DEL) Alkane: Cyclic11.819 Concentration Rank 29

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 11.819 | 8.40 ng/ul | 420658 | Naphthalene-d8   | 11.120 |

| Hit# | of 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|------|--------------------------------|-----|---------|-------------|------|
| 1    |      | Cyclohexane, hexyl-            | 168 | C12H24  | 004292-75-5 | 83   |
| 2    |      | Cyclohexane, (4-methylpentyl)- | 168 | C12H24  | 061142-20-9 | 80   |
| 3    |      | Cyclohexane, hexyl-            | 168 | C12H24  | 004292-75-5 | 72   |
| 4    |      | Cyclohexane, butyl-            | 140 | C10H20  | 001678-93-9 | 64   |
| 5    |      | Cyclohexane, (1-methylethyl)-  | 126 | C9H18   | 000696-29-7 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

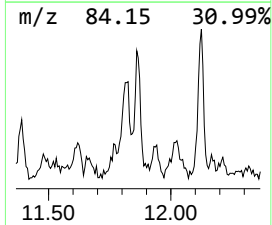
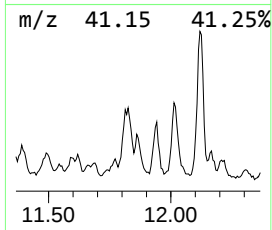
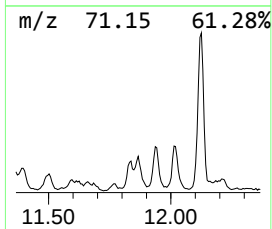
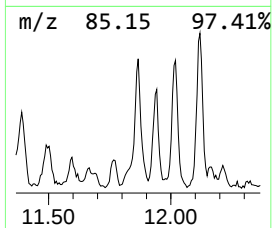
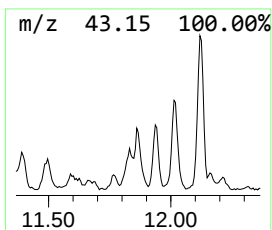
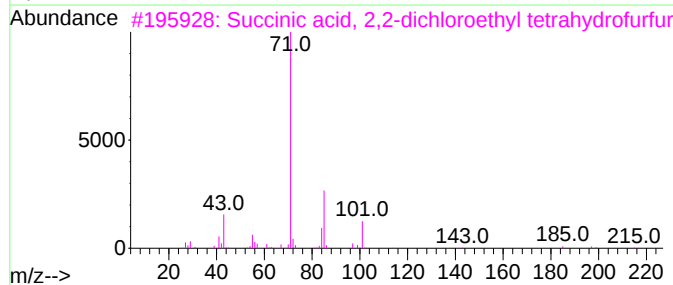
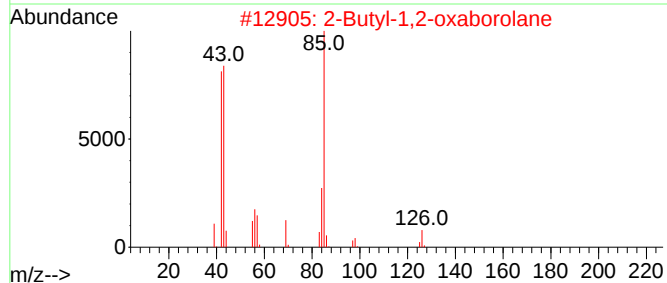
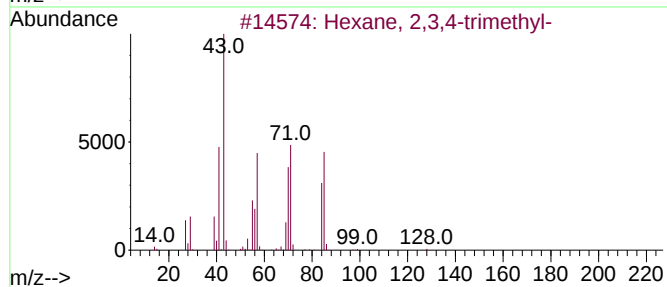
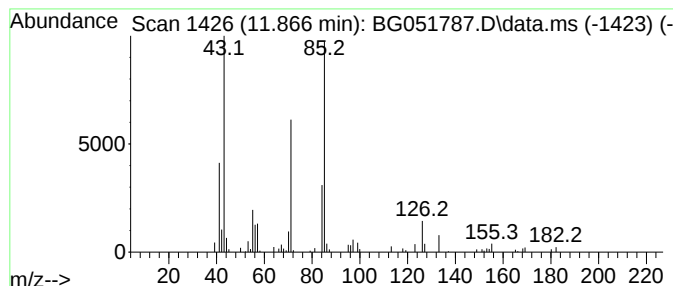
Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 49 (DEL) Alkane: Straight-Chai... Concentration Rank 47

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 11.866    | 4.38 ng/ul                          | 219057 | Naphthalene-d8   | 11.120       |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | Hexane, 2,3,4-trimethyl-            | 128    | C9H20            | 000921-47-1  | 50   |
| 2         | 2-Butyl-1,2-oxaborolane             | 126    | C7H15BO          | 005357-13-1  | 47   |
| 3         | Succinic acid, 2,2-dichloroethyl... | 298    | C11H16Cl2O5      | 1000390-71-6 | 47   |
| 4         | 1,3,2-Oxazaborolane, 2-butyl-       | 127    | C6H14BNO         | 031748-10-4  | 47   |
| 5         | Sulfurous acid, hexyl nonyl ester   | 292    | C15H32O3S        | 1000309-13-1 | 47   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

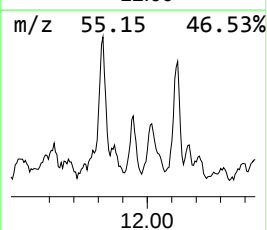
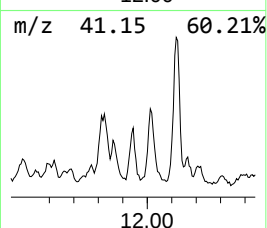
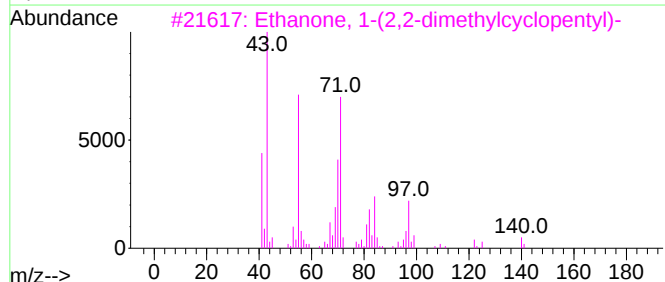
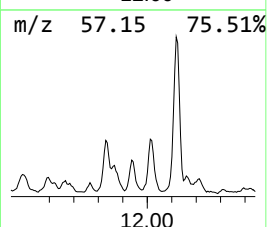
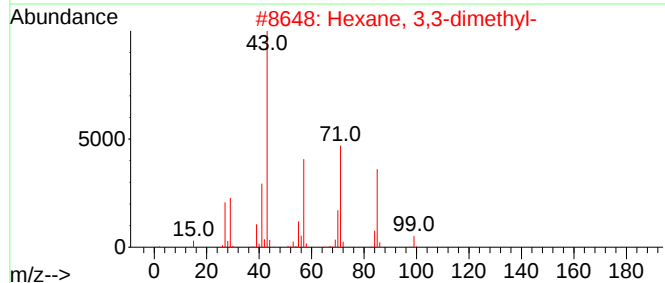
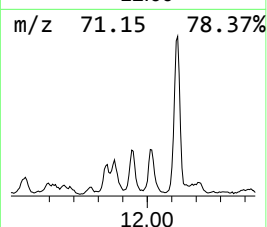
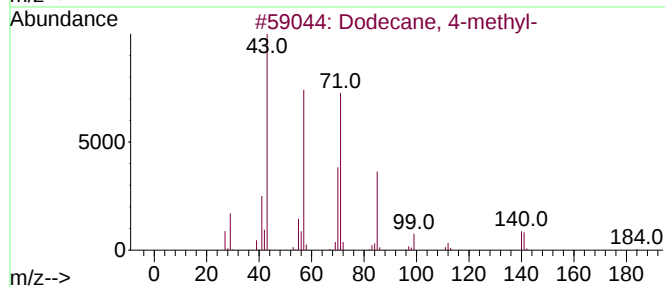
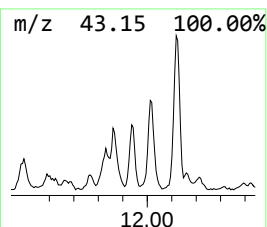
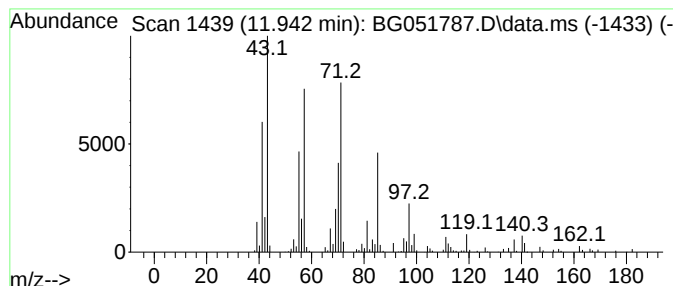
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 50 (DEL) Alkane: Straight-Chai... Concentration Rank 36

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|-------------|------|
| 11.942    | 6.44 ng/ul                          | 322564 | Naphthalene-d8   | 11.120      |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual |
| 1         | Dodecane, 4-methyl-                 | 184    | C13H28           | 006117-97-1 | 72   |
| 2         | Hexane, 3,3-dimethyl-               | 114    | C8H18            | 000563-16-6 | 64   |
| 3         | Ethanone, 1-(2,2-dimethylcyclope... | 140    | C9H16O           | 003664-75-3 | 64   |
| 4         | Dodecane, 4-methyl-                 | 184    | C13H28           | 006117-97-1 | 59   |
| 5         | Undecane, 4-methyl-                 | 170    | C12H26           | 002980-69-0 | 53   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

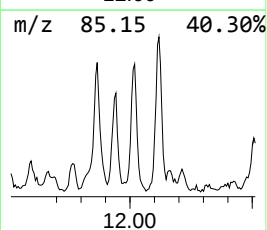
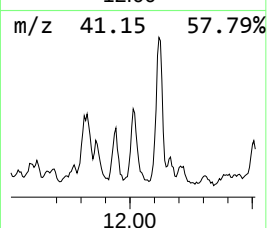
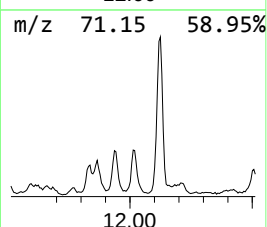
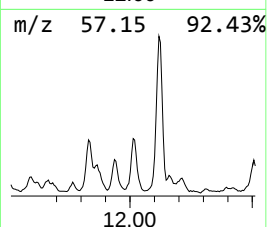
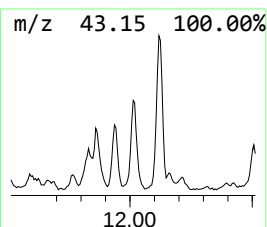
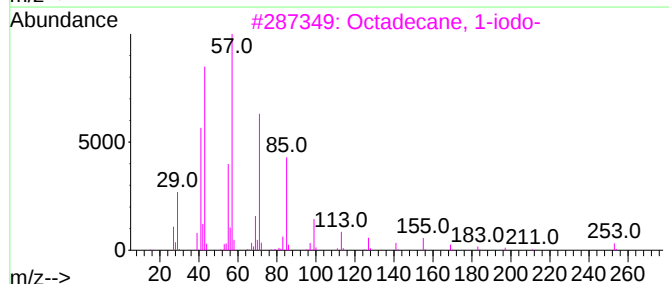
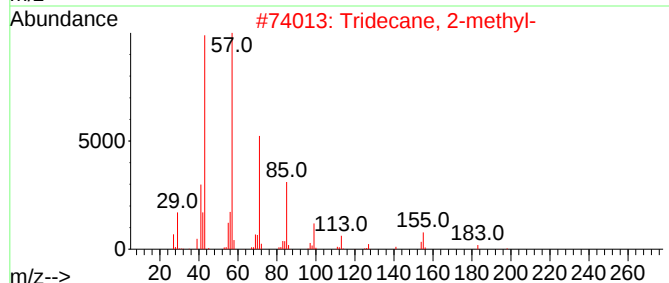
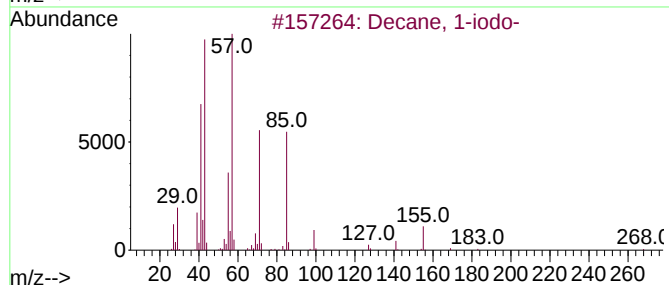
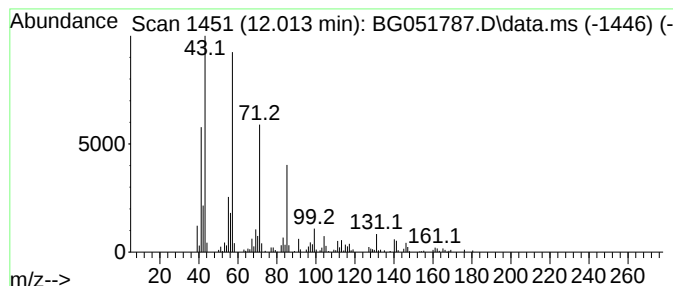
Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 51 Decane, 1-iodo- Concentration Rank 24

| R.T.   | EstConc    | Area                  | Relative to ISTD | R.T.            |
|--------|------------|-----------------------|------------------|-----------------|
| 12.013 | 9.51 ng/ul | 475949                | Naphthalene-d8   | 11.120          |
| Hit#   | of 5       | Tentative ID          | MW MolForm       | CAS# Qual       |
| 1      |            | Decane, 1-iodo-       | 268 C10H21I      | 002050-77-3 72  |
| 2      |            | Tridecane, 2-methyl-  | 198 C14H30       | 001560-96-9 64  |
| 3      |            | Octadecane, 1-iodo-   | 380 C18H37I      | 000629-93-6 64  |
| 4      |            | Decane, 3,8-dimethyl- | 170 C12H26       | 017312-55-9 64  |
| 5      |            | 1-Iodo-2-methylnonane | 268 C10H21I      | 1000101-47-9 59 |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

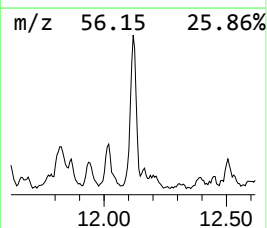
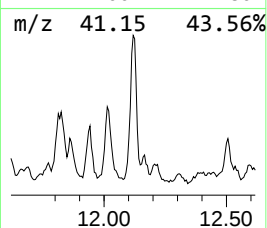
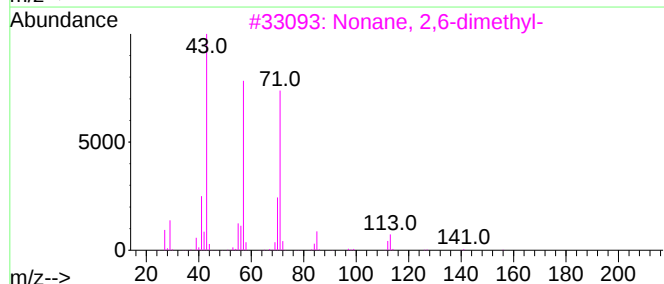
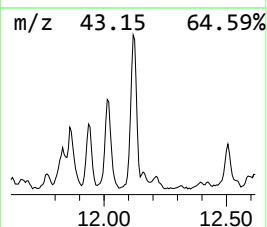
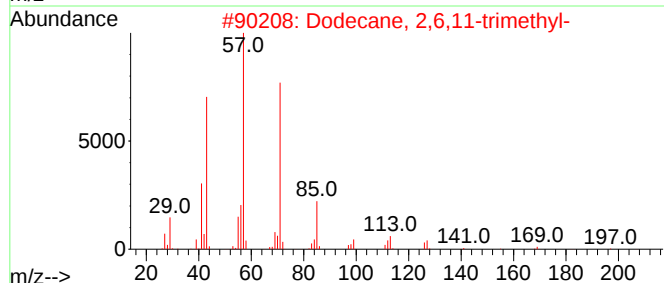
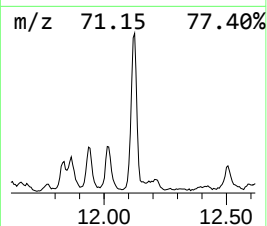
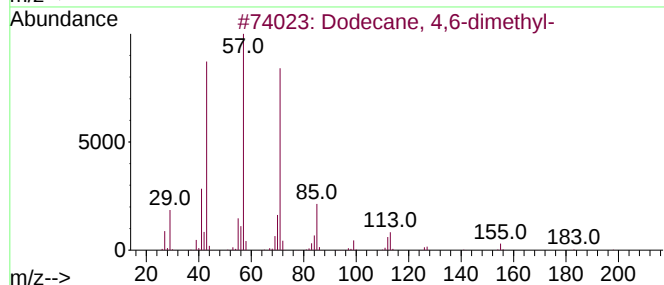
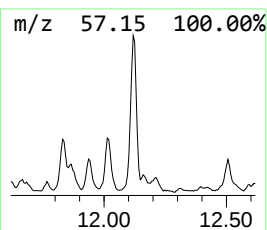
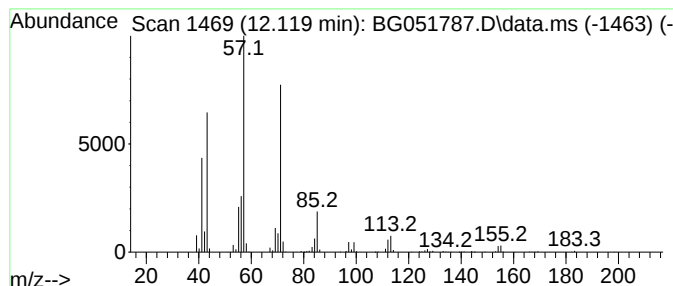
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 52 (DEL) Alkane: Straight-Chai... Concentration Rank 11

| R.T.      | EstConc                     | Area   | Relative to ISTD | R.T.        |      |
|-----------|-----------------------------|--------|------------------|-------------|------|
| 12.119    | 17.03 ng/ul                 | 852593 | Naphthalene-d8   | 11.120      |      |
| Hit# of 5 | Tentative ID                | MW     | MolForm          | CAS#        | Qual |
| 1         | Dodecane, 4,6-dimethyl-     | 198    | C14H30           | 061141-72-8 | 83   |
| 2         | Dodecane, 2,6,11-trimethyl- | 212    | C15H32           | 031295-56-4 | 83   |
| 3         | Nonane, 2,6-dimethyl-       | 156    | C11H24           | 017302-28-2 | 80   |
| 4         | Dodecane, 2,6,11-trimethyl- | 212    | C15H32           | 031295-56-4 | 78   |
| 5         | Octane, 2,3,7-trimethyl-    | 156    | C11H24           | 062016-34-6 | 72   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

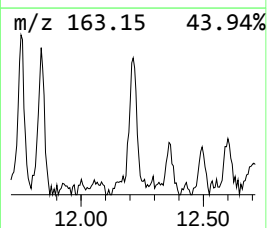
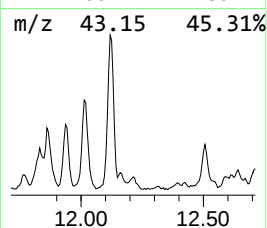
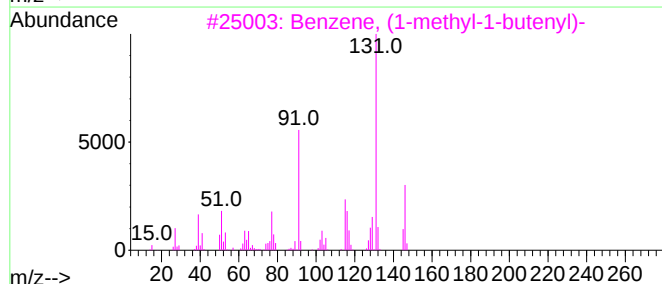
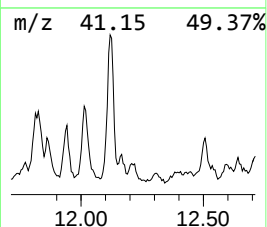
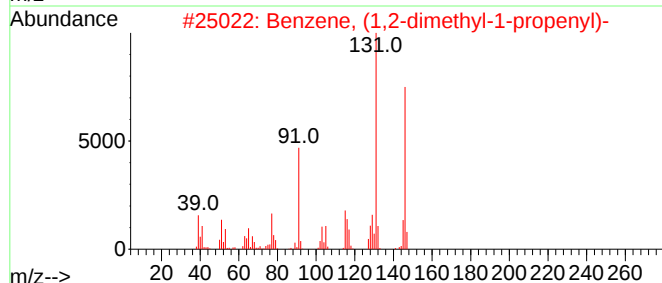
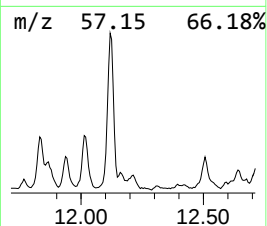
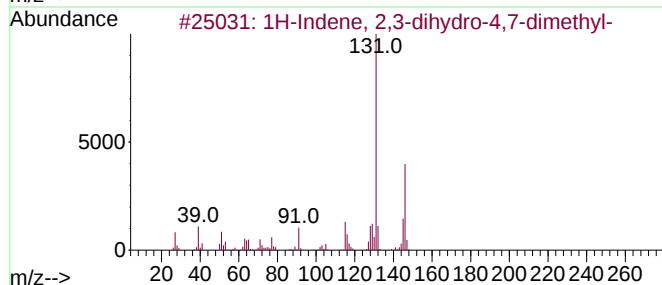
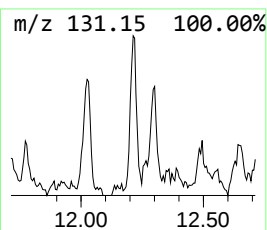
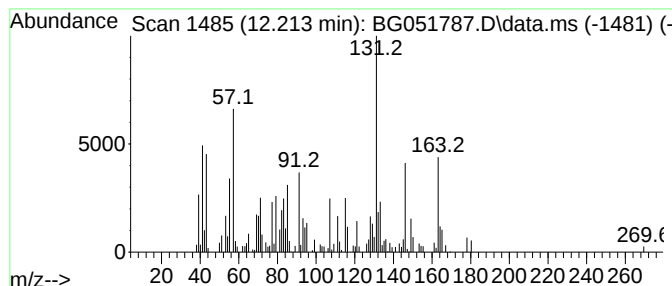
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 53 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 51

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 12.213 | 3.45 ng/ul | 172931 | Naphthalene-d8   | 11.120 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 | C11H14  | 006682-71-9 | 50   |
| 2    |    |   | Benzene, (1,2-dimethyl-1-propenyl)- | 146 | C11H14  | 000769-57-3 | 46   |
| 3    |    |   | Benzene, (1-methyl-1-butenyl)-      | 146 | C11H14  | 053172-84-2 | 45   |
| 4    |    |   | Benzene, (3-methyl-2-butenyl)-      | 146 | C11H14  | 004489-84-3 | 43   |
| 5    |    |   | 1H-Indene, 2,3-dihydro-4,7-dimet... | 146 | C11H14  | 006682-71-9 | 43   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

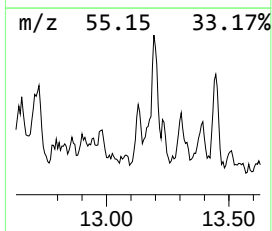
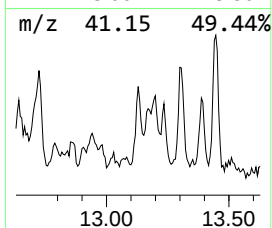
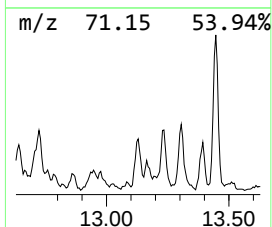
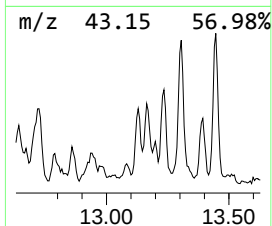
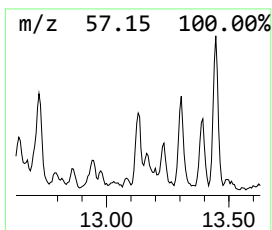
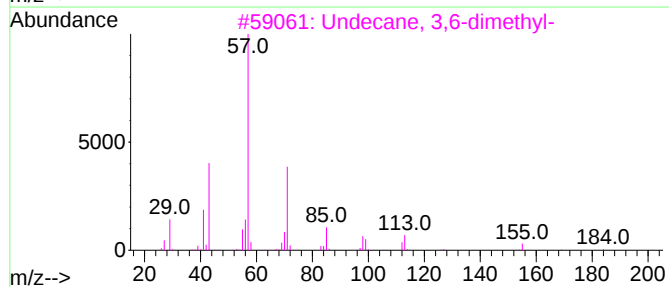
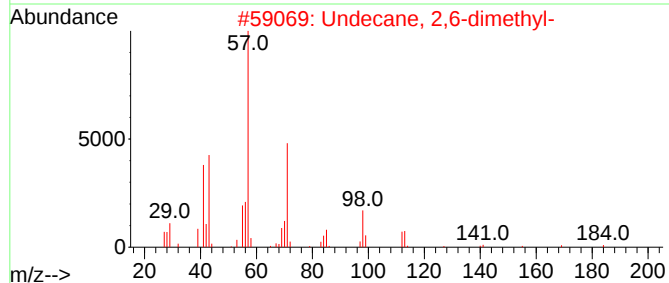
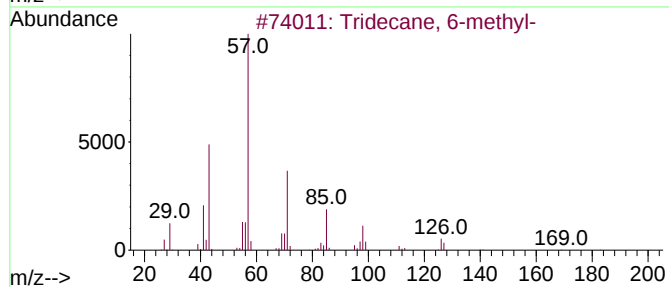
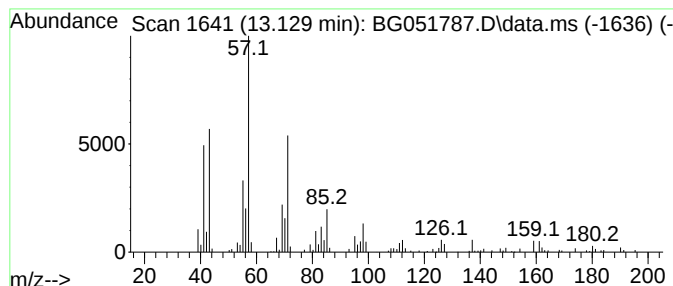
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 54 (DEL) Alkane: Straight-Chai... Concentration Rank 25

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 13.129 | 9.36 ng/ul | 194050 | Acenaphthene-d10 | 14.909 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Tridecane, 6-methyl-                | 198 | C14H30  | 013287-21-3 | 72   |
| 2    |    |   | Undecane, 2,6-dimethyl-             | 184 | C13H28  | 017301-23-4 | 70   |
| 3    |    |   | Undecane, 3,6-dimethyl-             | 184 | C13H28  | 017301-28-9 | 68   |
| 4    |    |   | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44  | 054833-48-6 | 59   |
| 5    |    |   | Decane, 2,3,5-trimethyl-            | 184 | C13H28  | 062238-11-3 | 50   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

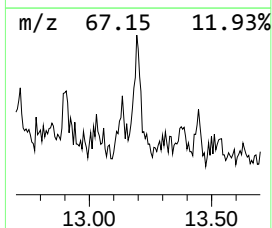
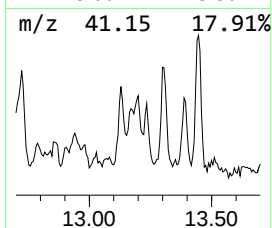
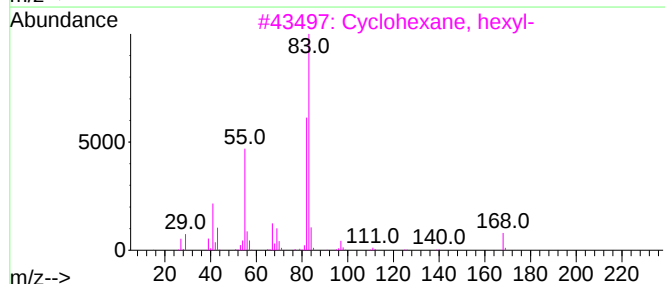
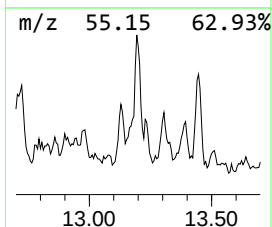
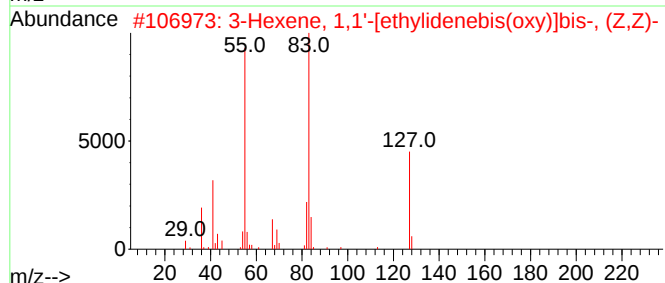
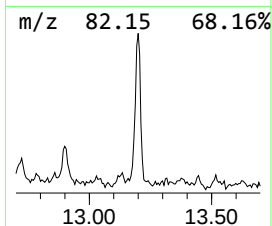
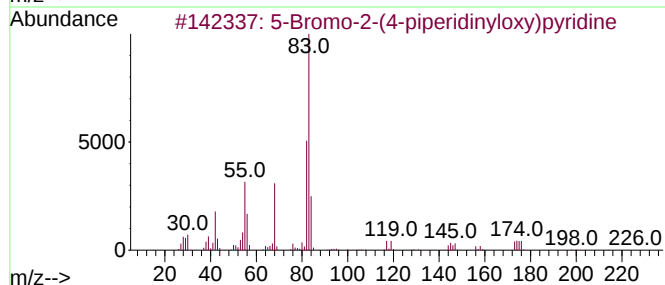
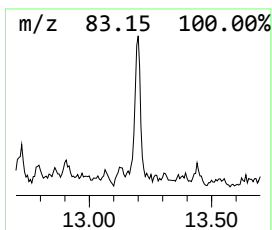
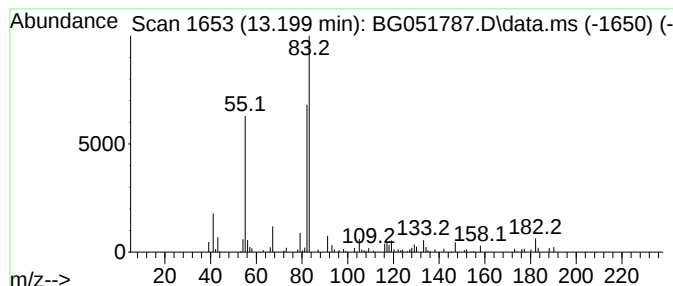
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 55 5-Bromo-2-(4-piperidinyloxy... Concentration Rank 27

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 13.200 | 8.56 ng/ul | 177370 | Acenaphthene-d10 | 14.909 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm     | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|-------------|-------------|------|
| 1    |    |   | 5-Bromo-2-(4-piperidinyloxy)pyri... | 256 | C10H13BrN2O | 194668-50-3 | 56   |
| 2    |    |   | 3-Hexene, 1,1'-[ethylidenebis(ox... | 226 | C14H26O2    | 063449-64-9 | 47   |
| 3    |    |   | Cyclohexane, hexyl-                 | 168 | C12H24      | 004292-75-5 | 45   |
| 4    |    |   | Cyclohexane, pentyl-                | 154 | C11H22      | 004292-92-6 | 45   |
| 5    |    |   | Cyclohexane, nonadecyl-             | 350 | C25H50      | 022349-03-7 | 45   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

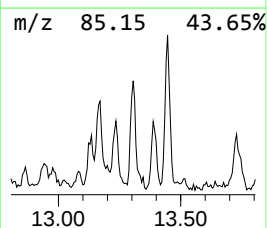
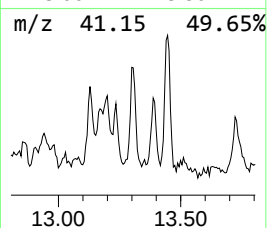
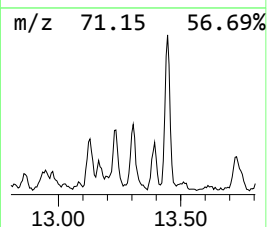
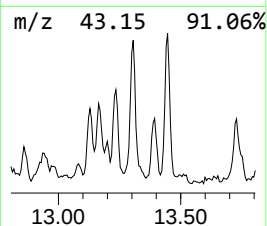
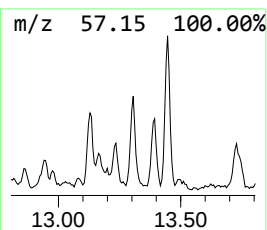
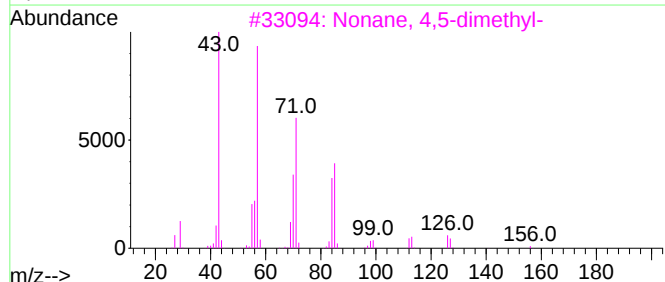
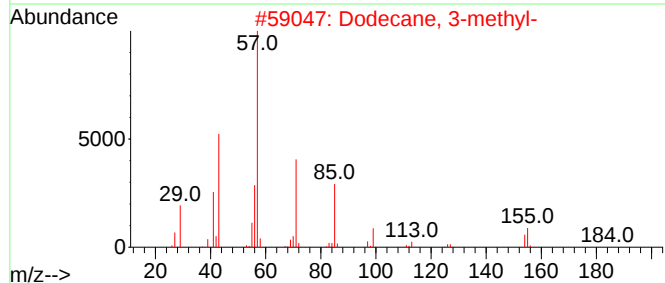
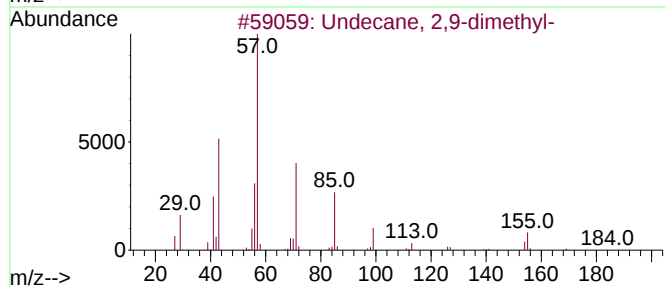
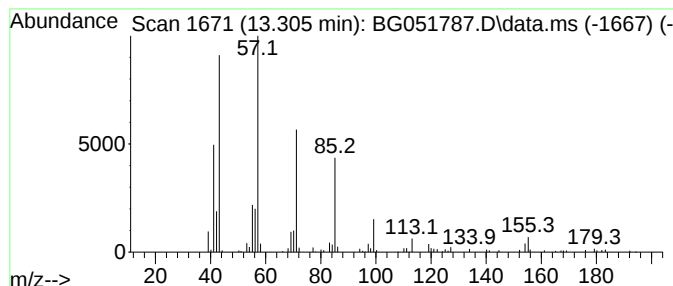
Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 56 (DEL) Alkane: Straight-Chai... Concentration Rank 20

| R.T.      | EstConc                      | Area   | Relative to ISTD | R.T.        |      |
|-----------|------------------------------|--------|------------------|-------------|------|
| 13.305    | 10.22 ng/ul                  | 211732 | Acenaphthene-d10 | 14.909      |      |
| Hit# of 5 | Tentative ID                 | MW     | MolForm          | CAS#        | Qual |
| 1         | Undecane, 2,9-dimethyl-      | 184    | C13H28           | 017301-26-7 | 90   |
| 2         | Dodecane, 3-methyl-          | 184    | C13H28           | 017312-57-1 | 90   |
| 3         | Nonane, 4,5-dimethyl-        | 156    | C11H24           | 017302-23-7 | 80   |
| 4         | Dodecane, 2-methyl-6-propyl- | 226    | C16H34           | 055045-08-4 | 72   |
| 5         | Tridecane                    | 184    | C13H28           | 000629-50-5 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

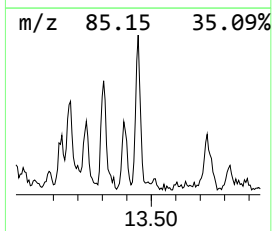
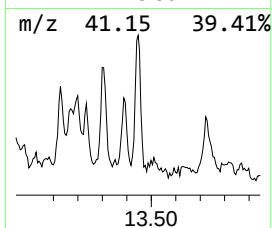
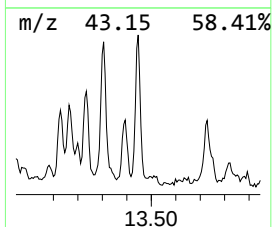
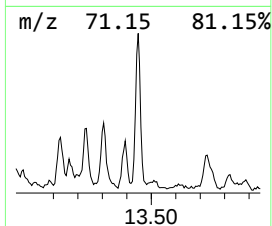
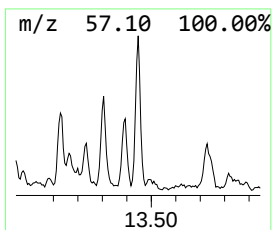
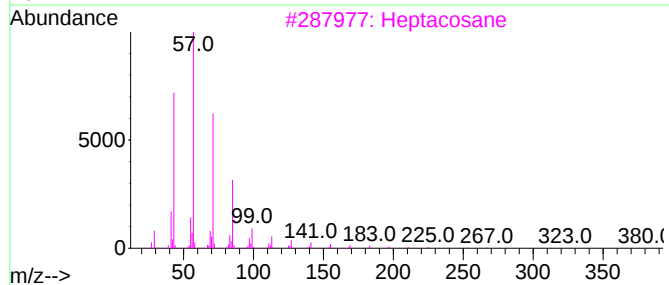
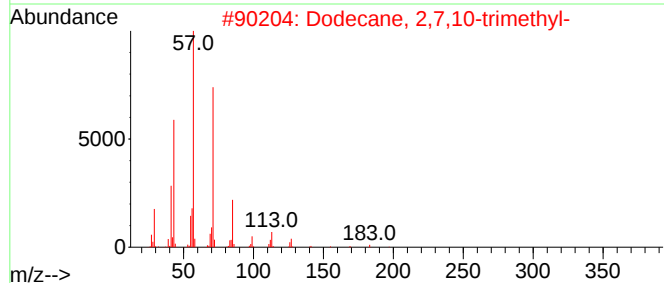
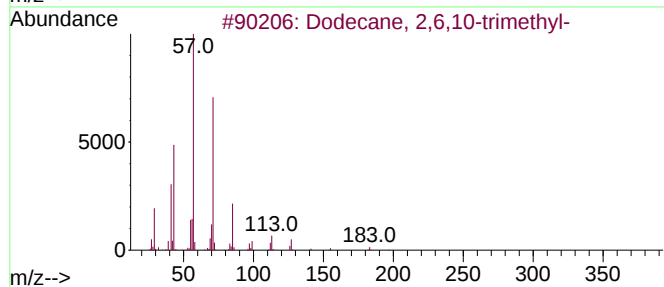
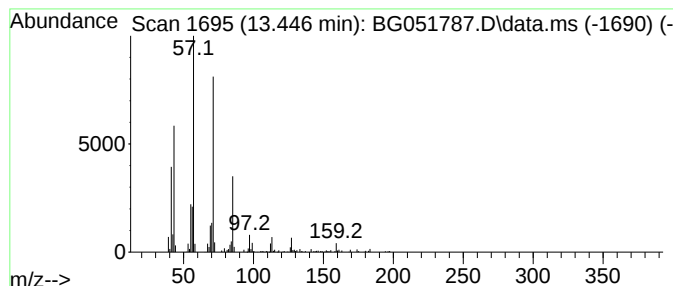
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 57 (DEL) Alkane: Straight-Chai... Concentration Rank 13

| R.T.      | EstConc                             | Area   | Relative to ISTD |          |              | R.T.   |
|-----------|-------------------------------------|--------|------------------|----------|--------------|--------|
| 13.446    | 15.90 ng/ul                         | 329582 | Acenaphthene-d10 |          |              | 14.909 |
| Hit# of 5 | Tentative ID                        |        | MW               | MolForm  | CAS#         | Qual   |
| 1         | Dodecane, 2,6,10-trimethyl-         |        | 212              | C15H32   | 003891-98-3  | 90     |
| 2         | Dodecane, 2,7,10-trimethyl-         |        | 212              | C15H32   | 074645-98-0  | 80     |
| 3         | Heptacosane                         |        | 380              | C27H56   | 000593-49-7  | 64     |
| 4         | Butane, 2,2-dimethyl-               |        | 86               | C6H14    | 000075-83-2  | 53     |
| 5         | Oxalic acid, bis(2-ethylhexyl) e... |        | 314              | C18H34O4 | 1000309-39-0 | 50     |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
Data File : BG051787.D  
Acq On : 23 Dec 2021 16:23  
Operator : CG/JU  
Sample : M5215-02  
Misc :  
ALS Vial : 38 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
BGLD4

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
Quant Title : SVOA CALIBRATION

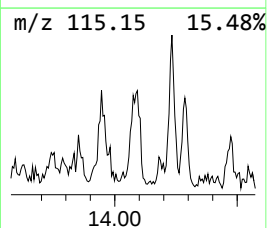
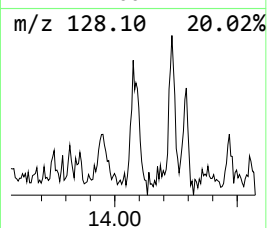
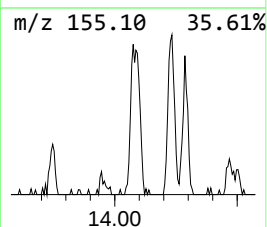
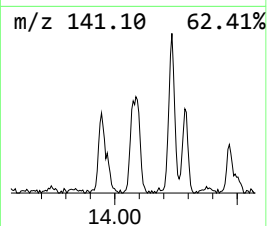
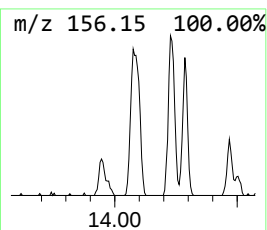
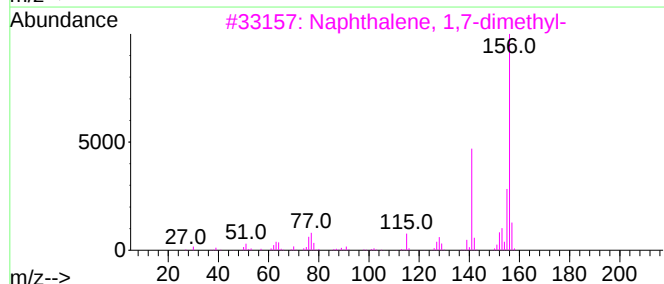
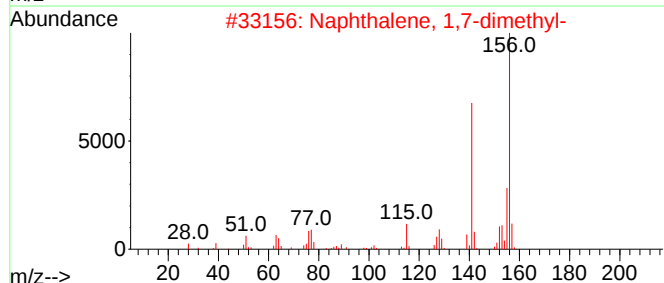
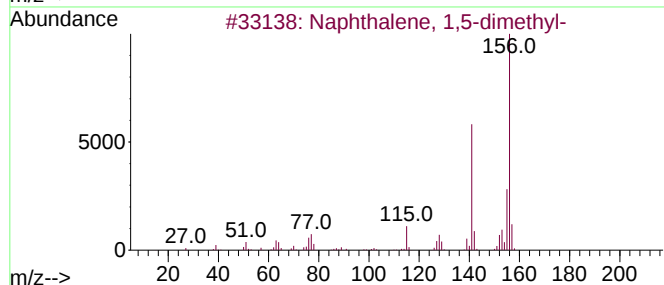
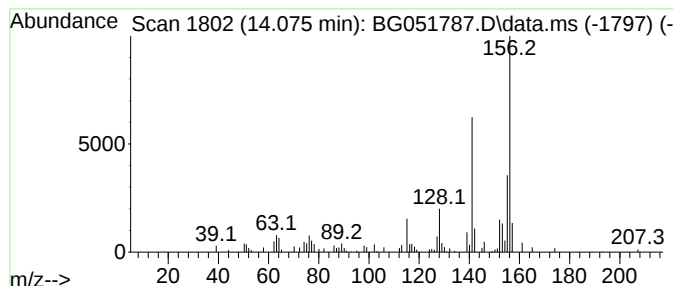
TIC Library : C:\DATABASE\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 58 Naphthalene, 1,5-dimethyl- Concentration Rank 28

| R.T.   | EstConc    | Area   | Relative to ISTD | R.T.   |
|--------|------------|--------|------------------|--------|
| 14.075 | 8.47 ng/ul | 175588 | Acenaphthene-d10 | 14.909 |

| Hit# | of 5 | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|------|----------------------------|-----|---------|-------------|------|
| 1    |      | Naphthalene, 1,5-dimethyl- | 156 | C12H12  | 000571-61-9 | 95   |
| 2    |      | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 94   |
| 3    |      | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 94   |
| 4    |      | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 94   |
| 5    |      | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 94   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_G\Data\BG122221\  
 Data File : BG051787.D  
 Acq On : 23 Dec 2021 16:23  
 Operator : CG/JU  
 Sample : M5215-02  
 Misc :  
 ALS Vial : 38 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 BGLD4

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_G\Methods\SFAM-EPA-BG121421.M  
 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L  
 TIC Integration Parameters: LSCINT.P

| TIC Top Hit name   | RT     | EstConc | Units | Response | --Internal Standard-- |        |         |      |
|--------------------|--------|---------|-------|----------|-----------------------|--------|---------|------|
|                    |        |         |       |          | #                     | RT     | Resp    | Conc |
| Ethylbenzene       | 5.745  | 20.3    | ng/ul | 4107870  | 1                     | 8.277  | 4038740 | 20.0 |
| p-Xylene           | 5.903  | 17.6    | ng/ul | 3546740  | 1                     | 8.277  | 4038740 | 20.0 |
| (DEL) Alkane: S... | 6.526  | 5.0     | ng/ul | 1001610  | 1                     | 8.277  | 4038740 | 20.0 |
| (DEL) Alkane: S... | 6.644  | 3.7     | ng/ul | 742963   | 1                     | 8.277  | 4038740 | 20.0 |
| 2,2-Dimethyl-3-... | 6.738  | 5.0     | ng/ul | 1004660  | 1                     | 8.277  | 4038740 | 20.0 |
| (DEL) Alkane: S... | 6.814  | 5.4     | ng/ul | 1095270  | 1                     | 8.277  | 4038740 | 20.0 |
| (DEL) Alkane: C... | 6.902  | 8.7     | ng/ul | 1747160  | 1                     | 8.277  | 4038740 | 20.0 |
| (DEL) Alkane: S... | 7.161  | 5.7     | ng/ul | 1140400  | 1                     | 8.277  | 4038740 | 20.0 |
| (DEL) Alkane: S... | 7.219  | 2.7     | ng/ul | 552764   | 1                     | 8.277  | 4038740 | 20.0 |
| (DEL) Alkane: S... | 7.255  | 10.7    | ng/ul | 2159970  | 1                     | 8.277  | 4038740 | 20.0 |
| (DEL) Alkane: S... | 7.313  | 7.8     | ng/ul | 1572270  | 1                     | 8.277  | 4038740 | 20.0 |
| Benzene, 1-ethy... | 7.666  | 2.9     | ng/ul | 582265   | 1                     | 8.277  | 4038740 | 20.0 |
| Mesitylene         | 7.930  | 29.7    | ng/ul | 5995630  | 1                     | 8.277  | 4038740 | 20.0 |
| Carbonic acid, ... | 8.106  | 3.1     | ng/ul | 618190   | 1                     | 8.277  | 4038740 | 20.0 |
| unknown-01         | 8.194  | 6.5     | ng/ul | 1312070  | 1                     | 8.277  | 4038740 | 20.0 |
| (DEL) Alkane: S... | 8.353  | 7.8     | ng/ul | 1568820  | 1                     | 8.277  | 4038740 | 20.0 |
| Benzene, 1,2,3-... | 8.406  | 12.8    | ng/ul | 2589570  | 1                     | 8.277  | 4038740 | 20.0 |
| 1-Iodo-2-methyl... | 8.471  | 6.2     | ng/ul | 1258790  | 1                     | 8.277  | 4038740 | 20.0 |
| (DEL) Alkane: S... | 8.535  | 6.3     | ng/ul | 1261730  | 1                     | 8.277  | 4038740 | 20.0 |
| (DEL) Alkane: C... | 8.582  | 8.2     | ng/ul | 1647530  | 1                     | 8.277  | 4038740 | 20.0 |
| Indane             | 8.664  | 4.8     | ng/ul | 958504   | 1                     | 8.277  | 4038740 | 20.0 |
| (DEL) Alkane: S... | 8.717  | 2.9     | ng/ul | 577022   | 1                     | 8.277  | 4038740 | 20.0 |
| (DEL) Alkane: S... | 8.817  | 9.8     | ng/ul | 1970700  | 1                     | 8.277  | 4038740 | 20.0 |
| 4-Methyl-3-hept... | 8.876  | 9.6     | ng/ul | 1935950  | 1                     | 8.277  | 4038740 | 20.0 |
| (DEL) Alkane: S... | 9.058  | 11.7    | ng/ul | 2362510  | 1                     | 8.277  | 4038740 | 20.0 |
| Naphthalene, de... | 9.140  | 3.4     | ng/ul | 676811   | 1                     | 8.277  | 4038740 | 20.0 |
| Benzene, 2-ethy... | 9.246  | 3.7     | ng/ul | 742508   | 1                     | 8.277  | 4038740 | 20.0 |
| o-Cymene           | 9.299  | 7.5     | ng/ul | 1514270  | 1                     | 8.277  | 4038740 | 20.0 |
| Benzene, 1-ethy... | 9.640  | 6.3     | ng/ul | 1278570  | 1                     | 8.277  | 4038740 | 20.0 |
| (DEL) Alkane: S... | 9.781  | 48.9    | ng/ul | 2446830  | 2                     | 11.120 | 1001180 | 20.0 |
| (DEL) Alkane: S... | 9.880  | 20.6    | ng/ul | 1030260  | 2                     | 11.120 | 1001180 | 20.0 |
| Benzene, 1-ethy... | 9.939  | 36.0    | ng/ul | 1801780  | 2                     | 11.120 | 1001180 | 20.0 |
| 1,3,8-p-Menthat... | 9.992  | 34.2    | ng/ul | 1711780  | 2                     | 11.120 | 1001180 | 20.0 |
| Pulegone           | 10.039 | 18.1    | ng/ul | 905674   | 2                     | 11.120 | 1001180 | 20.0 |
| (DEL) Alkane: C... | 10.233 | 20.1    | ng/ul | 1006810  | 2                     | 11.120 | 1001180 | 20.0 |
| Sulfurous acid,... | 10.444 | 16.4    | ng/ul | 818893   | 2                     | 11.120 | 1001180 | 20.0 |
| unknown-02         | 10.521 | 38.5    | ng/ul | 1925390  | 2                     | 11.120 | 1001180 | 20.0 |
| Benzene, 1-ethy... | 10.574 | 11.2    | ng/ul | 558973   | 2                     | 11.120 | 1001180 | 20.0 |
| (DEL) Alkane: S... | 10.632 | 14.9    | ng/ul | 743657   | 2                     | 11.120 | 1001180 | 20.0 |
| Benzene, 1-meth... | 10.809 | 10.1    | ng/ul | 505187   | 2                     | 11.120 | 1001180 | 20.0 |
| (DEL) Alkane: C... | 10.973 | 4.2     | ng/ul | 211378   | 2                     | 11.120 | 1001180 | 20.0 |
| unknown-03         | 11.008 | 11.8    | ng/ul | 591465   | 2                     | 11.120 | 1001180 | 20.0 |
| unknown-04         | 11.331 | 8.2     | ng/ul | 410757   | 2                     | 11.120 | 1001180 | 20.0 |
| unknown-05         | 11.390 | 5.5     | ng/ul | 273038   | 2                     | 11.120 | 1001180 | 20.0 |
| Benzene, 1,3-di... | 11.490 | 4.4     | ng/ul | 219344   | 2                     | 11.120 | 1001180 | 20.0 |
| (DEL) Alkane: C... | 11.819 | 8.4     | ng/ul | 420658   | 2                     | 11.120 | 1001180 | 20.0 |
| (DEL) Alkane: S... | 11.866 | 4.4     | ng/ul | 219057   | 2                     | 11.120 | 1001180 | 20.0 |
| (DEL) Alkane: S... | 11.942 | 6.4     | ng/ul | 322564   | 2                     | 11.120 | 1001180 | 20.0 |
| Decane, 1-iodo-    | 12.013 | 9.5     | ng/ul | 475949   | 2                     | 11.120 | 1001180 | 20.0 |
| (DEL) Alkane: S... | 12.119 | 17.0    | ng/ul | 852593   | 2                     | 11.120 | 1001180 | 20.0 |
| 1H-Indene, 2,3-... | 12.213 | 3.5     | ng/ul | 172931   | 2                     | 11.120 | 1001180 | 20.0 |
| (DEL) Alkane: S... | 13.129 | 9.4     | ng/ul | 194050   | 3                     | 14.909 | 414529  | 20.0 |

|                    |        |      |       |        |   |        |        |      |
|--------------------|--------|------|-------|--------|---|--------|--------|------|
| 5-Bromo-2-(4-pi... | 13.200 | 8.6  | ng/ul | 177370 | 3 | 14.909 | 414529 | 20.0 |
| (DEL) Alkane: S... | 13.305 | 10.2 | ng/ul | 211732 | 3 | 14.909 | 414529 | 20.0 |
| (DEL) Alkane: S... | 13.446 | 15.9 | ng/ul | 329582 | 3 | 14.909 | 414529 | 20.0 |
| Naphthalene, 1,... | 14.075 | 8.5  | ng/ul | 175588 | 3 | 14.909 | 414529 | 20.0 |

Instrument :

BNA\_G

ClientSampleId :

BGLD4