

Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF062716\  
 Data File : BF088466.D  
 Acq On : 27 Jun 2016 22:21  
 Operator : UM/SJ  
 Sample : H3630-03  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 B1409-TP01N-1224

Integration Parameters: rteint.p  
 Integrator: RTE  
 Smoothing : ON  
 Sampling : 1  
 Start Thrs: 0.2  
 Stop Thrs : 0  
 Filtering: 5  
 Min Area: 3 % of largest Peak  
 Max Peaks: 100  
 Peak Location: TOP

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

Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060716.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 1.690       | 7             | 9           | 17           | rBV      | 921832         | 1442800       | 15.08%          | 0.636%        |
| 2         | 2.204       | 43            | 54          | 58           | rVB      | 209064         | 570257        | 5.96%           | 0.251%        |
| 3         | 3.038       | 120           | 127         | 132          | rBV2     | 161213         | 460175        | 4.81%           | 0.203%        |
| 4         | 3.221       | 139           | 143         | 148          | rBV      | 100697         | 250351        | 2.62%           | 0.110%        |
| 5         | 3.438       | 153           | 162         | 165          | rBV      | 382574         | 1022407       | 10.69%          | 0.450%        |
| 6         | 3.507       | 165           | 168         | 173          | rVB      | 173334         | 345991        | 3.62%           | 0.152%        |
| 7         | 3.633       | 173           | 179         | 183          | rBV2     | 120479         | 334624        | 3.50%           | 0.147%        |
| 8         | 3.907       | 198           | 203         | 208          | rVB2     | 458312         | 981358        | 10.26%          | 0.432%        |
| 9         | 4.033       | 209           | 214         | 219          | rBV2     | 276315         | 509333        | 5.32%           | 0.224%        |
| 10        | 4.433       | 247           | 249         | 251          | rBV      | 315572         | 482545        | 5.04%           | 0.213%        |
| 11        | 4.536       | 255           | 258         | 260          | rVB      | 1133571        | 1546092       | 16.16%          | 0.681%        |
| 12        | 4.604       | 260           | 264         | 266          | rVB      | 856847         | 1074815       | 11.24%          | 0.473%        |
| 13        | 4.650       | 266           | 268         | 273          | rVB3     | 306826         | 527387        | 5.51%           | 0.232%        |
| 14        | 4.753       | 273           | 277         | 282          | rVB4     | 201394         | 348494        | 3.64%           | 0.154%        |
| 15        | 4.844       | 282           | 285         | 287          | rBV2     | 847456         | 1540975       | 16.11%          | 0.679%        |
| 16        | 4.890       | 287           | 289         | 292          | rVB4     | 187629         | 391324        | 4.09%           | 0.172%        |
| 17        | 4.947       | 292           | 294         | 295          | rBV      | 445568         | 576035        | 6.02%           | 0.254%        |
| 18        | 4.981       | 295           | 297         | 299          | rVB      | 525281         | 588384        | 6.15%           | 0.259%        |
| 19        | 5.027       | 299           | 301         | 302          | rBV      | 434582         | 460372        | 4.81%           | 0.203%        |
| 20        | 5.061       | 302           | 304         | 308          | rVB      | 855024         | 949347        | 9.92%           | 0.418%        |
| 21        | 5.119       | 308           | 309         | 310          | rBV      | 1093877        | 1079694       | 11.29%          | 0.476%        |
| 22        | 5.244       | 317           | 320         | 322          | rBV      | 801587         | 1211353       | 12.66%          | 0.534%        |
| 23        | 5.290       | 322           | 324         | 326          | rVB      | 1247865        | 1317625       | 13.78%          | 0.580%        |
| 24        | 5.336       | 326           | 328         | 332          | rVB3     | 1067069        | 1575466       | 16.47%          | 0.694%        |
| 25        | 5.416       | 332           | 335         | 338          | rBV      | 347532         | 520847        | 5.45%           | 0.229%        |
| 26        | 5.519       | 338           | 344         | 346          | rBV2     | 1272184        | 2873026       | 30.04%          | 1.266%        |
| 27        | 5.564       | 346           | 348         | 349          | rBV2     | 496739         | 838179        | 8.76%           | 0.369%        |
| 28        | 5.656       | 354           | 356         | 360          | rVV2     | 1496745        | 2958947       | 30.93%          | 1.303%        |
| 29        | 5.736       | 360           | 363         | 364          | rVV2     | 1032999        | 1846974       | 19.31%          | 0.814%        |
| 30        | 5.793       | 364           | 368         | 370          | rVV4     | 2758165        | 6409294       | 67.01%          | 2.823%        |
| 31        | 5.850       | 370           | 373         | 376          | rVV2     | 1757674        | 4253106       | 44.46%          | 1.874%        |
| 32        | 5.987       | 380           | 385         | 387          | rVV4     | 1796749        | 4124373       | 43.12%          | 1.817%        |
| 33        | 6.056       | 387           | 391         | 397          | rVV5     | 2795799        | 9274212       | 96.96%          | 4.085%        |
| 34        | 6.147       | 397           | 399         | 401          | rVV      | 1857546        | 3066710       | 32.06%          | 1.351%        |

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 ClientSampleId :  
 B1409-TP01N-1224

Integration Parameters: rteint.p  
 Integrator: RTE  
 Smoothing : ON  
 Sampling : 1  
 Start Thrs: 0.2  
 Stop Thrs : 0

Filtering: 5  
 Min Area: 3 % of largest Peak  
 Max Peaks: 100  
 Peak Location: TOP

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

Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060716.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |       |     |     |     |      |         |         |         |        |
|----|-------|-----|-----|-----|------|---------|---------|---------|--------|
| 35 | 6.216 | 401 | 405 | 406 | rVV2 | 2134958 | 4902801 | 51.26%  | 2.160% |
| 36 | 6.239 | 406 | 407 | 409 | rVV  | 1365327 | 2093896 | 21.89%  | 0.922% |
| 37 | 6.319 | 409 | 414 | 417 | rVV4 | 2278815 | 8588672 | 89.79%  | 3.783% |
| 38 | 6.399 | 419 | 421 | 423 | rVV3 | 1900987 | 3429693 | 35.86%  | 1.511% |
| 39 | 6.513 | 423 | 431 | 432 | rVV6 | 1606577 | 6875941 | 71.88%  | 3.029% |
| 40 | 6.593 | 435 | 438 | 440 | rVV  | 2895912 | 7159316 | 74.85%  | 3.154% |
| 41 | 6.639 | 440 | 442 | 445 | rVV2 | 1780733 | 4690096 | 49.03%  | 2.066% |
| 42 | 6.719 | 445 | 449 | 453 | rVV6 | 2872439 | 9565235 | 100.00% | 4.214% |
| 43 | 6.822 | 453 | 458 | 459 | rVV4 | 1695980 | 5502084 | 57.52%  | 2.424% |
| 44 | 6.856 | 459 | 461 | 462 | rVV2 | 1750608 | 2874047 | 30.05%  | 1.266% |
| 45 | 6.890 | 462 | 464 | 465 | rVV2 | 2058452 | 3427992 | 35.84%  | 1.510% |
| 46 | 6.913 | 465 | 466 | 468 | rVV  | 1800574 | 3051313 | 31.90%  | 1.344% |
| 47 | 6.947 | 468 | 469 | 473 | rVV4 | 2148933 | 5002131 | 52.29%  | 2.203% |
| 48 | 7.039 | 475 | 477 | 478 | rVV  | 954600  | 1874817 | 19.60%  | 0.826% |
| 49 | 7.107 | 479 | 483 | 484 | rVV  | 3020507 | 6256463 | 65.41%  | 2.756% |
| 50 | 7.130 | 484 | 485 | 487 | rVV2 | 2134900 | 3443813 | 36.00%  | 1.517% |
| 51 | 7.176 | 487 | 489 | 490 | rVV  | 1389902 | 2527072 | 26.42%  | 1.113% |
| 52 | 7.210 | 490 | 492 | 494 | rVV3 | 1994001 | 3566499 | 37.29%  | 1.571% |
| 53 | 7.256 | 494 | 496 | 497 | rVV2 | 1112174 | 2055317 | 21.49%  | 0.905% |
| 54 | 7.279 | 497 | 498 | 500 | rVV  | 1592826 | 2455471 | 25.67%  | 1.082% |
| 55 | 7.336 | 500 | 503 | 504 | rVV3 | 1820146 | 3651158 | 38.17%  | 1.608% |
| 56 | 7.359 | 504 | 505 | 508 | rVV2 | 2228075 | 3116333 | 32.58%  | 1.373% |
| 57 | 7.416 | 508 | 510 | 512 | rVV2 | 805988  | 1977732 | 20.68%  | 0.871% |
| 58 | 7.473 | 512 | 515 | 516 | rVV3 | 1849781 | 3768062 | 39.39%  | 1.660% |
| 59 | 7.507 | 516 | 518 | 520 | rVV3 | 1451855 | 3326475 | 34.78%  | 1.465% |
| 60 | 7.587 | 522 | 525 | 526 | rVV2 | 2114997 | 4110623 | 42.97%  | 1.811% |
| 61 | 7.610 | 526 | 527 | 529 | rVV2 | 1237767 | 1858798 | 19.43%  | 0.819% |
| 62 | 7.667 | 529 | 532 | 534 | rVV3 | 1127619 | 2771399 | 28.97%  | 1.221% |
| 63 | 7.713 | 534 | 536 | 540 | rVV  | 964527  | 2429726 | 25.40%  | 1.070% |
| 64 | 7.839 | 540 | 547 | 548 | rVV4 | 2162876 | 6157901 | 64.38%  | 2.713% |
| 65 | 7.862 | 548 | 549 | 553 | rVV4 | 1781689 | 4388091 | 45.88%  | 1.933% |
| 66 | 7.919 | 553 | 554 | 560 | rVV3 | 1714325 | 3792566 | 39.65%  | 1.671% |
| 67 | 8.022 | 560 | 563 | 564 | rVV2 | 870083  | 1833612 | 19.17%  | 0.808% |
| 68 | 8.136 | 567 | 573 | 575 | rVV6 | 1066108 | 4412588 | 46.13%  | 1.944% |
| 69 | 8.193 | 577 | 578 | 580 | rVV2 | 832406  | 1521193 | 15.90%  | 0.670% |
| 70 | 8.227 | 580 | 581 | 583 | rVV2 | 1536390 | 1802129 | 18.84%  | 0.794% |
| 71 | 8.273 | 583 | 585 | 589 | rVV3 | 1385259 | 3770048 | 39.41%  | 1.661% |

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Instrument :  
 BNA\_F  
 ClientSampleId :  
 B1409-TP01N-1224

Integration Parameters: rteint.p  
 Integrator: RTE  
 Smoothing : ON Filtering: 5  
 Sampling : 1 Min Area: 3 % 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:\HPCHEM1\BNA\_F\METHODS\8270-BF060716.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|     |        |     |     |     |      |         |         |        |        |
|-----|--------|-----|-----|-----|------|---------|---------|--------|--------|
| 72  | 8.342  | 589 | 591 | 593 | rVV2 | 768553  | 1758676 | 18.39% | 0.775% |
| 73  | 8.445  | 597 | 600 | 601 | rVV3 | 1093322 | 2077905 | 21.72% | 0.915% |
| 74  | 8.479  | 601 | 603 | 606 | rVV4 | 812442  | 2261304 | 23.64% | 0.996% |
| 75  | 8.547  | 606 | 609 | 612 | rVV2 | 1198174 | 2986098 | 31.22% | 1.315% |
| 76  | 8.605  | 612 | 614 | 615 | rVV2 | 387249  | 699565  | 7.31%  | 0.308% |
| 77  | 8.650  | 615 | 618 | 619 | rVV  | 1141278 | 1776197 | 18.57% | 0.782% |
| 78  | 8.673  | 619 | 620 | 622 | rVV2 | 354513  | 602023  | 6.29%  | 0.265% |
| 79  | 8.742  | 622 | 626 | 631 | rVV8 | 658216  | 2227807 | 23.29% | 0.981% |
| 80  | 8.867  | 634 | 637 | 638 | rVV  | 936677  | 1206568 | 12.61% | 0.532% |
| 81  | 8.913  | 638 | 641 | 643 | rVV  | 1519041 | 1756535 | 18.36% | 0.774% |
| 82  | 9.005  | 645 | 649 | 651 | rVV2 | 388837  | 760769  | 7.95%  | 0.335% |
| 83  | 9.050  | 651 | 653 | 656 | rVV3 | 209570  | 406382  | 4.25%  | 0.179% |
| 84  | 9.108  | 656 | 658 | 661 | rVB  | 218399  | 302936  | 3.17%  | 0.133% |
| 85  | 9.176  | 661 | 664 | 666 | rBV  | 411229  | 639107  | 6.68%  | 0.282% |
| 86  | 9.245  | 668 | 670 | 671 | rBV  | 567751  | 639638  | 6.69%  | 0.282% |
| 87  | 9.268  | 671 | 672 | 674 | rVV2 | 395318  | 512659  | 5.36%  | 0.226% |
| 88  | 9.313  | 674 | 676 | 679 | rVB  | 709254  | 814913  | 8.52%  | 0.359% |
| 89  | 9.519  | 693 | 694 | 696 | rBV  | 171846  | 240424  | 2.51%  | 0.106% |
| 90  | 9.576  | 696 | 699 | 701 | rBV3 | 430896  | 560466  | 5.86%  | 0.247% |
| 91  | 9.839  | 719 | 722 | 725 | rVB3 | 191749  | 397226  | 4.15%  | 0.175% |
| 92  | 9.976  | 732 | 734 | 737 | rBV4 | 120426  | 245051  | 2.56%  | 0.108% |
| 93  | 10.239 | 754 | 757 | 759 | rBV  | 282016  | 269793  | 2.82%  | 0.119% |
| 94  | 10.365 | 765 | 768 | 770 | rBV2 | 650218  | 699747  | 7.32%  | 0.308% |
| 95  | 10.502 | 777 | 780 | 783 | rBV2 | 339077  | 498999  | 5.22%  | 0.220% |
| 96  | 10.776 | 800 | 804 | 806 | rBV  | 140960  | 343821  | 3.59%  | 0.151% |
| 97  | 10.971 | 817 | 821 | 823 | rBV2 | 258263  | 480018  | 5.02%  | 0.211% |
| 98  | 11.051 | 827 | 828 | 829 | rBV  | 280827  | 260018  | 2.72%  | 0.115% |
| 99  | 12.114 | 919 | 921 | 924 | rBV2 | 222974  | 424710  | 4.44%  | 0.187% |
| 100 | 12.651 | 966 | 968 | 972 | rBV2 | 848360  | 1377448 | 14.40% | 0.607% |

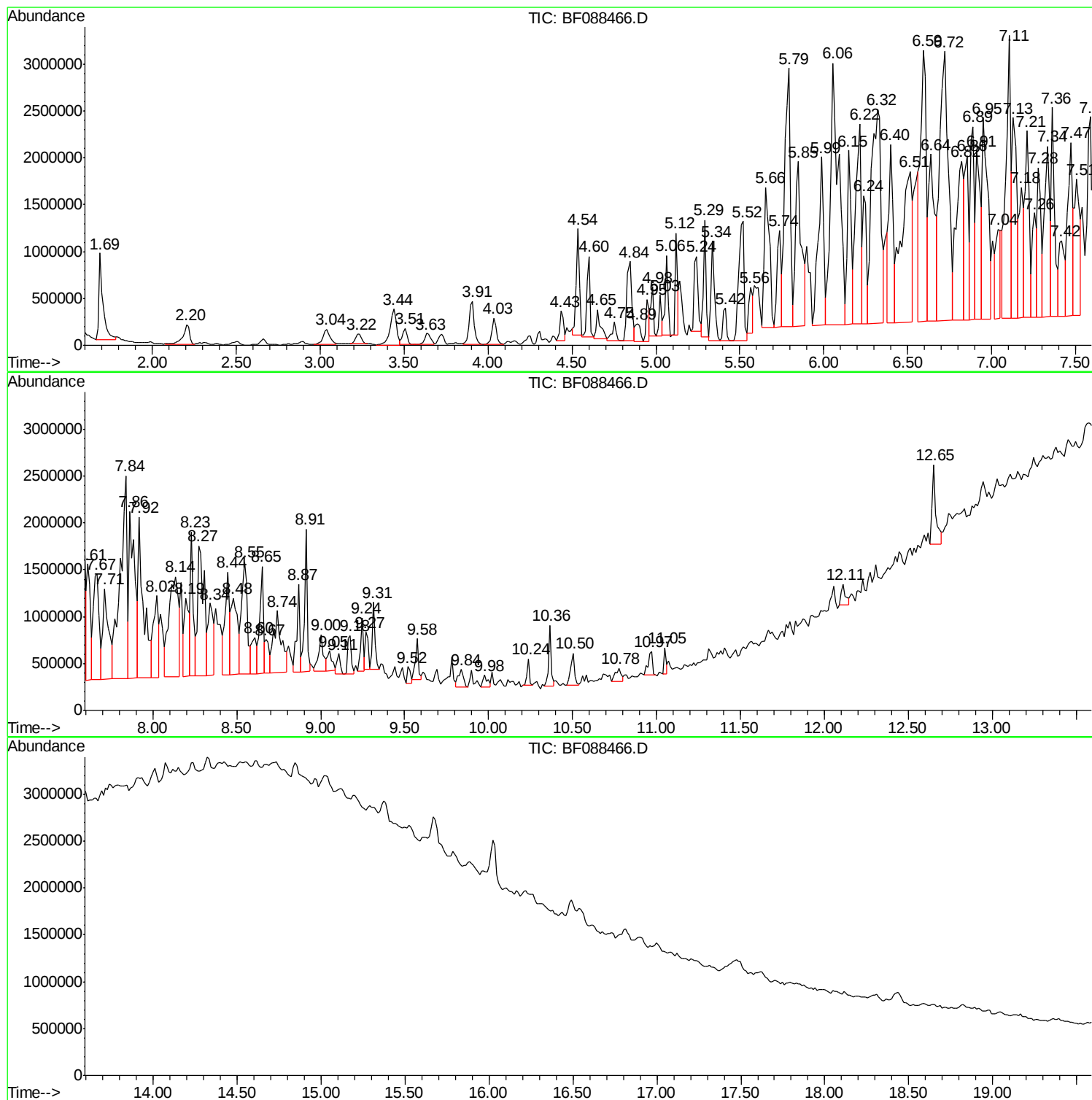
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Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
B1409-TP01N-1224

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF062716\  
Data File : BF088466.D  
Acq On : 27 Jun 2016 22:21  
Operator : UM/SJ  
Sample : H3630-03  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
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B1409-TP01N-1224

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

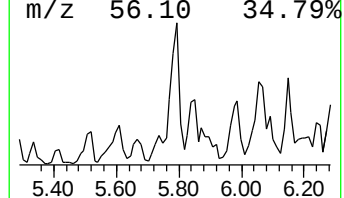
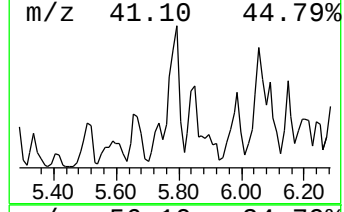
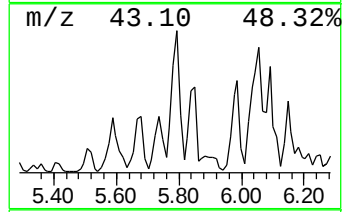
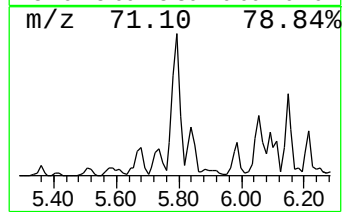
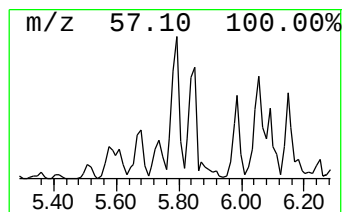
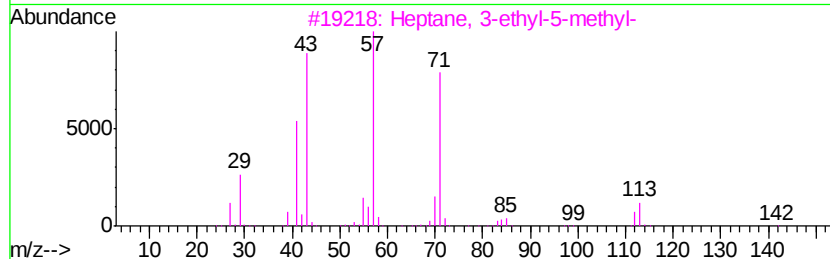
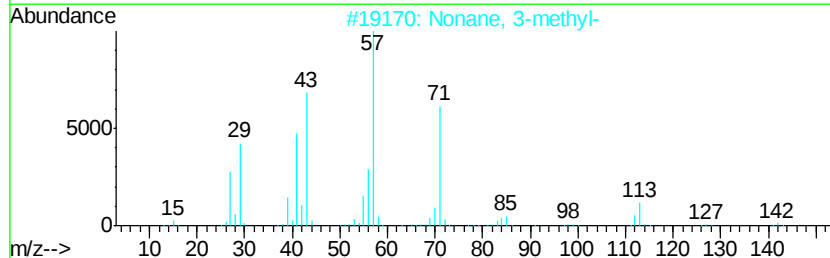
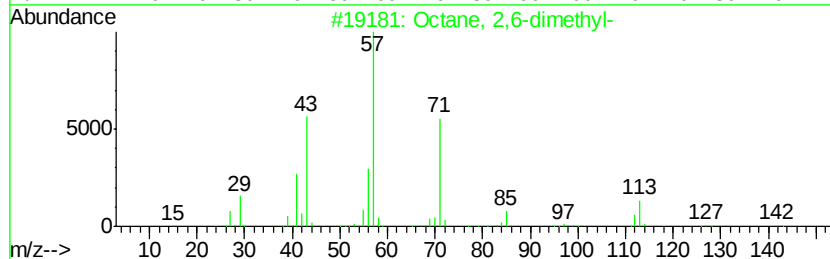
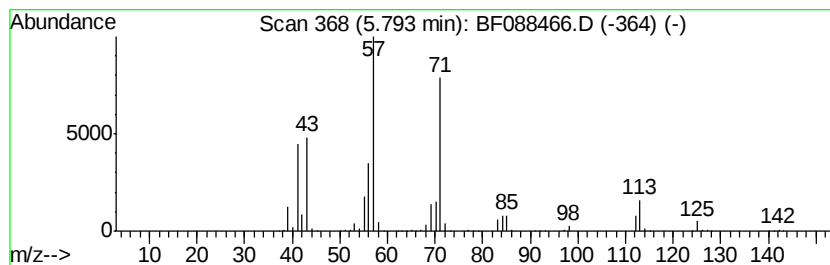
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 1 Octane, 2,6-dimethyl- Concentration Rank 12

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 5.79 | 18.64 ng | 6409290 | 1,4-Dichlorobenzene-d4 | 6.55 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Octane, 2,6-dimethyl-               | 142 | C10H22    | 002051-30-1  | 91   |
| 2    |    | Nonane, 3-methyl-                   | 142 | C10H22    | 005911-04-6  | 91   |
| 3    |    | Heptane, 3-ethyl-5-methyl-          | 142 | C10H22    | 052896-90-9  | 72   |
| 4    |    | Sulfurous acid, decyl 2-ethylhex... | 334 | C18H38O3S | 1000309-19-3 | 64   |
| 5    |    | Butane, 2,2-dimethyl-               | 86  | C6H14     | 000075-83-2  | 64   |



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ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

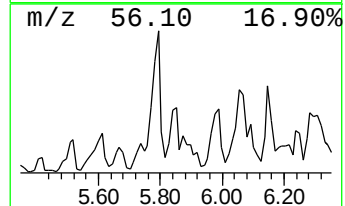
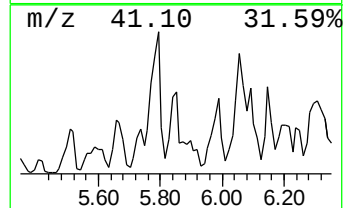
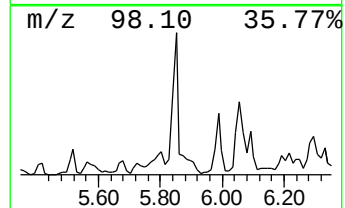
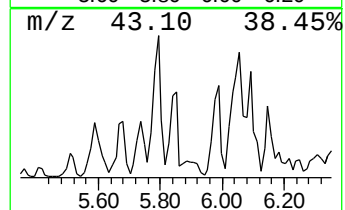
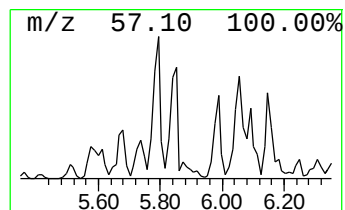
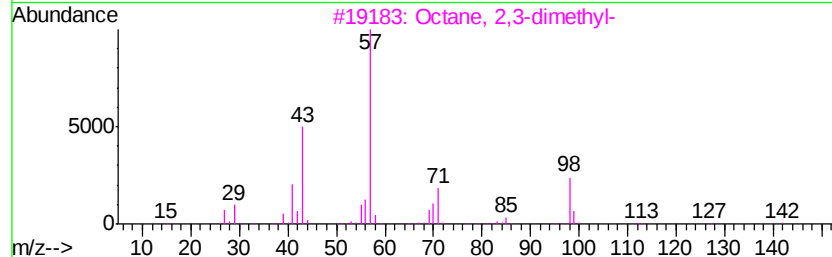
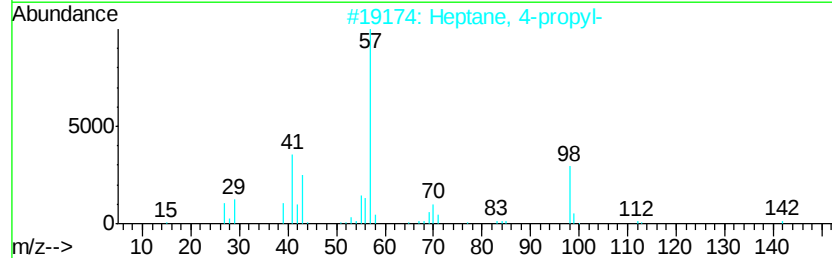
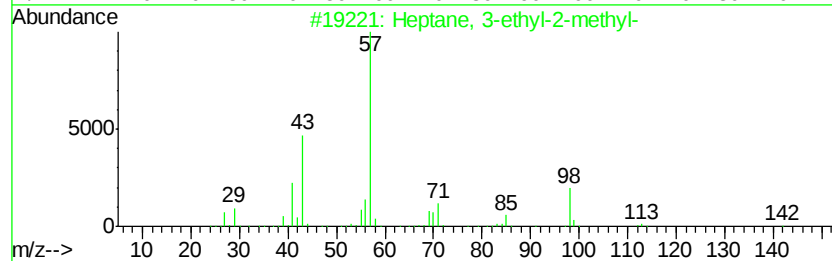
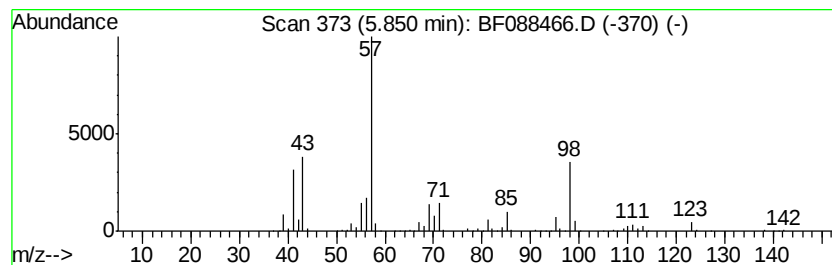
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 2 Heptane, 3-ethyl-2-methyl- Concentration Rank 20

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 5.85 | 12.37 ng | 4253110 | 1,4-Dichlorobenzene-d4 | 6.55 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | Heptane, 3-ethyl-2-methyl-          | 142 | C10H22    | 014676-29-0  | 59   |
| 2    |    | Heptane, 4-propyl-                  | 142 | C10H22    | 003178-29-8  | 50   |
| 3    |    | Octane, 2,3-dimethyl-               | 142 | C10H22    | 007146-60-3  | 49   |
| 4    |    | Ether, tert-butyl isopropylidene... | 154 | C10H18O   | 024524-56-9  | 47   |
| 5    |    | Sulfurous acid, 2-ethylhexyl hep... | 432 | C25H52O3S | 1000309-20-0 | 43   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF062716\  
Data File : BF088466.D  
Acq On : 27 Jun 2016 22:21  
Operator : UM/SJ  
Sample : H3630-03  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
B1409-TP01N-1224

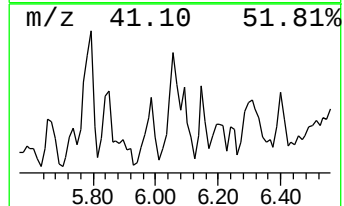
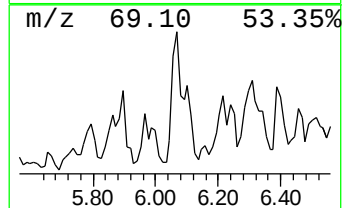
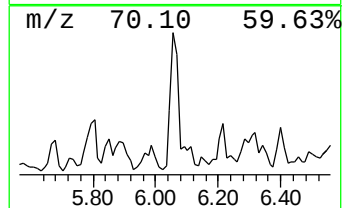
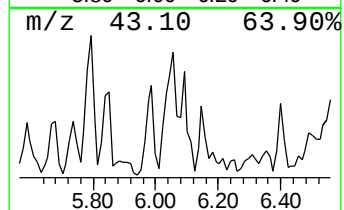
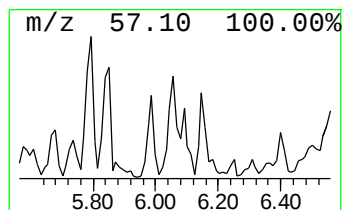
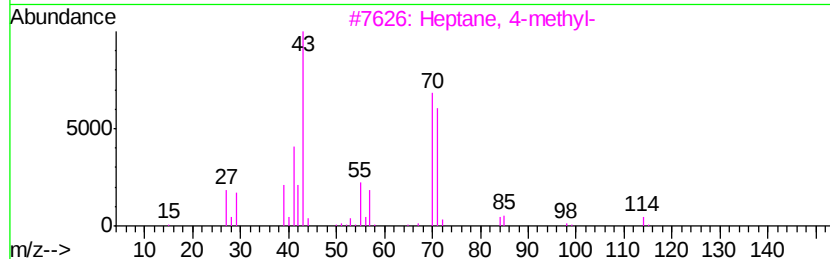
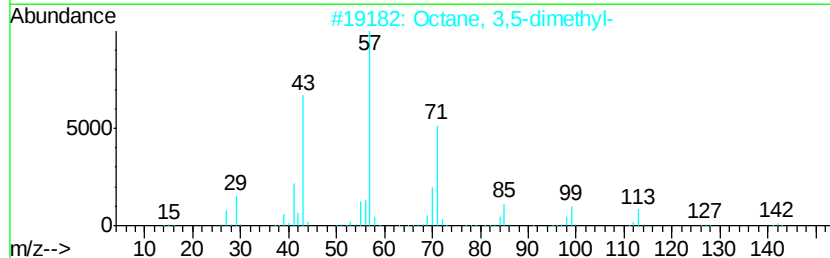
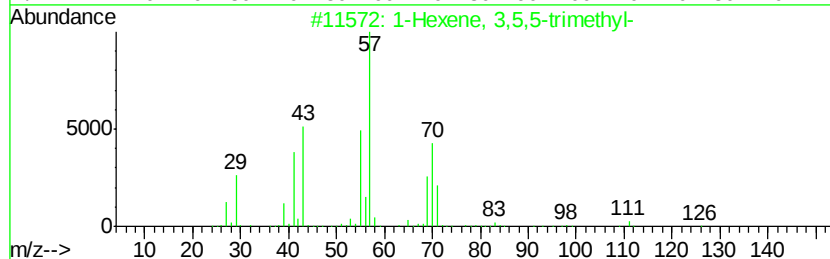
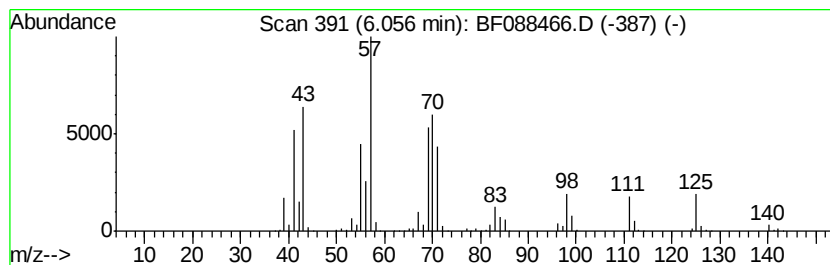
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 3 1-Hexene, 3,5,5-trimethyl- Concentration Rank 7

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.06 | 26.98 ng | 9274210 | 1,4-Dichlorobenzene-d4 | 6.55 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | 1-Hexene, 3,5,5-trimethyl- | 126 | C9H18   | 004316-65-8 | 52   |
| 2    |    | Octane, 3,5-dimethyl-      | 142 | C10H22  | 015869-93-9 | 43   |
| 3    |    | Heptane, 4-methyl-         | 114 | C8H18   | 000589-53-7 | 38   |
| 4    |    | Undecane, 4,5-dimethyl-    | 184 | C13H28  | 017312-79-7 | 38   |
| 5    |    | Nonane, 4-methyl-          | 142 | C10H22  | 017301-94-9 | 30   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF062716\  
Data File : BF088466.D  
Acq On : 27 Jun 2016 22:21  
Operator : UM/SJ  
Sample : H3630-03  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
B1409-TP01N-1224

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

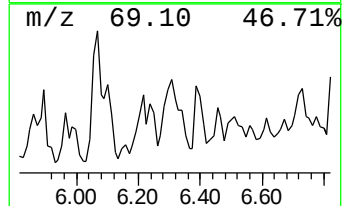
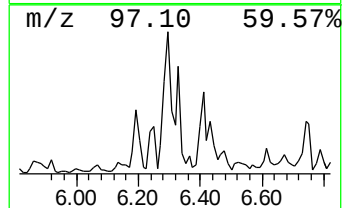
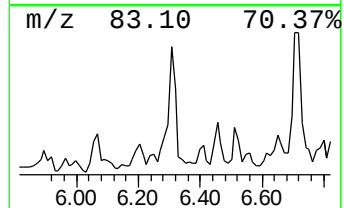
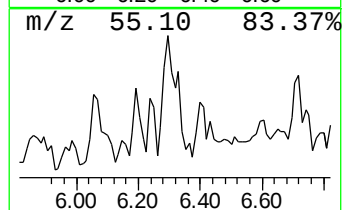
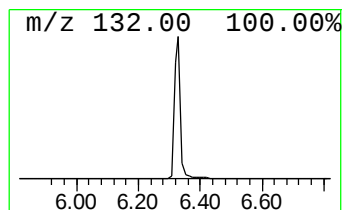
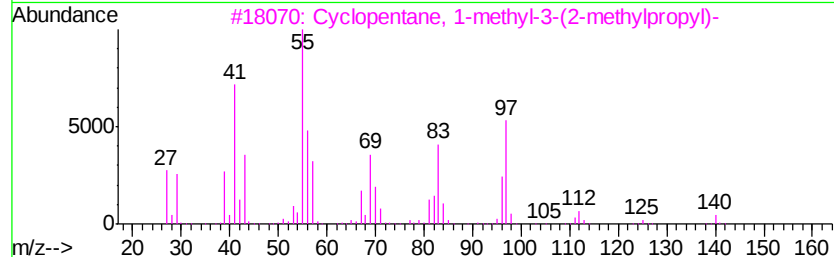
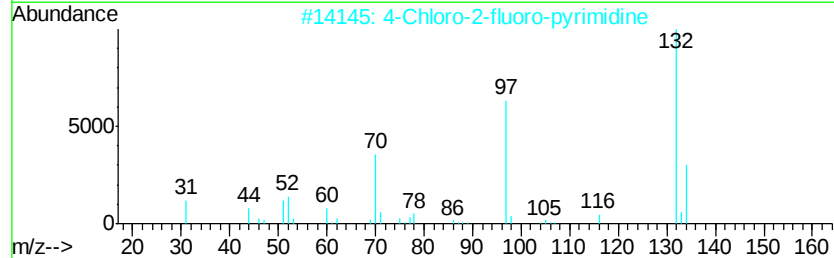
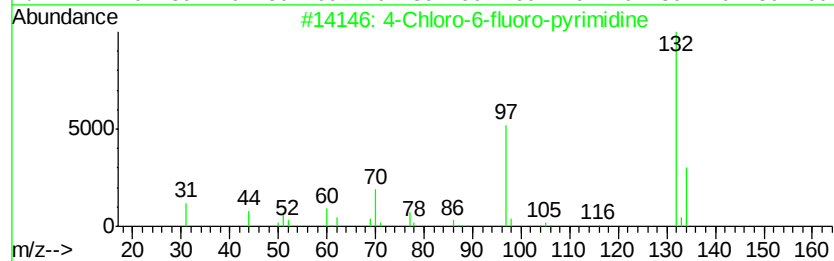
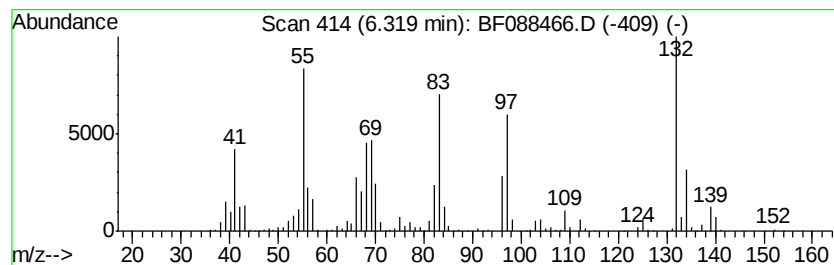
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 4 unknown6.32 Concentration Rank 9

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.32 | 24.98 ng | 8588670 | 1,4-Dichlorobenzene-d4 | 6.55 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|-------------------------------------|-----|-----------|-------------|------|
| 1    | 5  | 4-Chloro-6-fluoro-pyrimidine        | 132 | C4H2ClFN2 | 051422-01-6 | 27   |
| 2    |    | 4-Chloro-2-fluoro-pyrimidine        | 132 | C4H2ClFN2 | 051422-00-5 | 27   |
| 3    |    | Cyclopentane, 1-methyl-3-(2-meth... | 140 | C10H20    | 029053-04-1 | 15   |
| 4    |    | Cyclohexane, 1,2-dimethyl- (cis/... | 112 | C8H16     | 000583-57-3 | 11   |
| 5    |    | Cyclohexanemethanol                 | 114 | C7H14O    | 000100-49-2 | 10   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF062716\  
Data File : BF088466.D  
Acq On : 27 Jun 2016 22:21  
Operator : UM/SJ  
Sample : H3630-03  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
B1409-TP01N-1224

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

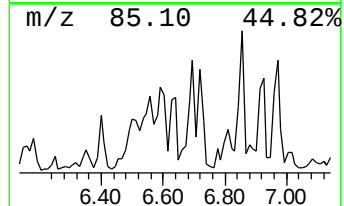
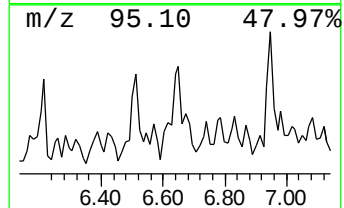
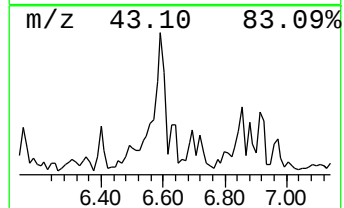
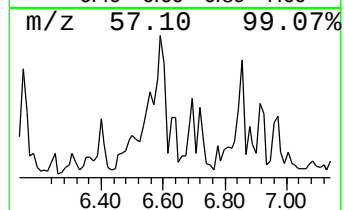
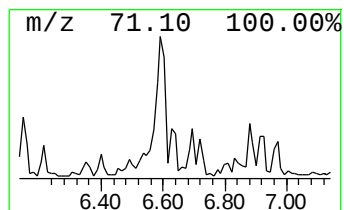
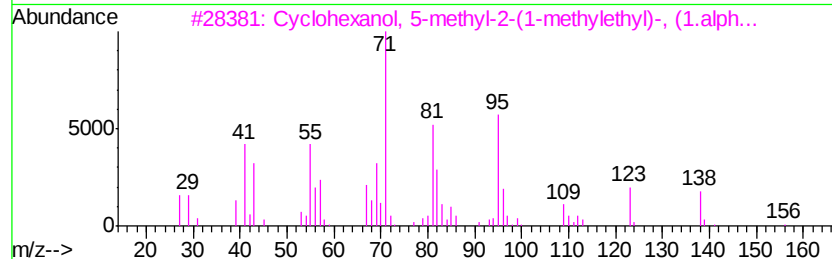
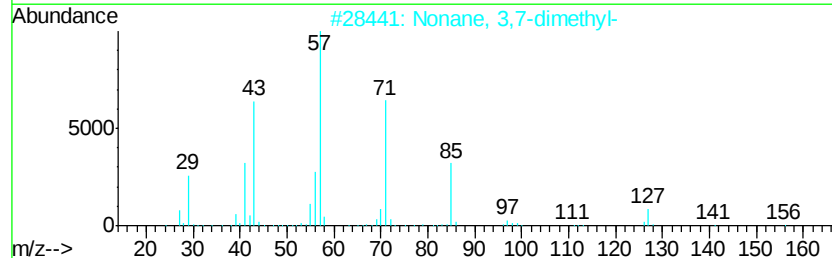
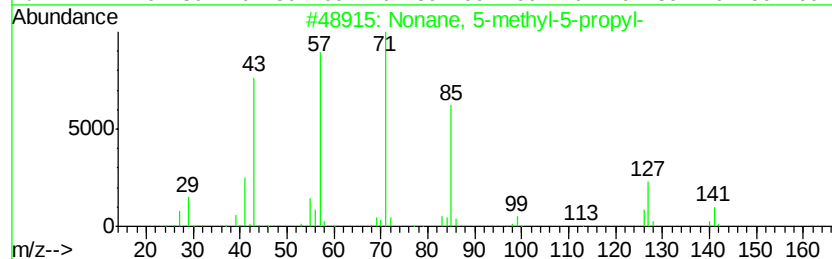
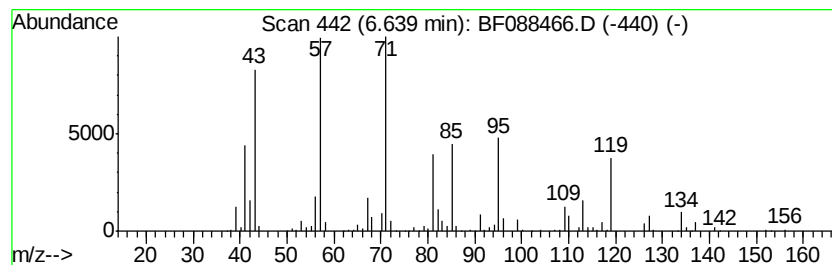
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 5 unknown6.64 Concentration Rank 18

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.64 | 13.64 ng | 4690100 | 1,4-Dichlorobenzene-d4 | 6.55 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Nonane, 5-methyl-5-propyl-          | 184 | C13H28  | 017312-75-3 | 47   |
| 2    |    | Nonane, 3,7-dimethyl-               | 156 | C11H24  | 017302-32-8 | 46   |
| 3    |    | Cyclohexanol, 5-methyl-2-(1-meth... | 156 | C10H20O | 000491-02-1 | 43   |
| 4    |    | Dodecane, 2,7,10-trimethyl-         | 212 | C15H32  | 074645-98-0 | 43   |
| 5    |    | 3-Bromooctane                       | 192 | C8H17Br | 000999-64-4 | 38   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF062716\  
Data File : BF088466.D  
Acq On : 27 Jun 2016 22:21  
Operator : UM/SJ  
Sample : H3630-03  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
B1409-TP01N-1224

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

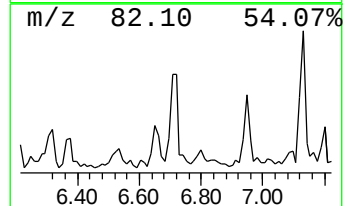
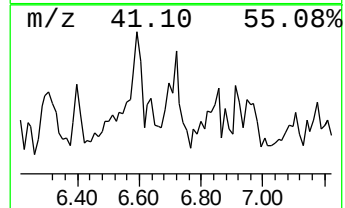
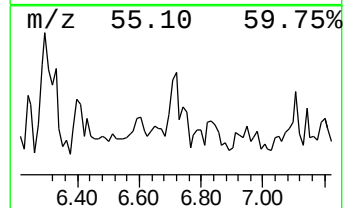
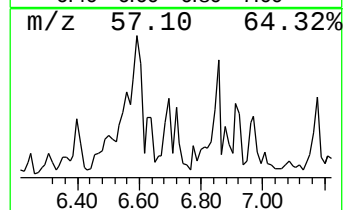
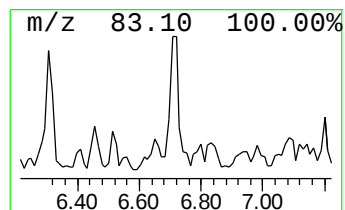
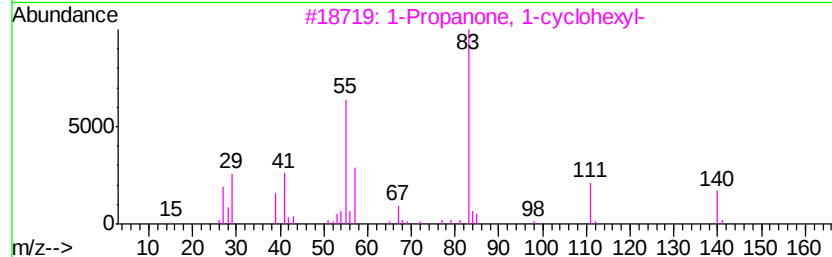
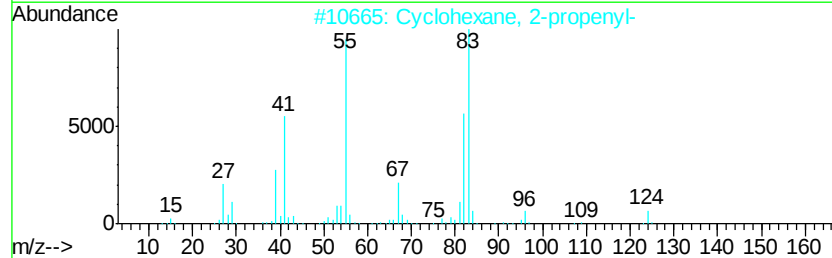
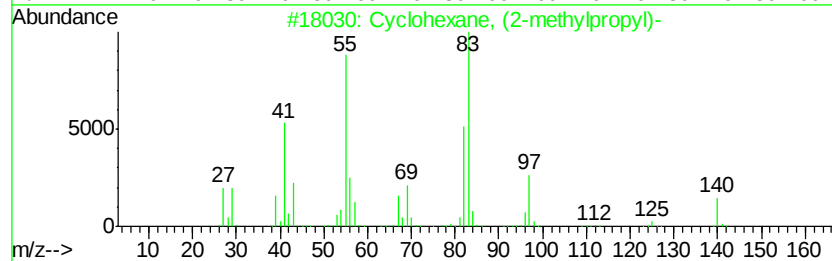
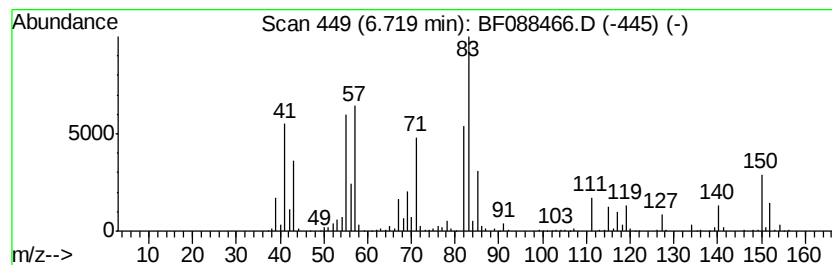
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TIC Integration Parameters: LSCINT.P

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Peak Number 6 unknown6.72 Concentration Rank 5

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.72 | 27.82 ng | 9565240 | 1,4-Dichlorobenzene-d4 | 6.55 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Cyclohexane, (2-methylpropyl)-      | 140 | C10H20   | 001678-98-4 | 38   |
| 2    |    | Cyclohexane, 2-propenyl-            | 124 | C9H16    | 002114-42-3 | 35   |
| 3    |    | 1-Propanone, 1-cyclohexyl-          | 140 | C9H16O   | 001123-86-0 | 30   |
| 4    |    | 1-Imidazol-1-yl-3-methylbut-2-en... | 150 | C8H10N2O | 061985-22-6 | 27   |
| 5    |    | 1-Butanone, 1-cyclohexyl-           | 154 | C10H18O  | 001462-27-7 | 27   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF062716\  
Data File : BF088466.D  
Acq On : 27 Jun 2016 22:21  
Operator : UM/SJ  
Sample : H3630-03  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
B1409-TP01N-1224

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

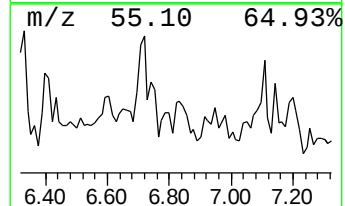
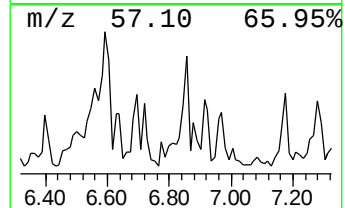
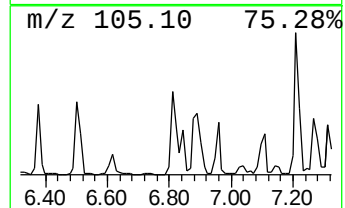
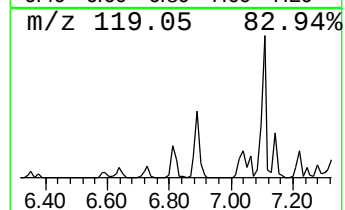
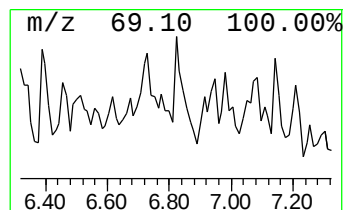
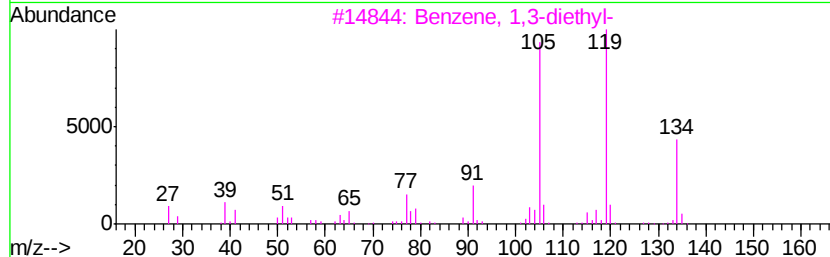
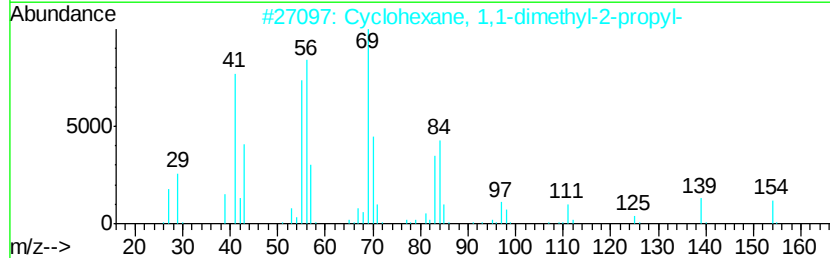
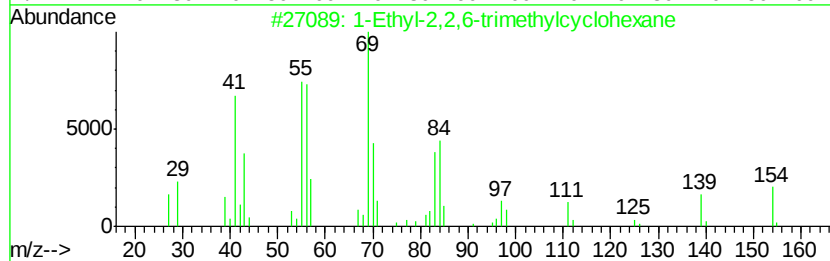
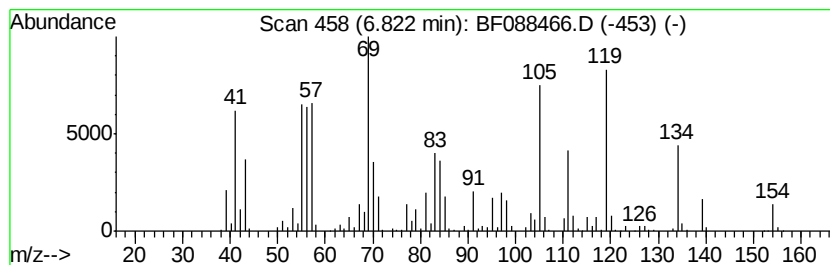
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TIC Integration Parameters: LSCINT.P

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Peak Number 7 1-Ethyl-2,2,6-trimethylcycl... Concentration Rank 13

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.82 | 16.00 ng | 5502080 | 1,4-Dichlorobenzene-d4 | 6.55 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 1-Ethyl-2,2,6-trimethylcyclohexane  | 154 | C11H22  | 071186-27-1 | 55   |
| 2    |    | Cyclohexane, 1,1-dimethyl-2-propyl- | 154 | C11H22  | 081983-71-3 | 55   |
| 3    |    | Benzene, 1,3-diethyl-               | 134 | C10H14  | 000141-93-5 | 48   |
| 4    |    | Benzene, 1,2-diethyl-               | 134 | C10H14  | 000135-01-3 | 48   |
| 5    |    | Benzene, 1,4-diethyl-               | 134 | C10H14  | 000105-05-5 | 47   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF062716\  
Data File : BF088466.D  
Acq On : 27 Jun 2016 22:21  
Operator : UM/SJ  
Sample : H3630-03  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
B1409-TP01N-1224

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

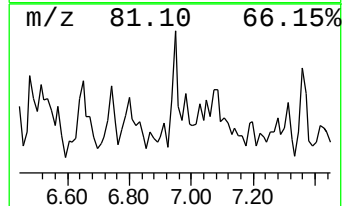
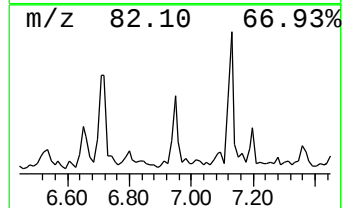
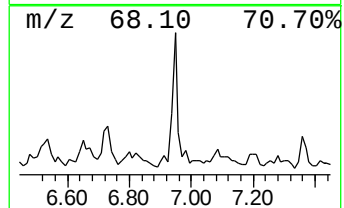
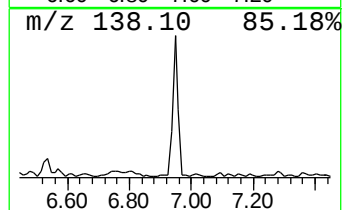
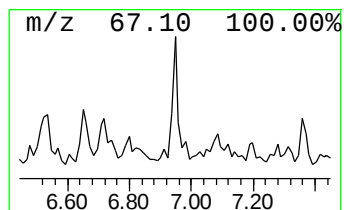
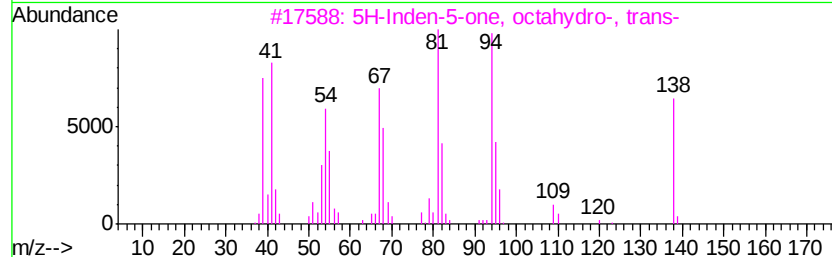
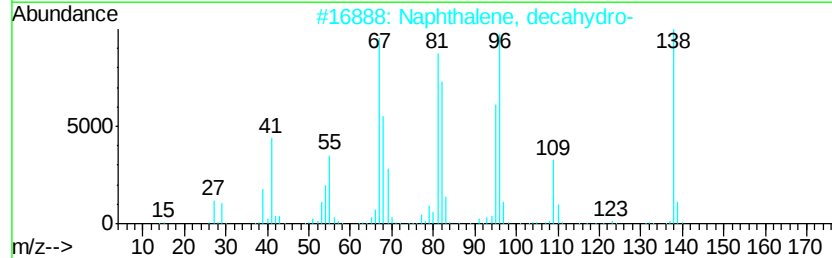
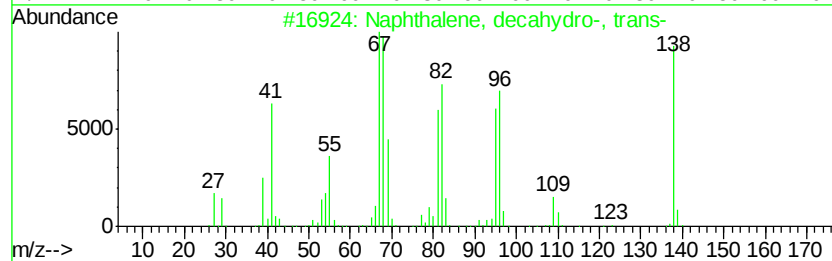
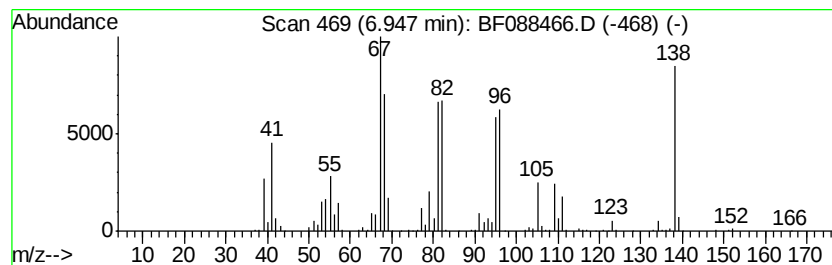
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TIC Integration Parameters: LSCINT.P

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Peak Number 8 Naphthalene, decahydro-, tr... Concentration Rank 14

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 6.95 | 14.55 ng | 5002130 | 1,4-Dichlorobenzene-d4 | 6.55 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, decahydro-, trans-     | 138 | C10H18  | 000493-02-7 | 95   |
| 2    |    | Naphthalene, decahydro-             | 138 | C10H18  | 000091-17-8 | 95   |
| 3    |    | 5H-Inden-5-one, octahydro-, trans-  | 138 | C9H14O  | 004668-81-9 | 72   |
| 4    |    | Bicyclo[3.1.0]hexan-2-one, 5-(1-... | 138 | C9H14O  | 000513-20-2 | 72   |
| 5    |    | Cyclohexene, 1-butyl-               | 138 | C10H18  | 003282-53-9 | 68   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF062716\  
Data File : BF088466.D  
Acq On : 27 Jun 2016 22:21  
Operator : UM/SJ  
Sample : H3630-03  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
B1409-TP01N-1224

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

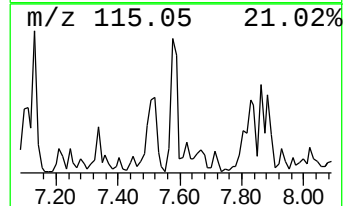
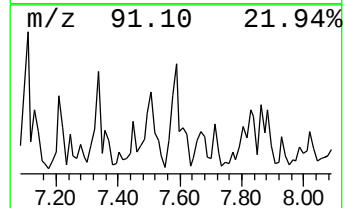
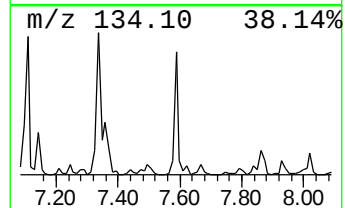
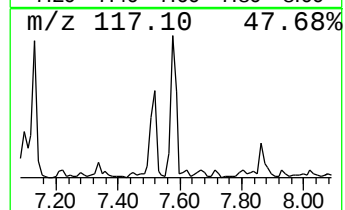
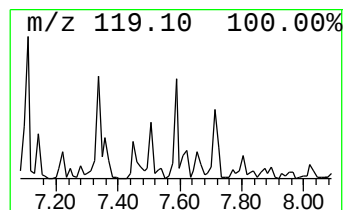
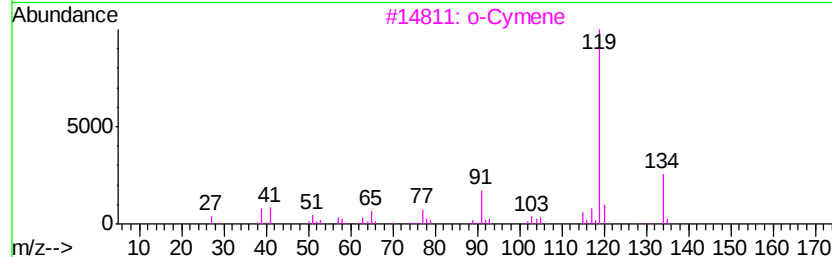
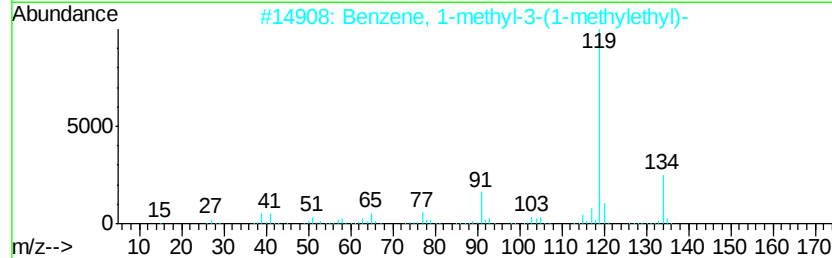
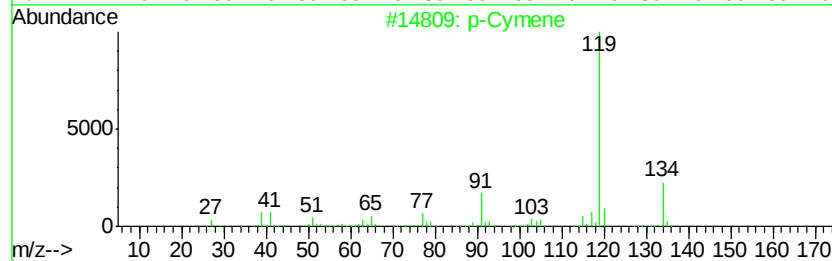
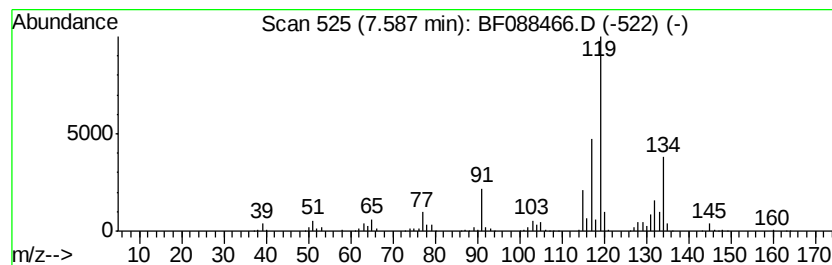
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 p-Cymene Concentration Rank 19

| R.T. | EstConc  | Area    | Relative to ISTD | R.T. |
|------|----------|---------|------------------|------|
| 7.59 | 13.35 ng | 4110620 | Napthalene-d8    | 7.84 |

| Hit# | of | Tentative ID                         | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------------|-----|---------|-------------|------|
| 1    | 5  | p-Cymene                             | 134 | C10H14  | 000099-87-6 | 64   |
| 2    |    | Benzene, 1-methyl-3-(1-methylethyl-) | 134 | C10H14  | 000535-77-3 | 64   |
| 3    |    | o-Cymene                             | 134 | C10H14  | 000527-84-4 | 64   |
| 4    |    | Benzene, 1,2,3,4-tetramethyl-        | 134 | C10H14  | 000488-23-3 | 64   |
| 5    |    | Benzene, 1,2,4,5-tetramethyl-        | 134 | C10H14  | 000095-93-2 | 64   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF062716\  
Data File : BF088466.D  
Acq On : 27 Jun 2016 22:21  
Operator : UM/SJ  
Sample : H3630-03  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
B1409-TP01N-1224

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

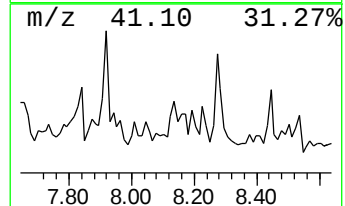
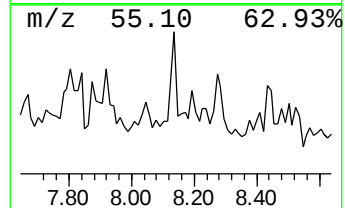
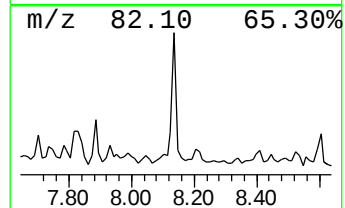
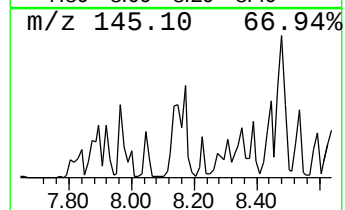
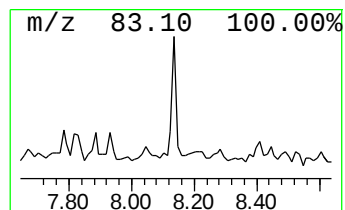
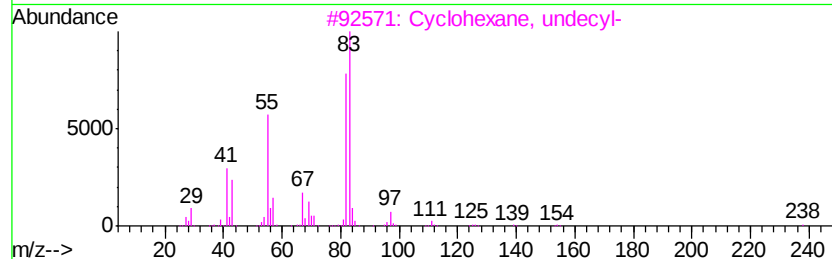
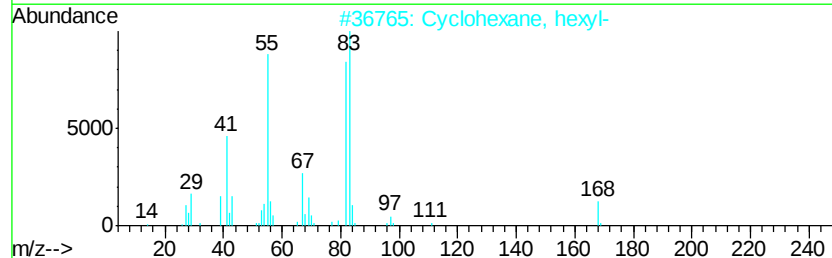
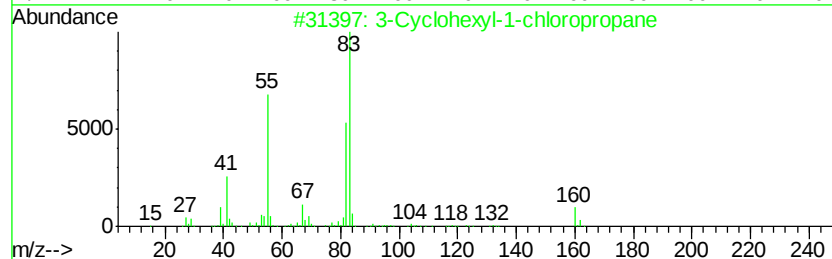
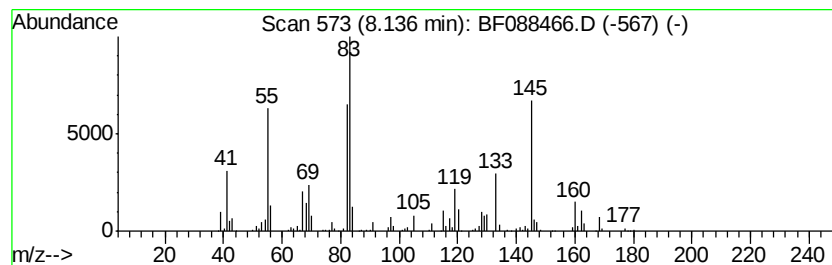
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TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 unknown8.14 Concentration Rank 16

| R.T. | EstConc  | Area    | Relative to ISTD | R.T. |
|------|----------|---------|------------------|------|
| 8.14 | 14.33 ng | 4412590 | Napthalene-d8    | 7.84 |

| Hit# | of | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|-------------------------------------|-----|---------|-------------|------|
| 1    | 5  | 3-Cyclohexyl-1-chloropropane        | 160 | C9H17Cl | 001124-62-5 | 46   |
| 2    |    | Cyclohexane, hexyl-                 | 168 | C12H24  | 004292-75-5 | 45   |
| 3    |    | Cyclohexane, undecyl-               | 238 | C17H34  | 054105-66-7 | 43   |
| 4    |    | Cyclohexane, (4-methylpentyl)-      | 168 | C12H24  | 061142-20-9 | 43   |
| 5    |    | 1H-Indene, 2,3-dihydro-1,4,7-tri... | 160 | C12H16  | 054340-87-3 | 40   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF062716\  
Data File : BF088466.D  
Acq On : 27 Jun 2016 22:21  
Operator : UM/SJ  
Sample : H3630-03  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
B1409-TP01N-1224

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

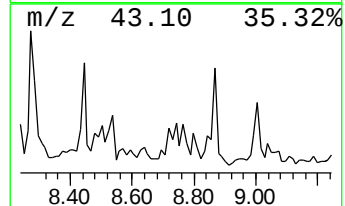
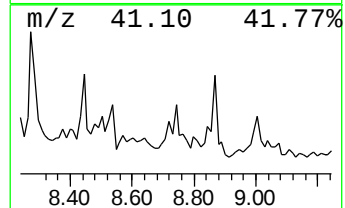
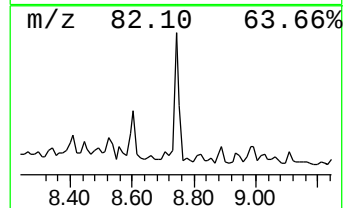
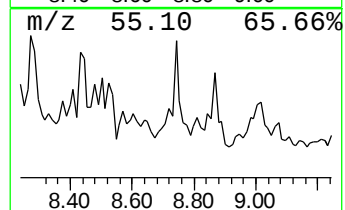
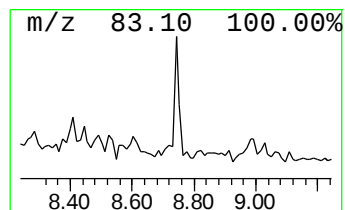
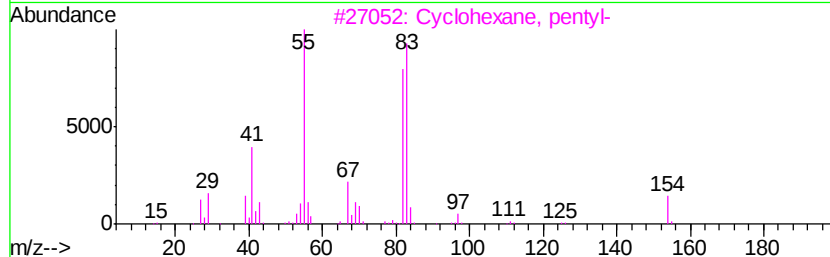
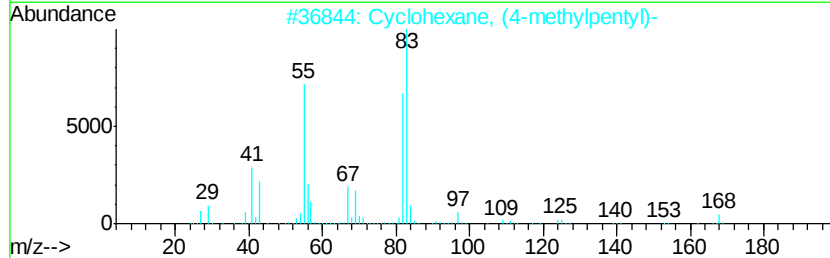
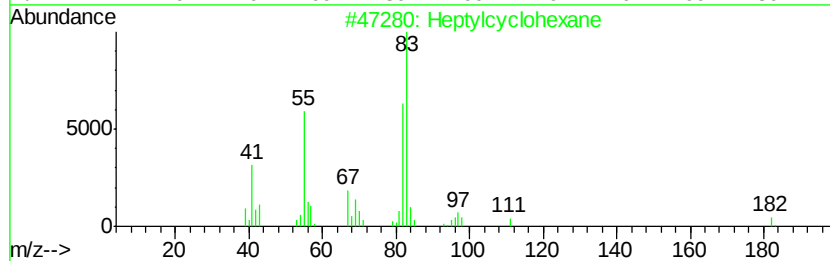
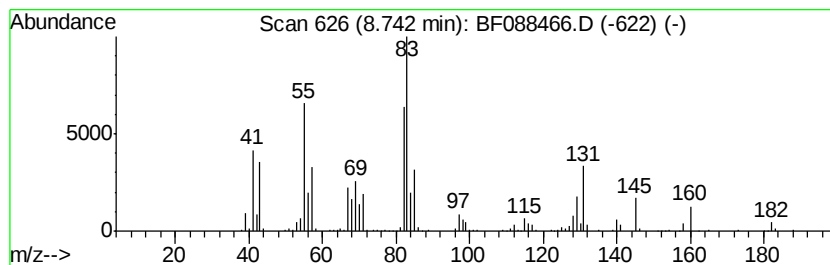
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 11 Heptylcyclohexane Concentration Rank 1

| R.T. | EstConc  | Area    | Relative to ISTD | R.T. |
|------|----------|---------|------------------|------|
| 8.74 | 79.50 ng | 2227810 | Acenaphthene-d10 | 9.58 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | Heptylcyclohexane              | 182 | C13H26  | 005617-41-4 | 52   |
| 2    |    | Cyclohexane, (4-methylpentyl)- | 168 | C12H24  | 061142-20-9 | 50   |
| 3    |    | Cyclohexane, pentyl-           | 154 | C11H22  | 004292-92-6 | 47   |
| 4    |    | Cyclohexane, (1-methylethyl)-  | 126 | C9H18   | 000696-29-7 | 47   |
| 5    |    | Cyclohexane, butyl-            | 140 | C10H20  | 001678-93-9 | 47   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF062716\  
Data File : BF088466.D  
Acq On : 27 Jun 2016 22:21  
Operator : UM/SJ  
Sample : H3630-03  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
B1409-TP01N-1224

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

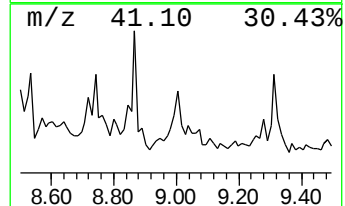
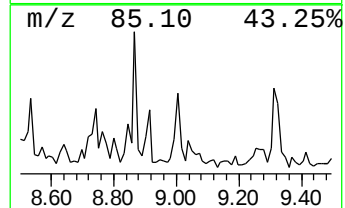
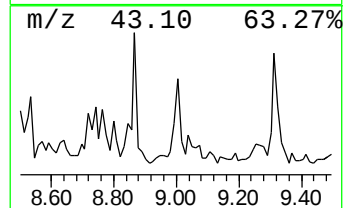
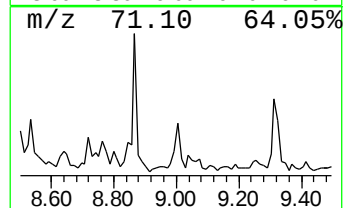
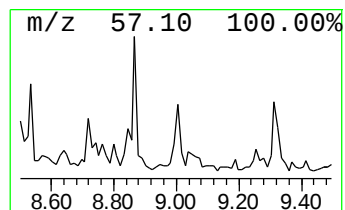
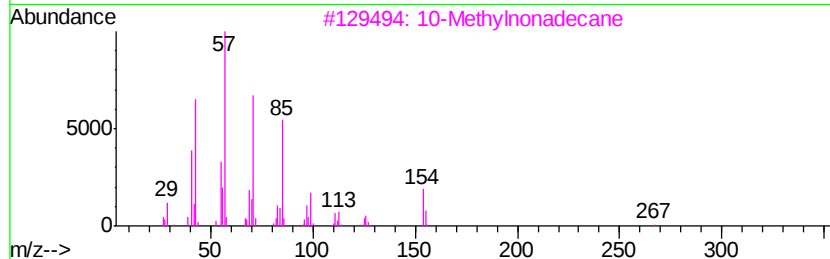
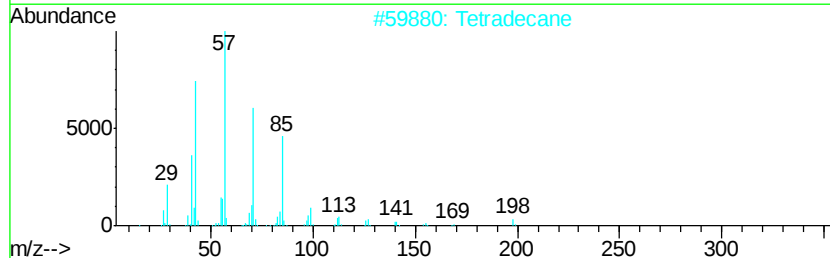
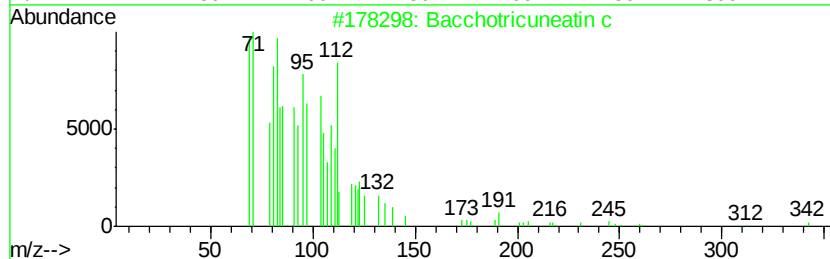
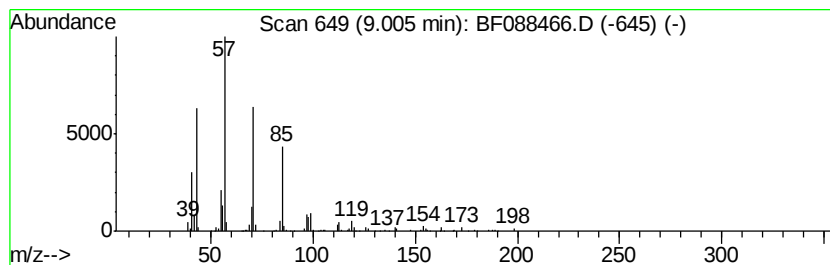
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 12 Bacchotricuneatin c Concentration Rank 6

| R.T. | EstConc  | Area   | Relative to ISTD | R.T. |
|------|----------|--------|------------------|------|
| 9.00 | 27.15 ng | 760769 | Acenaphthene-d10 | 9.58 |

| Hit# | of | Tentative ID        | MW  | MolForm  | CAS#        | Qual |
|------|----|---------------------|-----|----------|-------------|------|
| 1    | 5  | Bacchotricuneatin c | 342 | C20H22O5 | 066563-30-2 | 96   |
| 2    |    | Tetradecane         | 198 | C14H30   | 000629-59-4 | 86   |
| 3    |    | 10-Methylnonadecane | 282 | C20H42   | 056862-62-5 | 78   |
| 4    |    | Dodecane, 2-methyl- | 184 | C13H28   | 001560-97-0 | 72   |
| 5    |    | Nonadecane          | 268 | C19H40   | 000629-92-5 | 64   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF062716\  
Data File : BF088466.D  
Acq On : 27 Jun 2016 22:21  
Operator : UM/SJ  
Sample : H3630-03  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
B1409-TP01N-1224

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

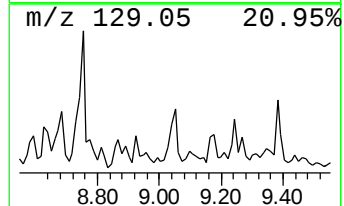
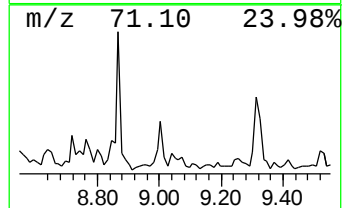
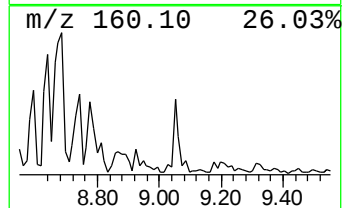
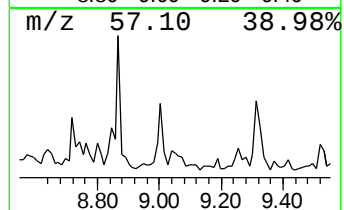
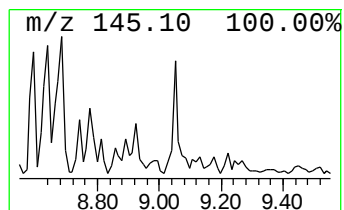
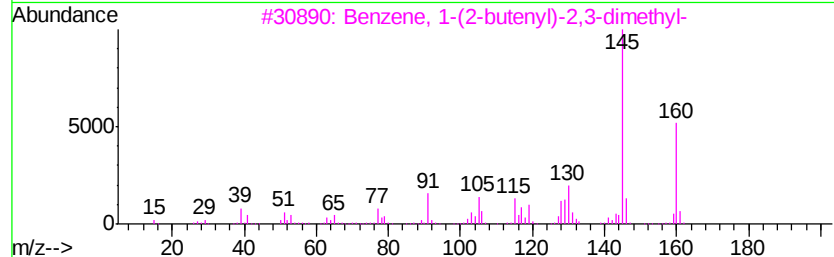
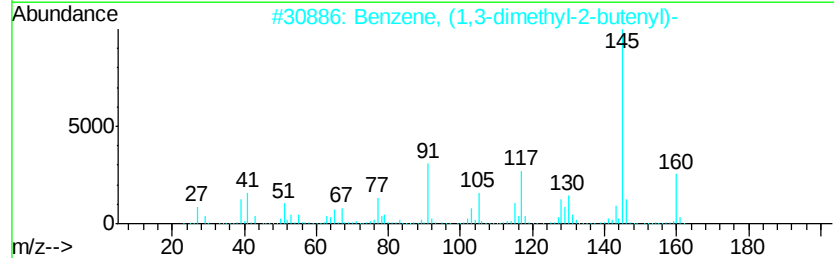
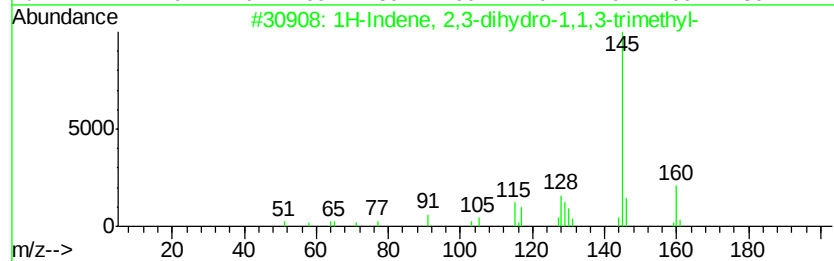
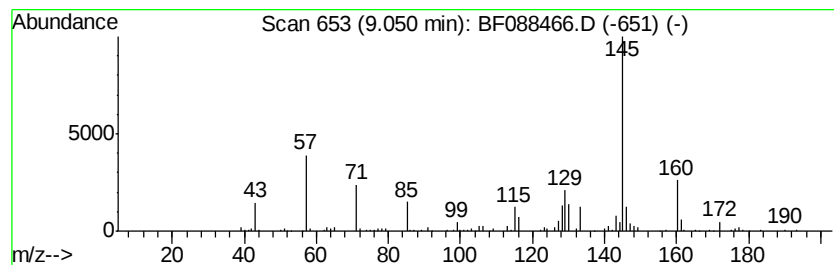
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 13 1H-Indene, 2,3-dihydro-1,1,... Concentration Rank 15

| R.T. | EstConc  | Area   | Relative to ISTD | R.T. |
|------|----------|--------|------------------|------|
| 9.05 | 14.50 ng | 406382 | Acenaphthene-d10 | 9.58 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | 1H-Indene, 2,3-dihydro-1,1,3-tri... | 160 | C12H16  | 002613-76-5 | 70   |
| 2    |    |   | Benzene, (1,3-dimethyl-2-butenyl)-  | 160 | C12H16  | 050704-01-3 | 70   |
| 3    |    |   | Benzene, 1-(2-butenyl)-2,3-dimet... | 160 | C12H16  | 054340-85-1 | 62   |
| 4    |    |   | Benzene, 4-(2-butenyl)-1,2-dimet... | 160 | C12H16  | 054340-86-2 | 62   |
| 5    |    |   | 1H-Indene, 2,3-dihydro-1,1,5-tri... | 160 | C12H16  | 040650-41-7 | 62   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF062716\  
Data File : BF088466.D  
Acq On : 27 Jun 2016 22:21  
Operator : UM/SJ  
Sample : H3630-03  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
B1409-TP01N-1224

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

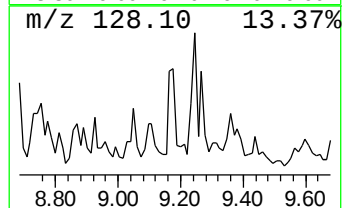
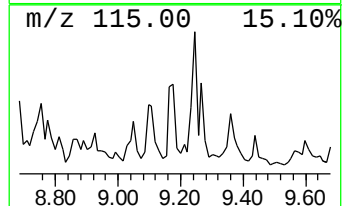
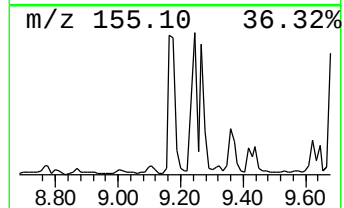
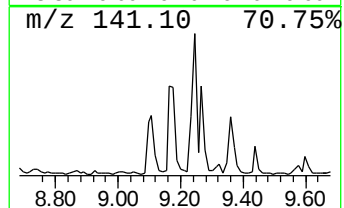
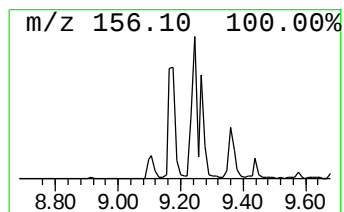
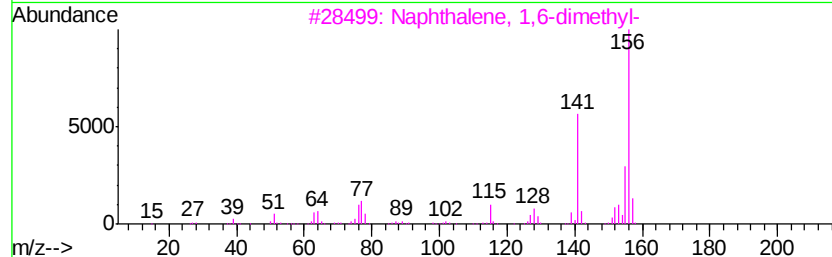
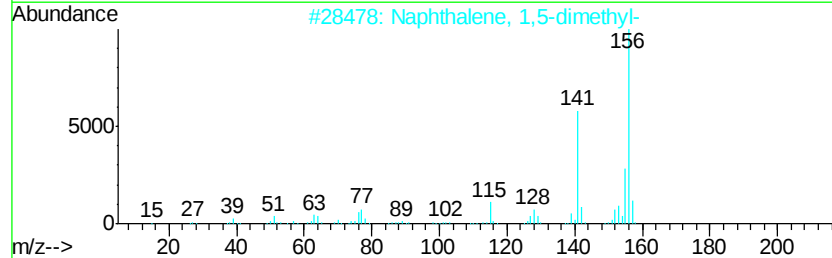
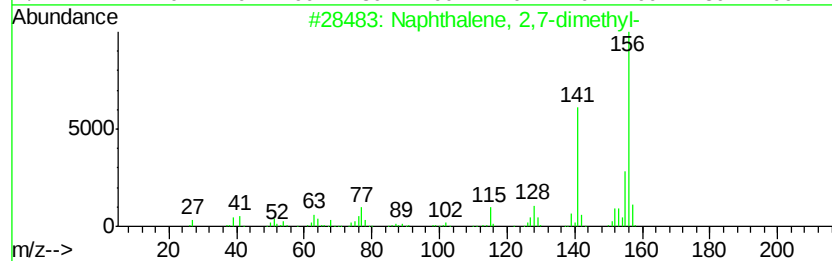
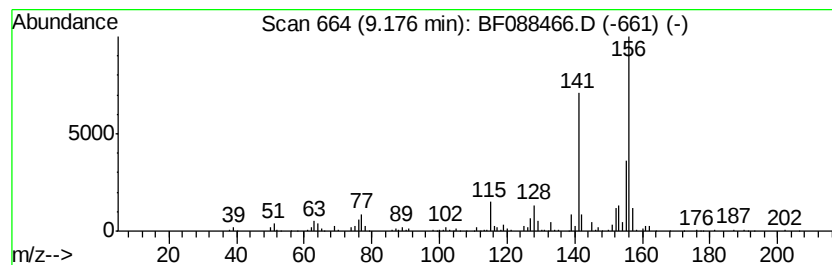
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 Naphthalene, 2,7-dimethyl- Concentration Rank 11

| R.T. | EstConc  | Area   | Relative to ISTD | R.T. |
|------|----------|--------|------------------|------|
| 9.18 | 22.81 ng | 639107 | Acenaphthene-d10 | 9.58 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 2,7-dimethyl- | 156 | C12H12  | 000582-16-1 | 97   |
| 2    |    | Naphthalene, 1,5-dimethyl- | 156 | C12H12  | 000571-61-9 | 97   |
| 3    |    | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 96   |
| 4    |    | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 96   |
| 5    |    | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 96   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF062716\  
Data File : BF088466.D  
Acq On : 27 Jun 2016 22:21  
Operator : UM/SJ  
Sample : H3630-03  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
B1409-TP01N-1224

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

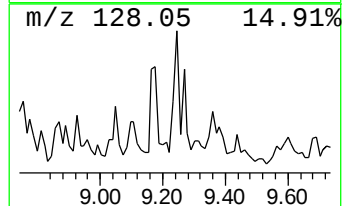
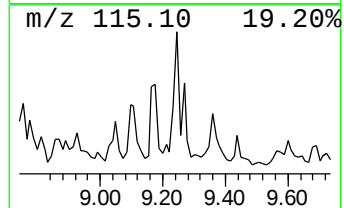
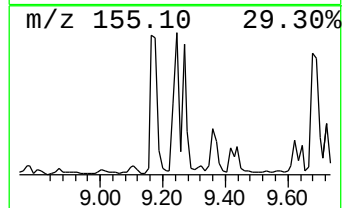
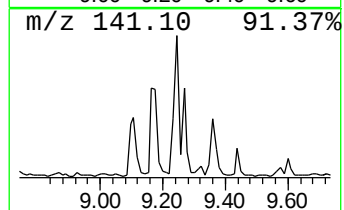
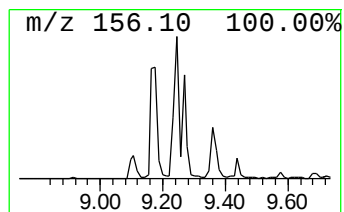
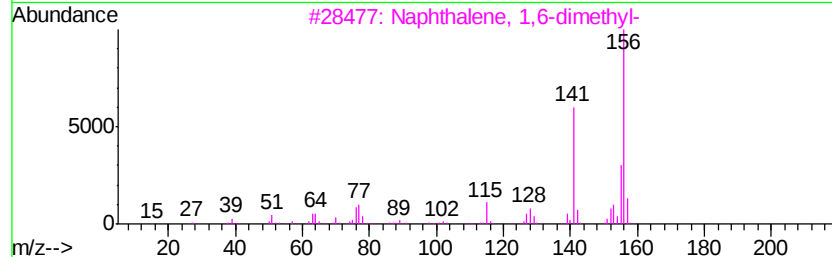
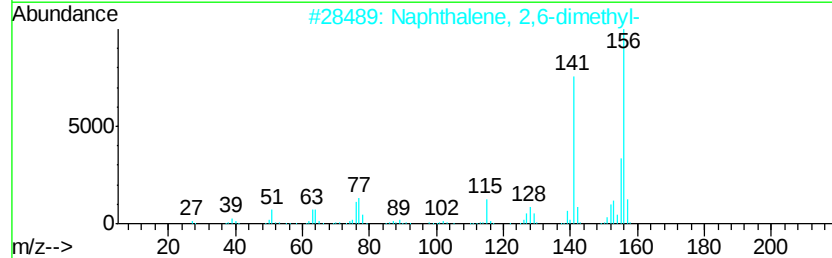
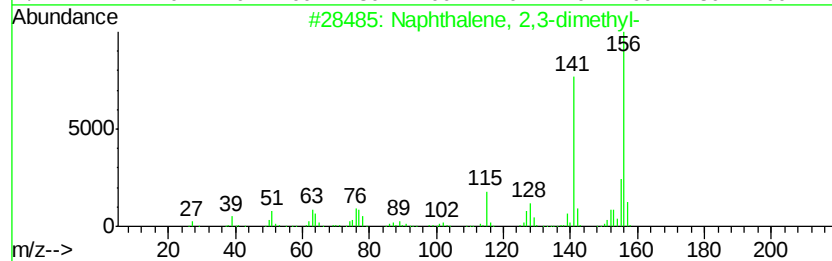
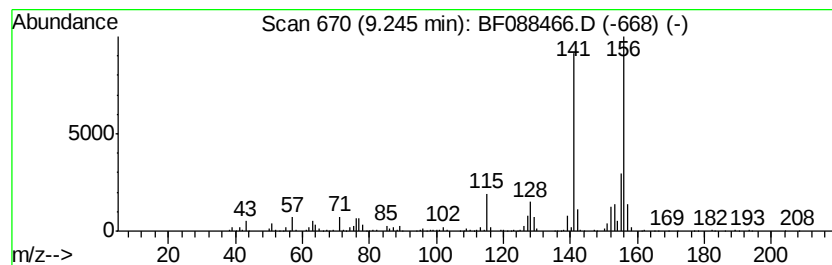
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 15 Naphthalene, 2,3-dimethyl- Concentration Rank 10

| R.T. | EstConc  | Area   | Relative to ISTD | R.T. |
|------|----------|--------|------------------|------|
| 9.24 | 22.83 ng | 639638 | Acenaphthene-d10 | 9.58 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, 2,3-dimethyl- | 156 | C12H12  | 000581-40-8 | 97   |
| 2    |    | Naphthalene, 2,6-dimethyl- | 156 | C12H12  | 000581-42-0 | 97   |
| 3    |    | Naphthalene, 1,6-dimethyl- | 156 | C12H12  | 000575-43-9 | 97   |
| 4    |    | Naphthalene, 1,4-dimethyl- | 156 | C12H12  | 000571-58-4 | 97   |
| 5    |    | Naphthalene, 1,7-dimethyl- | 156 | C12H12  | 000575-37-1 | 97   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF062716\  
Data File : BF088466.D  
Acq On : 27 Jun 2016 22:21  
Operator : UM/SJ  
Sample : H3630-03  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
B1409-TP01N-1224

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

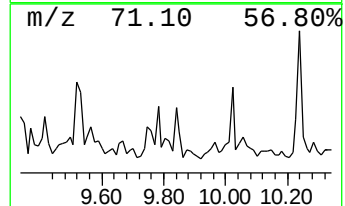
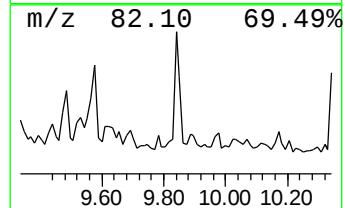
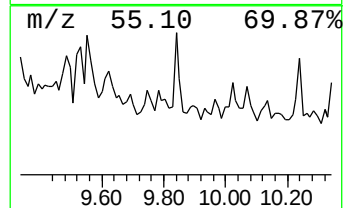
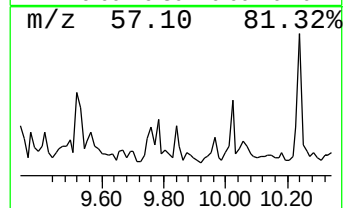
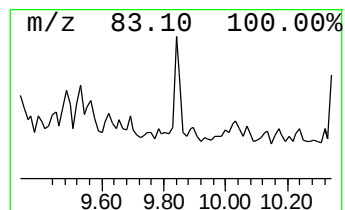
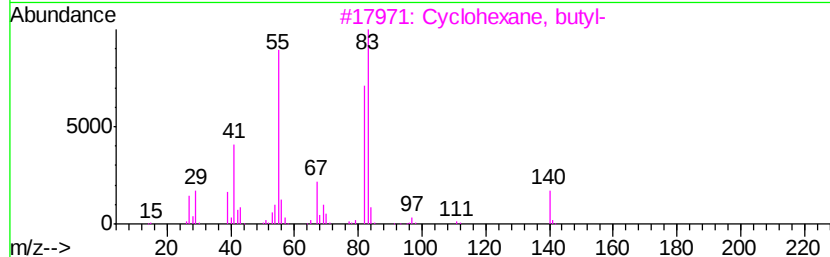
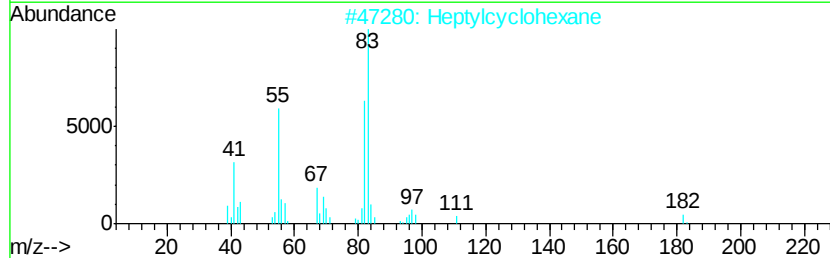
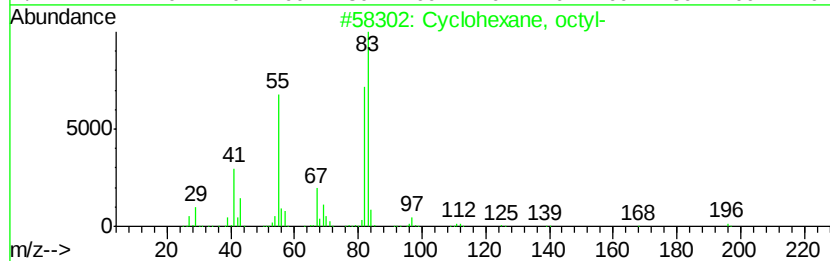
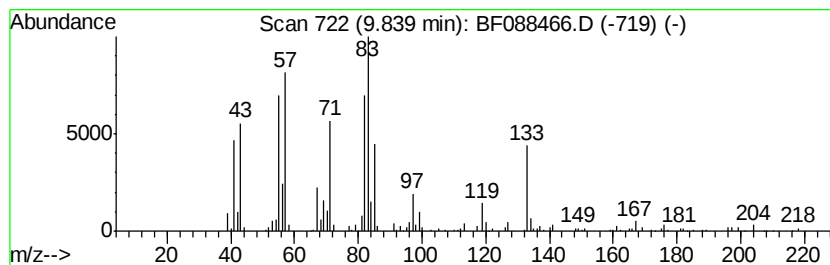
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 16 Cyclohexane, octyl- Concentration Rank 17

| R.T. | EstConc  | Area   | Relative to ISTD | R.T. |
|------|----------|--------|------------------|------|
| 9.84 | 14.17 ng | 397226 | Acenaphthene-d10 | 9.58 |

| Hit# | of | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------|-----|---------|-------------|------|
| 1    | 5  | Cyclohexane, octyl-  | 196 | C14H28  | 001795-15-9 | 50   |
| 2    |    | Heptylcyclohexane    | 182 | C13H26  | 005617-41-4 | 43   |
| 3    |    | Cyclohexane, butyl-  | 140 | C10H20  | 001678-93-9 | 38   |
| 4    |    | Cyclohexane, propyl- | 126 | C9H18   | 001678-92-8 | 38   |
| 5    |    | Cyclohexane, pentyl- | 154 | C11H22  | 004292-92-6 | 35   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF062716\  
Data File : BF088466.D  
Acq On : 27 Jun 2016 22:21  
Operator : UM/SJ  
Sample : H3630-03  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

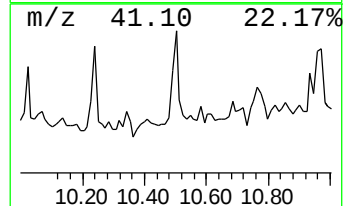
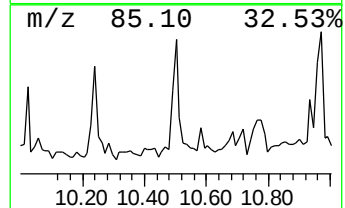
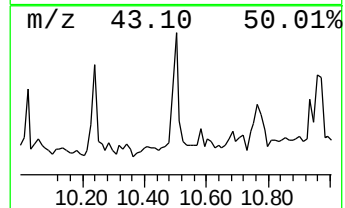
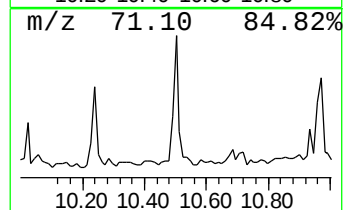
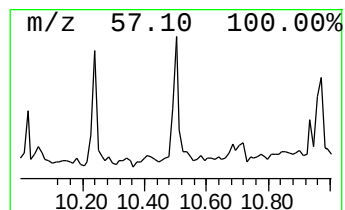
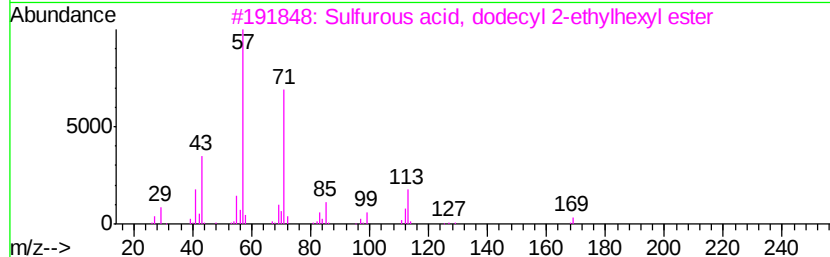
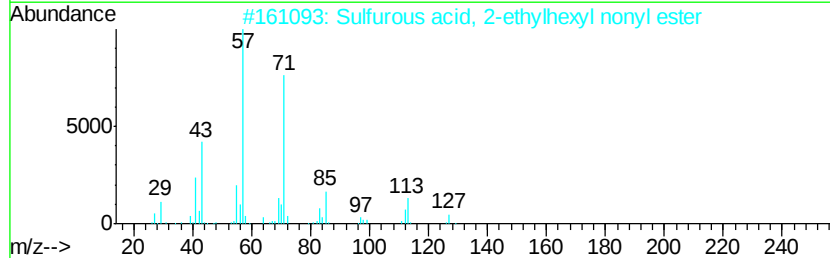
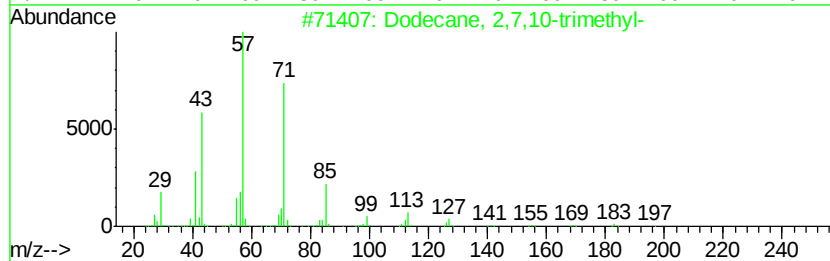
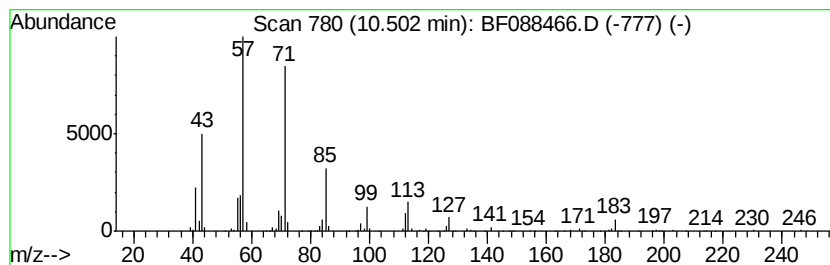
Instrument :  
BNA\_F  
ClientSampleId :  
B1409-TP01N-1224

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 17 Dodecane, 2,7,10-trimethyl- Concentration Rank 2

| R.T.    | EstConc         | Area                | Relative to ISTD | R.T.      |              |      |
|---------|-----------------|---------------------|------------------|-----------|--------------|------|
| 10.50   | 38.38 ng        | 498999              | Phenanthrene-d10 | 11.05     |              |      |
| Hit# of | 5               | Tentative ID        | MW               | MolForm   | CAS#         | Qual |
| 1       | Dodecane,       | 2,7,10-trimethyl-   | 212              | C15H32    | 074645-98-0  | 86   |
| 2       | Sulfurous acid, | 2-ethylhexyl non... | 320              | C17H36O3S | 1000309-19-2 | 80   |
| 3       | Sulfurous acid, | dodecyl 2-ethylh... | 362              | C20H42O3S | 1000309-19-5 | 72   |
| 4       | Dodecane,       | 4,6-dimethyl-       | 198              | C14H30    | 061141-72-8  | 72   |
| 5       | Sulfurous acid, | 2-ethylhexyl und... | 348              | C19H40O3S | 1000309-19-4 | 72   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF062716\  
Data File : BF088466.D  
Acq On : 27 Jun 2016 22:21  
Operator : UM/SJ  
Sample : H3630-03  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
B1409-TP01N-1224

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

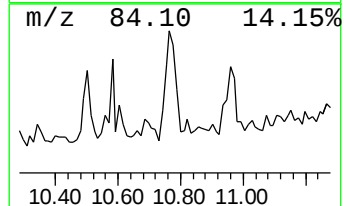
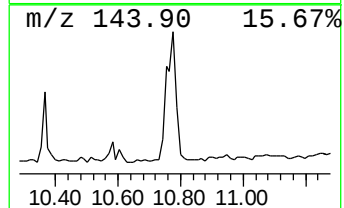
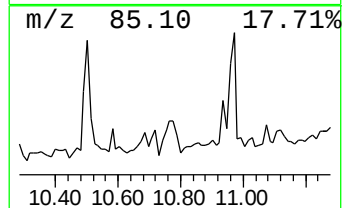
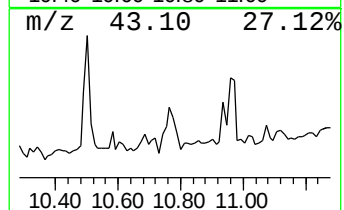
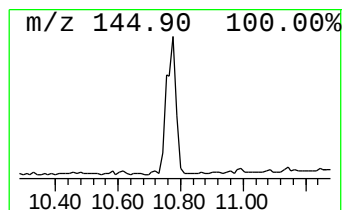
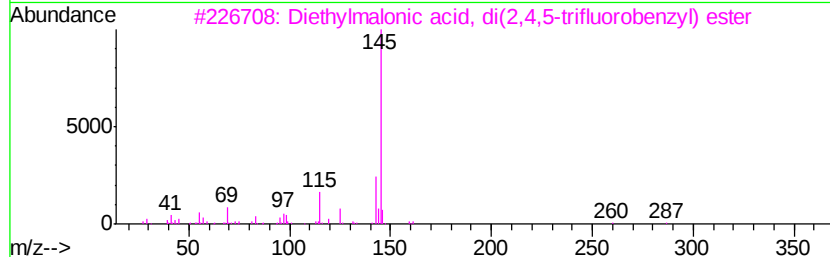
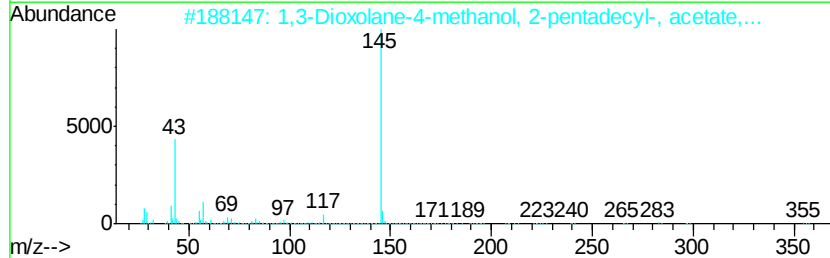
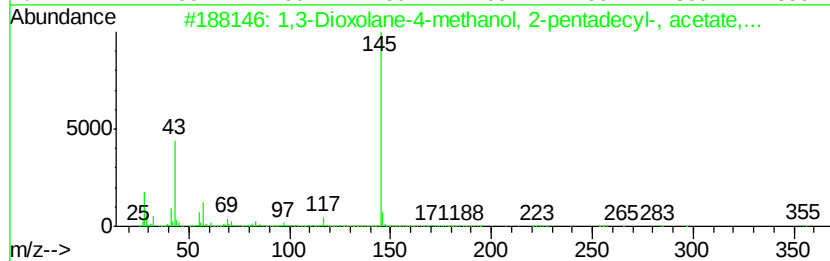
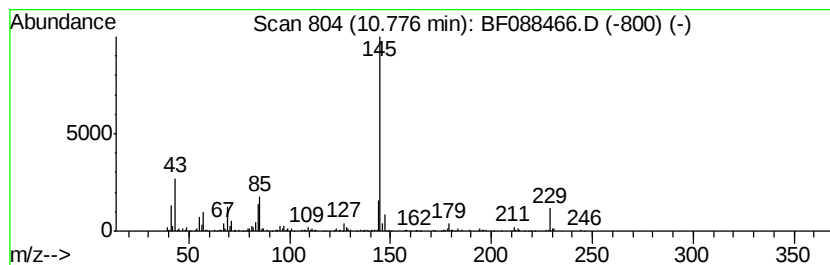
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 18 unknown10.78 Concentration Rank 8

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 10.78 | 26.45 ng | 343821 | Phenanthrene-d10 | 11.05 |

| Hit# | of | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|-------------------------------------|-----|------------|--------------|------|
| 1    | 5  | 1,3-Dioxolane-4-methanol, 2-pent... | 356 | C21H40O4   | 030889-29-3  | 47   |
| 2    |    | 1,3-Dioxolane-4-methanol, 2-pent... | 356 | C21H40O4   | 030889-32-8  | 47   |
| 3    |    | Diethylmalonic acid, di(2,4,5-tr... | 448 | C21H18F6O4 | 1000369-27-0 | 38   |
| 4    |    | 1,3-Dioxan-5-ol, 2-pentadecyl-, ... | 356 | C21H40O4   | 030889-23-7  | 38   |
| 5    |    | 1,3-Dioxan-5-ol, 2-pentadecyl-, ... | 356 | C21H40O4   | 030889-26-0  | 25   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF062716\  
Data File : BF088466.D  
Acq On : 27 Jun 2016 22:21  
Operator : UM/SJ  
Sample : H3630-03  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
B1409-TP01N-1224

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

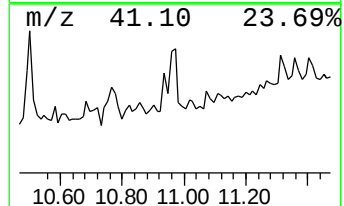
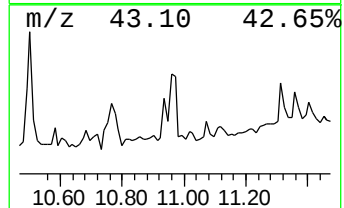
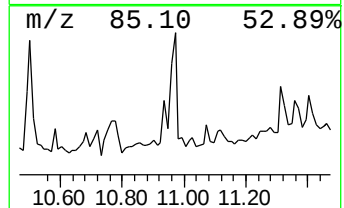
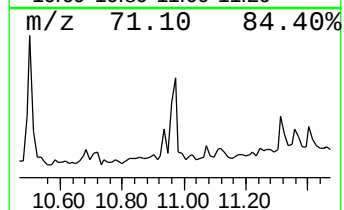
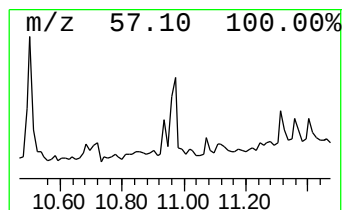
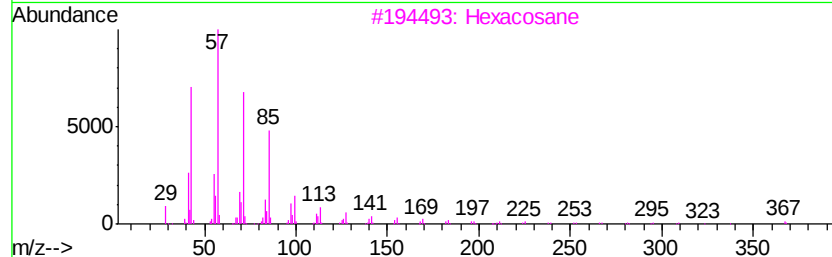
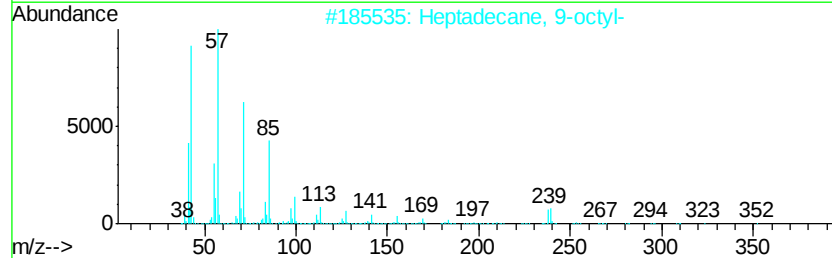
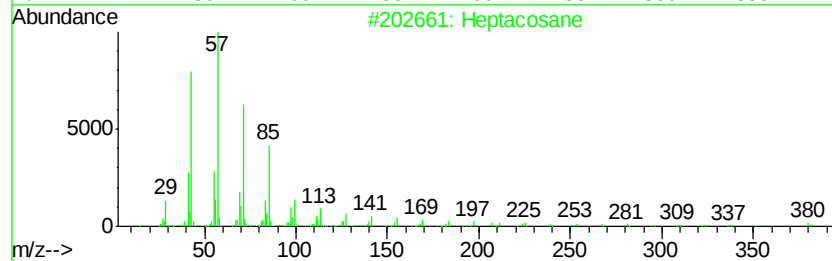
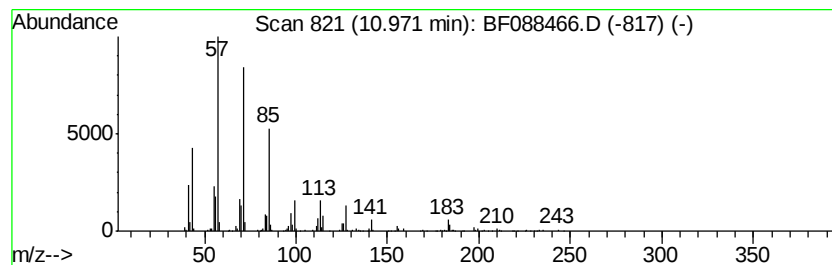
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 Heptacosane Concentration Rank 3

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 10.97 | 36.92 ng | 480018 | Phenanthrene-d10 | 11.05 |

| Hit# | of | Tentative ID           | MW  | MolForm | CAS#        | Qual |
|------|----|------------------------|-----|---------|-------------|------|
| 1    | 5  | Heptacosane            | 380 | C27H56  | 000593-49-7 | 90   |
| 2    |    | Heptadecane, 9-octyl-  | 352 | C25H52  | 007225-64-1 | 90   |
| 3    |    | Hexacosane             | 366 | C26H54  | 000630-01-3 | 87   |
| 4    |    | Heptadecane, 8-methyl- | 254 | C18H38  | 013287-23-5 | 86   |
| 5    |    | Heneicosane            | 296 | C21H44  | 000629-94-7 | 86   |



Data Path : Z:\HPCHEM1\BNA F\DATA\BF062716\  
Data File : BF088466.D  
Acq On : 27 Jun 2016 22:21  
Operator : UM/SJ  
Sample : H3630-03  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
B1409-TP01N-1224

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

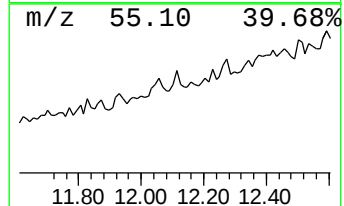
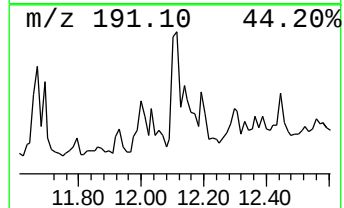
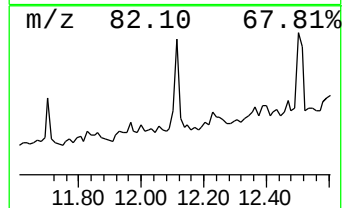
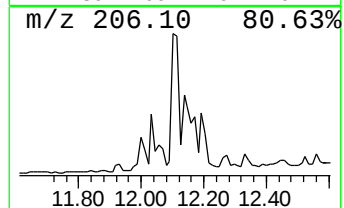
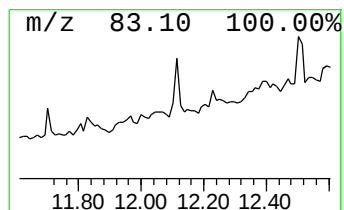
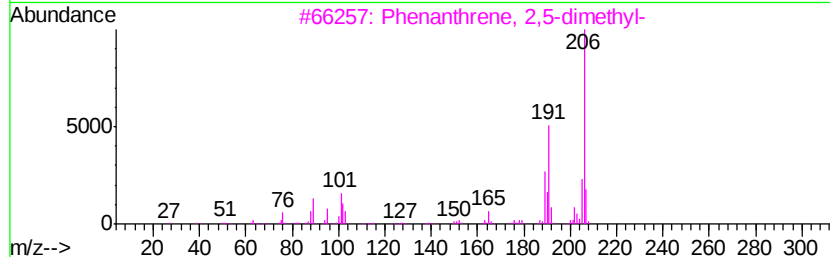
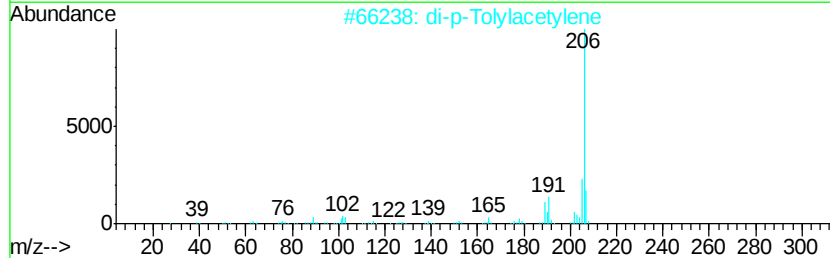
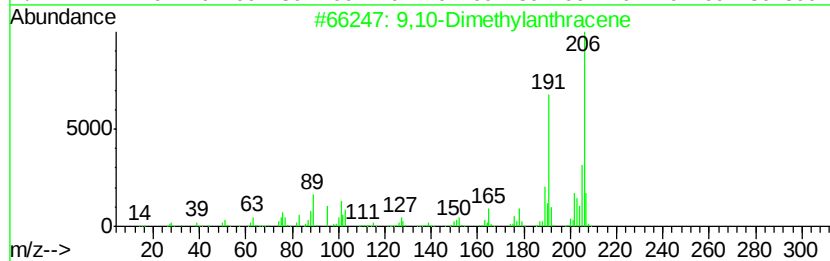
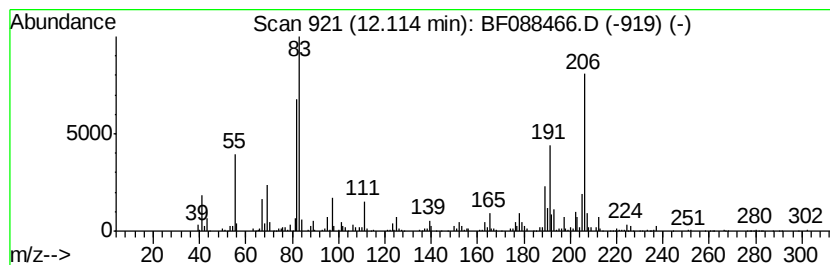
TIC Library : C:\Database\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 20 9,10-Dimethylantracene Concentration Rank 4

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 12.11 | 32.67 ng | 424710 | Phenanthrene-d10 | 11.05 |

| Hit# | of | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------|-----|---------|-------------|------|
| 1    | 5  | 9,10-Dimethylantracene      | 206 | C16H14  | 000781-43-1 | 64   |
| 2    |    | di-p-Tolylacetylene         | 206 | C16H14  | 002789-88-0 | 64   |
| 3    |    | Phenanthrene, 2,5-dimethyl- | 206 | C16H14  | 003674-66-6 | 50   |
| 4    |    | Phenanthrene, 2,3-dimethyl- | 206 | C16H14  | 003674-65-5 | 50   |
| 5    |    | Phenanthrene, 1,7-dimethyl- | 206 | C16H14  | 000483-87-4 | 41   |



Data Path : Z:\HPCHEM1\BNA\_F\DATA\BF062716\  
 Data File : BF088466.D  
 Acq On : 27 Jun 2016 22:21  
 Operator : UM/SJ  
 Sample : H3630-03  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 B1409-TP01N-1224

Quant Method : Z:\HPCHEM1\BNA\_F\METHODS\8270-BF060716.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |         |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|---------|------|
|                      |       |         |       |          | #                     | RT    | Resp    | Conc |
| Octane, 2,6-dimet... | 5.79  | 18.6    | ng    | 6409290  | 1                     | 6.55  | 6875940 | 20.0 |
| Heptane, 3-ethyl-... | 5.85  | 12.4    | ng    | 4253110  | 1                     | 6.55  | 6875940 | 20.0 |
| 1-Hexene, 3,5,5-t... | 6.06  | 27.0    | ng    | 9274210  | 1                     | 6.55  | 6875940 | 20.0 |
| unknown6.32          | 6.32  | 25.0    | ng    | 8588670  | 1                     | 6.55  | 6875940 | 20.0 |
| unknown6.64          | 6.64  | 13.6    | ng    | 4690100  | 1                     | 6.55  | 6875940 | 20.0 |
| unknown6.72          | 6.72  | 27.8    | ng    | 9565240  | 1                     | 6.55  | 6875940 | 20.0 |
| 1-Ethyl-2,2,6-tri... | 6.82  | 16.0    | ng    | 5502080  | 1                     | 6.55  | 6875940 | 20.0 |
| Naphthalene, deca... | 6.95  | 14.6    | ng    | 5002130  | 1                     | 6.55  | 6875940 | 20.0 |
| p-Cymene             | 7.59  | 13.4    | ng    | 4110620  | 2                     | 7.84  | 6157900 | 20.0 |
| unknown8.14          | 8.14  | 14.3    | ng    | 4412590  | 2                     | 7.84  | 6157900 | 20.0 |
| Heptylcyclohexane    | 8.74  | 79.5    | ng    | 2227810  | 3                     | 9.58  | 560466  | 20.0 |
| Bacchotricuneatin c  | 9.00  | 27.1    | ng    | 760769   | 3                     | 9.58  | 560466  | 20.0 |
| 1H-Indene, 2,3-di... | 9.05  | 14.5    | ng    | 406382   | 3                     | 9.58  | 560466  | 20.0 |
| Naphthalene, 2,7-... | 9.18  | 22.8    | ng    | 639107   | 3                     | 9.58  | 560466  | 20.0 |
| Naphthalene, 2,3-... | 9.24  | 22.8    | ng    | 639638   | 3                     | 9.58  | 560466  | 20.0 |
| Cyclohexane, octyl-  | 9.84  | 14.2    | ng    | 397226   | 3                     | 9.58  | 560466  | 20.0 |
| Dodecane, 2,7,10-... | 10.50 | 38.4    | ng    | 498999   | 4                     | 11.05 | 260018  | 20.0 |
| unknown10.78         | 10.78 | 26.4    | ng    | 343821   | 4                     | 11.05 | 260018  | 20.0 |
| Heptacosane          | 10.97 | 36.9    | ng    | 480018   | 4                     | 11.05 | 260018  | 20.0 |
| 9,10-Dimethylanth... | 12.11 | 32.7    | ng    | 424710   | 4                     | 11.05 | 260018  | 20.0 |