

Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF102919\  
 Data File : BF117580.D  
 Acq On : 29 Oct 2019 10:30  
 Operator : JU  
 Sample : K5563-07  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 TP-1

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:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF101819.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         | 4.928       | 478           | 483         | 496          | rBV      | 108929         | 162425        | 18.28%          | 0.951%        |
| 2         | 5.357       | 553           | 556         | 575          | rBV      | 657469         | 805161        | 90.63%          | 4.712%        |
| 3         | 5.516       | 580           | 583         | 589          | rVB2     | 20543          | 26229         | 2.95%           | 0.154%        |
| 4         | 5.969       | 658           | 660         | 663          | rBV      | 21974          | 23758         | 2.67%           | 0.139%        |
| 5         | 6.251       | 705           | 708         | 711          | rBV3     | 40426          | 41829         | 4.71%           | 0.245%        |
| 6         | 6.328       | 717           | 721         | 724          | rBV2     | 23106          | 27492         | 3.09%           | 0.161%        |
| 7         | 6.386       | 727           | 731         | 742          | rBV      | 660943         | 882241        | 99.30%          | 5.164%        |
| 8         | 6.504       | 748           | 751         | 756          | rVV      | 972104         | 888441        | 100.00%         | 5.200%        |
| 9         | 6.551       | 756           | 759         | 766          | rVB3     | 93122          | 124745        | 14.04%          | 0.730%        |
| 10        | 6.692       | 779           | 783         | 785          | rVV3     | 14549          | 20807         | 2.34%           | 0.122%        |
| 11        | 6.728       | 785           | 789         | 795          | rVB      | 337817         | 352713        | 39.70%          | 2.064%        |
| 12        | 6.781       | 795           | 798         | 801          | rBV      | 54262          | 57577         | 6.48%           | 0.337%        |
| 13        | 6.880       | 811           | 815         | 822          | rVB      | 704062         | 687082        | 77.34%          | 4.021%        |
| 14        | 6.957       | 822           | 828         | 829          | rBV7     | 18516          | 24978         | 2.81%           | 0.146%        |
| 15        | 6.975       | 829           | 831         | 833          | rBV      | 37695          | 32634         | 3.67%           | 0.191%        |
| 16        | 7.004       | 833           | 836         | 838          | rVB2     | 53335          | 56559         | 6.37%           | 0.331%        |
| 17        | 7.045       | 838           | 843         | 845          | rBV2     | 110151         | 100191        | 11.28%          | 0.586%        |
| 18        | 7.069       | 845           | 847         | 850          | rVB2     | 24282          | 22362         | 2.52%           | 0.131%        |
| 19        | 7.116       | 852           | 855         | 858          | rBV2     | 109033         | 96844         | 10.90%          | 0.567%        |
| 20        | 7.192       | 863           | 868         | 870          | rBV2     | 69652          | 91950         | 10.35%          | 0.538%        |
| 21        | 7.216       | 870           | 872         | 874          | rVV      | 58582          | 48042         | 5.41%           | 0.281%        |
| 22        | 7.257       | 874           | 879         | 882          | rVV      | 147369         | 168447        | 18.96%          | 0.986%        |
| 23        | 7.292       | 882           | 885         | 893          | rVB      | 528451         | 592872        | 66.73%          | 3.470%        |
| 24        | 7.369       | 893           | 898         | 902          | rVB5     | 113462         | 151766        | 17.08%          | 0.888%        |
| 25        | 7.410       | 902           | 905         | 907          | rBV      | 64018          | 68084         | 7.66%           | 0.398%        |
| 26        | 7.439       | 907           | 910         | 914          | rVB2     | 64823          | 74414         | 8.38%           | 0.436%        |
| 27        | 7.475       | 914           | 916         | 918          | rBV2     | 58803          | 59290         | 6.67%           | 0.347%        |
| 28        | 7.498       | 918           | 920         | 922          | rVV      | 130859         | 104803        | 11.80%          | 0.613%        |
| 29        | 7.528       | 922           | 925         | 928          | rVB2     | 160834         | 150409        | 16.93%          | 0.880%        |
| 30        | 7.610       | 934           | 939         | 942          | rBV4     | 70425          | 115932        | 13.05%          | 0.679%        |
| 31        | 7.657       | 942           | 947         | 950          | rVV3     | 156146         | 205541        | 23.14%          | 1.203%        |
| 32        | 7.692       | 950           | 953         | 956          | rVB2     | 47902          | 41957         | 4.72%           | 0.246%        |
| 33        | 7.751       | 956           | 963         | 965          | rBV4     | 136075         | 188292        | 21.19%          | 1.102%        |
| 34        | 7.775       | 965           | 967         | 971          | rVB2     | 108873         | 102796        | 11.57%          | 0.602%        |

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 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 TP-1

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:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF101819.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |        |        |        |        |
|----|--------|------|------|------|------|--------|--------|--------|--------|
| 35 | 7.816  | 971  | 974  | 978  | rBV2 | 82076  | 118798 | 13.37% | 0.695% |
| 36 | 7.875  | 982  | 984  | 988  | rVB  | 95336  | 92529  | 10.41% | 0.542% |
| 37 | 7.910  | 988  | 990  | 992  | rBV  | 25618  | 20472  | 2.30%  | 0.120% |
| 38 | 7.957  | 992  | 998  | 1001 | rBV5 | 69203  | 118432 | 13.33% | 0.693% |
| 39 | 8.010  | 1001 | 1007 | 1009 | rVV  | 446855 | 584785 | 65.82% | 3.423% |
| 40 | 8.057  | 1013 | 1015 | 1020 | rVV4 | 108775 | 159745 | 17.98% | 0.935% |
| 41 | 8.098  | 1020 | 1022 | 1025 | rVB  | 68449  | 57286  | 6.45%  | 0.335% |
| 42 | 8.133  | 1025 | 1028 | 1030 | rVB2 | 62280  | 58253  | 6.56%  | 0.341% |
| 43 | 8.163  | 1030 | 1033 | 1035 | rBV2 | 42782  | 37078  | 4.17%  | 0.217% |
| 44 | 8.186  | 1035 | 1037 | 1040 | rVB  | 45734  | 33933  | 3.82%  | 0.199% |
| 45 | 8.298  | 1052 | 1056 | 1060 | rVV5 | 47145  | 68264  | 7.68%  | 0.400% |
| 46 | 8.386  | 1068 | 1071 | 1074 | rVB4 | 31924  | 35262  | 3.97%  | 0.206% |
| 47 | 9.080  | 1184 | 1189 | 1200 | rBV  | 938026 | 835666 | 94.06% | 4.891% |
| 48 | 9.480  | 1253 | 1257 | 1265 | rVB  | 138966 | 136374 | 15.35% | 0.798% |
| 49 | 9.757  | 1300 | 1304 | 1309 | rBV  | 493798 | 450286 | 50.68% | 2.635% |
| 50 | 10.192 | 1375 | 1378 | 1383 | rVB2 | 18312  | 20832  | 2.34%  | 0.122% |
| 51 | 10.310 | 1394 | 1398 | 1403 | rVB2 | 17631  | 21660  | 2.44%  | 0.127% |
| 52 | 10.427 | 1414 | 1418 | 1423 | rBV  | 134763 | 138415 | 15.58% | 0.810% |
| 53 | 10.557 | 1436 | 1440 | 1449 | rVV  | 464431 | 477835 | 53.78% | 2.797% |
| 54 | 10.627 | 1449 | 1452 | 1456 | rVV  | 124125 | 136392 | 15.35% | 0.798% |
| 55 | 10.674 | 1458 | 1460 | 1465 | rVB  | 26519  | 24697  | 2.78%  | 0.145% |
| 56 | 11.245 | 1554 | 1557 | 1560 | rBV  | 420856 | 407979 | 45.92% | 2.388% |
| 57 | 11.268 | 1560 | 1561 | 1566 | rVV  | 152154 | 124267 | 13.99% | 0.727% |
| 58 | 11.327 | 1568 | 1571 | 1574 | rBV  | 34541  | 30552  | 3.44%  | 0.179% |
| 59 | 11.492 | 1594 | 1599 | 1602 | rBV  | 16020  | 20834  | 2.35%  | 0.122% |
| 60 | 11.774 | 1644 | 1647 | 1654 | rVB3 | 119066 | 139052 | 15.65% | 0.814% |
| 61 | 11.863 | 1658 | 1662 | 1665 | rBV  | 36645  | 37138  | 4.18%  | 0.217% |
| 62 | 12.068 | 1694 | 1697 | 1707 | rVB  | 111473 | 155411 | 17.49% | 0.910% |
| 63 | 12.298 | 1733 | 1736 | 1738 | rVV  | 24366  | 26997  | 3.04%  | 0.158% |
| 64 | 12.321 | 1738 | 1740 | 1751 | rVV2 | 121845 | 188701 | 21.24% | 1.104% |
| 65 | 12.398 | 1751 | 1753 | 1760 | rVB  | 54038  | 62026  | 6.98%  | 0.363% |
| 66 | 12.462 | 1760 | 1764 | 1767 | rBV  | 294877 | 256179 | 28.83% | 1.499% |
| 67 | 12.492 | 1767 | 1769 | 1773 | rVV3 | 32367  | 34626  | 3.90%  | 0.203% |
| 68 | 12.557 | 1777 | 1780 | 1782 | rVV  | 34192  | 34544  | 3.89%  | 0.202% |
| 69 | 12.610 | 1786 | 1789 | 1793 | rVV4 | 36941  | 53822  | 6.06%  | 0.315% |
| 70 | 12.692 | 1797 | 1803 | 1809 | rVV  | 245717 | 268302 | 30.20% | 1.570% |
| 71 | 12.827 | 1822 | 1826 | 1829 | rBV  | 599173 | 512514 | 57.69% | 3.000% |

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Instrument :  
 BNA\_F  
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 TP-1

Integration Parameters: rteint.p  
 Integrator: RTE  
 Smoothing : ON  
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 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:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF101819.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|     |        |      |      |      |      |        |        |        |        |
|-----|--------|------|------|------|------|--------|--------|--------|--------|
| 72  | 12.915 | 1829 | 1841 | 1848 | rVV3 | 130620 | 432955 | 48.73% | 2.534% |
| 73  | 13.027 | 1857 | 1860 | 1862 | rVV2 | 34494  | 36163  | 4.07%  | 0.212% |
| 74  | 13.051 | 1862 | 1864 | 1868 | rVB  | 24744  | 24218  | 2.73%  | 0.142% |
| 75  | 13.398 | 1918 | 1923 | 1926 | rBV2 | 34082  | 42337  | 4.77%  | 0.248% |
| 76  | 13.545 | 1944 | 1948 | 1954 | rBV  | 22715  | 32556  | 3.66%  | 0.191% |
| 77  | 13.645 | 1961 | 1965 | 1971 | rBV5 | 31220  | 60478  | 6.81%  | 0.354% |
| 78  | 13.727 | 1971 | 1979 | 1982 | rBV3 | 96774  | 155904 | 17.55% | 0.912% |
| 79  | 13.857 | 1993 | 2001 | 2003 | rBV  | 160731 | 308002 | 34.67% | 1.803% |
| 80  | 13.892 | 2003 | 2007 | 2010 | rVV2 | 320793 | 432256 | 48.65% | 2.530% |
| 81  | 13.915 | 2010 | 2011 | 2017 | rVB  | 100736 | 93518  | 10.53% | 0.547% |
| 82  | 13.998 | 2023 | 2025 | 2028 | rVB2 | 22527  | 19552  | 2.20%  | 0.114% |
| 83  | 14.039 | 2028 | 2032 | 2039 | rBV  | 107702 | 124097 | 13.97% | 0.726% |
| 84  | 14.286 | 2069 | 2074 | 2076 | rBV2 | 31333  | 50023  | 5.63%  | 0.293% |
| 85  | 14.345 | 2080 | 2084 | 2089 | rVB  | 212068 | 197925 | 22.28% | 1.158% |
| 86  | 14.509 | 2107 | 2112 | 2120 | rVB  | 106996 | 195192 | 21.97% | 1.142% |
| 87  | 14.639 | 2127 | 2134 | 2138 | rBV  | 296170 | 344341 | 38.76% | 2.015% |
| 88  | 14.921 | 2178 | 2182 | 2184 | rBV2 | 117224 | 144420 | 16.26% | 0.845% |
| 89  | 14.956 | 2184 | 2188 | 2197 | rVB2 | 325628 | 411430 | 46.31% | 2.408% |
| 90  | 15.027 | 2197 | 2200 | 2206 | rVB  | 33083  | 43878  | 4.94%  | 0.257% |
| 91  | 15.115 | 2210 | 2215 | 2220 | rBV4 | 35599  | 65394  | 7.36%  | 0.383% |
| 92  | 15.209 | 2227 | 2231 | 2234 | rVB  | 68789  | 77188  | 8.69%  | 0.452% |
| 93  | 15.268 | 2238 | 2241 | 2244 | rBV  | 89615  | 109210 | 12.29% | 0.639% |
| 94  | 15.315 | 2244 | 2249 | 2255 | rVB2 | 250282 | 520970 | 58.64% | 3.049% |
| 95  | 15.715 | 2311 | 2317 | 2323 | rBV  | 142854 | 220330 | 24.80% | 1.290% |
| 96  | 15.951 | 2353 | 2357 | 2362 | rBV7 | 15160  | 26045  | 2.93%  | 0.152% |
| 97  | 16.186 | 2391 | 2397 | 2402 | rVB2 | 66103  | 107175 | 12.06% | 0.627% |
| 98  | 16.745 | 2485 | 2492 | 2500 | rVB3 | 57730  | 139884 | 15.74% | 0.819% |
| 99  | 16.909 | 2514 | 2520 | 2524 | rBV8 | 13344  | 24333  | 2.74%  | 0.142% |
| 100 | 17.174 | 2560 | 2565 | 2571 | rVB2 | 33587  | 67458  | 7.59%  | 0.395% |

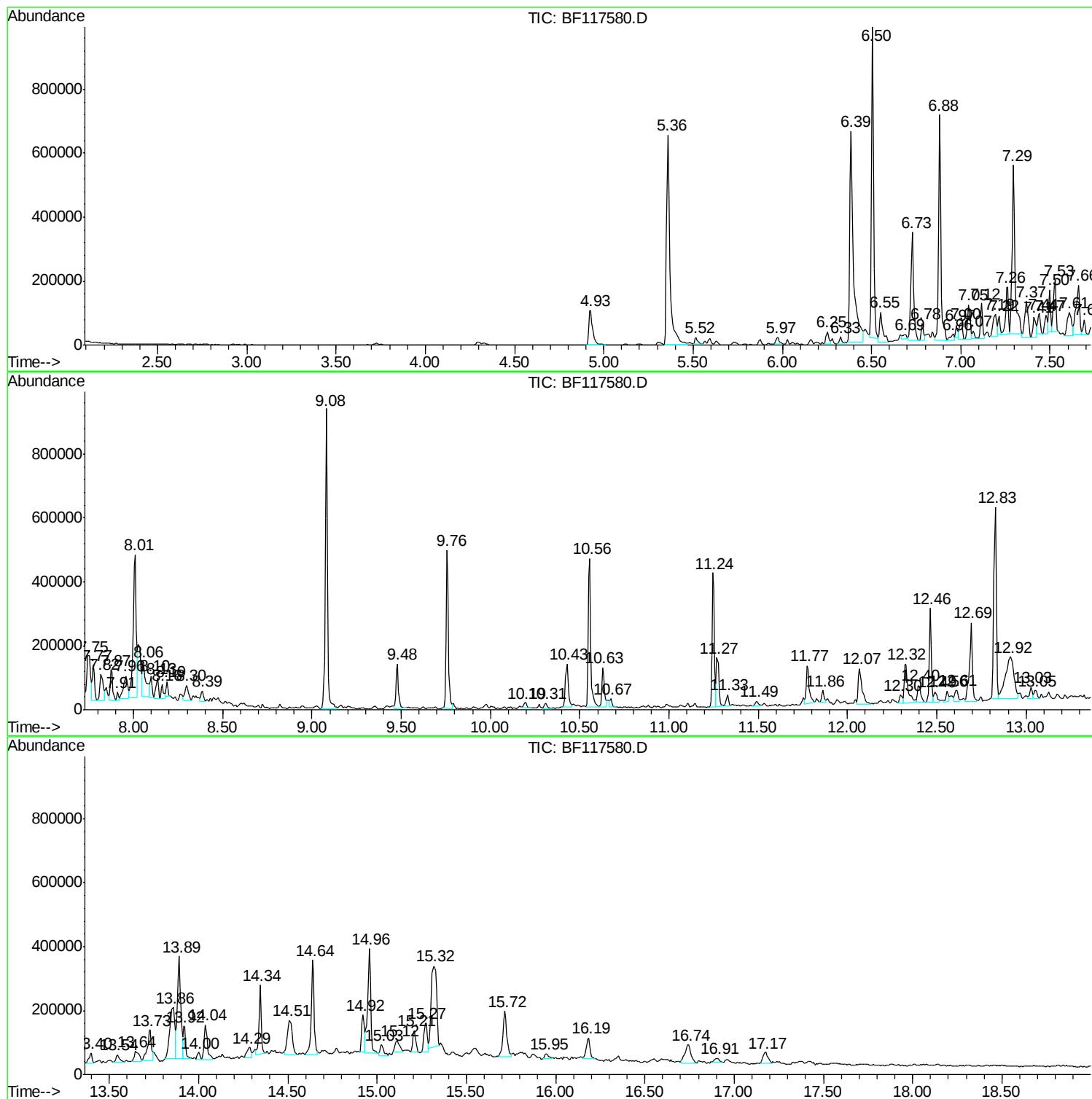
Sum of corrected areas: 17085833

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Data File : BF117580.D  
Acq On : 29 Oct 2019 10:30  
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Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampled :  
TP-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102919\  
Data File : BF117580.D  
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Sample : K5563-07  
Misc :  
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Instrument :  
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ClientSampleId :  
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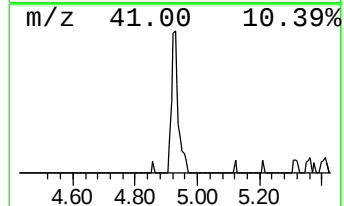
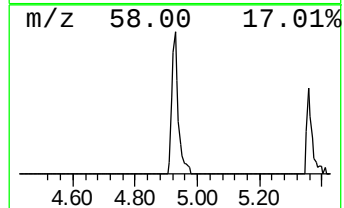
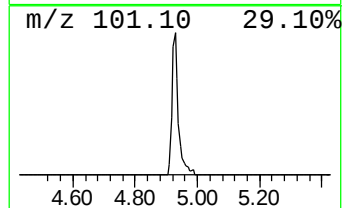
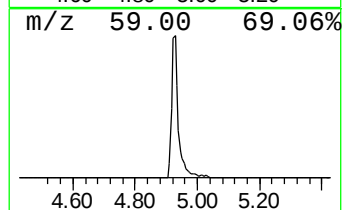
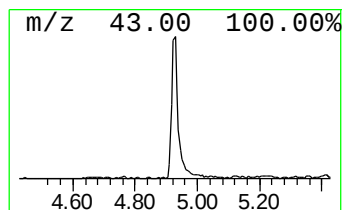
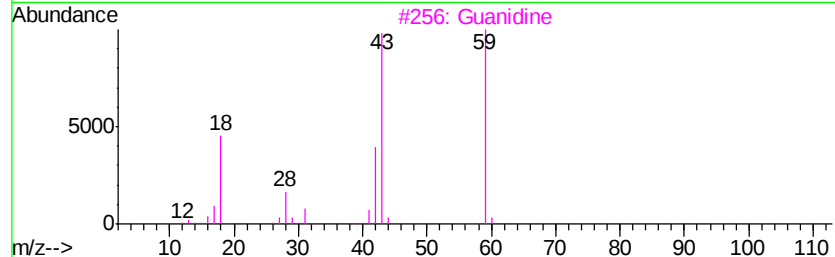
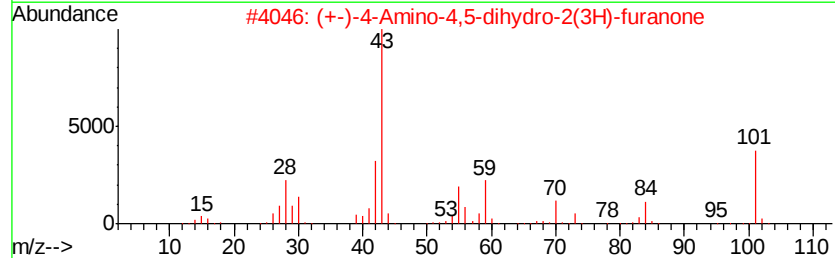
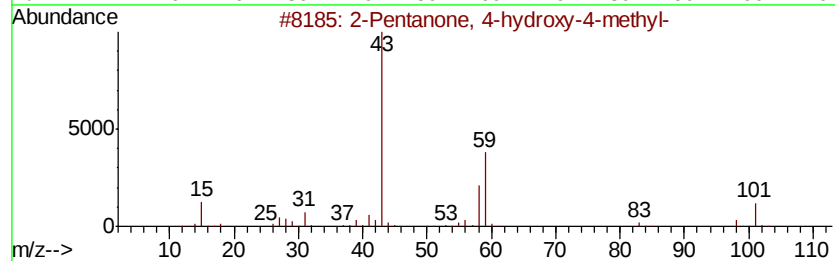
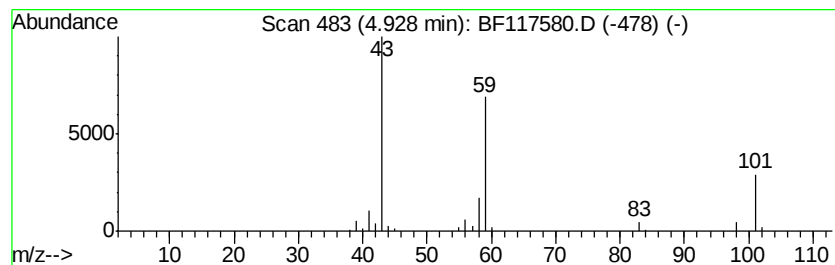
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 6

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 4.93 | 9.21 ng | 162425 | 1,4-Dichlorobenzene-d4 | 6.73 |

| Hit# | of | 5 | Tentative ID                           | MW  | MolForm  | CAS#        | Qual |
|------|----|---|----------------------------------------|-----|----------|-------------|------|
| 1    |    |   | 2-Pentanone, 4-hydroxy-4-methyl-       | 116 | C6H12O2  | 000123-42-2 | 56   |
| 2    |    |   | (+)-4-Amino-4,5-dihydro-2(3H)-furanone | 101 | C4H7NO2  | 016504-58-8 | 47   |
| 3    |    |   | Guanidine                              | 59  | CH5N3    | 000113-00-8 | 43   |
| 4    |    |   | Butyl aldoxime, 3-methyl-, syn-        | 101 | C5H11NO  | 005780-40-5 | 35   |
| 5    |    |   | Acetic acid, cyano-, 1,1-dimethyl-     | 141 | C7H11NO2 | 001116-98-9 | 17   |



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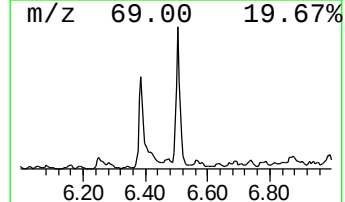
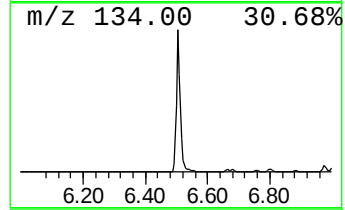
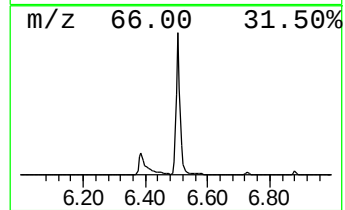
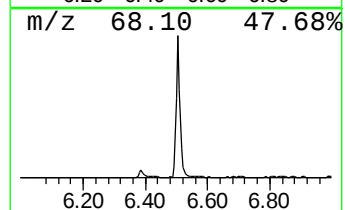
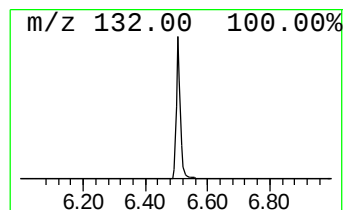
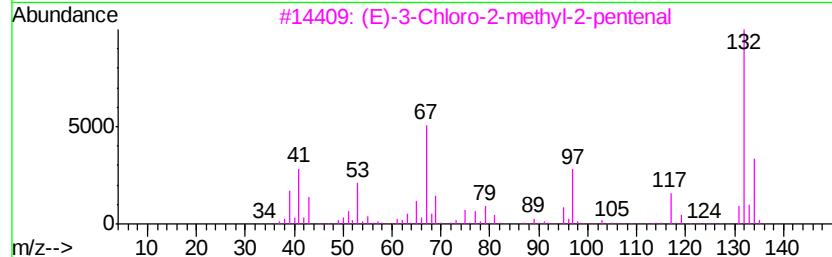
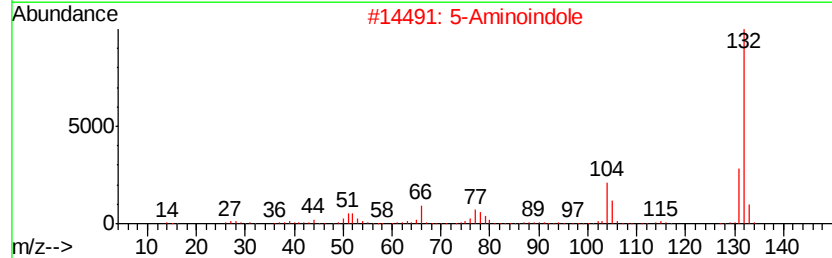
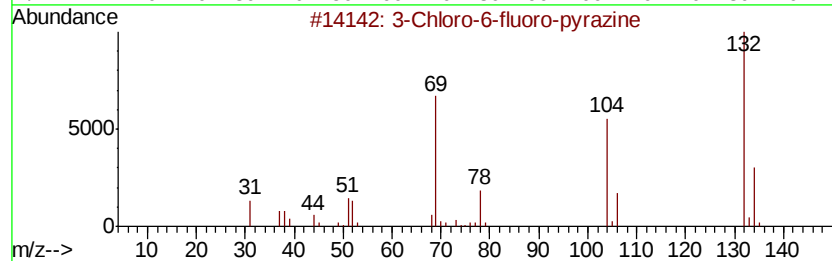
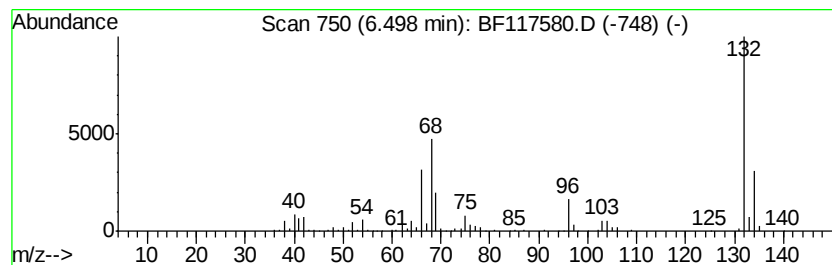
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 2 unknown6.50 Concentration Rank 1

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 6.50 | 50.38 ng | 888441 | 1,4-Dichlorobenzene-d4 | 6.73 |

| Hit# | of                                  | Tentative ID | MW        | MolForm      | CAS# | Qual |
|------|-------------------------------------|--------------|-----------|--------------|------|------|
| 1    | 3-Chloro-6-fluoro-pyrazine          | 132          | C4H2ClFN2 | 1000146-10-7 | 27   |      |
| 2    | 5-Aminoindole                       | 132          | C8H8N2    | 005192-03-0  | 10   |      |
| 3    | (E)-3-Chloro-2-methyl-2-pentenal    | 132          | C6H9ClO   | 031357-76-3  | 9    |      |
| 4    | Bicyclo[4.2.0]octa-1,3,5-triene,... | 132          | C10H12    | 028749-81-7  | 9    |      |
| 5    | 5-Fluoro-2-chloropyrimidine         | 132          | C4H2ClFN2 | 062802-42-0  | 9    |      |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102919\  
Data File : BF117580.D  
Acq On : 29 Oct 2019 10:30  
Operator : JU  
Sample : K5563-07  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-1

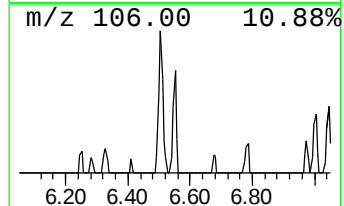
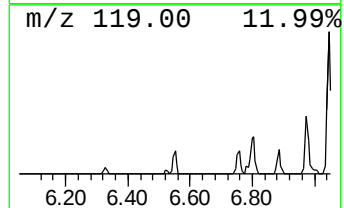
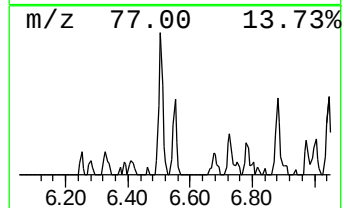
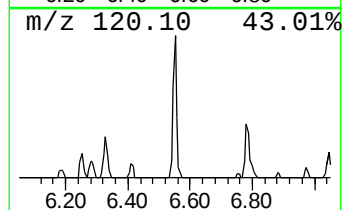
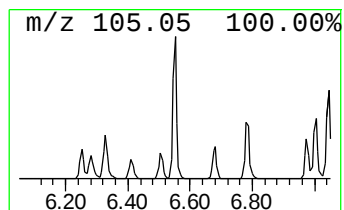
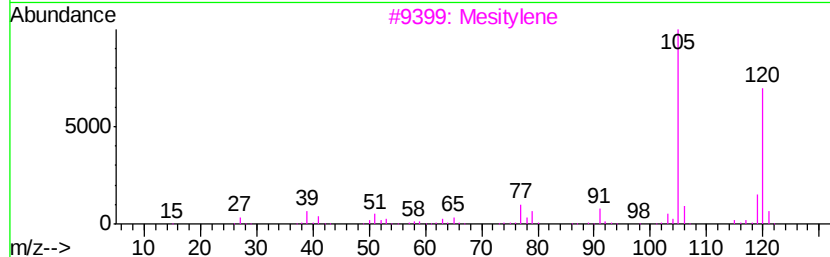
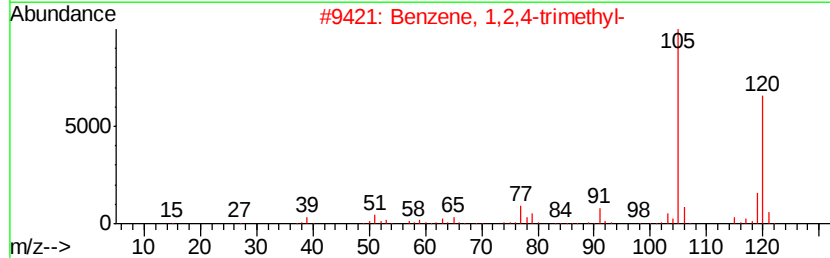
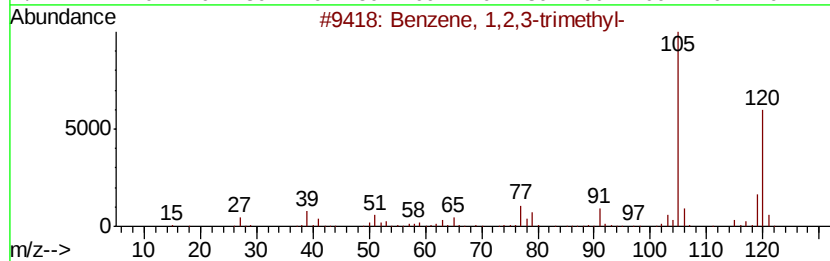
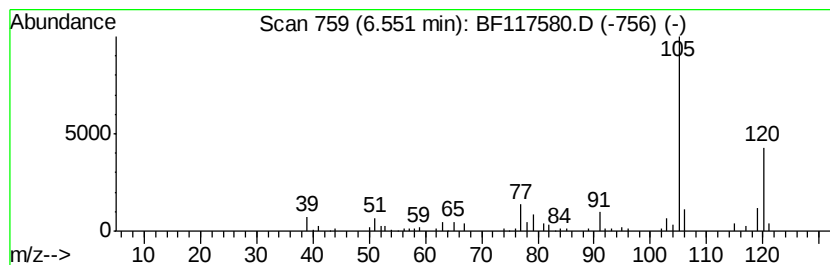
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 6.55 | 7.07 ng | 124745 | 1,4-Dichlorobenzene-d4 | 6.73 |

| Hit# | of | Tentative ID               | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2,3-trimethyl-  | 120 | C9H12   | 000526-73-8 | 97   |
| 2    |    | Benzene, 1,2,4-trimethyl-  | 120 | C9H12   | 000095-63-6 | 91   |
| 3    |    | Mesitylene                 | 120 | C9H12   | 000108-67-8 | 91   |
| 4    |    | Benzene, 1-ethyl-3-methyl- | 120 | C9H12   | 000620-14-4 | 91   |
| 5    |    | Benzene, 1-ethyl-4-methyl- | 120 | C9H12   | 000622-96-8 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102919\  
Data File : BF117580.D  
Acq On : 29 Oct 2019 10:30  
Operator : JU  
Sample : K5563-07  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampled :  
TP-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

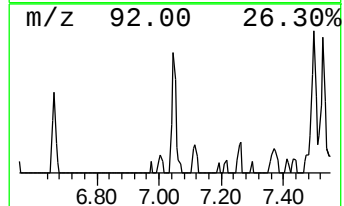
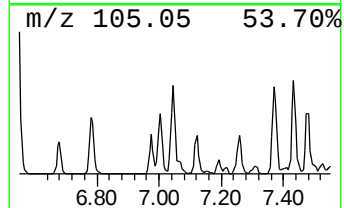
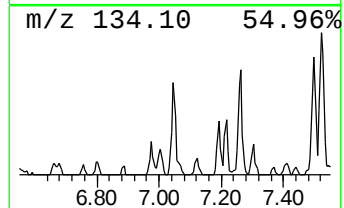
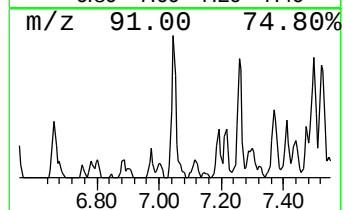
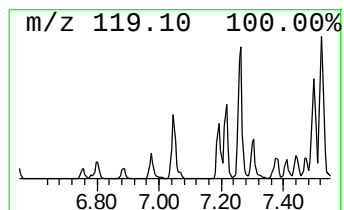
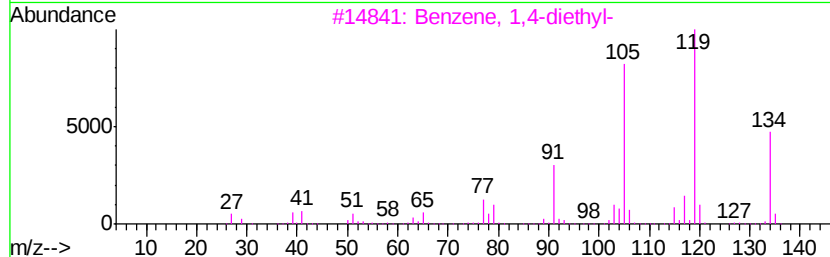
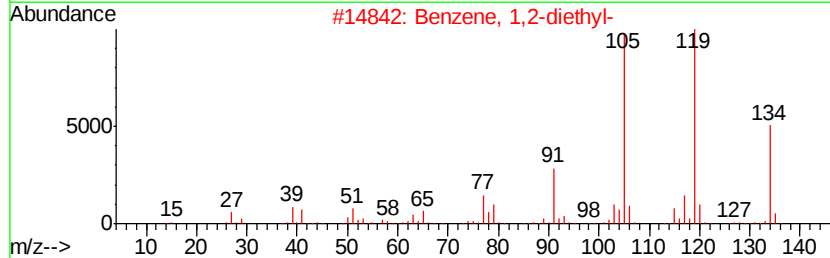
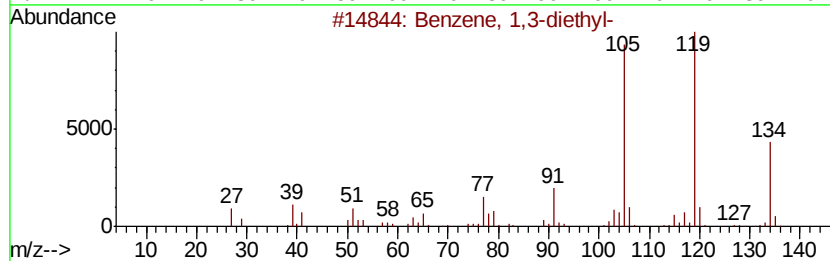
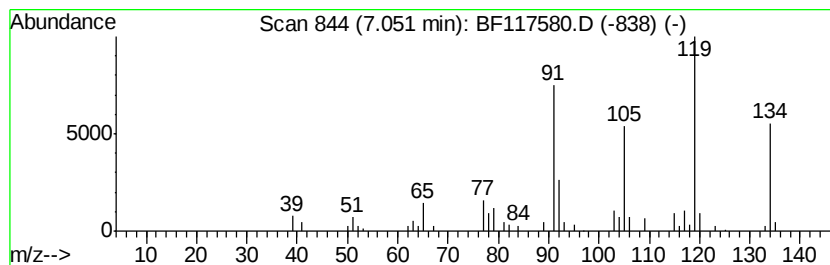
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TIC Integration Parameters: LSCINT.P

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Peak Number 5 Benzene, 1,3-diethyl- Concentration Rank 19

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 7.05 | 5.68 ng | 100191 | 1,4-Dichlorobenzene-d4 | 6.73 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,3-diethyl-          | 134 | C10H14  | 000141-93-5 | 96   |
| 2    |    | Benzene, 1,2-diethyl-          | 134 | C10H14  | 000135-01-3 | 95   |
| 3    |    | Benzene, 1,4-diethyl-          | 134 | C10H14  | 000105-05-5 | 95   |
| 4    |    | Benzene, 4-ethyl-1,2-dimethyl- | 134 | C10H14  | 000934-80-5 | 93   |
| 5    |    | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14  | 000933-98-2 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102919\  
Data File : BF117580.D  
Acq On : 29 Oct 2019 10:30  
Operator : JU  
Sample : K5563-07  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

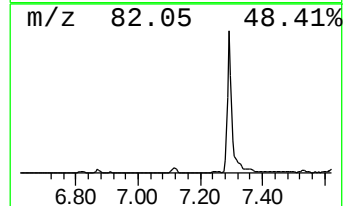
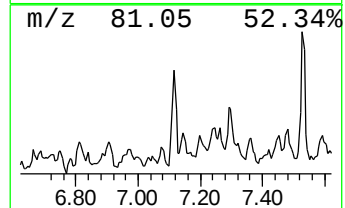
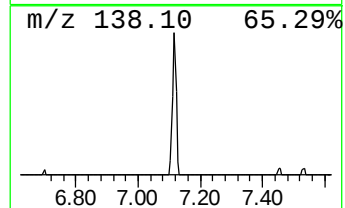
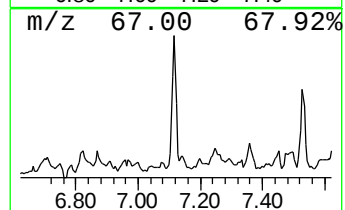
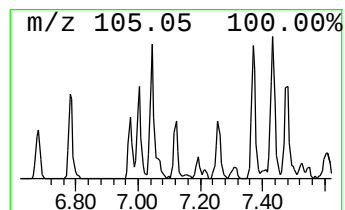
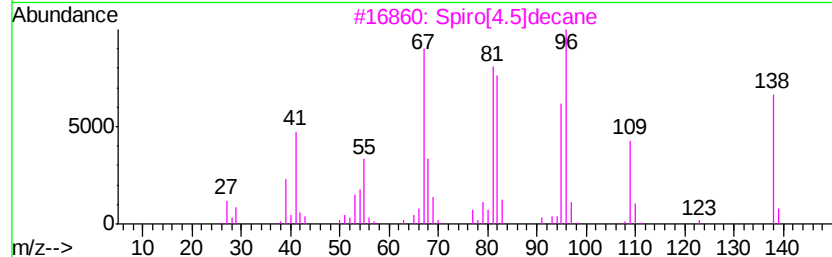
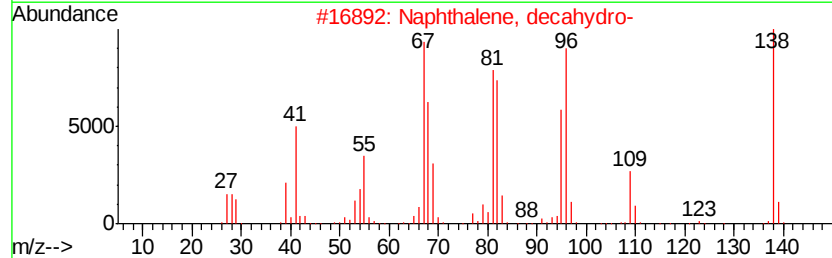
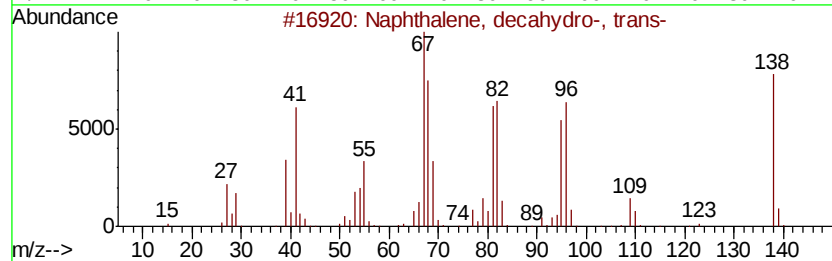
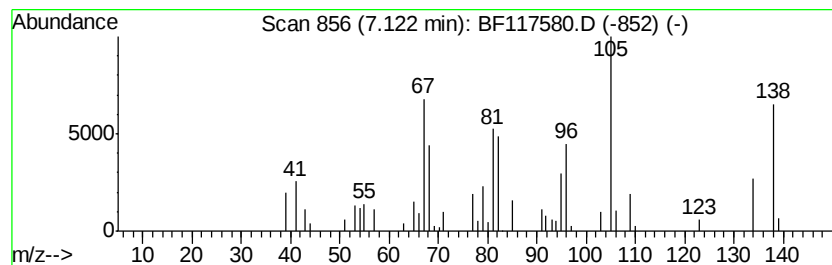
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 6 Naphthalene, decahydro-, tr... Concentration Rank 20

| R.T. | EstConc | Area  | Relative to ISTD       | R.T. |
|------|---------|-------|------------------------|------|
| 7.12 | 5.49 ng | 96844 | 1,4-Dichlorobenzene-d4 | 6.73 |

| Hit# | of | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------------|-----|---------|-------------|------|
| 1    | 5  | Naphthalene, decahydro-, trans-  | 138 | C10H18  | 000493-02-7 | 96   |
| 2    |    | Naphthalene, decahydro-          | 138 | C10H18  | 000091-17-8 | 94   |
| 3    |    | Spiro[4.5]decane                 | 138 | C10H18  | 000176-63-6 | 93   |
| 4    |    | Naphthalene, decahydro-, cis-    | 138 | C10H18  | 000493-01-6 | 81   |
| 5    |    | 5H-Inden-5-one, octahydro-, cis- | 138 | C9H14O  | 004668-91-1 | 80   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102919\  
Data File : BF117580.D  
Acq On : 29 Oct 2019 10:30  
Operator : JU  
Sample : K5563-07  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

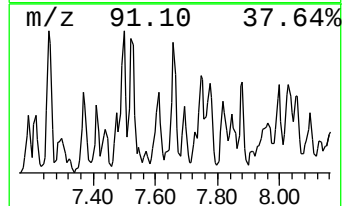
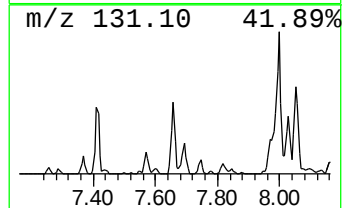
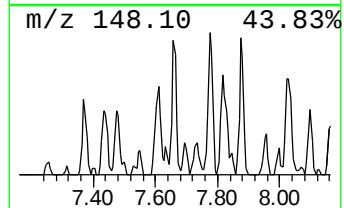
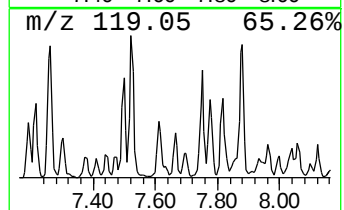
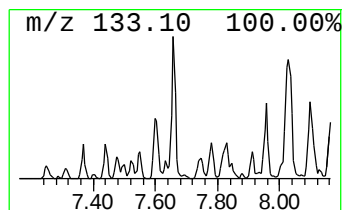
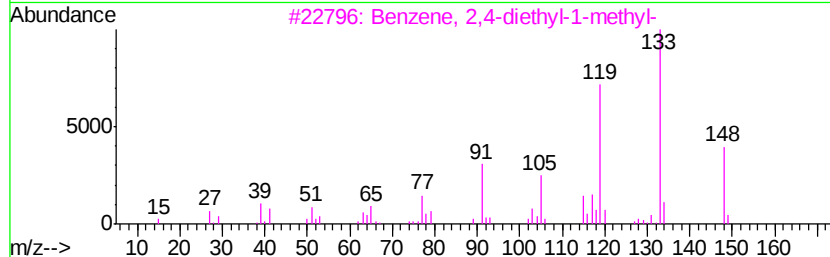
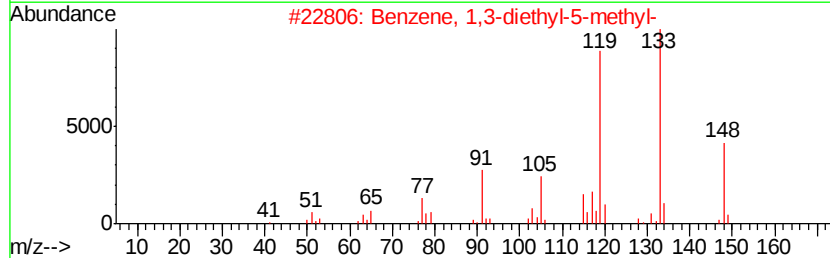
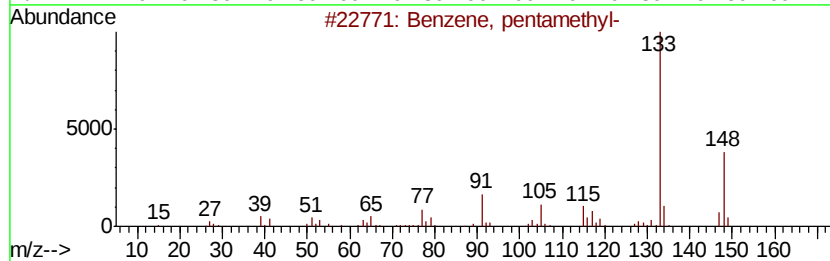
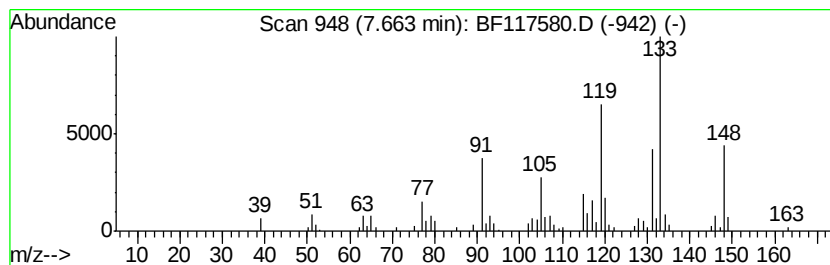
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 7 Benzene, pentamethyl- Concentration Rank 13

| R.T. | EstConc | Area   | Relative to ISTD | R.T. |
|------|---------|--------|------------------|------|
| 7.66 | 7.03 ng | 205541 | Naphthalene-d8   | 8.01 |

| Hit# | of | Tentative ID                         | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, pentamethyl-                | 148 | C11H16  | 000700-12-9 | 70   |
| 2    |    | Benzene, 1,3-diethyl-5-methyl-       | 148 | C11H16  | 002050-24-0 | 64   |
| 3    |    | Benzene, 2,4-diethyl-1-methyl-       | 148 | C11H16  | 001758-85-6 | 64   |
| 4    |    | Benzene, 1,3-dimethyl-5-(1-methyl... | 148 | C11H16  | 004706-90-5 | 58   |
| 5    |    | Benzene, 1-ethyl-4-(1-methylethyl)-  | 148 | C11H16  | 004218-48-8 | 58   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102919\  
Data File : BF117580.D  
Acq On : 29 Oct 2019 10:30  
Operator : JU  
Sample : K5563-07  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

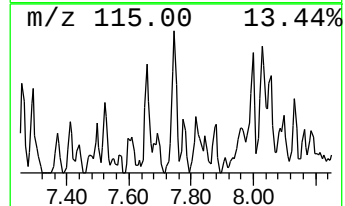
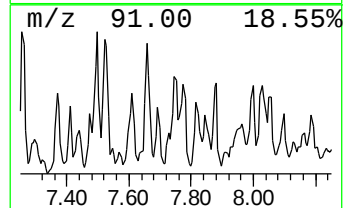
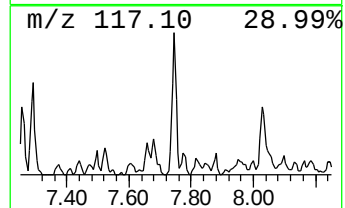
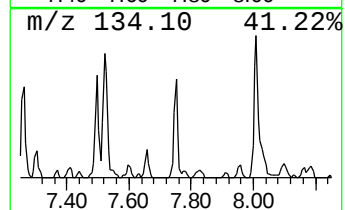
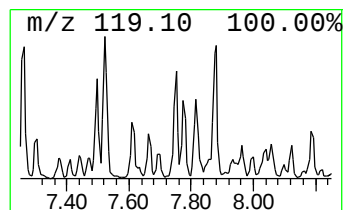
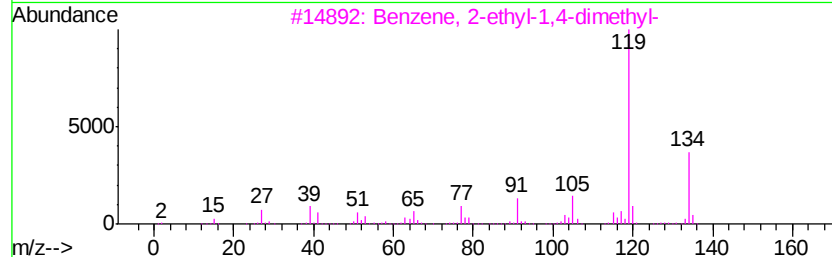
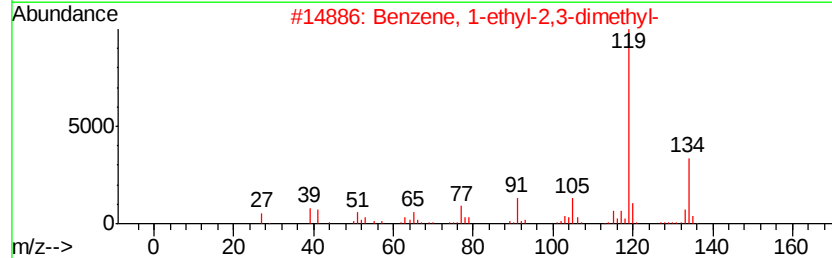
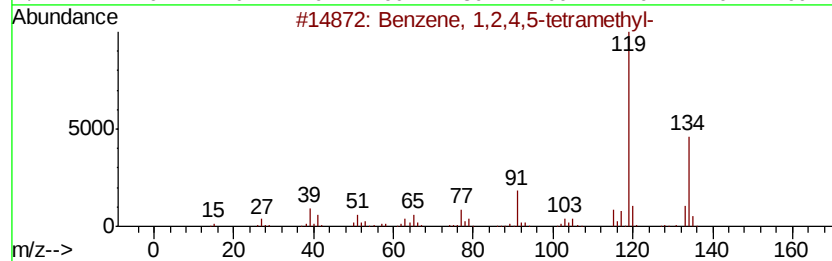
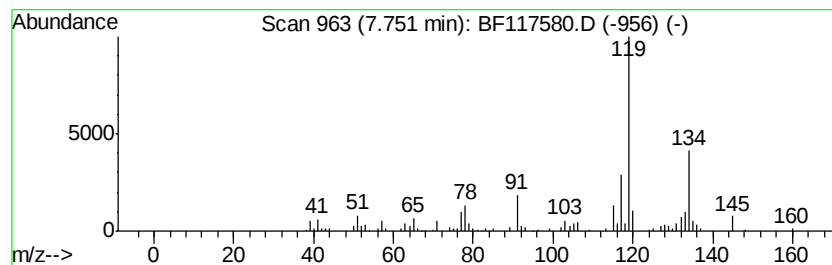
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 16

| R.T. | EstConc | Area   | Relative to ISTD | R.T. |
|------|---------|--------|------------------|------|
| 7.75 | 6.44 ng | 188292 | Naphthalene-d8   | 8.01 |

| Hit# | of | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|--------------------------------|-----|---------|-------------|------|
| 1    | 5  | Benzene, 1,2,4,5-tetramethyl-  | 134 | C10H14  | 000095-93-2 | 95   |
| 2    |    | Benzene, 1-ethyl-2,3-dimethyl- | 134 | C10H14  | 000933-98-2 | 81   |
| 3    |    | Benzene, 2-ethyl-1,4-dimethyl- | 134 | C10H14  | 001758-88-9 | 81   |
| 4    |    | o-Cymene                       | 134 | C10H14  | 000527-84-4 | 81   |
| 5    |    | Benzene, 1,2,3,5-tetramethyl-  | 134 | C10H14  | 000527-53-7 | 81   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102919\  
Data File : BF117580.D  
Acq On : 29 Oct 2019 10:30  
Operator : JU  
Sample : K5563-07  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

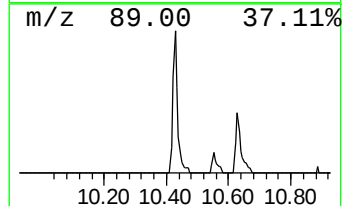
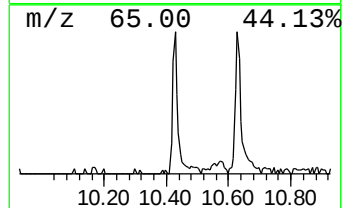
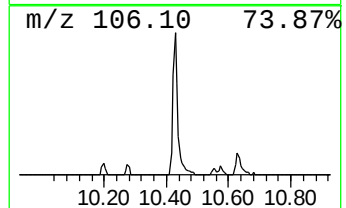
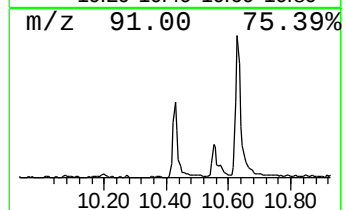
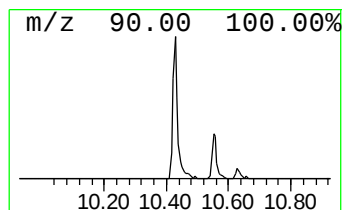
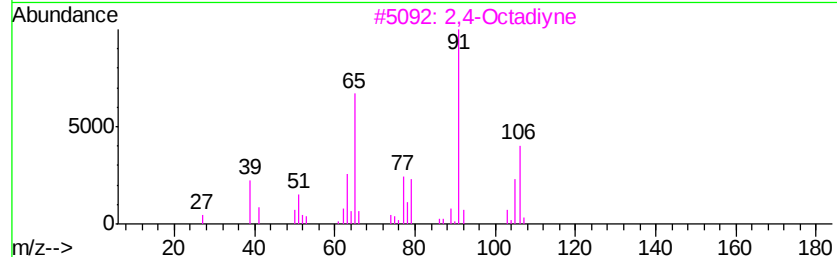
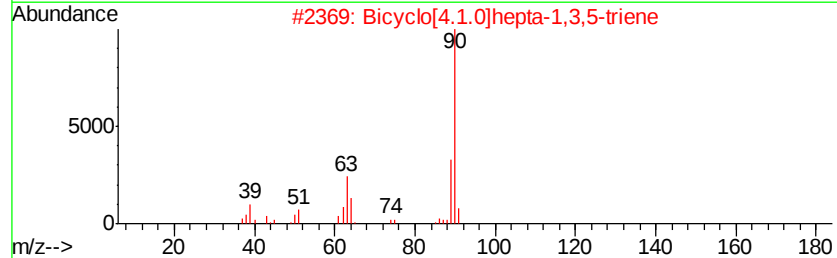
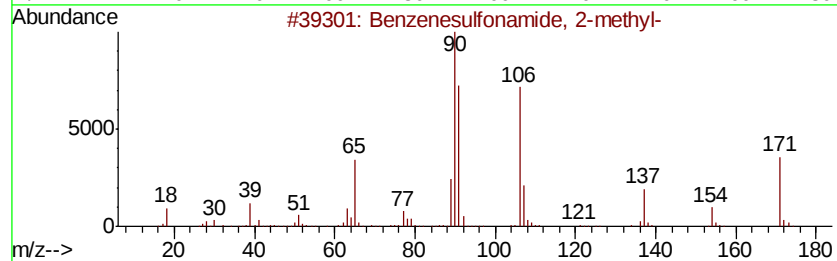
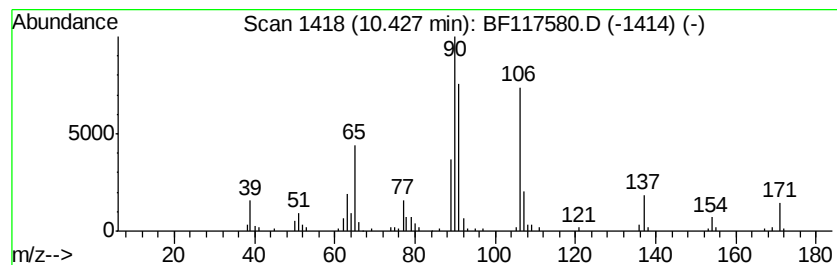
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 Benzenesulfonamide, 2-methyl- Concentration Rank 17

| R.T.  | EstConc | Area   | Relative to ISTD | R.T. |
|-------|---------|--------|------------------|------|
| 10.43 | 6.15 ng | 138415 | Acenaphthene-d10 | 9.76 |

| Hit# | of | 5 | Tentative ID                       | MW  | MolForm    | CAS#        | Qual |
|------|----|---|------------------------------------|-----|------------|-------------|------|
| 1    |    |   | Benzenesulfonamide, 2-methyl-      | 171 | C7H9NO2S   | 000088-19-7 | 97   |
| 2    |    |   | Bicyclo[4.1.0]hepta-1,3,5-triene   | 90  | C7H6       | 004646-69-9 | 43   |
| 3    |    |   | 2,4-Octadiyne                      | 106 | C8H10      | 002809-71-4 | 38   |
| 4    |    |   | Benzene, 1-methyl-4-nitro-         | 137 | C7H7NO2    | 000099-99-0 | 35   |
| 5    |    |   | 4-(2-Aminoethyl)benzenesulfonamide | 200 | C8H12N2O2S | 035303-76-5 | 28   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102919\  
Data File : BF117580.D  
Acq On : 29 Oct 2019 10:30  
Operator : JU  
Sample : K5563-07  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

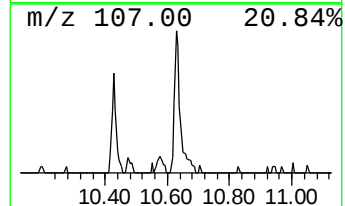
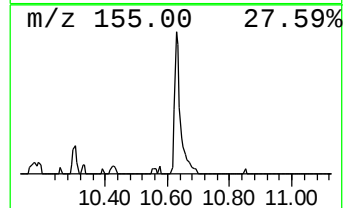
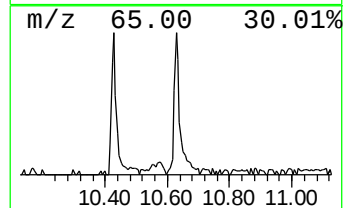
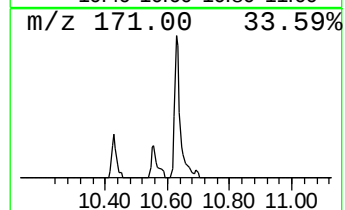
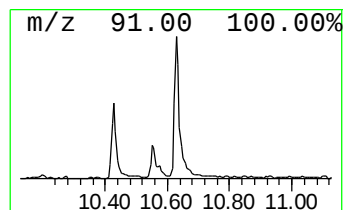
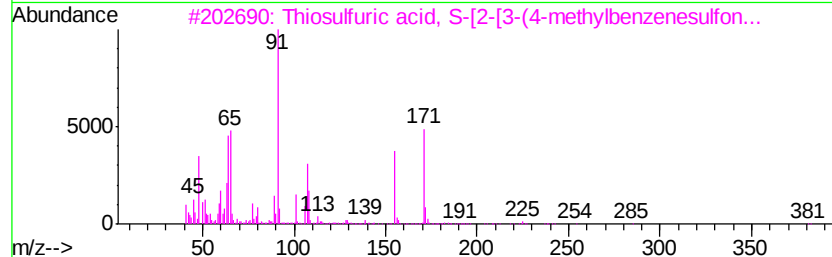
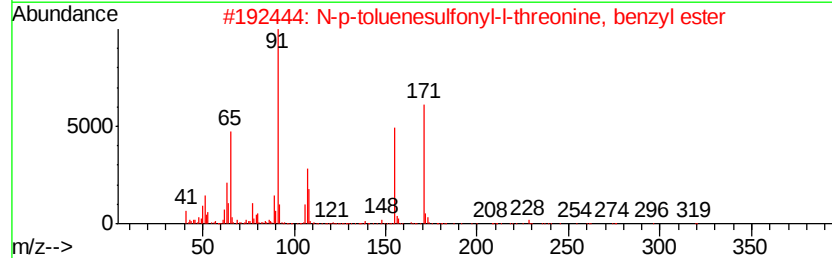
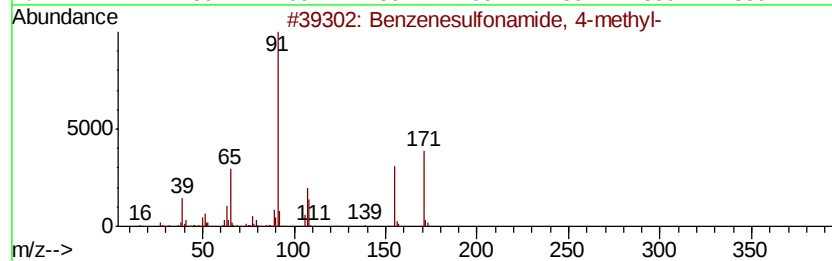
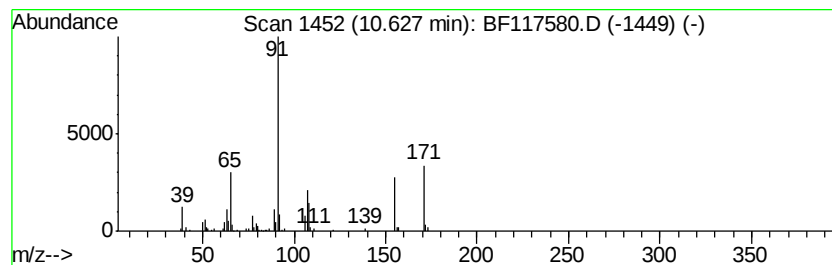
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 10 Benzenesulfonamide, 4-methyl- Concentration Rank 15

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 10.63 | 6.69 ng | 136392 | Phenanthrene-d10 | 11.24 |

| Hit# | of | 5 | Tentative ID                         | MW  | MolForm      | CAS#         | Qual |
|------|----|---|--------------------------------------|-----|--------------|--------------|------|
| 1    |    |   | Benzenesulfonamide, 4-methyl-        | 171 | C7H9NO2S     | 000070-55-3  | 97   |
| 2    |    |   | N-p-toluenesulfonyl-L-threonine,...  | 363 | C18H21NO5S   | 1000130-34-4 | 83   |
| 3    |    |   | Thiosulfuric acid, S-[2-[3-(4-me...  | 381 | C12H19N3O5S3 | 1000255-60-3 | 78   |
| 4    |    |   | O-p-toluenesulfonyl-L-N-acetylser... | 315 | C13H17NO6S   | 1000126-44-8 | 53   |
| 5    |    |   | Benzenemethanesulfonamide            | 171 | C7H9NO2S     | 004563-33-1  | 43   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102919\  
Data File : BF117580.D  
Acq On : 29 Oct 2019 10:30  
Operator : JU  
Sample : K5563-07  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

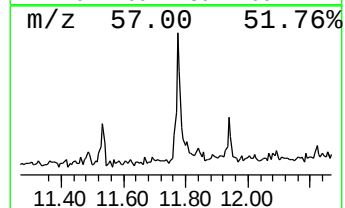
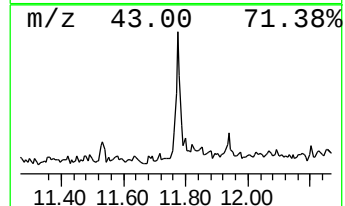
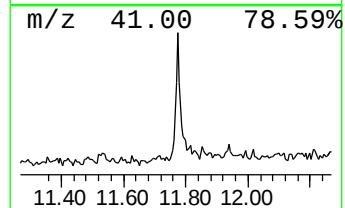
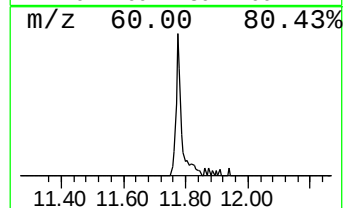
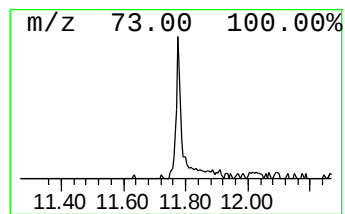
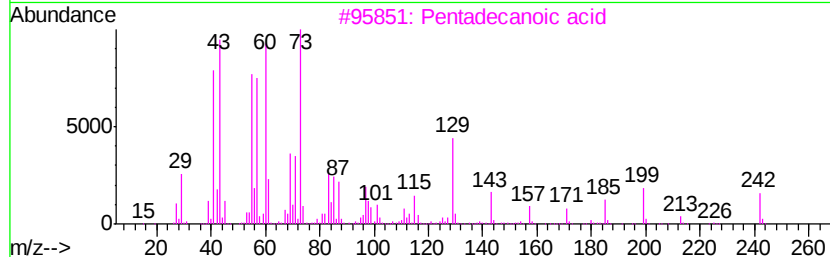
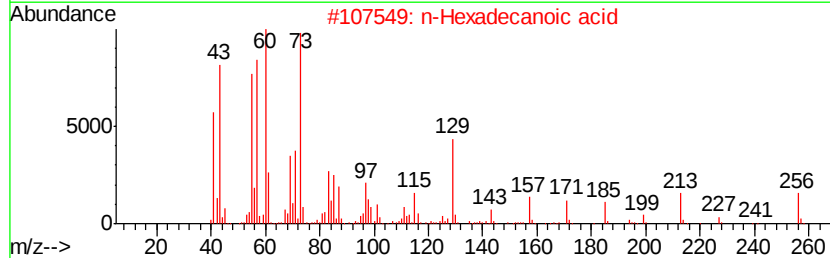
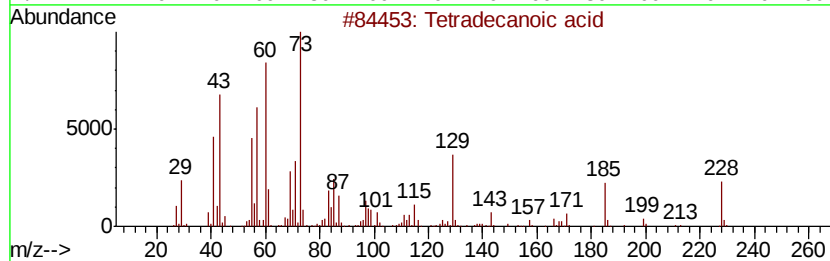
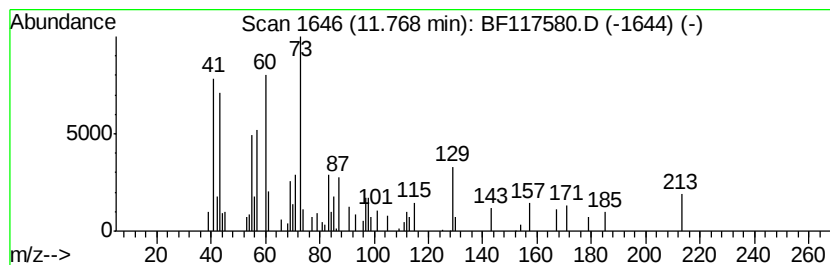
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 Tetradecanoic acid Concentration Rank 14

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 11.77 | 6.82 ng | 139052 | Phenanthrene-d10 | 11.24 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm    | CAS#        | Qual |
|------|----|---|--------------------------------|-----|------------|-------------|------|
| 1    |    |   | Tetradecanoic acid             | 228 | C14H28O2   | 000544-63-8 | 80   |
| 2    |    |   | n-Hexadecanoic acid            | 256 | C16H32O2   | 000057-10-3 | 76   |
| 3    |    |   | Pentadecanoic acid             | 242 | C15H30O2   | 001002-84-2 | 68   |
| 4    |    |   | Decanoic acid, silver(1+) salt | 278 | C10H19AgO2 | 013126-67-5 | 53   |
| 5    |    |   | Nonanoic acid                  | 158 | C9H18O2    | 000112-05-0 | 38   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102919\  
Data File : BF117580.D  
Acq On : 29 Oct 2019 10:30  
Operator : JU  
Sample : K5563-07  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

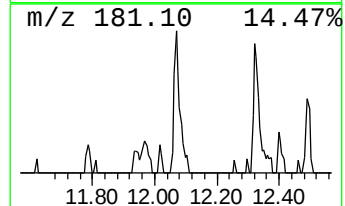
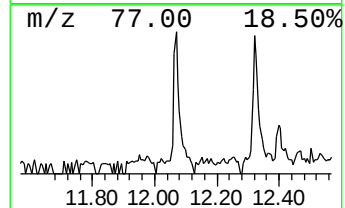
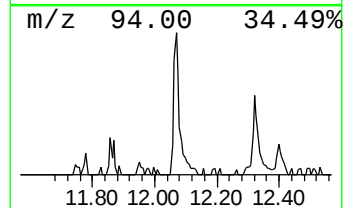
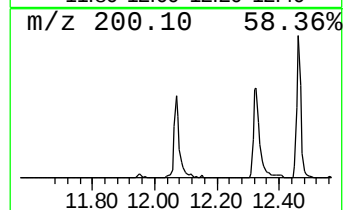
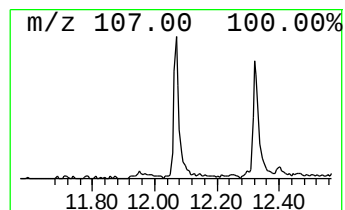
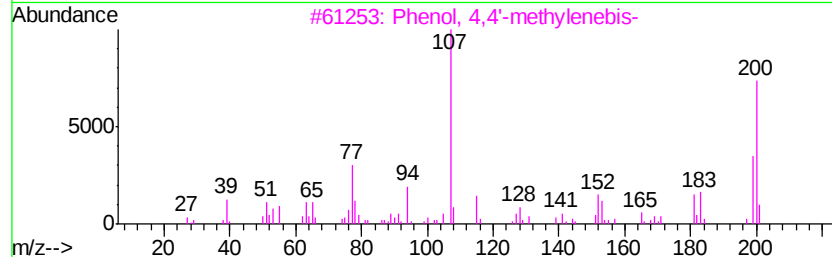
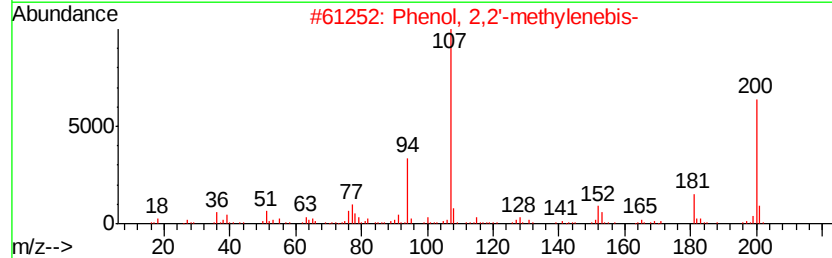
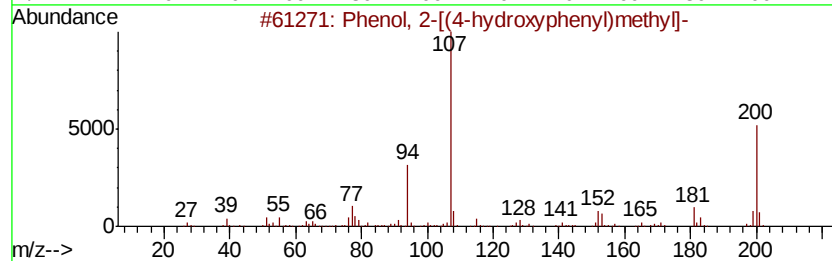
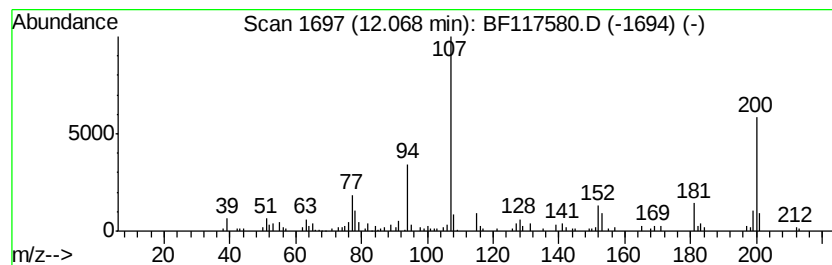
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 12 Phenol, 2-[(4-hydroxyphenyl)... Concentration Rank 10

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.07 | 7.62 ng | 155411 | Phenanthrene-d10 | 11.24 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Phenol, 2-[(4-hydroxyphenyl)meth... | 200 | C13H12O2 | 002467-03-0 | 97   |
| 2    |    | Phenol, 2,2'-methylenebis-          | 200 | C13H12O2 | 002467-02-9 | 94   |
| 3    |    | Phenol, 4,4'-methylenebis-          | 200 | C13H12O2 | 000620-92-8 | 72   |
| 4    |    | Ethyl 2-hydroxybenzyl sulfone       | 200 | C9H12O3S | 053380-27-1 | 47   |
| 5    |    | 4-(2-Methoxyethyl)phenol            | 152 | C9H12O2  | 056718-71-9 | 38   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102919\  
Data File : BF117580.D  
Acq On : 29 Oct 2019 10:30  
Operator : JU  
Sample : K5563-07  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

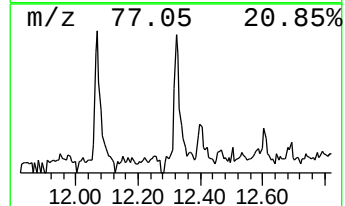
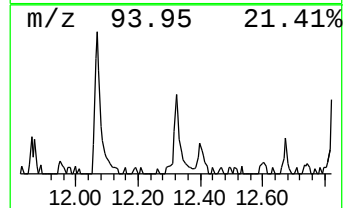
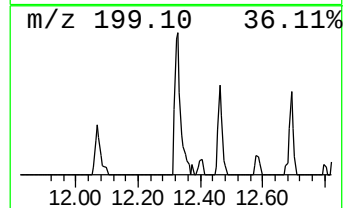
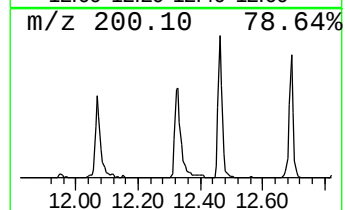
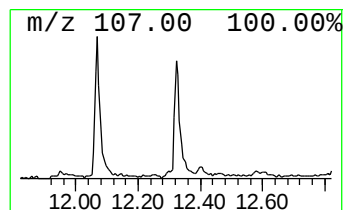
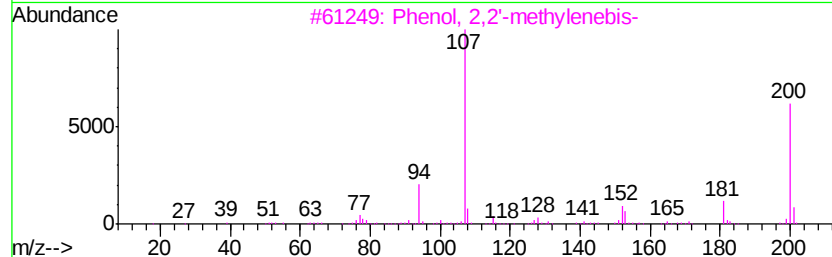
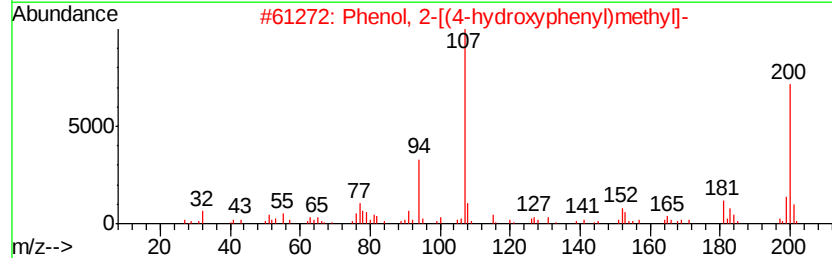
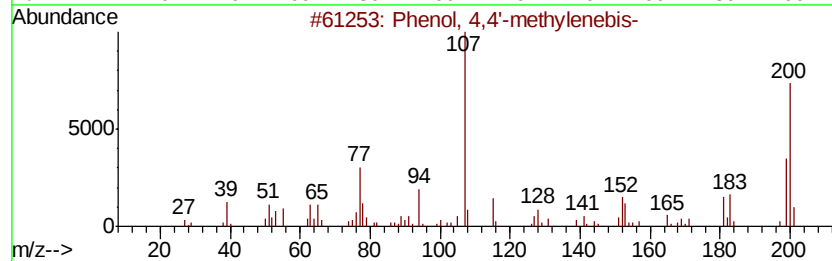
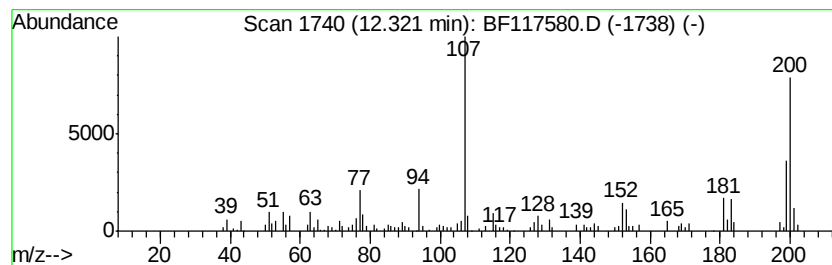
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 13 Phenol, 4,4'-methylenebis- Concentration Rank 5

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 12.32 | 9.25 ng | 188701 | Phenanthrene-d10 | 11.24 |

| Hit# | of | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|-------------------------------------|-----|----------|-------------|------|
| 1    | 5  | Phenol, 4,4'-methylenebis-          | 200 | C13H12O2 | 000620-92-8 | 94   |
| 2    |    | Phenol, 2-[(4-hydroxyphenyl)meth... | 200 | C13H12O2 | 002467-03-0 | 91   |
| 3    |    | Phenol, 2,2'-methylenebis-          | 200 | C13H12O2 | 002467-02-9 | 87   |
| 4    |    | 1,3-Benzenediol, 4-(phenylmethyl)-  | 200 | C13H12O2 | 002284-30-2 | 60   |
| 5    |    | Benzenemethanol, 3-phenoxy-         | 200 | C13H12O2 | 013826-35-2 | 49   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102919\  
Data File : BF117580.D  
Acq On : 29 Oct 2019 10:30  
Operator : JU  
Sample : K5563-07  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

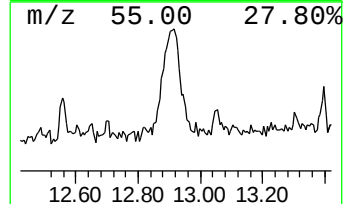
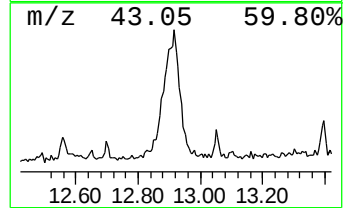
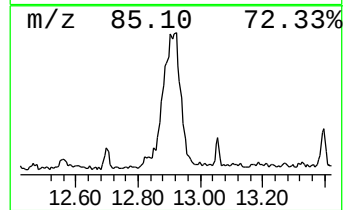
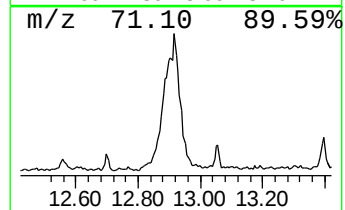
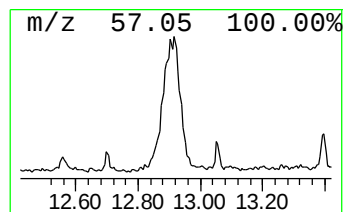
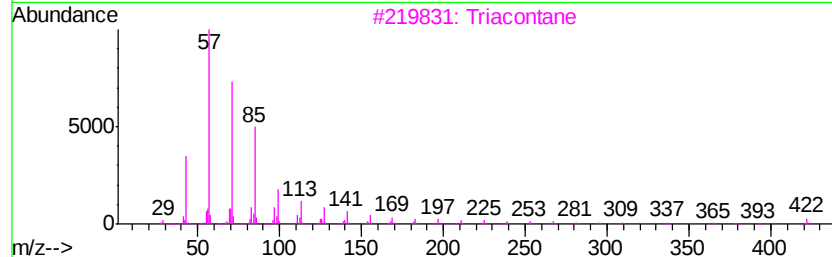
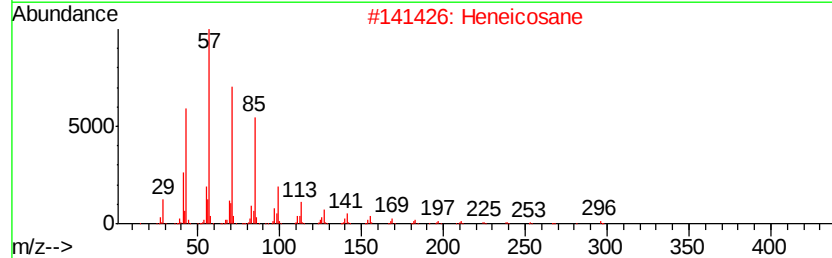
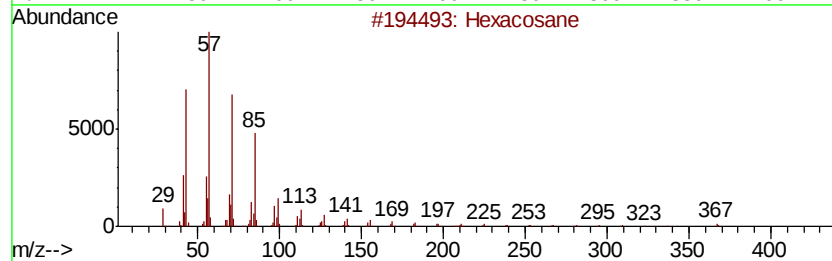
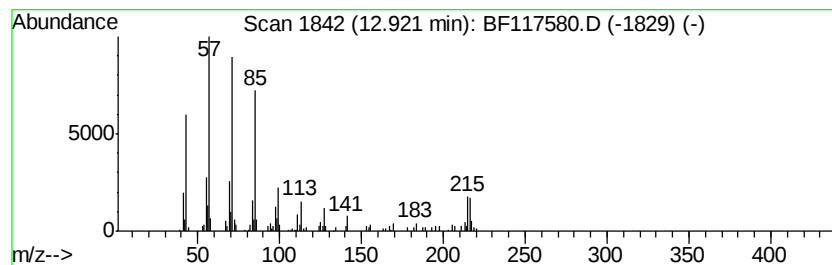
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*  
Peak Number 14 Hexacosane Concentration Rank 3

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 12.92 | 20.03 ng | 432955 | Chrysene-d12     | 13.89 |

| Hit# | of | Tentative ID                     | MW  | MolForm | CAS#        | Qual |
|------|----|----------------------------------|-----|---------|-------------|------|
| 1    | 5  | Hexacosane                       | 366 | C26H54  | 000630-01-3 | 87   |
| 2    |    | Heneicosane                      | 296 | C21H44  | 000629-94-7 | 83   |
| 3    |    | Triacontane                      | 422 | C30H62  | 000638-68-6 | 83   |
| 4    |    | Hentriacontane                   | 437 | C31H64  | 000630-04-6 | 81   |
| 5    |    | Heneicosane, 11-(1-ethylpropyl)- | 366 | C26H54  | 055282-11-6 | 80   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102919\  
Data File : BF117580.D  
Acq On : 29 Oct 2019 10:30  
Operator : JU  
Sample : K5563-07  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

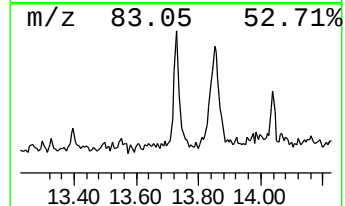
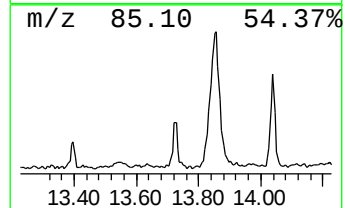
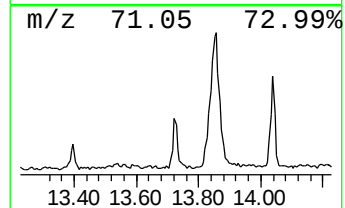
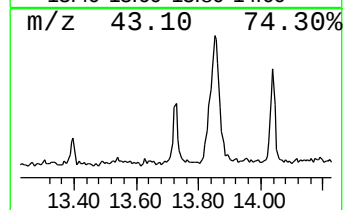
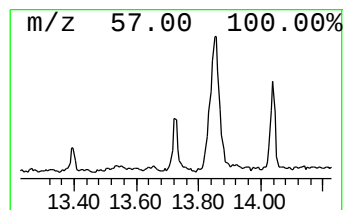
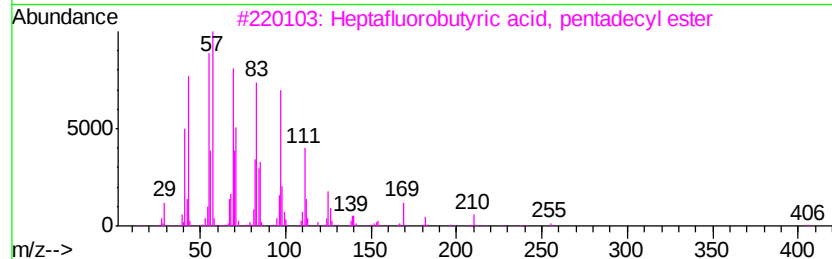
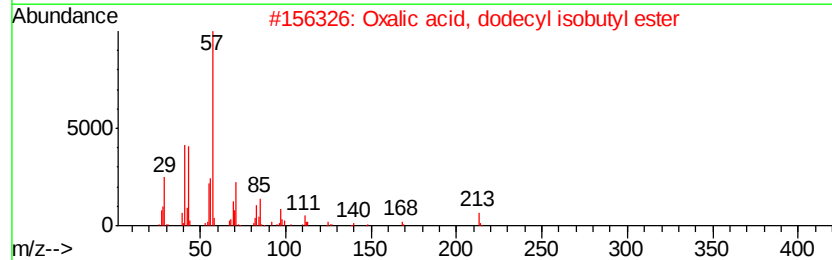
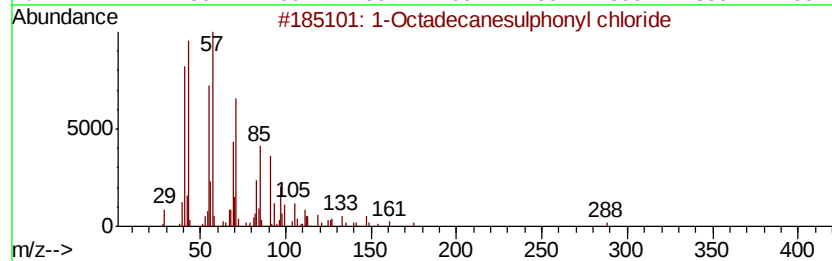
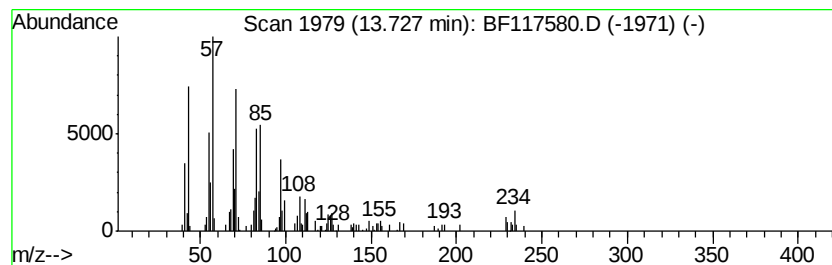
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 15 1-Octadecanesulphonyl chloride Concentration Rank 11

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 13.73 | 7.21 ng | 155904 | Chrysene-d12     | 13.89 |

| Hit# | of | Tentative ID                              | MW  | MolForm     | CAS#         | Qual |
|------|----|-------------------------------------------|-----|-------------|--------------|------|
| 1    | 5  | 1-Octadecanesulphonyl chloride            | 352 | C18H37ClO2S | 1000342-70-4 | 58   |
| 2    |    | Oxalic acid, dodecyl isobutyl ester       | 314 | C18H34O4    | 1000309-37-7 | 58   |
| 3    |    | Heptafluorobutyric acid, pentadecyl ester | 424 | C19H31F7O2  | 959261-23-5  | 49   |
| 4    |    | Trifluoroacetic acid, pentadecyl ester    | 324 | C17H31F3O2  | 959010-23-2  | 49   |
| 5    |    | 1-Decanol, 2-octyl-                       | 270 | C18H38O     | 045235-48-1  | 46   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102919\  
Data File : BF117580.D  
Acq On : 29 Oct 2019 10:30  
Operator : JU  
Sample : K5563-07  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

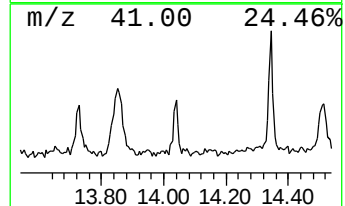
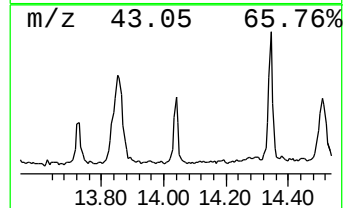
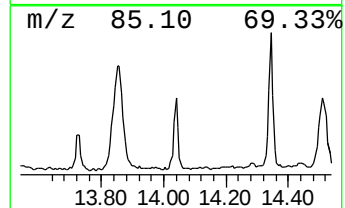
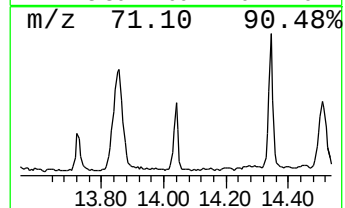
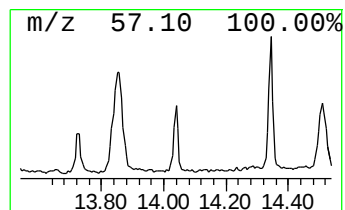
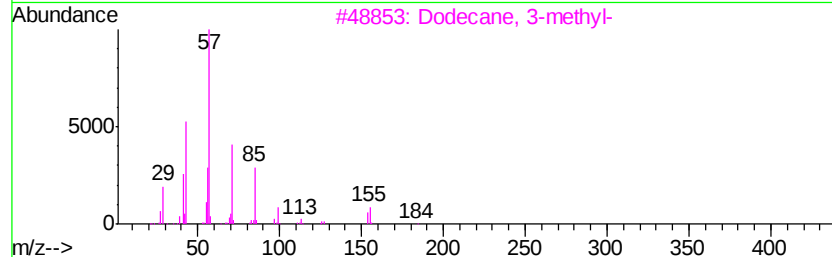
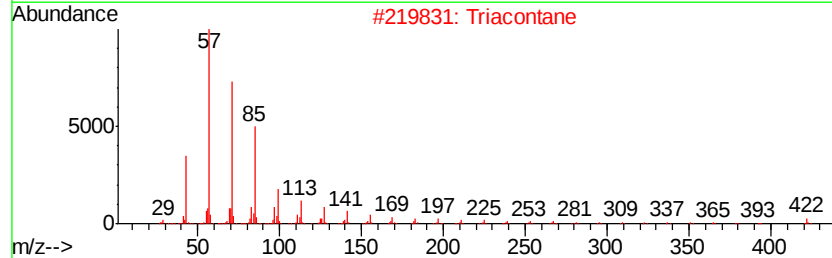
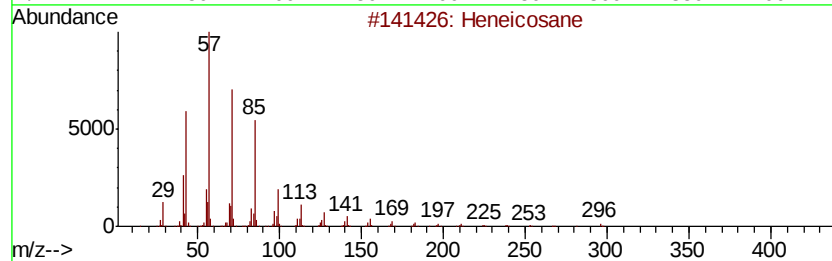
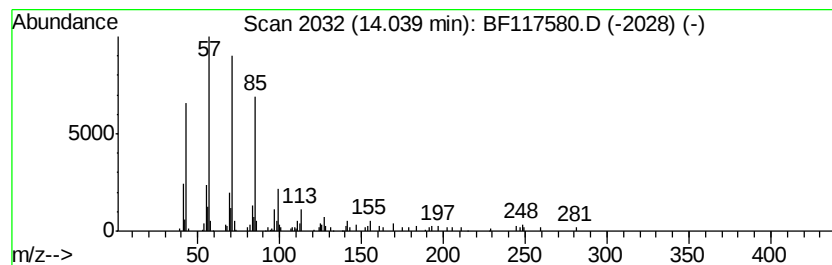
Instrument :  
BNA\_F  
ClientSampleId :  
TP-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.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 16 Heneicosane Concentration Rank 18

| R.T.      | EstConc                          | Area   | Relative to ISTD | R.T.         |      |
|-----------|----------------------------------|--------|------------------|--------------|------|
| 14.04     | 5.74 ng                          | 124097 | Chrysene-d12     | 13.89        |      |
| Hit# of 5 | Tentative ID                     | MW     | MolForm          | CAS#         | Qual |
| 1         | Heneicosane                      | 296    | C21H44           | 000629-94-7  | 91   |
| 2         | Triacontane                      | 422    | C30H62           | 000638-68-6  | 91   |
| 3         | Dodecane, 3-methyl-              | 184    | C13H28           | 017312-57-1  | 90   |
| 4         | Heneicosane, 11-(1-ethylpropyl)- | 366    | C26H54           | 055282-11-6  | 86   |
| 5         | 5-Ethyl-5-methylnonadecane       | 310    | C22H46           | 1000360-42-7 | 86   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102919\  
Data File : BF117580.D  
Acq On : 29 Oct 2019 10:30  
Operator : JU  
Sample : K5563-07  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

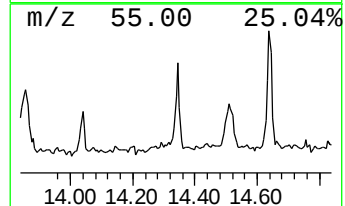
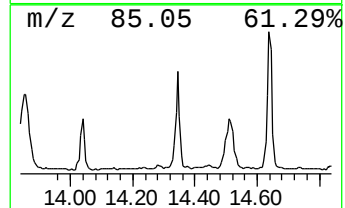
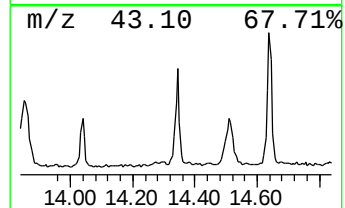
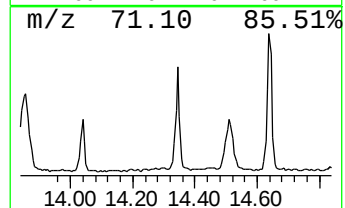
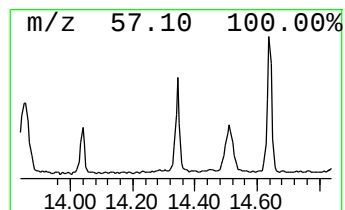
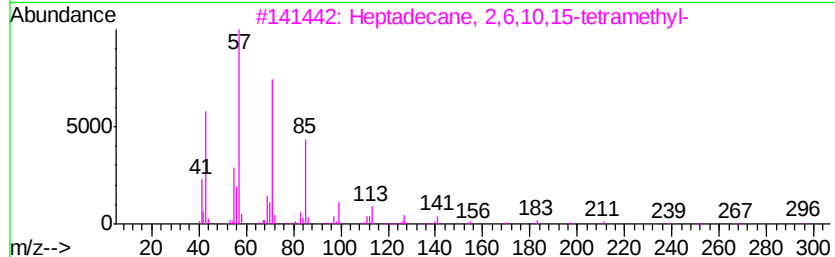
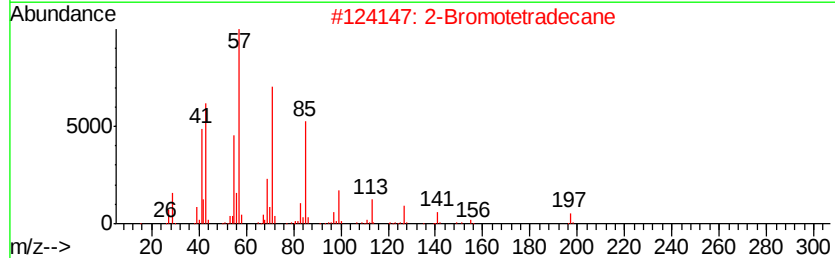
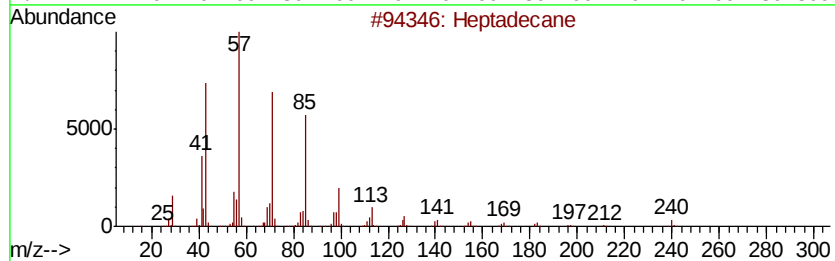
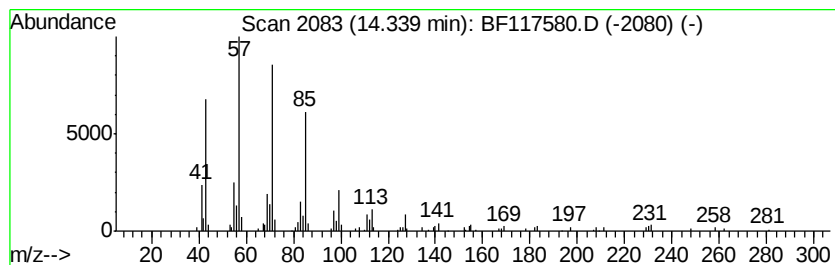
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 17 Heptadecane Concentration Rank 7

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 14.34 | 9.16 ng | 197925 | Chrysene-d12     | 13.89 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm  | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|----------|-------------|------|
| 1    |    |   | Heptadecane                         | 240 | C17H36   | 000629-78-7 | 91   |
| 2    |    |   | 2-Bromotetradecane                  | 276 | C14H29Br | 074036-95-6 | 90   |
| 3    |    |   | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44   | 054833-48-6 | 87   |
| 4    |    |   | Heneicosane, 11-(1-ethylpropyl)-    | 366 | C26H54   | 055282-11-6 | 86   |
| 5    |    |   | Octadecane, 1-iodo-                 | 380 | C18H37I  | 000629-93-6 | 86   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102919\  
Data File : BF117580.D  
Acq On : 29 Oct 2019 10:30  
Operator : JU  
Sample : K5563-07  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

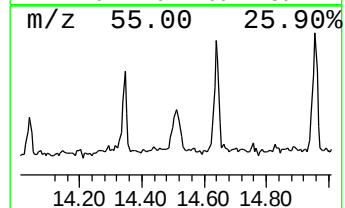
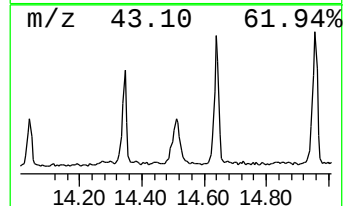
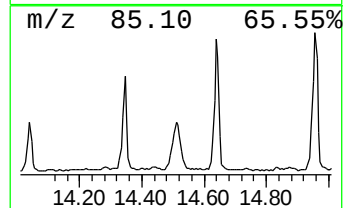
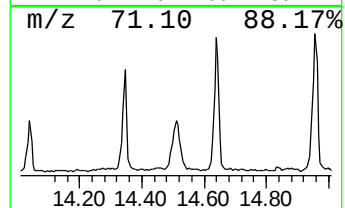
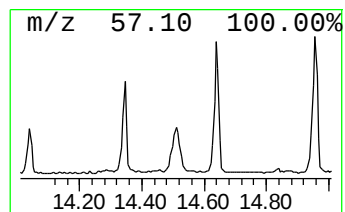
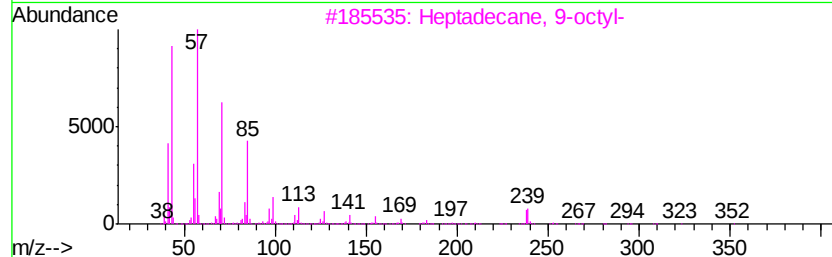
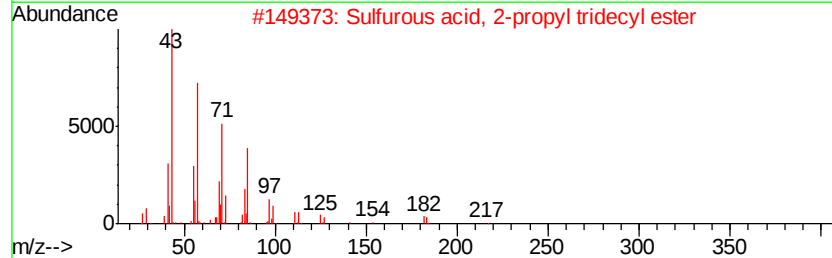
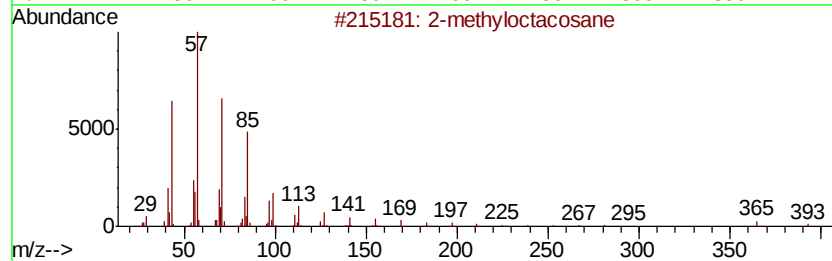
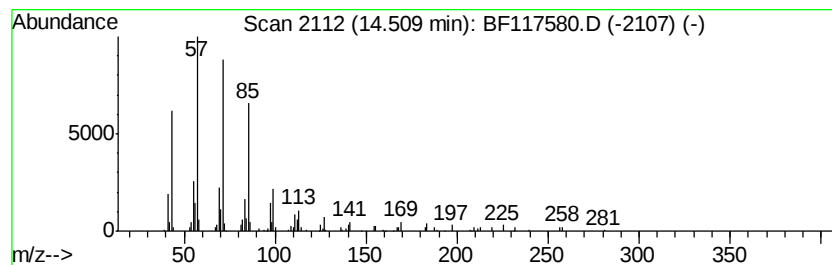
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 2-methyloctacosane Concentration Rank 8

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 14.51 | 9.03 ng | 195192 | Chrysene-d12     | 13.89 |

| Hit# | of | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|-------------------------------------|-----|-----------|--------------|------|
| 1    | 5  | 2-methyloctacosane                  | 408 | C29H60    | 1000376-72-8 | 91   |
| 2    |    | Sulfurous acid, 2-propyl tridecy... | 306 | C16H34O3S | 1000309-12-4 | 86   |
| 3    |    | Heptadecane, 9-octyl-               | 352 | C25H52    | 007225-64-1  | 86   |
| 4    |    | Heneicosane, 11-(1-ethylpropyl)-    | 366 | C26H54    | 055282-11-6  | 86   |
| 5    |    | Heptadecane                         | 240 | C17H36    | 000629-78-7  | 86   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102919\  
Data File : BF117580.D  
Acq On : 29 Oct 2019 10:30  
Operator : JU  
Sample : K5563-07  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

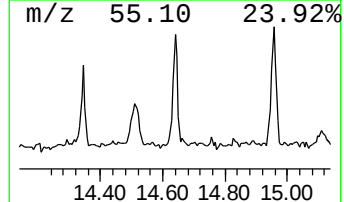
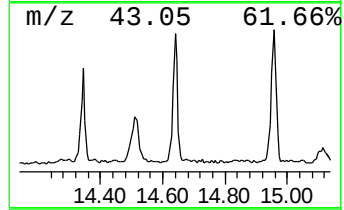
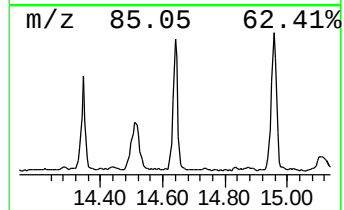
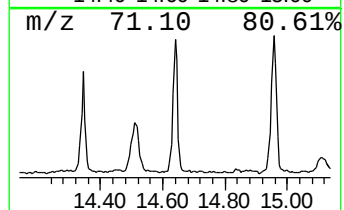
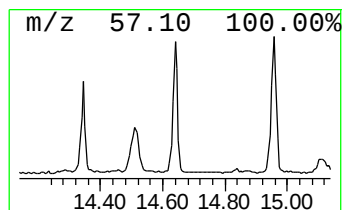
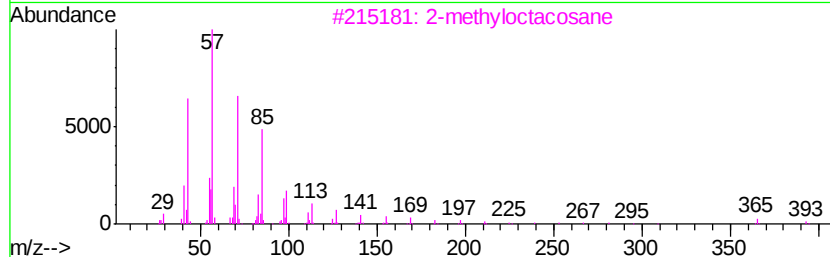
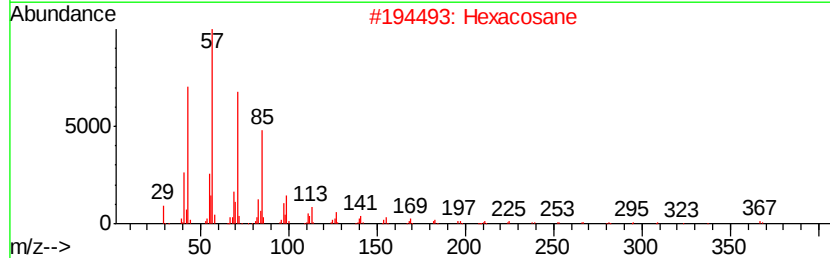
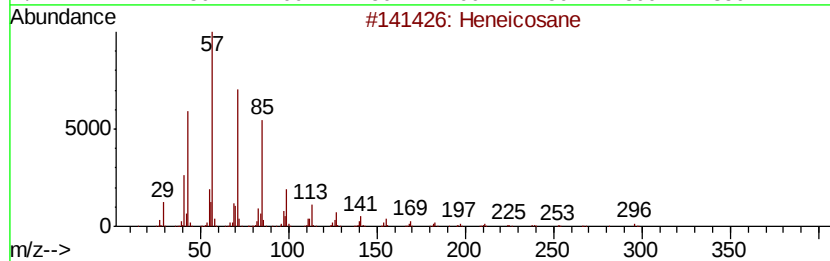
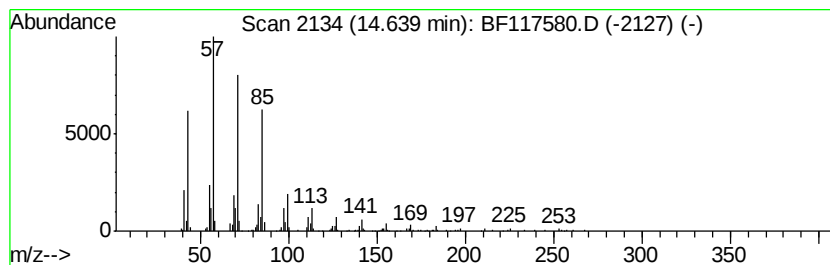
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 19 Octadecane Concentration Rank 4

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 14.64 | 13.22 ng | 344341 | Perylene-d12     | 15.33 |

| Hit# | of | Tentative ID       | MW  | MolForm | CAS#         | Qual |
|------|----|--------------------|-----|---------|--------------|------|
| 1    | 5  | Heneicosane        | 296 | C21H44  | 000629-94-7  | 91   |
| 2    |    | Hexacosane         | 366 | C26H54  | 000630-01-3  | 91   |
| 3    |    | 2-methyloctacosane | 408 | C29H60  | 1000376-72-8 | 91   |
| 4    |    | Octadecane         | 254 | C18H38  | 000593-45-3  | 87   |
| 5    |    | Heptadecane        | 240 | C17H36  | 000629-78-7  | 86   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102919\  
Data File : BF117580.D  
Acq On : 29 Oct 2019 10:30  
Operator : JU  
Sample : K5563-07  
Misc :  
ALS Vial : 3 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
TP-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

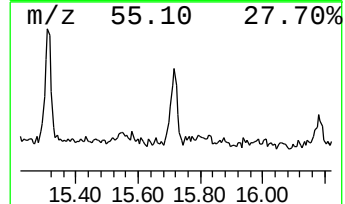
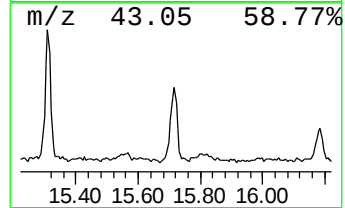
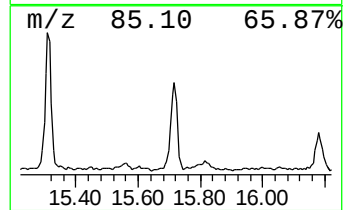
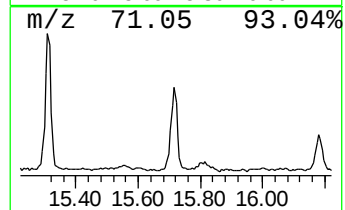
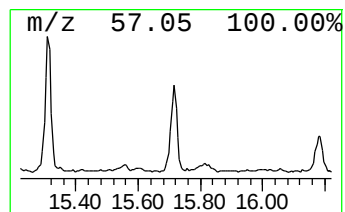
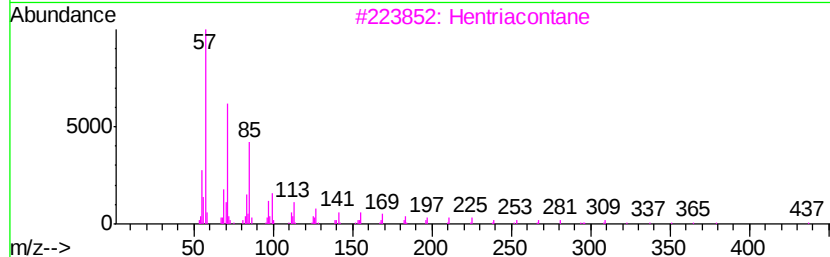
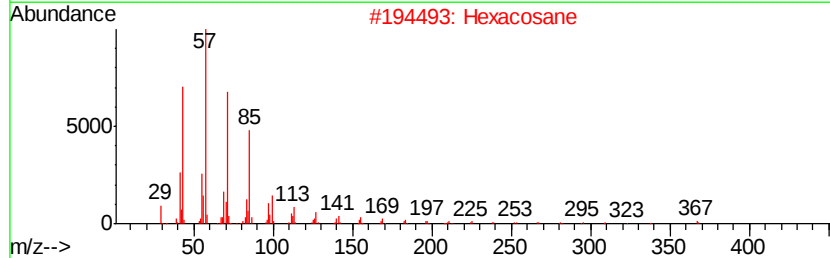
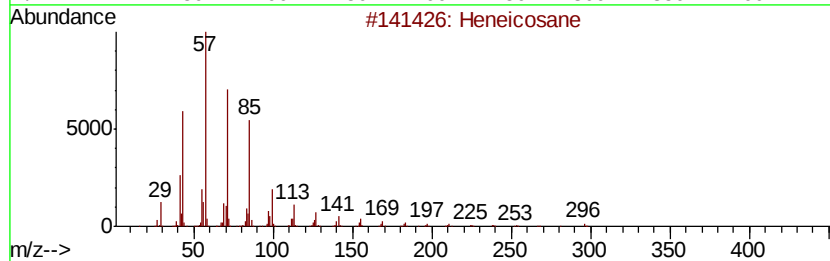
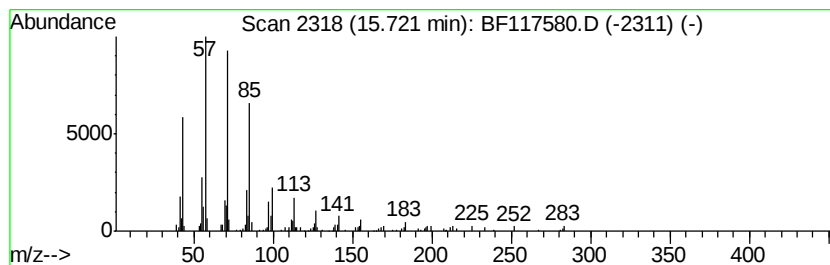
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 Triacontane Concentration Rank 9

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 15.72 | 8.46 ng | 220330 | Perylene-d12     | 15.33 |

| Hit# | of | Tentative ID          | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------|-----|---------|-------------|------|
| 1    | 5  | Heneicosane           | 296 | C21H44  | 000629-94-7 | 91   |
| 2    |    | Hexacosane            | 366 | C26H54  | 000630-01-3 | 91   |
| 3    |    | Hentriacontane        | 437 | C31H64  | 000630-04-6 | 91   |
| 4    |    | Triacontane           | 422 | C30H62  | 000638-68-6 | 91   |
| 5    |    | Heptadecane, 9-octyl- | 352 | C25H52  | 007225-64-1 | 91   |



Data Path : Z:\SVOASRV\HPCHEM1\BNA\_F\DATA\BF102919\  
 Data File : BF117580.D  
 Acq On : 29 Oct 2019 10:30  
 Operator : JU  
 Sample : K5563-07  
 Misc :  
 ALS Vial : 3 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 TP-1

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF101819.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 |
| 2-Pentanone, 4-hy... | 4.93  | 9.2 ng  |       | 162425   | 1                     | 6.73  | 352713 | 20.0 |
| unknown6.50          | 6.50  | 50.4 ng |       | 888441   | 1                     | 6.73  | 352713 | 20.0 |
| Benzene, 1,2,3-tr... | 6.55  | 7.1 ng  |       | 124745   | 1                     | 6.73  | 352713 | 20.0 |
| Benzene, 1,3-diet... | 7.05  | 5.7 ng  |       | 100191   | 1                     | 6.73  | 352713 | 20.0 |
| Naphthalene, deca... | 7.12  | 5.5 ng  |       | 96844    | 1                     | 6.73  | 352713 | 20.0 |
| Benzene, pentamet... | 7.66  | 7.0 ng  |       | 205541   | 2                     | 8.01  | 584785 | 20.0 |
| Benzene, 1,2,4,5-... | 7.75  | 6.4 ng  |       | 188292   | 2                     | 8.01  | 584785 | 20.0 |
| Benzenesulfonamid... | 10.43 | 6.2 ng  |       | 138415   | 3                     | 9.76  | 450286 | 20.0 |
| Benzenesulfonamid... | 10.63 | 6.7 ng  |       | 136392   | 4                     | 11.24 | 407979 | 20.0 |
| Tetradecanoic acid   | 11.77 | 6.8 ng  |       | 139052   | 4                     | 11.24 | 407979 | 20.0 |
| Phenol, 2-[(4-hyd... | 12.07 | 7.6 ng  |       | 155411   | 4                     | 11.24 | 407979 | 20.0 |
| Phenol, 4,4'-meth... | 12.32 | 9.3 ng  |       | 188701   | 4                     | 11.24 | 407979 | 20.0 |
| Hexacosane           | 12.92 | 20.0 ng |       | 432955   | 5                     | 13.89 | 432256 | 20.0 |
| 1-Octadecanesulph... | 13.73 | 7.2 ng  |       | 155904   | 5                     | 13.89 | 432256 | 20.0 |
| Heneicosane          | 14.04 | 5.7 ng  |       | 124097   | 5                     | 13.89 | 432256 | 20.0 |
| Heptadecane          | 14.34 | 9.2 ng  |       | 197925   | 5                     | 13.89 | 432256 | 20.0 |
| 2-methyloctacosane   | 14.51 | 9.0 ng  |       | 195192   | 5                     | 13.89 | 432256 | 20.0 |
| Octadecane           | 14.64 | 13.2 ng |       | 344341   | 6                     | 15.33 | 520970 | 20.0 |
| triacontane          | 15.72 | 8.5 ng  |       | 220330   | 6                     | 15.33 | 520970 | 20.0 |