

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101624\  
 Data File : BF139826.D  
 Acq On : 16 Oct 2024 12:42  
 Operator : RC/JU  
 Sample : P4395-01  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 F05308-SOLID

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:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF101524.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF139826.D\data.ms

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 3.910       | 304           | 309         | 316          | rVB      | 17822          | 28855         | 1.49%           | 0.123%        |
| 2         | 5.063       | 499           | 505         | 511          | rBV      | 86646          | 131400        | 6.76%           | 0.558%        |
| 3         | 5.263       | 533           | 539         | 546          | rBV      | 17425          | 29024         | 1.49%           | 0.123%        |
| 4         | 5.469       | 559           | 574         | 583          | rBV      | 1302145        | 1756838       | 90.43%          | 7.461%        |
| 5         | 5.616       | 595           | 599         | 604          | rBV2     | 19499          | 32247         | 1.66%           | 0.137%        |
| 6         | 5.704       | 610           | 614         | 620          | rBV4     | 24706          | 46684         | 2.40%           | 0.198%        |
| 7         | 5.769       | 620           | 625         | 630          | rVB3     | 22065          | 40827         | 2.10%           | 0.173%        |
| 8         | 5.863       | 635           | 641         | 646          | rBV3     | 82097          | 148490        | 7.64%           | 0.631%        |
| 9         | 5.916       | 646           | 650         | 654          | rBV2     | 34459          | 58933         | 3.03%           | 0.250%        |
| 10        | 6.004       | 661           | 665         | 669          | rVB2     | 66462          | 89890         | 4.63%           | 0.382%        |
| 11        | 6.051       | 669           | 673         | 676          | rBV      | 63156          | 93300         | 4.80%           | 0.396%        |
| 12        | 6.098       | 676           | 681         | 687          | rVV5     | 119322         | 253836        | 13.07%          | 1.078%        |
| 13        | 6.151       | 687           | 690         | 693          | rVV      | 57365          | 63477         | 3.27%           | 0.270%        |
| 14        | 6.187       | 693           | 696         | 699          | rVV3     | 41387          | 60721         | 3.13%           | 0.258%        |
| 15        | 6.240       | 703           | 705         | 708          | rVB2     | 39451          | 39333         | 2.02%           | 0.167%        |
| 16        | 6.287       | 708           | 713         | 716          | rBV2     | 95268          | 147751        | 7.61%           | 0.627%        |
| 17        | 6.322       | 716           | 719         | 725          | rVB6     | 40314          | 70915         | 3.65%           | 0.301%        |
| 18        | 6.381       | 725           | 729         | 738          | rBV3     | 167942         | 396525        | 20.41%          | 1.684%        |
| 19        | 6.487       | 738           | 747         | 751          | rBV      | 1288160        | 1942699       | 100.00%         | 8.250%        |
| 20        | 6.522       | 751           | 753         | 756          | rVB3     | 98477          | 96880         | 4.99%           | 0.411%        |
| 21        | 6.598       | 761           | 766         | 769          | rBV4     | 109231         | 199151        | 10.25%          | 0.846%        |
| 22        | 6.628       | 769           | 771         | 776          | rVB5     | 128983         | 163127        | 8.40%           | 0.693%        |
| 23        | 6.681       | 776           | 780         | 782          | rBV2     | 112389         | 161633        | 8.32%           | 0.686%        |
| 24        | 6.704       | 782           | 784         | 788          | rVB2     | 102593         | 114771        | 5.91%           | 0.487%        |
| 25        | 6.769       | 788           | 795         | 798          | rBV7     | 76979          | 171262        | 8.82%           | 0.727%        |
| 26        | 6.816       | 799           | 803         | 805          | rBV4     | 71763          | 98746         | 5.08%           | 0.419%        |
| 27        | 6.851       | 805           | 809         | 814          | rVB2     | 482681         | 645120        | 33.21%          | 2.740%        |
| 28        | 6.910       | 815           | 819         | 821          | rBV2     | 109793         | 141876        | 7.30%           | 0.603%        |
| 29        | 6.940       | 821           | 824         | 825          | rBV2     | 55310          | 50566         | 2.60%           | 0.215%        |
| 30        | 6.993       | 830           | 833         | 836          | rVV2     | 156960         | 225479        | 11.61%          | 0.958%        |
| 31        | 7.028       | 836           | 839         | 843          | rVV2     | 93877          | 121980        | 6.28%           | 0.518%        |
| 32        | 7.075       | 844           | 847         | 850          | rVV3     | 53239          | 69659         | 3.59%           | 0.296%        |
| 33        | 7.122       | 850           | 855         | 859          | rVB5     | 129505         | 216658        | 11.15%          | 0.920%        |
| 34        | 7.163       | 859           | 862         | 865          | rBV4     | 43165          | 64317         | 3.31%           | 0.273%        |
| 35        | 7.245       | 872           | 876         | 878          | rBV2     | 115651         | 147440        | 7.59%           | 0.626%        |
| 36        | 7.328       | 887           | 890         | 893          | rBV3     | 33986          | 42019         | 2.16%           | 0.178%        |

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 ALS Vial : 6 Sample Multiplier: 1

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 BNA\_F  
 ClientSampleId :  
 F05308-SOLID

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:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF101524.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |        |        |
|----|--------|------|------|------|------|---------|---------|--------|--------|
| 37 | 7.381  | 893  | 899  | 901  | rVV4 | 133351  | 264287  | 13.60% | 1.122% |
| 38 | 7.416  | 901  | 905  | 909  | rVV  | 847842  | 1148423 | 59.11% | 4.877% |
| 39 | 7.487  | 913  | 917  | 922  | rVB4 | 148441  | 244813  | 12.60% | 1.040% |
| 40 | 7.540  | 922  | 926  | 930  | rBV3 | 56632   | 104241  | 5.37%  | 0.443% |
| 41 | 7.628  | 938  | 941  | 943  | rBV3 | 58557   | 74859   | 3.85%  | 0.318% |
| 42 | 7.657  | 943  | 946  | 953  | rVB4 | 163692  | 249992  | 12.87% | 1.062% |
| 43 | 7.728  | 953  | 958  | 962  | rBV6 | 89255   | 158616  | 8.16%  | 0.674% |
| 44 | 7.769  | 962  | 965  | 973  | rVB7 | 103252  | 142597  | 7.34%  | 0.606% |
| 45 | 7.881  | 980  | 984  | 991  | rVB7 | 36947   | 79574   | 4.10%  | 0.338% |
| 46 | 7.951  | 992  | 996  | 1001 | rBV4 | 51190   | 85768   | 4.41%  | 0.364% |
| 47 | 8.004  | 1001 | 1005 | 1008 | rBV3 | 58199   | 85689   | 4.41%  | 0.364% |
| 48 | 8.081  | 1015 | 1018 | 1020 | rBV2 | 33516   | 46872   | 2.41%  | 0.199% |
| 49 | 8.134  | 1021 | 1027 | 1032 | rVV  | 579179  | 856748  | 44.10% | 3.638% |
| 50 | 8.187  | 1033 | 1036 | 1039 | rVV  | 117733  | 142510  | 7.34%  | 0.605% |
| 51 | 8.216  | 1039 | 1041 | 1049 | rVB6 | 57477   | 94526   | 4.87%  | 0.401% |
| 52 | 8.281  | 1049 | 1052 | 1054 | rBV3 | 29778   | 39061   | 2.01%  | 0.166% |
| 53 | 8.416  | 1073 | 1075 | 1081 | rVB6 | 68709   | 101131  | 5.21%  | 0.429% |
| 54 | 8.475  | 1082 | 1085 | 1088 | rBV3 | 47051   | 51422   | 2.65%  | 0.218% |
| 55 | 8.516  | 1089 | 1092 | 1094 | rVB3 | 36529   | 38490   | 1.98%  | 0.163% |
| 56 | 8.551  | 1094 | 1098 | 1103 | rBV  | 243362  | 425094  | 21.88% | 1.805% |
| 57 | 8.634  | 1109 | 1112 | 1115 | rBV2 | 44826   | 58468   | 3.01%  | 0.248% |
| 58 | 8.669  | 1115 | 1118 | 1123 | rVV5 | 54035   | 90351   | 4.65%  | 0.384% |
| 59 | 8.910  | 1156 | 1159 | 1167 | rBV7 | 49303   | 118510  | 6.10%  | 0.503% |
| 60 | 9.010  | 1173 | 1176 | 1178 | rBV4 | 48144   | 66046   | 3.40%  | 0.280% |
| 61 | 9.151  | 1196 | 1200 | 1204 | rVB  | 217983  | 274351  | 14.12% | 1.165% |
| 62 | 9.204  | 1204 | 1209 | 1214 | rVB  | 1334920 | 1723365 | 88.71% | 7.319% |
| 63 | 9.286  | 1219 | 1223 | 1227 | rBV3 | 124796  | 196170  | 10.10% | 0.833% |
| 64 | 9.475  | 1251 | 1255 | 1259 | rVB  | 91424   | 117037  | 6.02%  | 0.497% |
| 65 | 9.545  | 1264 | 1267 | 1271 | rVV  | 97361   | 137900  | 7.10%  | 0.586% |
| 66 | 9.598  | 1271 | 1276 | 1281 | rVB2 | 306098  | 471227  | 24.26% | 2.001% |
| 67 | 9.745  | 1298 | 1301 | 1305 | rBV4 | 98682   | 136689  | 7.04%  | 0.580% |
| 68 | 9.792  | 1306 | 1309 | 1313 | rVB  | 136012  | 167194  | 8.61%  | 0.710% |
| 69 | 9.886  | 1320 | 1325 | 1329 | rVB  | 526479  | 661263  | 34.04% | 2.808% |
| 70 | 9.992  | 1339 | 1343 | 1347 | rVB2 | 84684   | 118871  | 6.12%  | 0.505% |
| 71 | 10.081 | 1355 | 1358 | 1363 | rVB3 | 86714   | 124988  | 6.43%  | 0.531% |
| 72 | 10.128 | 1363 | 1366 | 1375 | rVB4 | 103878  | 216560  | 11.15% | 0.920% |
| 73 | 10.204 | 1375 | 1379 | 1381 | rBV  | 92988   | 134487  | 6.92%  | 0.571% |
| 74 | 10.286 | 1390 | 1393 | 1400 | rVB4 | 78934   | 146988  | 7.57%  | 0.624% |
| 75 | 10.598 | 1441 | 1446 | 1450 | rVB  | 318381  | 412832  | 21.25% | 1.753% |
| 76 | 10.675 | 1454 | 1459 | 1464 | rVV  | 702194  | 878426  | 45.22% | 3.730% |

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 Sample : P4395-01  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 F05308-SOLID

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:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF101524.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |        |         |        |        |
|----|--------|------|------|------|------|--------|---------|--------|--------|
| 77 | 10.792 | 1476 | 1479 | 1486 | rVB  | 141397 | 207027  | 10.66% | 0.879% |
| 78 | 10.869 | 1489 | 1492 | 1495 | rVV2 | 49810  | 69143   | 3.56%  | 0.294% |
| 79 | 10.904 | 1495 | 1498 | 1502 | rVV  | 71715  | 105285  | 5.42%  | 0.447% |
| 80 | 10.975 | 1506 | 1510 | 1513 | rVV2 | 121789 | 180583  | 9.30%  | 0.767% |
| 81 | 11.010 | 1513 | 1516 | 1527 | rVB5 | 77999  | 182655  | 9.40%  | 0.776% |
| 82 | 11.263 | 1555 | 1559 | 1566 | rVB4 | 85645  | 144540  | 7.44%  | 0.614% |
| 83 | 11.369 | 1573 | 1577 | 1585 | rBV  | 454654 | 636638  | 32.77% | 2.704% |
| 84 | 11.480 | 1592 | 1596 | 1602 | rBV4 | 45834  | 90143   | 4.64%  | 0.383% |
| 85 | 11.786 | 1645 | 1648 | 1651 | rBV  | 22932  | 28753   | 1.48%  | 0.122% |
| 86 | 11.892 | 1660 | 1666 | 1671 | rBV3 | 125084 | 218410  | 11.24% | 0.928% |
| 87 | 12.345 | 1741 | 1743 | 1751 | rVB2 | 35415  | 60942   | 3.14%  | 0.259% |
| 88 | 12.422 | 1752 | 1756 | 1759 | rBV  | 66071  | 86159   | 4.44%  | 0.366% |
| 89 | 12.586 | 1780 | 1784 | 1788 | rBV  | 91101  | 114839  | 5.91%  | 0.488% |
| 90 | 12.816 | 1818 | 1823 | 1829 | rBV2 | 102146 | 183781  | 9.46%  | 0.780% |
| 91 | 12.869 | 1829 | 1832 | 1836 | rVB  | 46806  | 57767   | 2.97%  | 0.245% |
| 92 | 12.951 | 1842 | 1846 | 1854 | rVB  | 791936 | 1058200 | 54.47% | 4.494% |
| 93 | 14.010 | 2019 | 2026 | 2037 | rBV2 | 396411 | 693907  | 35.72% | 2.947% |
| 94 | 15.415 | 2262 | 2265 | 2270 | rVB  | 48423  | 72427   | 3.73%  | 0.308% |
| 95 | 15.474 | 2270 | 2275 | 2288 | rVB  | 311211 | 507209  | 26.11% | 2.154% |

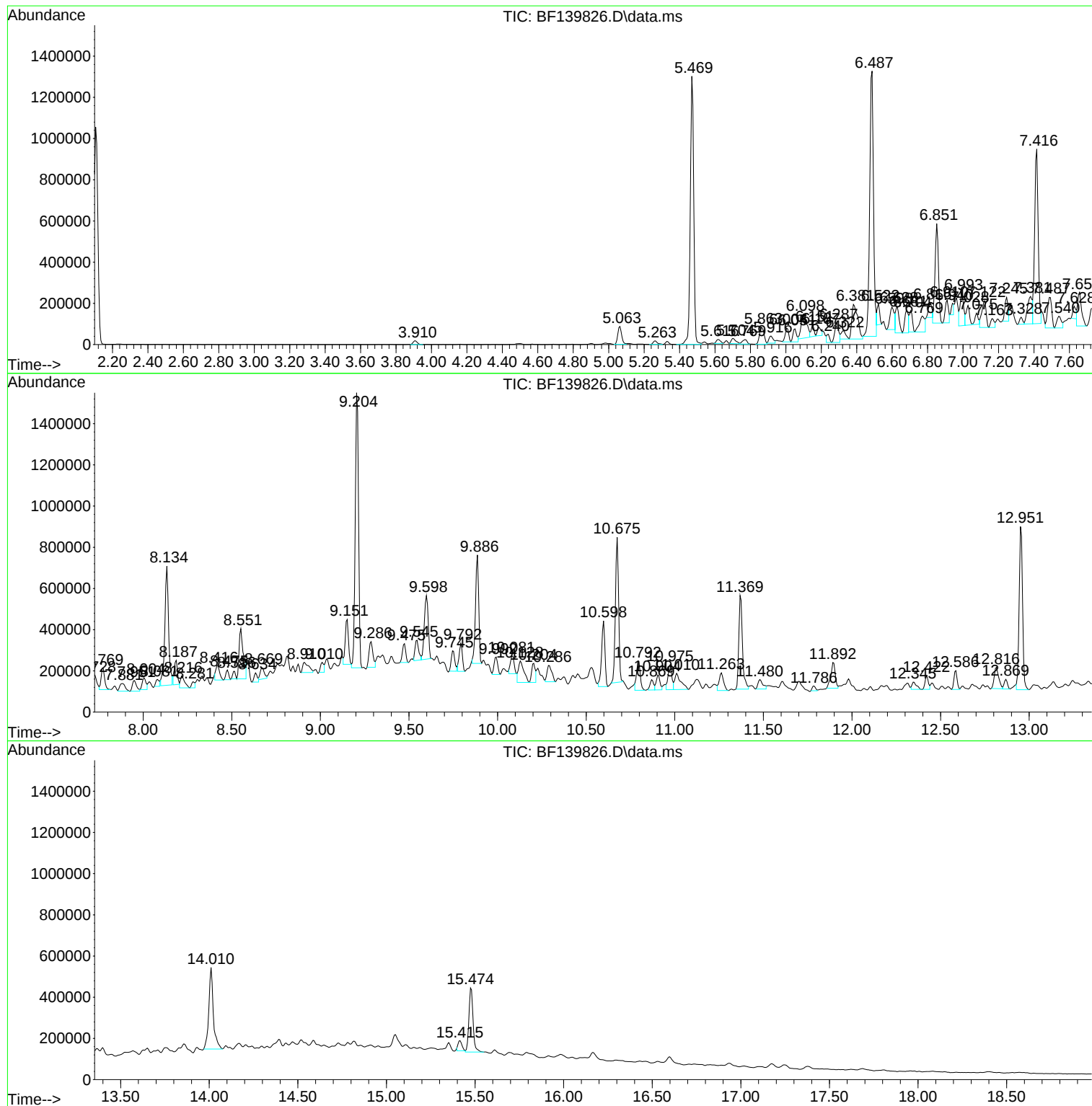
Sum of corrected areas: 23547273

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101624\  
Data File : BF139826.D  
Acq On : 16 Oct 2024 12:42  
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Sample : P4395-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
F05308-SOLID

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF101524.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101624\  
Data File : BF139826.D  
Acq On : 16 Oct 2024 12:42  
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Sample : P4395-01  
Misc :  
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Instrument :  
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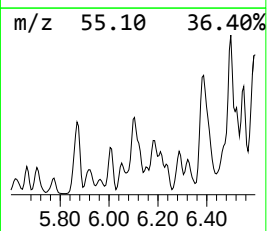
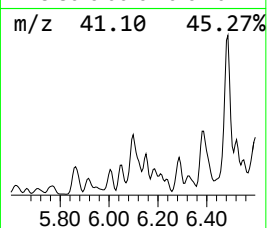
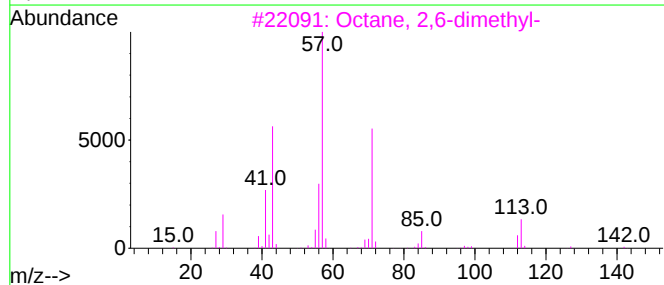
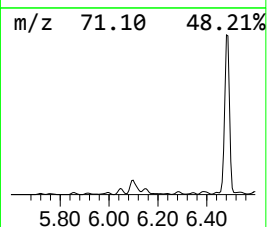
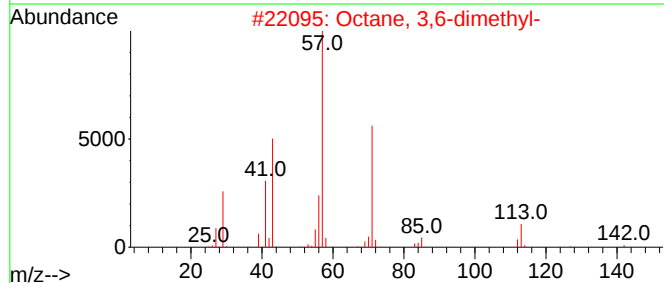
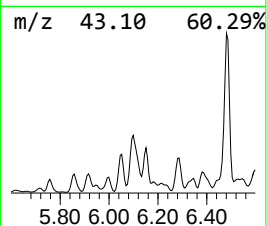
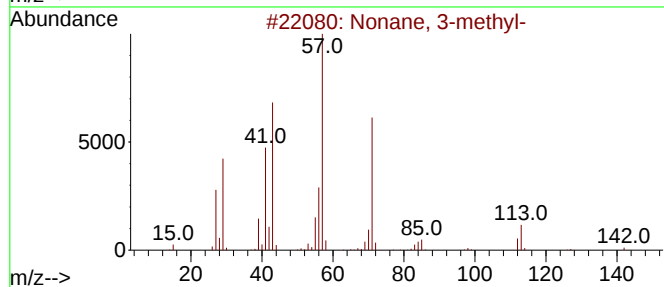
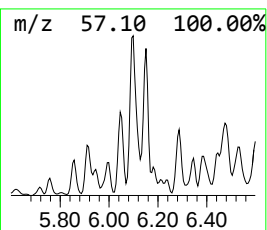
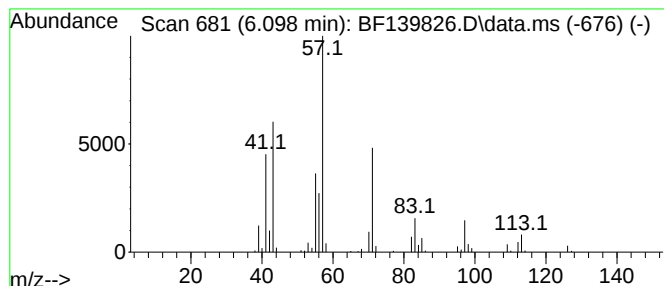
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF101524.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 1 Nonane, 3-methyl- Concentration Rank 6

| R.T.  | EstConc | Area   | Relative to ISTD       | R.T.  |
|-------|---------|--------|------------------------|-------|
| 6.098 | 7.87 ng | 253836 | 1,4-Dichlorobenzene-d4 | 6.851 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Nonane, 3-methyl-                   | 142 | C10H22    | 005911-04-6  | 64   |
| 2    |    |   | Octane, 3,6-dimethyl-               | 142 | C10H22    | 015869-94-0  | 59   |
| 3    |    |   | Octane, 2,6-dimethyl-               | 142 | C10H22    | 002051-30-1  | 59   |
| 4    |    |   | Sulfurous acid, 2-ethylhexyl oct... | 446 | C26H54O3S | 1000309-20-1 | 53   |
| 5    |    |   | Sulfurous acid, 2-ethylhexyl hep... | 432 | C25H52O3S | 1000309-20-0 | 53   |



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ClientSampleId :  
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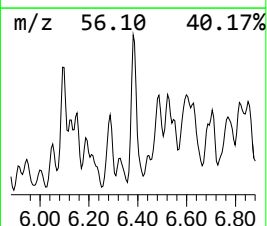
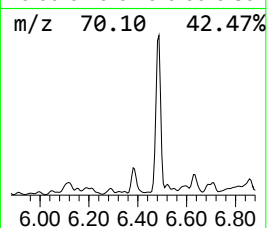
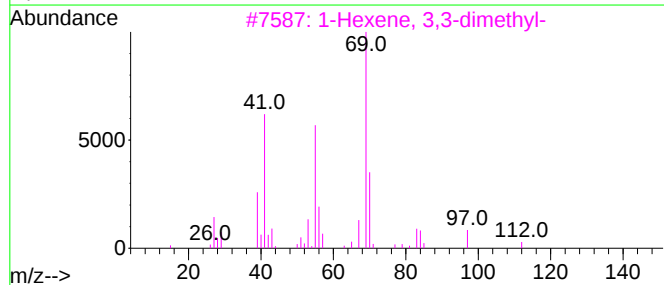
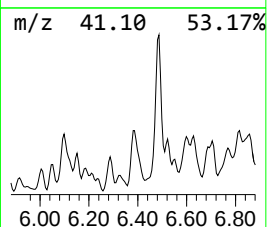
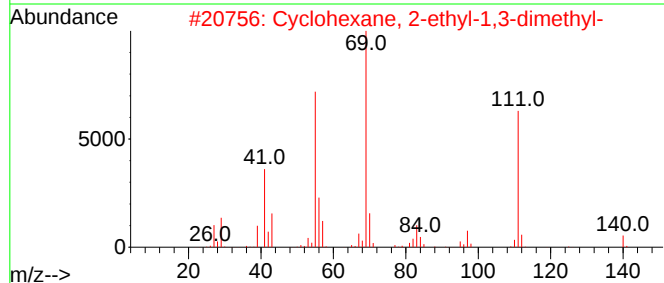
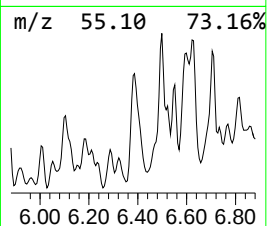
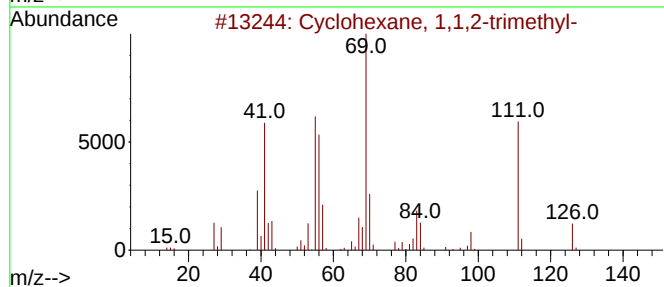
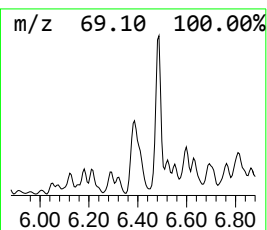
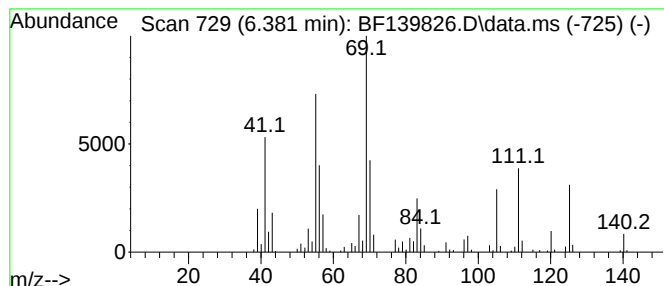
Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF101524.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

| R.T.  | EstConc  | Area   | Relative to ISTD       | R.T.  |
|-------|----------|--------|------------------------|-------|
| 6.381 | 12.29 ng | 396525 | 1,4-Dichlorobenzene-d4 | 6.851 |

| Hit# | of | 5 | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|----|---|------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, 1,1,2-trimethyl-      | 126 | C9H18   | 007094-26-0 | 58   |
| 2    |    |   | Cyclohexane, 2-ethyl-1,3-dimethyl- | 140 | C10H20  | 007045-67-2 | 47   |
| 3    |    |   | 1-Hexene, 3,3-dimethyl-            | 112 | C8H16   | 003404-77-1 | 46   |
| 4    |    |   | 3-Methyl-2-hexene                  | 98  | C7H14   | 017618-77-8 | 43   |
| 5    |    |   | 3-Octene, (E)-                     | 112 | C8H16   | 014919-01-8 | 41   |



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Data File : BF139826.D  
Acq On : 16 Oct 2024 12:42  
Operator : RC/JU  
Sample : P4395-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
F05308-SOLID

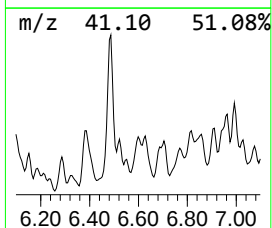
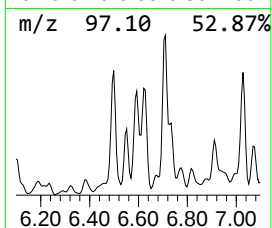
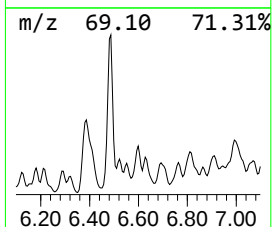
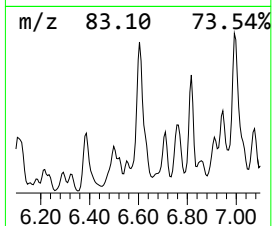
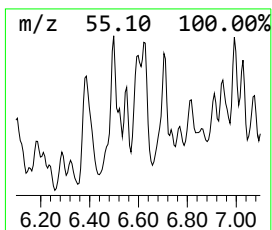
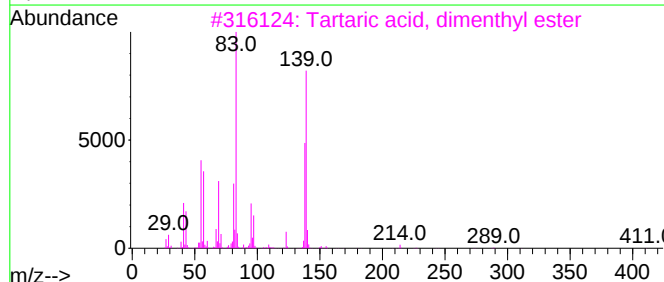
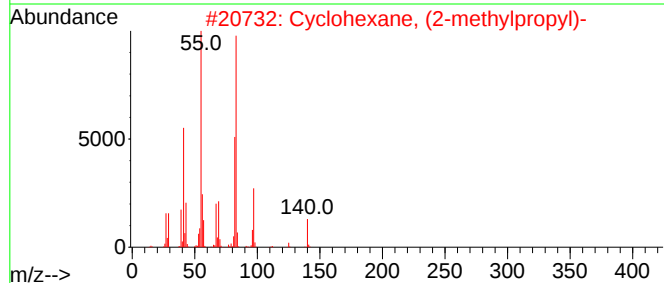
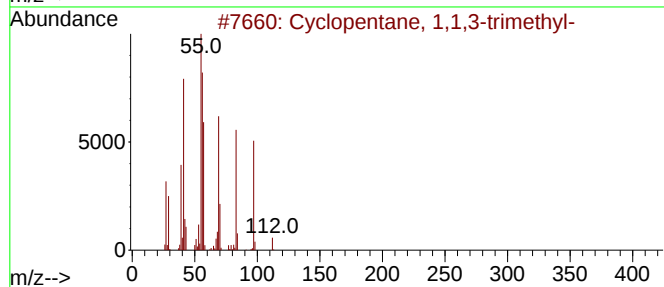
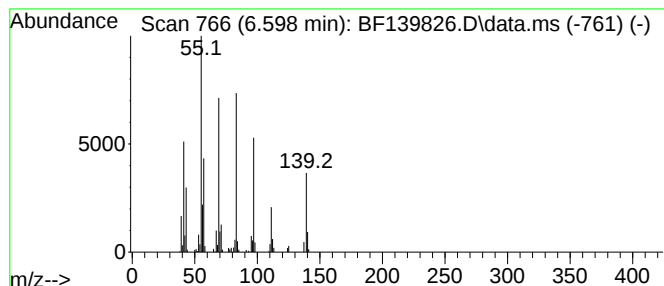
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 3 Cyclopentane, 1,1,3-trimethyl- Concentration Rank 13

| R.T.      | EstConc                        | Area   | Relative to ISTD       |          | R.T.         |      |
|-----------|--------------------------------|--------|------------------------|----------|--------------|------|
| 6.598     | 6.17 ng                        | 199151 | 1,4-Dichlorobenzene-d4 |          | 6.851        |      |
| Hit# of 5 | Tentative ID                   |        | MW                     | MolForm  | CAS#         | Qual |
| 1         | Cyclopentane, 1,1,3-trimethyl- |        | 112                    | C8H16    | 004516-69-2  | 64   |
| 2         | Cyclohexane, (2-methylpropyl)- |        | 140                    | C10H20   | 001678-98-4  | 49   |
| 3         | Tartaric acid, dimethyl ester  |        | 426                    | C24H42O6 | 1000149-89-9 | 43   |
| 4         | 1-Pentene, 2,3,3-trimethyl-    |        | 112                    | C8H16    | 000560-23-6  | 38   |
| 5         | 1-Pentyn-3-ol                  |        | 84                     | C5H8O    | 004187-86-4  | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101624\  
Data File : BF139826.D  
Acq On : 16 Oct 2024 12:42  
Operator : RC/JU  
Sample : P4395-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
F05308-SOLID

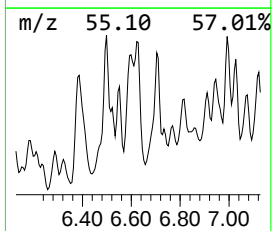
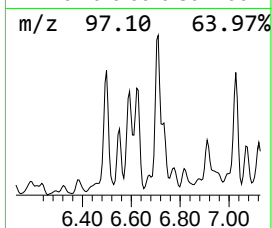
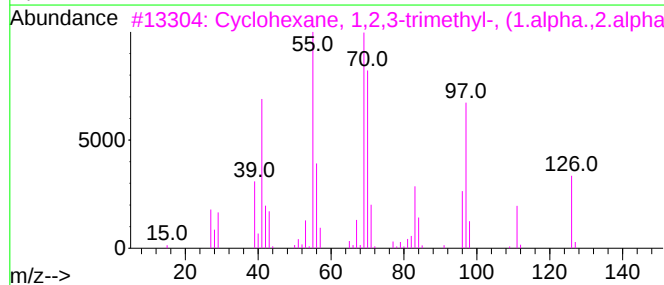
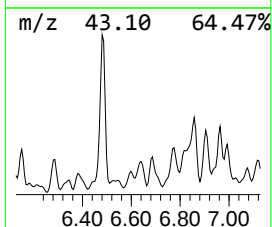
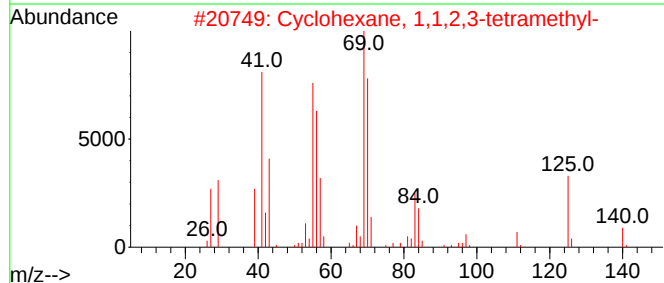
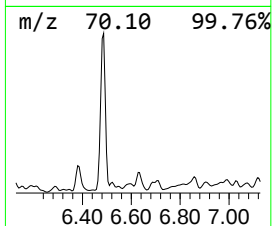
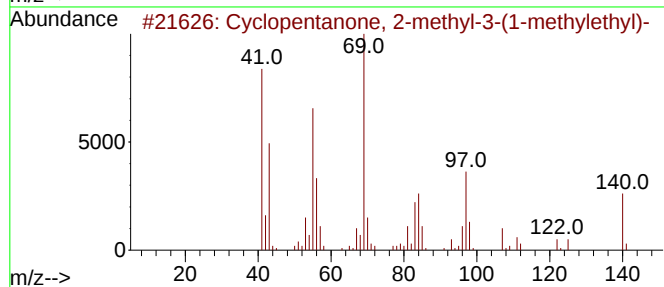
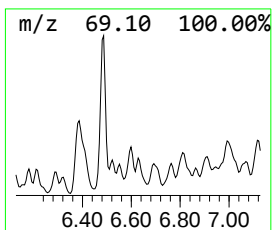
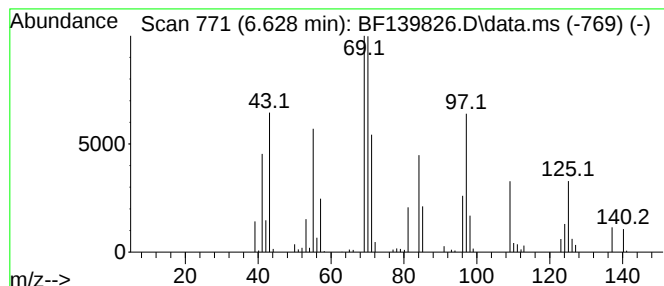
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 4 unknown6.628 Concentration Rank 19

| R.T.  | EstConc | Area   | Relative to ISTD       | R.T.  |
|-------|---------|--------|------------------------|-------|
| 6.628 | 5.06 ng | 163127 | 1,4-Dichlorobenzene-d4 | 6.851 |

| Hit# | of | 5 | Tentative ID                         | MW  | MolForm     | CAS#         | Qual |
|------|----|---|--------------------------------------|-----|-------------|--------------|------|
| 1    |    |   | Cyclopentanone, 2-methyl-3-(1-me...  | 140 | C9H16O      | 054549-81-4  | 43   |
| 2    |    |   | Cyclohexane, 1,1,2,3-tetramethyl-    | 140 | C10H20      | 006783-92-2  | 43   |
| 3    |    |   | Cyclohexane, 1,2,3-trimethyl-, (...) | 126 | C9H18       | 001839-88-9  | 40   |
| 4    |    |   | Dichloroacetic acid, 6-ethyl-3-o...  | 268 | C12H22Cl2O2 | 1000282-56-6 | 37   |
| 5    |    |   | 1-Hexene, 4-ethyl-                   | 112 | C8H16       | 016746-85-3  | 22   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101624\  
Data File : BF139826.D  
Acq On : 16 Oct 2024 12:42  
Operator : RC/JU  
Sample : P4395-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
F05308-SOLID

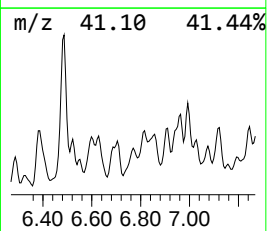
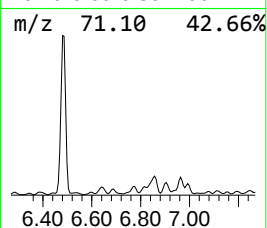
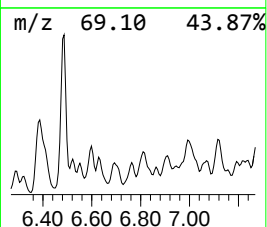
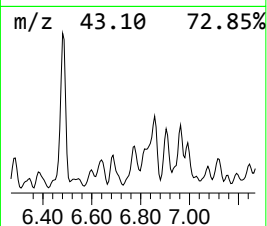
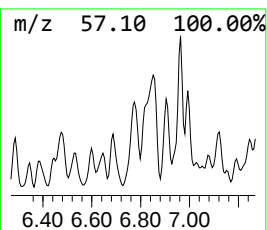
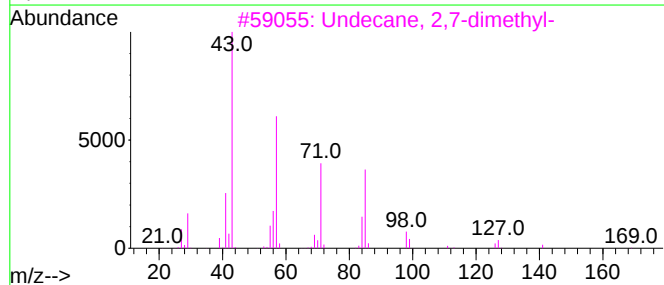
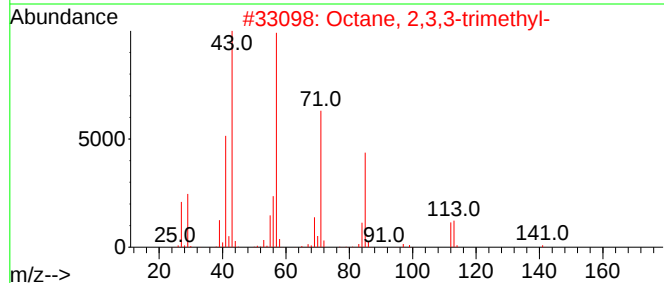
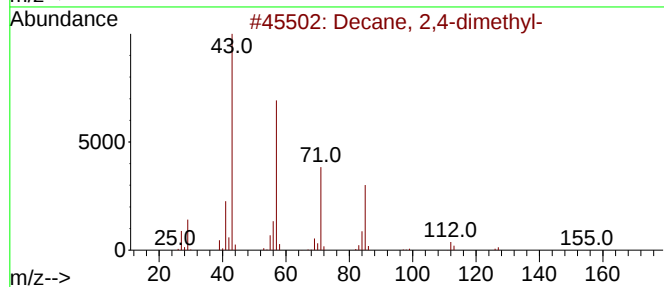
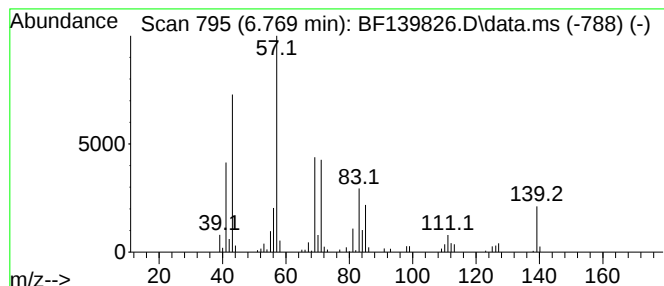
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 5 Decane, 2,4-dimethyl- Concentration Rank 18

| R.T.  | EstConc | Area   | Relative to ISTD       | R.T.  |
|-------|---------|--------|------------------------|-------|
| 6.769 | 5.31 ng | 171262 | 1,4-Dichlorobenzene-d4 | 6.851 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Decane, 2,4-dimethyl-               | 170 | C12H26    | 002801-84-5  | 52   |
| 2    |    |   | Octane, 2,3,3-trimethyl-            | 156 | C11H24    | 062016-30-2  | 47   |
| 3    |    |   | Undecane, 2,7-dimethyl-             | 184 | C13H28    | 017301-24-5  | 47   |
| 4    |    |   | Octane, 3-ethyl-2,7-dimethyl-       | 170 | C12H26    | 062183-55-5  | 43   |
| 5    |    |   | Sulfurous acid, 2-ethylhexyl hex... | 278 | C14H30O3S | 1000309-20-2 | 43   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101624\  
Data File : BF139826.D  
Acq On : 16 Oct 2024 12:42  
Operator : RC/JU  
Sample : P4395-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
F05308-SOLID

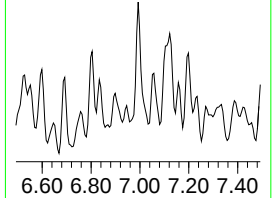
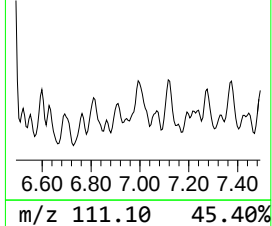
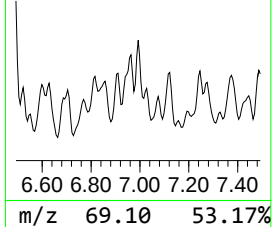
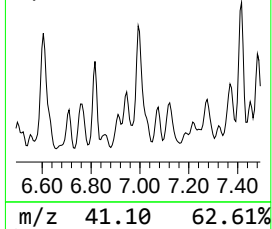
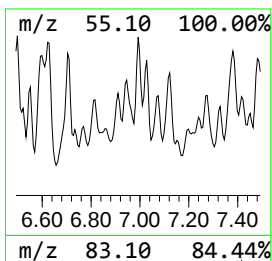
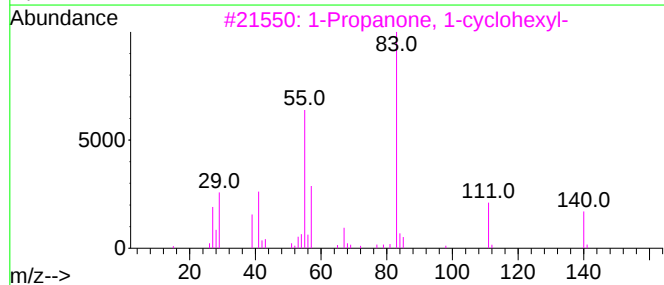
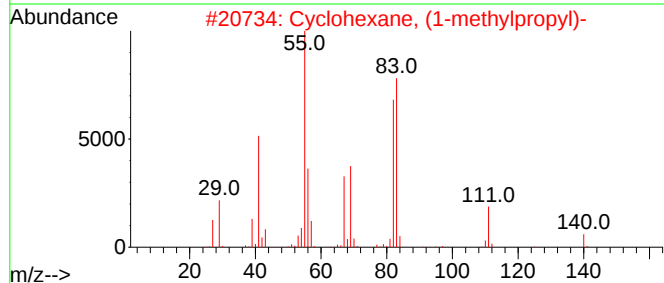
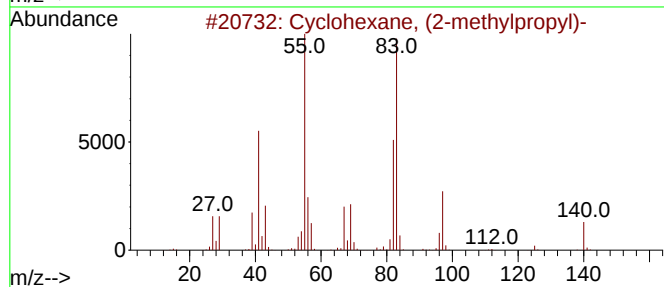
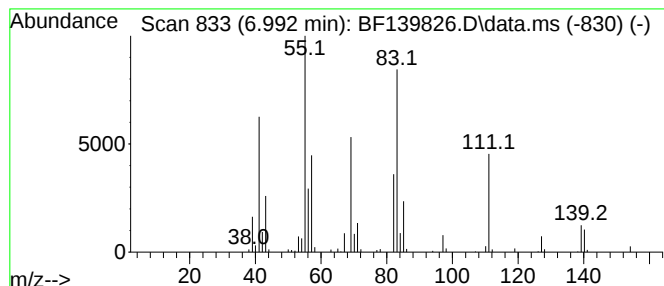
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 6 unknown6.992 Concentration Rank 8

| R.T.  | EstConc | Area   | Relative to ISTD       | R.T.  |
|-------|---------|--------|------------------------|-------|
| 6.992 | 6.99 ng | 225479 | 1,4-Dichlorobenzene-d4 | 6.851 |

| Hit# | of | 5 | Tentative ID                   | MW  | MolForm | CAS#        | Qual |
|------|----|---|--------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, (2-methylpropyl)- | 140 | C10H20  | 001678-98-4 | 49   |
| 2    |    |   | Cyclohexane, (1-methylpropyl)- | 140 | C10H20  | 007058-01-7 | 47   |
| 3    |    |   | 1-Propanone, 1-cyclohexyl-     | 140 | C9H16O  | 001123-86-0 | 43   |
| 4    |    |   | Cyclohexane, ethyl-            | 112 | C8H16   | 001678-91-7 | 43   |
| 5    |    |   | Cyclohexane, butyl-            | 140 | C10H20  | 001678-93-9 | 43   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101624\  
Data File : BF139826.D  
Acq On : 16 Oct 2024 12:42  
Operator : RC/JU  
Sample : P4395-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
F05308-SOLID

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

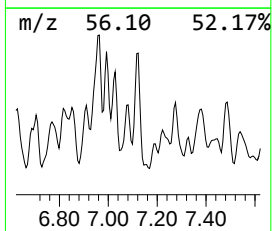
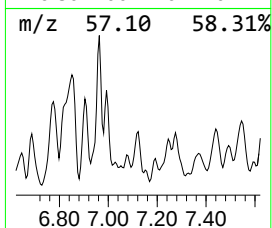
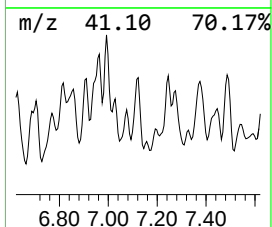
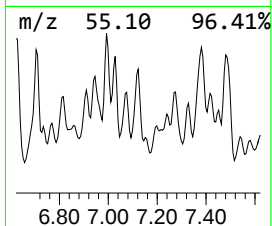
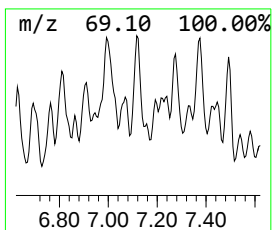
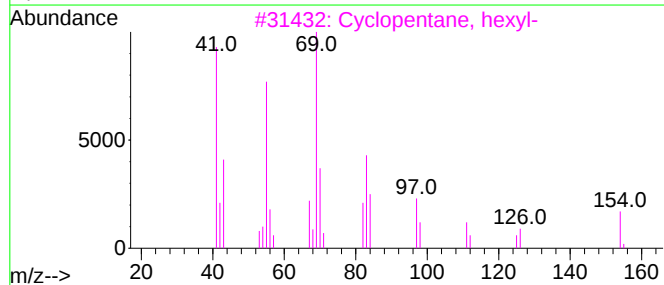
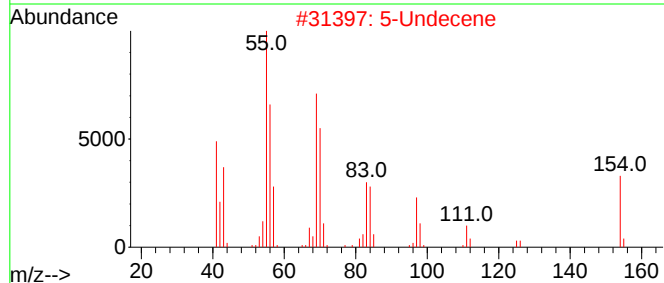
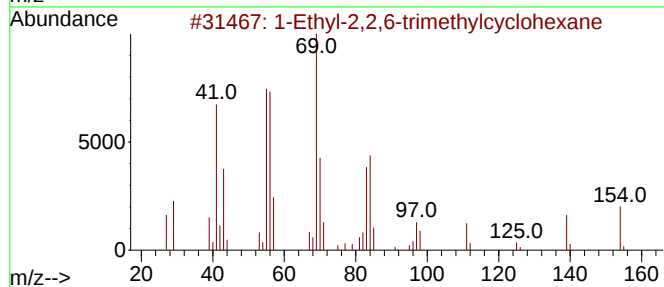
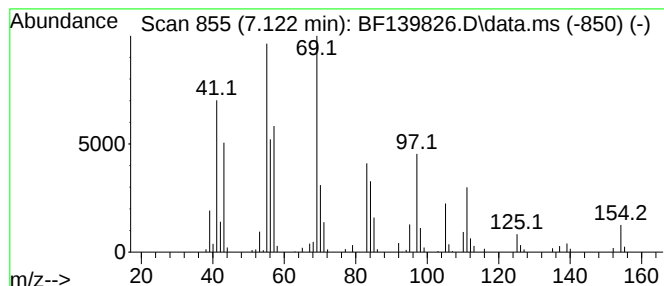
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TIC Integration Parameters: LSCINT.P

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Peak Number 7 1-Ethyl-2,2,6-trimethylcycl... Concentration Rank 10

| R.T.  | EstConc | Area   | Relative to ISTD       |  |  | R.T.  |
|-------|---------|--------|------------------------|--|--|-------|
| 7.122 | 6.72 ng | 216658 | 1,4-Dichlorobenzene-d4 |  |  | 6.851 |

| Hit# | of 5 | Tentative ID                       | MW  | MolForm | CAS#        | Qual |
|------|------|------------------------------------|-----|---------|-------------|------|
| 1    |      | 1-Ethyl-2,2,6-trimethylcyclohexane | 154 | C11H22  | 071186-27-1 | 76   |
| 2    |      | 5-Undecene                         | 154 | C11H22  | 004941-53-1 | 55   |
| 3    |      | Cyclopentane, hexyl-               | 154 | C11H22  | 004457-00-5 | 49   |
| 4    |      | 1-Tridecene                        | 182 | C13H26  | 002437-56-1 | 46   |
| 5    |      | Nonane, 2-methyl-3-methylene-      | 154 | C11H22  | 055499-08-6 | 43   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101624\  
Data File : BF139826.D  
Acq On : 16 Oct 2024 12:42  
Operator : RC/JU  
Sample : P4395-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
F05308-SOLID

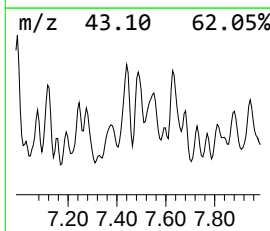
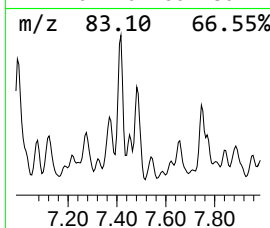
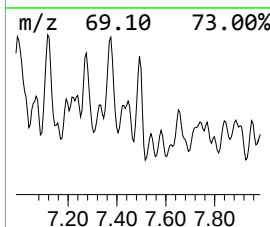
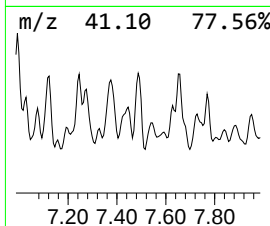
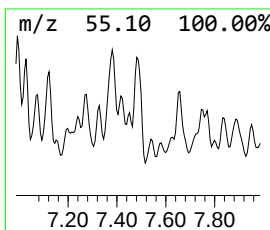
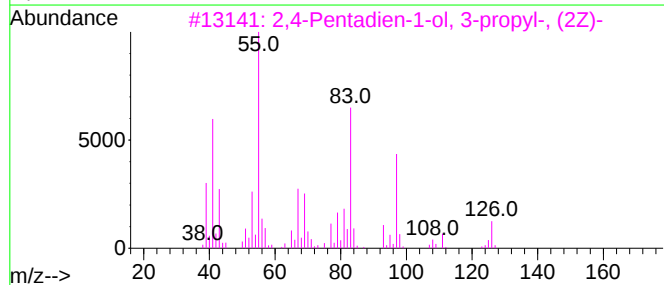
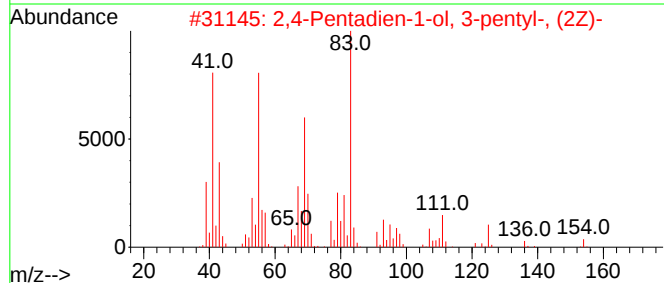
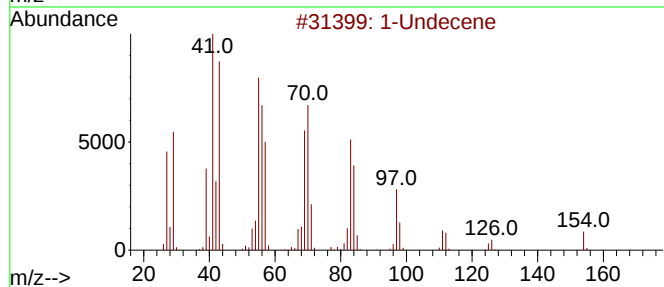
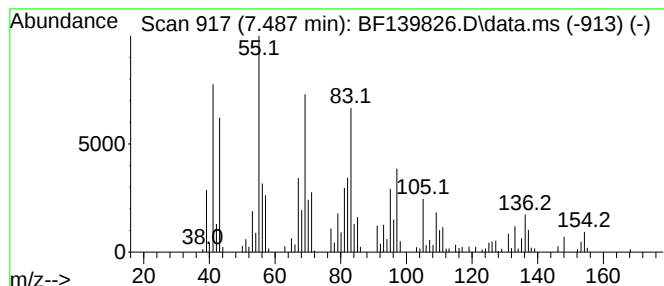
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 8 unknown7.487 Concentration Rank 7

| R.T.  | EstConc | Area   | Relative to ISTD       | R.T.  |
|-------|---------|--------|------------------------|-------|
| 7.487 | 7.59 ng | 244813 | 1,4-Dichlorobenzene-d4 | 6.851 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|---------|--------------|------|
| 1    |    |   | 1-Undecene                          | 154 | C11H22  | 000821-95-4  | 46   |
| 2    |    |   | 2,4-Pentadien-1-ol, 3-pentyl-, (... | 154 | C10H18O | 1010142-19-7 | 38   |
| 3    |    |   | 2,4-Pentadien-1-ol, 3-propyl-, (... | 126 | C8H14O  | 1000142-17-9 | 38   |
| 4    |    |   | 1-Eicosanol                         | 298 | C20H42O | 000629-96-9  | 35   |
| 5    |    |   | 6-Octenal, 3,7-dimethyl-, (S)-      | 154 | C10H18O | 005949-05-3  | 30   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101624\  
Data File : BF139826.D  
Acq On : 16 Oct 2024 12:42  
Operator : RC/JU  
Sample : P4395-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
F05308-SOLID

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF101524.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

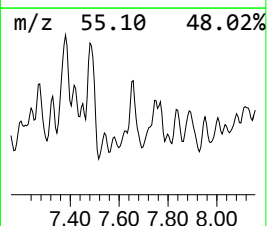
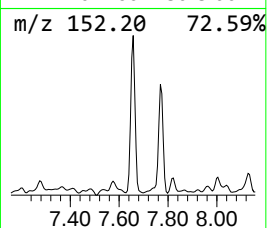
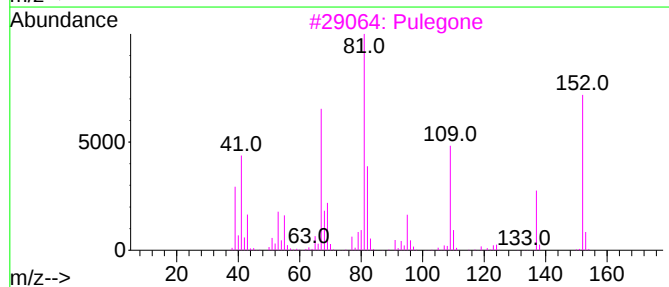
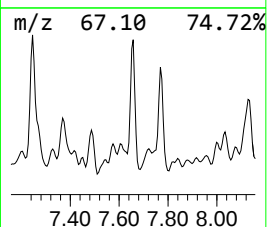
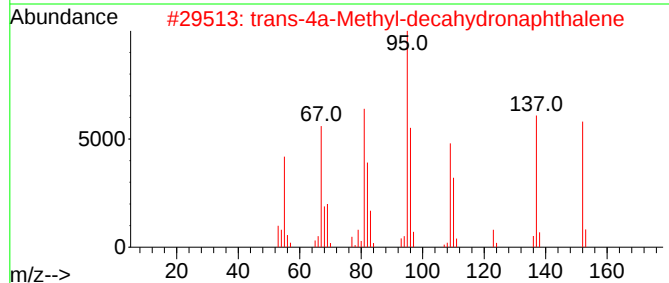
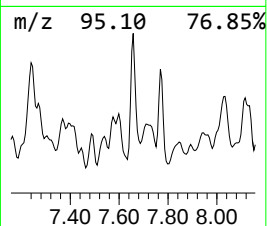
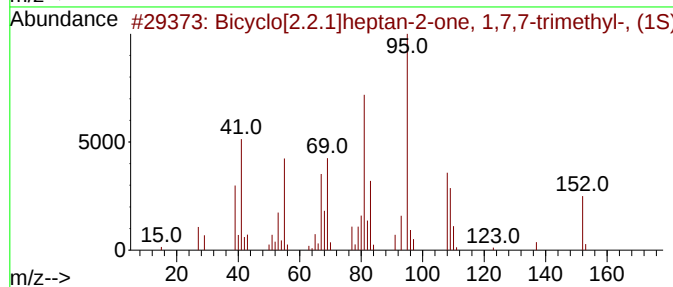
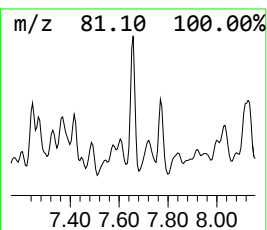
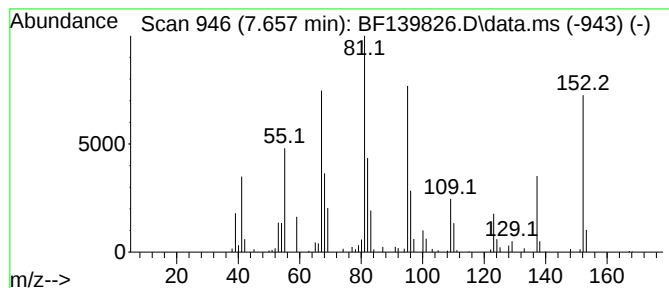
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Peak Number 9 Bicyclo[2.2.1]heptan-2-one,... Concentration Rank 15

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 7.657 | 5.84 ng | 249992 | Naphthalene-d8   | 8.134 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------------|-----|---------|-------------|------|
| 1    |      | Bicyclo[2.2.1]heptan-2-one, 1,7,... | 152 | C10H16O | 000464-48-2 | 72   |
| 2    |      | trans-4a-Methyl-decahydronaphtha... | 152 | C11H20  | 002547-27-5 | 70   |
| 3    |      | Pulegone                            | 152 | C10H16O | 000089-82-7 | 70   |
| 4    |      | Naphthalene, decahydro-2-methyl-    | 152 | C11H20  | 002958-76-1 | 62   |
| 5    |      | (+)-Dihydrocarvone                  | 152 | C10H16O | 005524-05-0 | 58   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101624\  
Data File : BF139826.D  
Acq On : 16 Oct 2024 12:42  
Operator : RC/JU  
Sample : P4395-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
F05308-SOLID

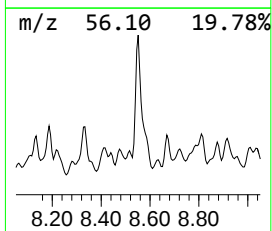
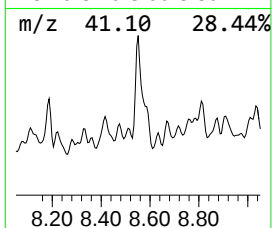
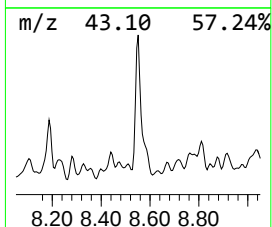
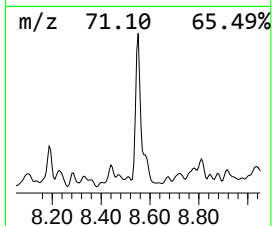
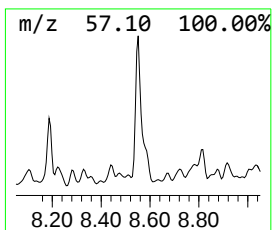
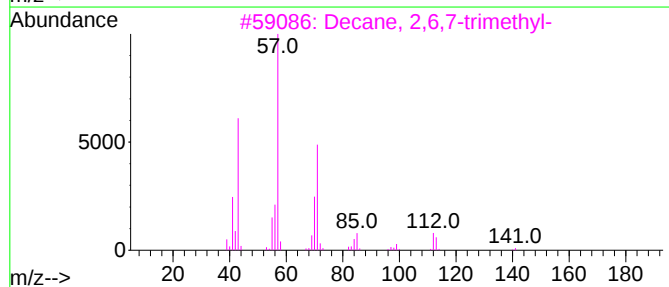
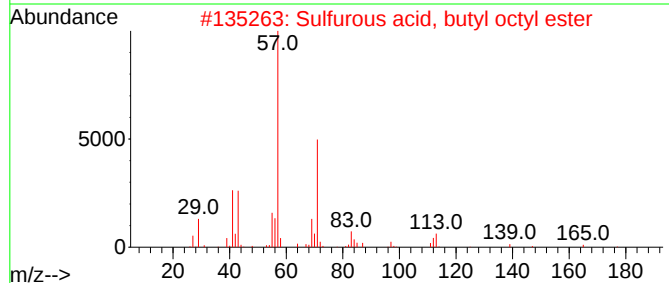
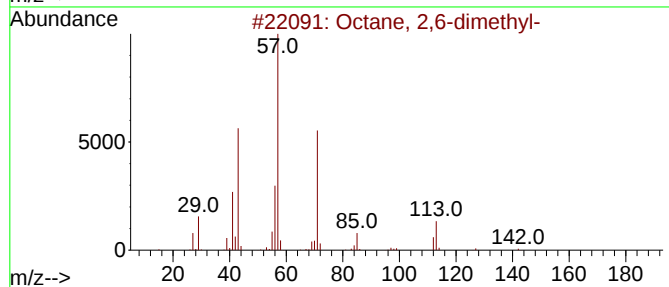
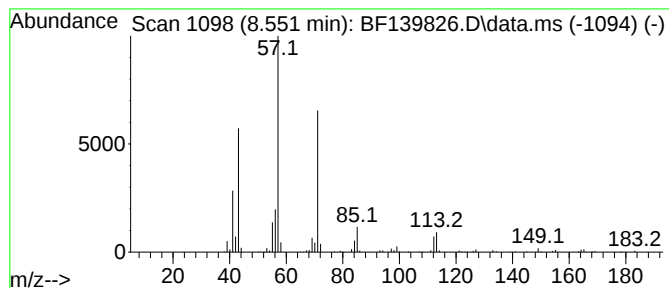
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 10 Octane, 2,6-dimethyl- Concentration Rank 4

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 8.551 | 9.92 ng | 425094 | Naphthalene-d8   | 8.134 |

| Hit# | of 5 | Tentative ID                      | MW  | MolForm   | CAS#         | Qual |
|------|------|-----------------------------------|-----|-----------|--------------|------|
| 1    |      | Octane, 2,6-dimethyl-             | 142 | C10H22    | 002051-30-1  | 78   |
| 2    |      | Sulfurous acid, butyl octyl ester | 250 | C12H26O3S | 1000309-17-5 | 78   |
| 3    |      | Decane, 2,6,7-trimethyl-          | 184 | C13H28    | 062108-25-2  | 72   |
| 4    |      | Octane, 3,6-dimethyl-             | 142 | C10H22    | 015869-94-0  | 72   |
| 5    |      | Octane, 2,3,7-trimethyl-          | 156 | C11H24    | 062016-34-6  | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101624\  
Data File : BF139826.D  
Acq On : 16 Oct 2024 12:42  
Operator : RC/JU  
Sample : P4395-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
F05308-SOLID

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF101524.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

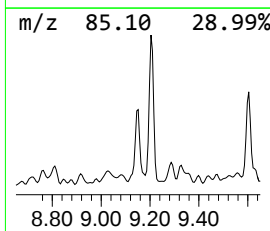
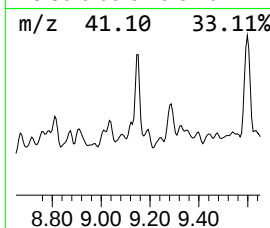
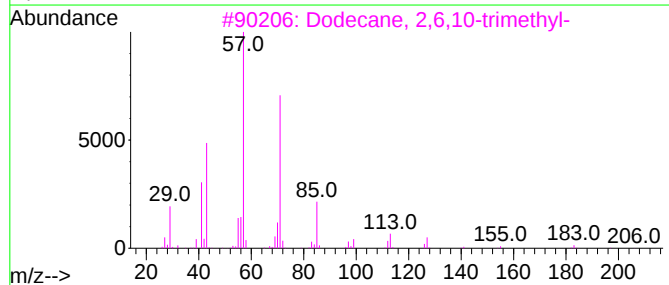
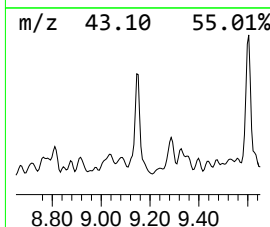
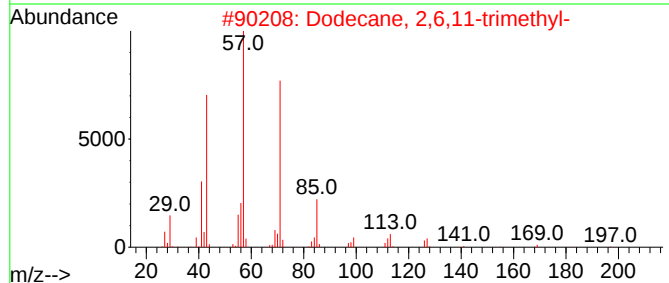
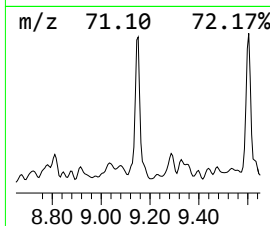
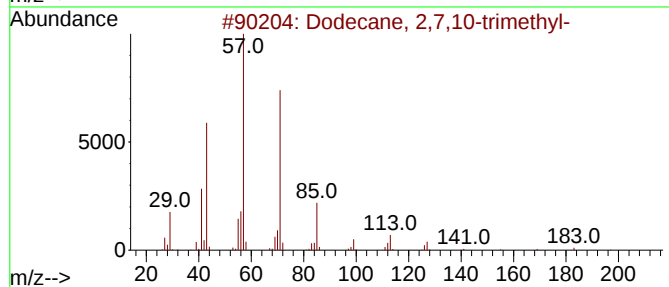
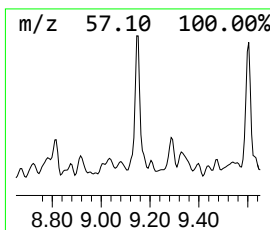
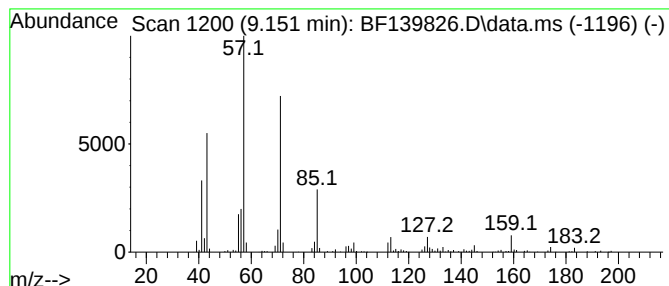
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TIC Integration Parameters: LSCINT.P

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Peak Number 11 Dodecane, 2,7,10-trimethyl- Concentration Rank 5

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 9.151 | 8.30 ng | 274351 | Acenaphthene-d10 | 9.886 |

| Hit# | of 5 | Tentative ID                | MW  | MolForm | CAS#        | Qual |
|------|------|-----------------------------|-----|---------|-------------|------|
| 1    |      | Dodecane, 2,7,10-trimethyl- | 212 | C15H32  | 074645-98-0 | 91   |
| 2    |      | Dodecane, 2,6,11-trimethyl- | 212 | C15H32  | 031295-56-4 | 90   |
| 3    |      | Dodecane, 2,6,10-trimethyl- | 212 | C15H32  | 003891-98-3 | 64   |
| 4    |      | Undecane, 3,5-dimethyl-     | 184 | C13H28  | 017312-81-1 | 64   |
| 5    |      | Nonane, 3,7-dimethyl-       | 156 | C11H24  | 017302-32-8 | 64   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101624\  
Data File : BF139826.D  
Acq On : 16 Oct 2024 12:42  
Operator : RC/JU  
Sample : P4395-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
F05308-SOLID

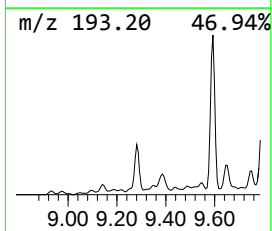
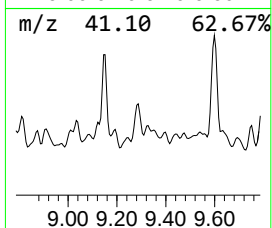
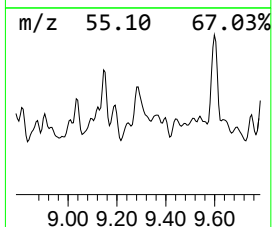
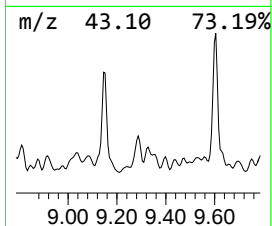
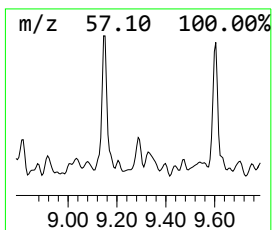
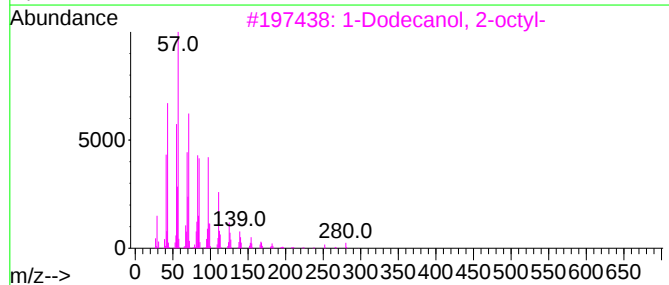
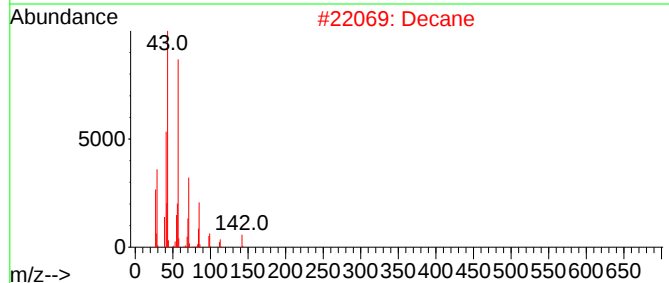
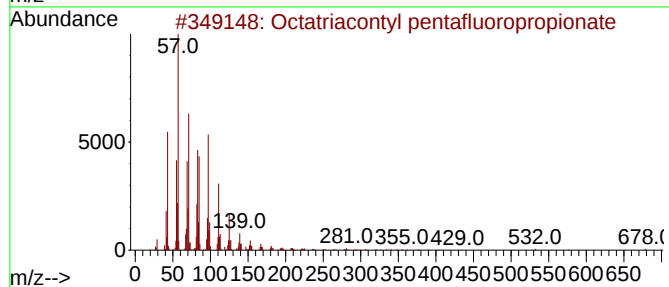
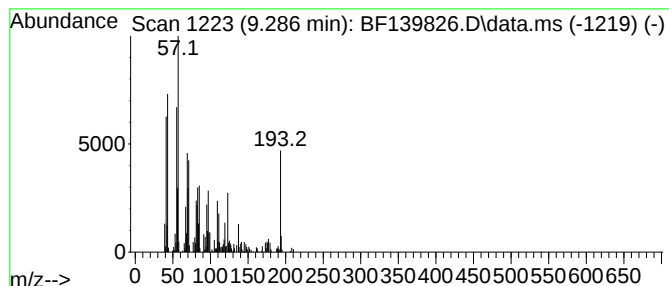
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 12 unknown9.286 Concentration Rank 14

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 9.286 | 5.93 ng | 196170 | Acenaphthene-d10 | 9.886 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | Octatriacontyl pentafluoropropio... | 697 | C41H77F5O2 | 1000351-89-1 | 38   |
| 2    |    |   | Decane                              | 142 | C10H22     | 000124-18-5  | 35   |
| 3    |    |   | 1-Dodecanol, 2-octyl-               | 298 | C20H42O    | 005333-42-6  | 30   |
| 4    |    |   | Tetratriacontyl heptafluorobutyrate | 691 | C38H69F7O2 | 1000351-84-1 | 30   |
| 5    |    |   | Hexatriacontyl pentafluoropropio... | 669 | C39H73F5O2 | 1000351-89-0 | 30   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101624\  
Data File : BF139826.D  
Acq On : 16 Oct 2024 12:42  
Operator : RC/JU  
Sample : P4395-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
F05308-SOLID

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF101524.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

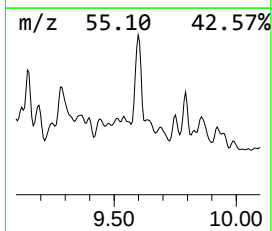
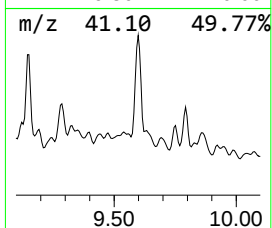
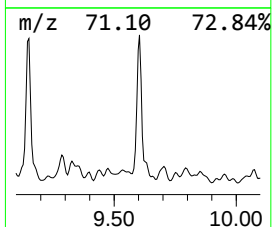
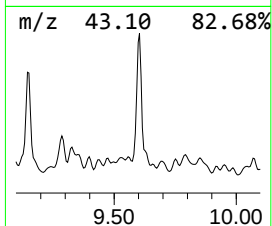
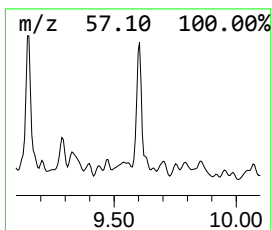
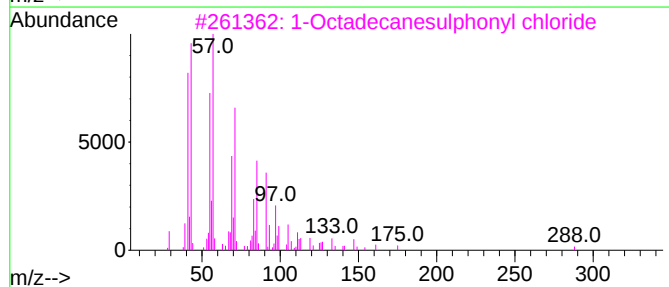
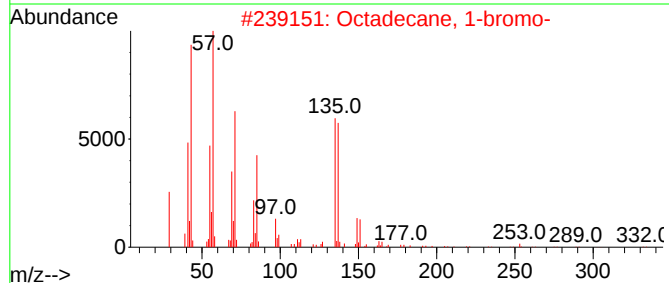
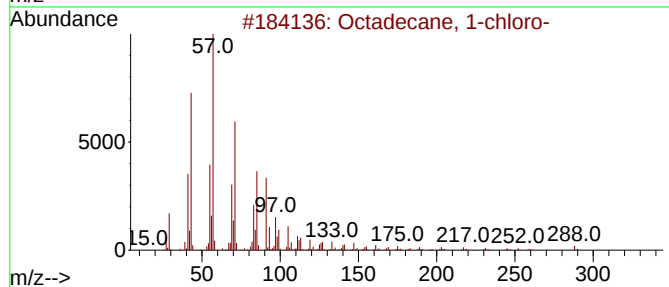
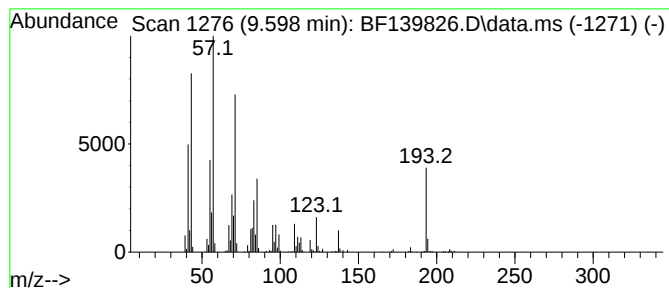
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TIC Integration Parameters: LSCINT.P

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Peak Number 13 unknown9.598 Concentration Rank 1

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 9.598 | 14.25 ng | 471227 | Acenaphthene-d10 | 9.886 |

| Hit# | of 5 | Tentative ID                   | MW  | MolForm     | CAS#         | Qual |
|------|------|--------------------------------|-----|-------------|--------------|------|
| 1    |      | Octadecane, 1-chloro-          | 288 | C18H37Cl    | 003386-33-2  | 43   |
| 2    |      | Octadecane, 1-bromo-           | 332 | C18H37Br    | 000112-89-0  | 43   |
| 3    |      | 1-Octadecanesulphonyl chloride | 352 | C18H37ClO2S | 1000342-70-4 | 38   |
| 4    |      | Tetradecane, 1-chloro-         | 232 | C14H29Cl    | 002425-54-9  | 38   |
| 5    |      | Dotriacontane                  | 451 | C32H66      | 000544-85-4  | 35   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101624\  
Data File : BF139826.D  
Acq On : 16 Oct 2024 12:42  
Operator : RC/JU  
Sample : P4395-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
F05308-SOLID

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF101524.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

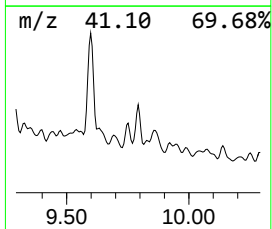
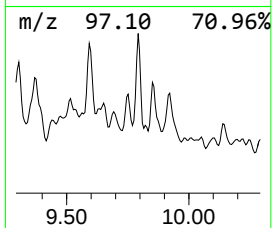
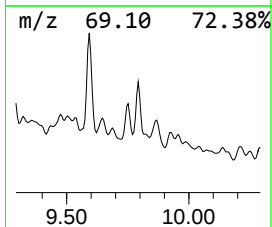
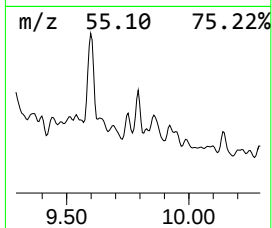
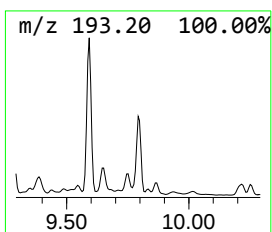
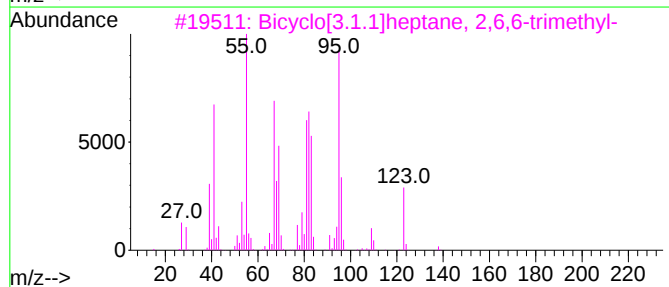
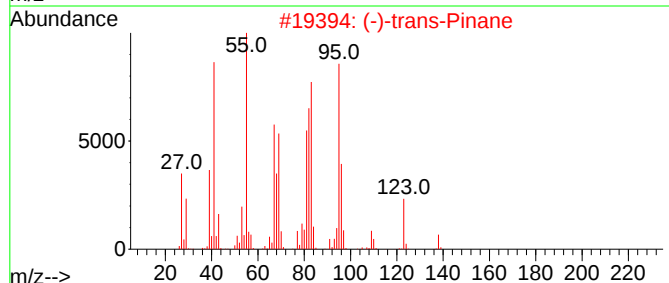
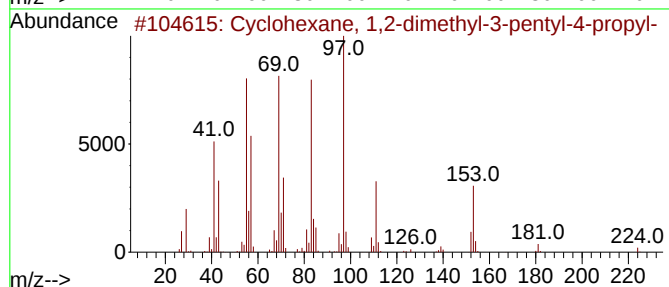
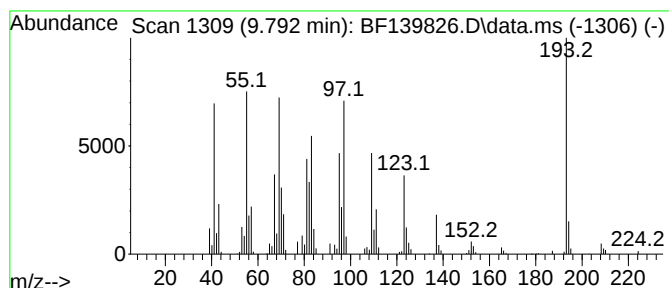
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TIC Integration Parameters: LSCINT.P

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Peak Number 14 unknown9.792 Concentration Rank 20

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 9.792 | 5.06 ng | 167194 | Acenaphthene-d10 | 9.886 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Cyclohexane, 1,2-dimethyl-3-pent... | 224 | C16H32  | 062376-17-4 | 38   |
| 2    |    |   | (-)-trans-Pinane                    | 138 | C10H18  | 033626-25-4 | 38   |
| 3    |    |   | Bicyclo[3.1.1]heptane, 2,6,6-tri... | 138 | C10H18  | 000473-55-2 | 30   |
| 4    |    |   | Cyclododecanone, 2-methylene-       | 194 | C13H22O | 003045-76-9 | 25   |
| 5    |    |   | [1,1'-Biphenyl]-4-acetonitrile      | 193 | C14H11N | 031603-77-7 | 22   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101624\  
Data File : BF139826.D  
Acq On : 16 Oct 2024 12:42  
Operator : RC/JU  
Sample : P4395-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
F05308-SOLID

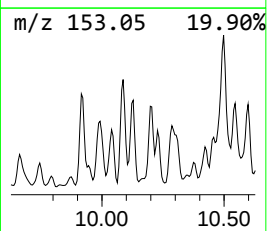
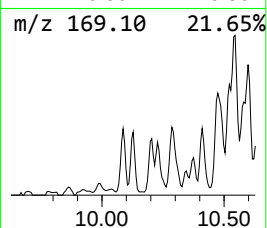
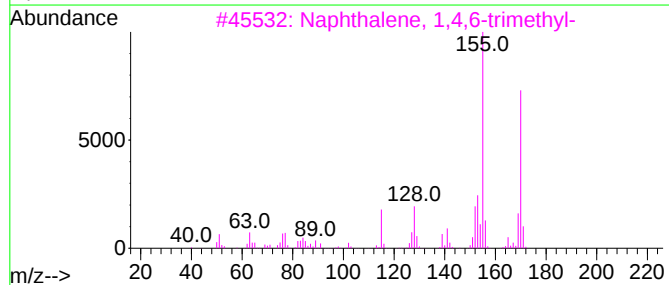
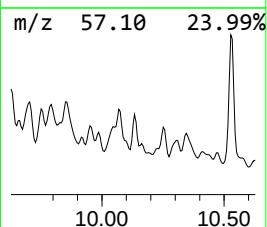
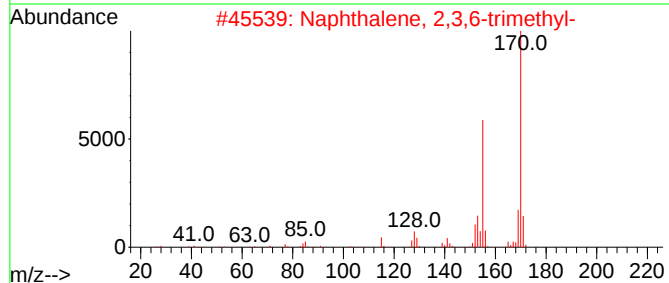
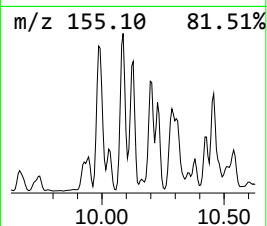
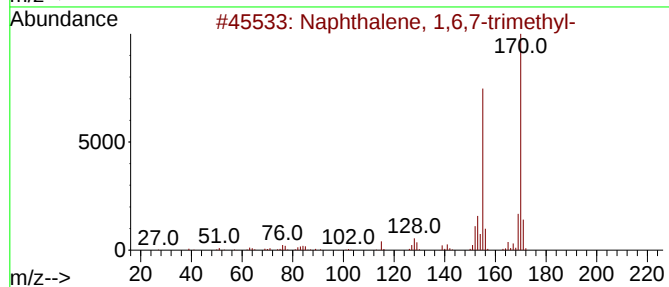
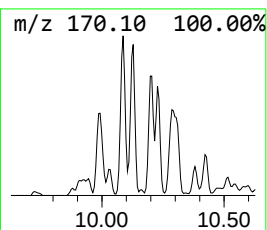
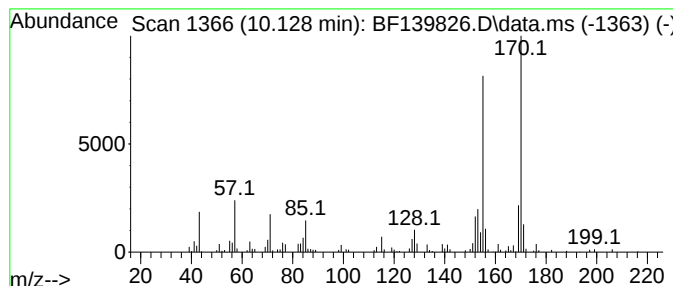
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 15 Naphthalene, 1,6,7-trimethyl- Concentration Rank 11

| R.T.      | EstConc                       | Area   | Relative to ISTD |         |             | R.T.  |
|-----------|-------------------------------|--------|------------------|---------|-------------|-------|
| 10.128    | 6.55 ng                       | 216560 | Acenaphthene-d10 |         |             | 9.886 |
| Hit# of 5 | Tentative ID                  |        | MW               | MolForm | CAS#        | Qual  |
| 1         | Naphthalene, 1,6,7-trimethyl- |        | 170              | C13H14  | 002245-38-7 | 97    |
| 2         | Naphthalene, 2,3,6-trimethyl- |        | 170              | C13H14  | 000829-26-5 | 94    |
| 3         | Naphthalene, 1,4,6-trimethyl- |        | 170              | C13H14  | 002131-42-2 | 94    |
| 4         | Naphthalene, 1,4,5-trimethyl- |        | 170              | C13H14  | 002131-41-1 | 91    |
| 5         | 4,6,8-Trimethylazulene        |        | 170              | C13H14  | 000941-81-1 | 87    |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101624\  
Data File : BF139826.D  
Acq On : 16 Oct 2024 12:42  
Operator : RC/JU  
Sample : P4395-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

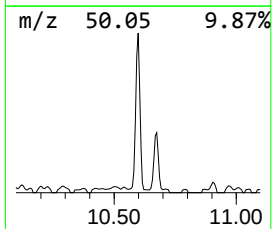
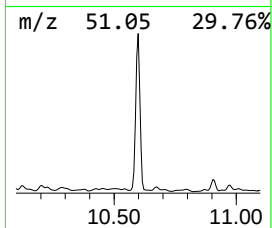
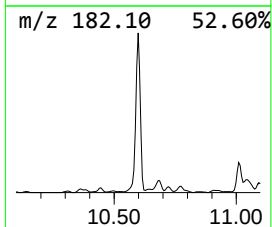
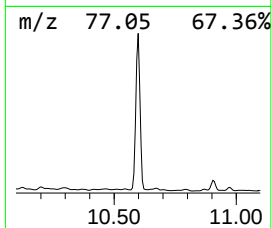
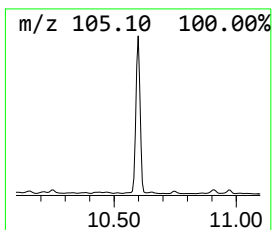
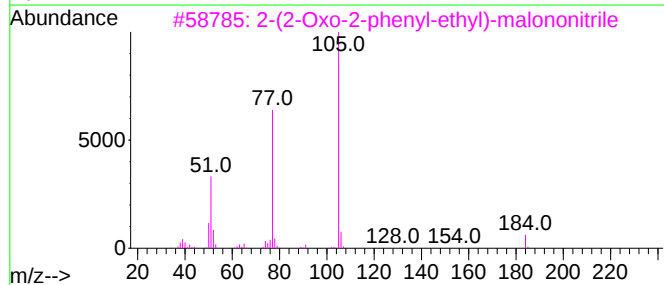
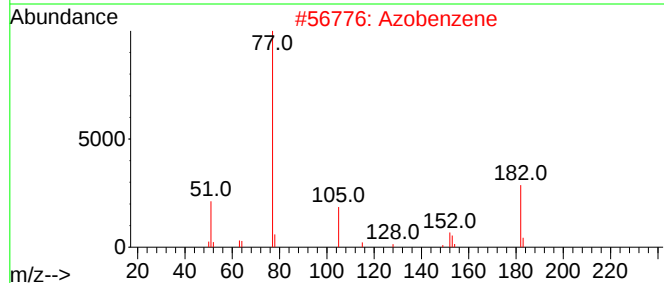
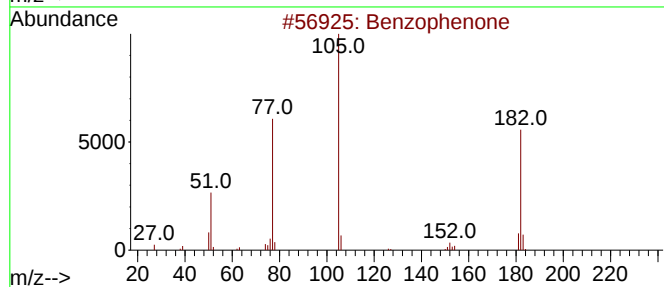
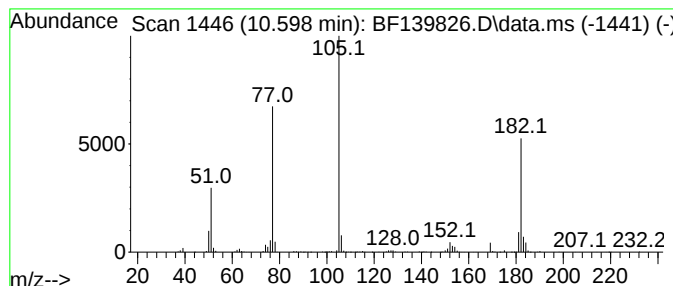
Instrument :  
BNA\_F  
ClientSampleId :  
F05308-SOLID

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF101524.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 16 Benzophenone Concentration Rank 2

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 10.598    | 12.49 ng                            | 412832 | Acenaphthene-d10 | 9.886        |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | Benzophenone                        | 182    | C13H10O          | 000119-61-9  | 96   |
| 2         | Azobenzene                          | 182    | C12H10N2         | 000103-33-3  | 59   |
| 3         | 2-(2-Oxo-2-phenyl-ethyl)-malonon... | 184    | C11H8N2O         | 1000296-76-9 | 55   |
| 4         | Disulfide, dibenzoyl                | 274    | C14H10O2S2       | 000644-32-6  | 46   |
| 5         | Benzoylformic acid                  | 150    | C8H6O3           | 000611-73-4  | 46   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101624\  
Data File : BF139826.D  
Acq On : 16 Oct 2024 12:42  
Operator : RC/JU  
Sample : P4395-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
F05308-SOLID

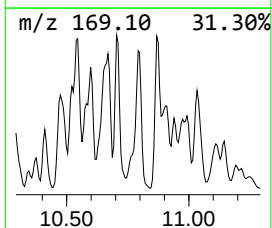
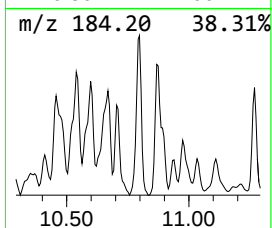
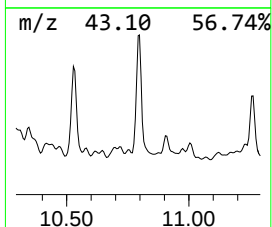
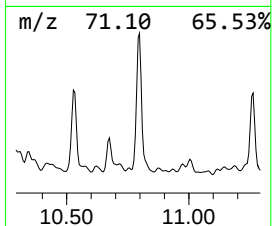
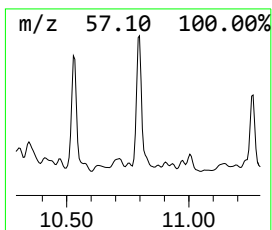
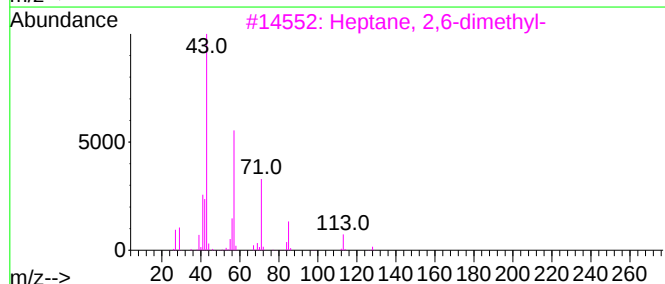
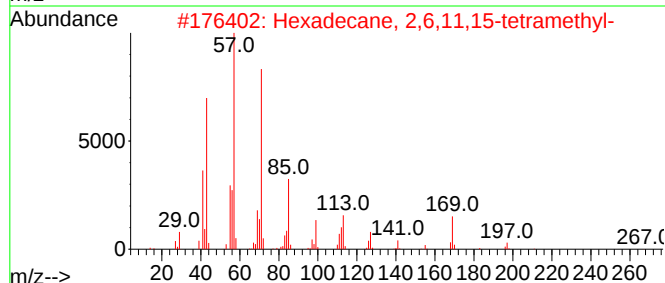
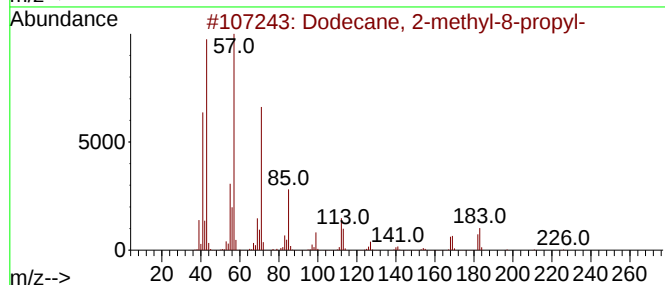
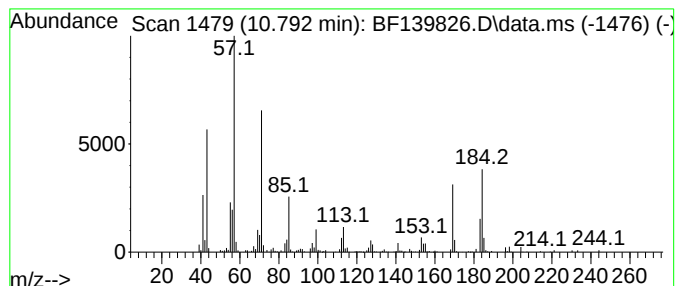
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 17 Dodecane, 2-methyl-8-propyl- Concentration Rank 12

| R.T.      | EstConc                            | Area   | Relative to ISTD | R.T.        |      |
|-----------|------------------------------------|--------|------------------|-------------|------|
| 10.792    | 6.50 ng                            | 207027 | Phenanthrene-d10 | 11.369      |      |
| Hit# of 5 | Tentative ID                       | MW     | MolForm          | CAS#        | Qual |
| 1         | Dodecane, 2-methyl-8-propyl-       | 226    | C16H34           | 055045-07-3 | 53   |
| 2         | Hexadecane, 2,6,11,15-tetramethyl- | 282    | C20H42           | 000504-44-9 | 53   |
| 3         | Heptane, 2,6-dimethyl-             | 128    | C9H20            | 001072-05-5 | 46   |
| 4         | Octadecane, 2,6-dimethyl-          | 282    | C20H42           | 075163-97-2 | 46   |
| 5         | Tridecane, 7-methyl-               | 198    | C14H30           | 026730-14-3 | 45   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101624\  
Data File : BF139826.D  
Acq On : 16 Oct 2024 12:42  
Operator : RC/JU  
Sample : P4395-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
F05308-SOLID

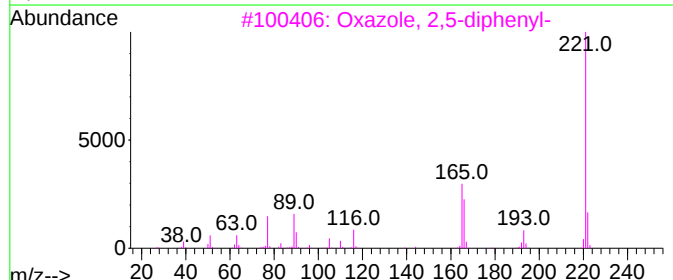
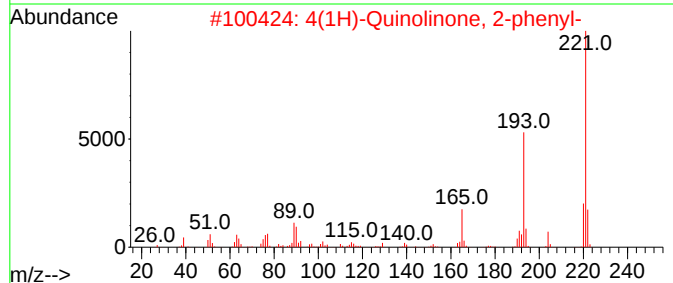
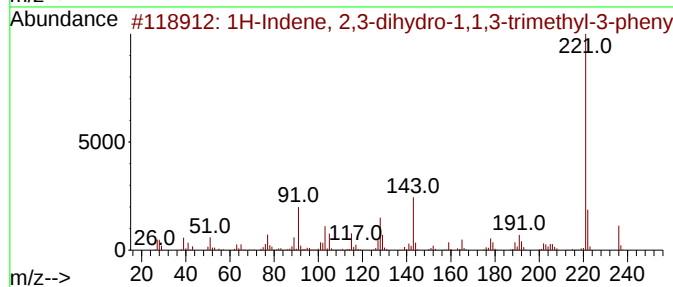
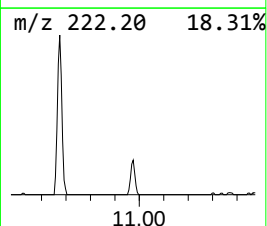
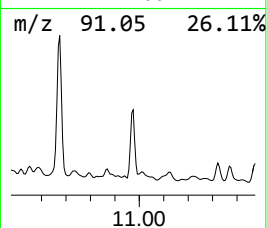
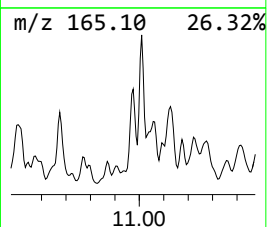
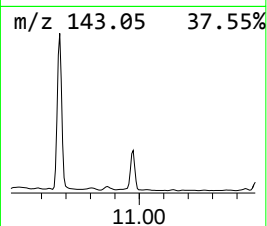
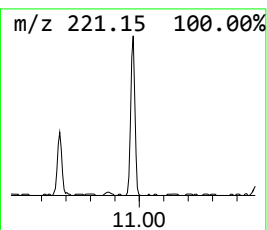
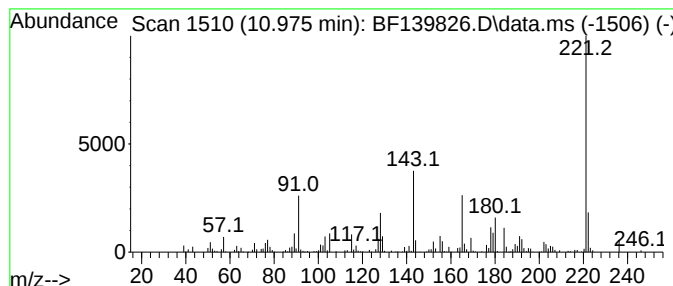
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 18 1H-Indene, 2,3-dihydro-1,1,... Concentration Rank 17

| R.T.      | EstConc                             | Area   | Relative to ISTD |          | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|----------|-------------|------|
| 10.975    | 5.67 ng                             | 180583 | Phenanthrene-d10 |          | 11.369      |      |
| Hit# of 5 | Tentative ID                        |        | MW               | MolForm  | CAS#        | Qual |
| 1         | 1H-Indene, 2,3-dihydro-1,1,3-tri... |        | 236              | C18H20   | 003910-35-8 | 89   |
| 2         | 4(1H)-Quinolinone, 2-phenyl-        |        | 221              | C15H11NO | 014802-18-7 | 53   |
| 3         | Oxazole, 2,5-diphenyl-              |        | 221              | C15H11NO | 000092-71-7 | 52   |
| 4         | 1H-Quinolin-2-one, 3-phenyl-        |        | 221              | C15H11NO | 038035-81-3 | 49   |
| 5         | 4-Quinolinol, 2-phenyl-             |        | 221              | C15H11NO | 001144-20-3 | 49   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101624\  
Data File : BF139826.D  
Acq On : 16 Oct 2024 12:42  
Operator : RC/JU  
Sample : P4395-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
F05308-SOLID

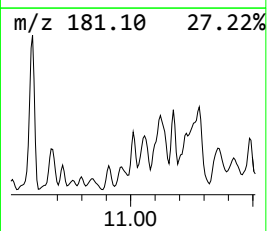
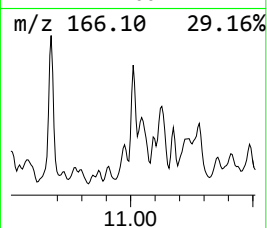
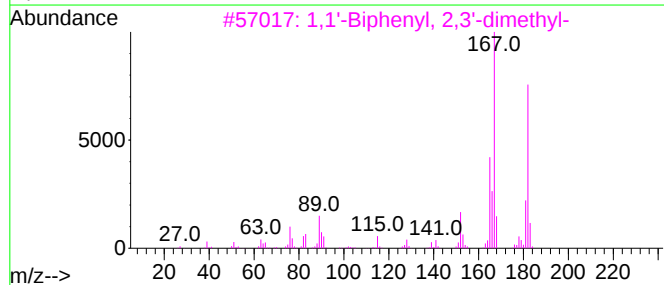
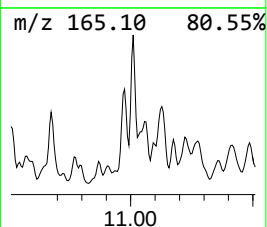
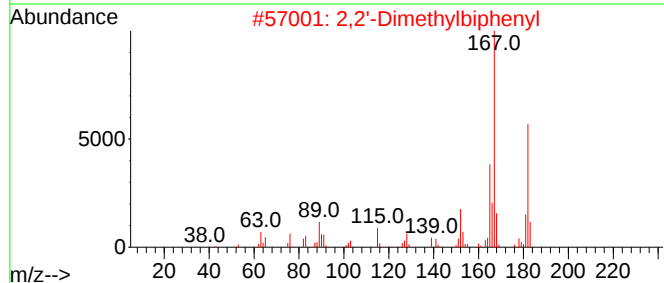
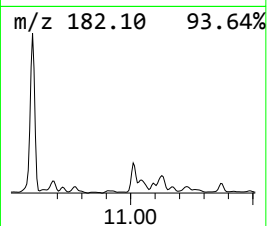
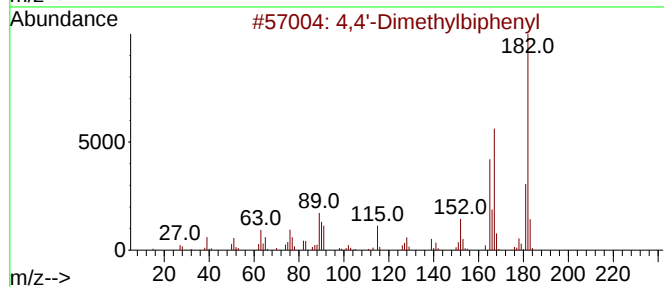
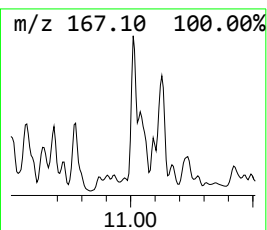
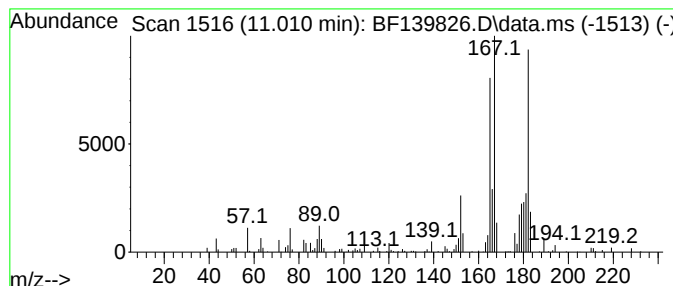
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 19 4,4'-Dimethylbiphenyl Concentration Rank 16

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.        |      |
|-----------|-------------------------------------|--------|------------------|-------------|------|
| 11.010    | 5.74 ng                             | 182655 | Phenanthrene-d10 | 11.369      |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#        | Qual |
| 1         | 4,4'-Dimethylbiphenyl               | 182    | C14H14           | 000613-33-2 | 91   |
| 2         | 2,2'-Dimethylbiphenyl               | 182    | C14H14           | 000605-39-0 | 83   |
| 3         | 1,1'-Biphenyl, 2,3'-dimethyl-       | 182    | C14H14           | 000611-43-8 | 76   |
| 4         | 1,1'-Biphenyl, 3,4'-dimethyl-       | 182    | C14H14           | 007383-90-6 | 72   |
| 5         | 1H,2H,3H-Cyclopenta[a]naphthalen... | 182    | C14H14           | 001624-22-2 | 68   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101624\  
Data File : BF139826.D  
Acq On : 16 Oct 2024 12:42  
Operator : RC/JU  
Sample : P4395-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
F05308-SOLID

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF101524.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

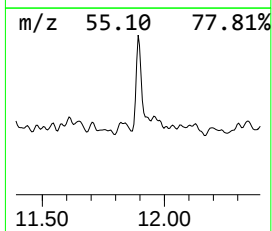
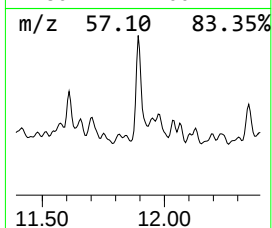
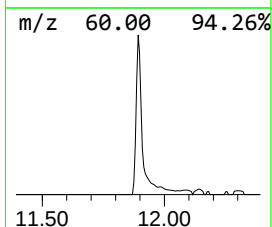
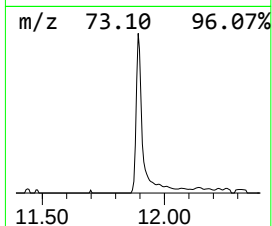
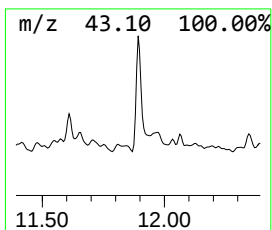
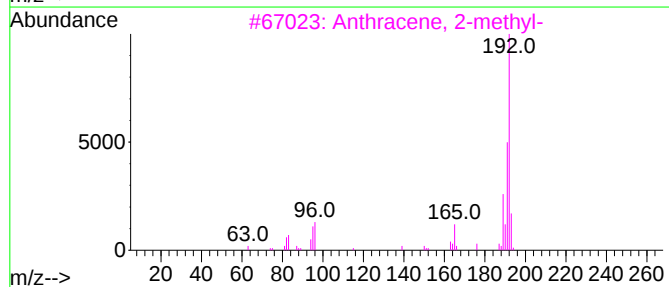
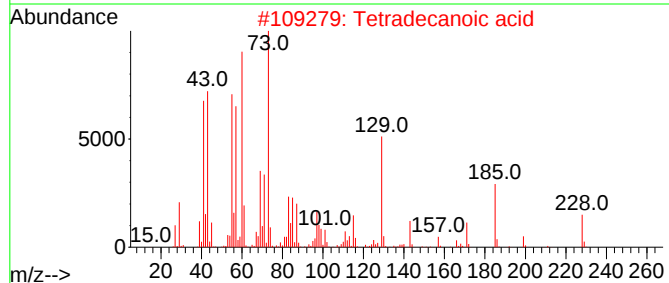
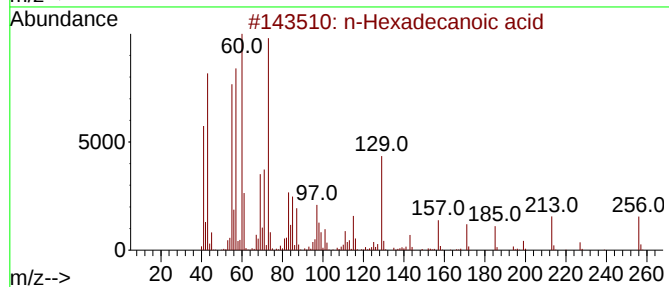
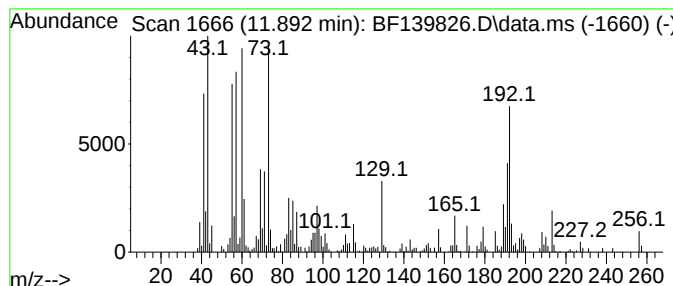
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 20 n-Hexadecanoic acid Concentration Rank 9

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 11.892 | 6.86 ng | 218410 | Phenanthrene-d10 | 11.369 |

| Hit# | of 5 | Tentative ID          | MW  | MolForm  | CAS#        | Qual |
|------|------|-----------------------|-----|----------|-------------|------|
| 1    |      | n-Hexadecanoic acid   | 256 | C16H32O2 | 000057-10-3 | 97   |
| 2    |      | Tetradecanoic acid    | 228 | C14H28O2 | 000544-63-8 | 89   |
| 3    |      | Anthracene, 2-methyl- | 192 | C15H12   | 000613-12-7 | 78   |
| 4    |      | Tridecanoic acid      | 214 | C13H26O2 | 000638-53-9 | 64   |
| 5    |      | 1H-Indene, 2-phenyl-  | 192 | C15H12   | 004505-48-0 | 56   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101624\  
Data File : BF139826.D  
Acq On : 16 Oct 2024 12:42  
Operator : RC/JU  
Sample : P4395-01  
Misc :  
ALS Vial : 6 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
F05308-SOLID

Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF101524.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

| TIC Top Hit name   | RT     | EstConc | Units | Response | --Internal Standard-- |        |        |      |
|--------------------|--------|---------|-------|----------|-----------------------|--------|--------|------|
|                    |        |         |       |          | #                     | RT     | Resp   | Conc |
| Nonane, 3-methyl-  | 6.098  | 7.9     | ng    | 253836   | 1                     | 6.851  | 645120 | 20.0 |
| Cyclohexane, 1,... | 6.381  | 12.3    | ng    | 396525   | 1                     | 6.851  | 645120 | 20.0 |
| Cyclopentane, 1... | 6.598  | 6.2     | ng    | 199151   | 1                     | 6.851  | 645120 | 20.0 |
| unknown6.628       | 6.628  | 5.1     | ng    | 163127   | 1                     | 6.851  | 645120 | 20.0 |
| Decane, 2,4-dim... | 6.769  | 5.3     | ng    | 171262   | 1                     | 6.851  | 645120 | 20.0 |
| unknown6.992       | 6.992  | 7.0     | ng    | 225479   | 1                     | 6.851  | 645120 | 20.0 |
| 1-Ethyl-2,2,6-t... | 7.122  | 6.7     | ng    | 216658   | 1                     | 6.851  | 645120 | 20.0 |
| unknown7.487       | 7.487  | 7.6     | ng    | 244813   | 1                     | 6.851  | 645120 | 20.0 |
| Bicyclo[2.2.1]h... | 7.657  | 5.8     | ng    | 249992   | 2                     | 8.134  | 856748 | 20.0 |
| Octane, 2,6-dim... | 8.551  | 9.9     | ng    | 425094   | 2                     | 8.134  | 856748 | 20.0 |
| Dodecane, 2,7,1... | 9.151  | 8.3     | ng    | 274351   | 3                     | 9.886  | 661263 | 20.0 |
| unknown9.286       | 9.286  | 5.9     | ng    | 196170   | 3                     | 9.886  | 661263 | 20.0 |
| unknown9.598       | 9.598  | 14.3    | ng    | 471227   | 3                     | 9.886  | 661263 | 20.0 |
| unknown9.792       | 9.792  | 5.1     | ng    | 167194   | 3                     | 9.886  | 661263 | 20.0 |
| Naphthalene, 1,... | 10.128 | 6.5     | ng    | 216560   | 3                     | 9.886  | 661263 | 20.0 |
| Benzophenone       | 10.598 | 12.5    | ng    | 412832   | 3                     | 9.886  | 661263 | 20.0 |
| Dodecane, 2-met... | 10.792 | 6.5     | ng    | 207027   | 4                     | 11.369 | 636638 | 20.0 |
| 1H-Indene, 2,3-... | 10.975 | 5.7     | ng    | 180583   | 4                     | 11.369 | 636638 | 20.0 |
| 4,4'-Dimethylbi... | 11.010 | 5.7     | ng    | 182655   | 4                     | 11.369 | 636638 | 20.0 |
| n-Hexadecanoic ... | 11.892 | 6.9     | ng    | 218410   | 4                     | 11.369 | 636638 | 20.0 |