

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101321\  
 Data File : BF125825.D  
 Acq On : 13 Oct 2021 14:14  
 Operator : JU/CG  
 Sample : M4127-05  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 CHRT-20426

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-BF100721.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF125825.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       | 305           | 310         | 316          | rVB      | 103606         | 139620        | 3.06%           | 0.230%        |
| 2         | 5.457       | 566           | 573         | 576          | rBV      | 3710129        | 3720022       | 81.53%          | 6.123%        |
| 3         | 6.469       | 740           | 745         | 748          | rBV      | 3596660        | 3662377       | 80.27%          | 6.028%        |
| 4         | 6.845       | 805           | 809         | 812          | rBV      | 1396899        | 1168945       | 25.62%          | 1.924%        |
| 5         | 7.022       | 836           | 839         | 843          | rVB      | 126275         | 105609        | 2.31%           | 0.174%        |
| 6         | 7.404       | 894           | 904         | 907          | rBV      | 2652950        | 2335755       | 51.19%          | 3.844%        |
| 7         | 7.486       | 914           | 918         | 922          | rVB      | 104742         | 117995        | 2.59%           | 0.194%        |
| 8         | 8.028       | 1006          | 1010        | 1015         | rBV      | 786424         | 685909        | 15.03%          | 1.129%        |
| 9         | 8.122       | 1022          | 1026        | 1033         | rBV      | 1843767        | 1739570       | 38.13%          | 2.863%        |
| 10        | 8.422       | 1069          | 1077        | 1080         | rBV7     | 103574         | 172526        | 3.78%           | 0.284%        |
| 11        | 8.510       | 1088          | 1092        | 1096         | rBV5     | 52853          | 75291         | 1.65%           | 0.124%        |
| 12        | 8.592       | 1104          | 1106        | 1109         | rVB3     | 140024         | 114682        | 2.51%           | 0.189%        |
| 13        | 8.675       | 1115          | 1120        | 1124         | rBV6     | 77028          | 152676        | 3.35%           | 0.251%        |
| 14        | 8.757       | 1129          | 1134        | 1135         | rBV5     | 74920          | 121556        | 2.66%           | 0.200%        |
| 15        | 8.816       | 1140          | 1144        | 1146         | rVV4     | 117441         | 123178        | 2.70%           | 0.203%        |
| 16        | 8.875       | 1151          | 1154        | 1157         | rBV4     | 168050         | 224716        | 4.93%           | 0.370%        |
| 17        | 8.969       | 1168          | 1170        | 1174         | rVB3     | 70664          | 72512         | 1.59%           | 0.119%        |
| 18        | 9.004       | 1174          | 1176        | 1177         | rBV2     | 105240         | 92492         | 2.03%           | 0.152%        |
| 19        | 9.086       | 1184          | 1190        | 1193         | rBV6     | 171982         | 304135        | 6.67%           | 0.501%        |
| 20        | 9.122       | 1193          | 1196        | 1198         | rVV2     | 335141         | 263031        | 5.76%           | 0.433%        |
| 21        | 9.151       | 1198          | 1201        | 1204         | rBV5     | 273755         | 289061        | 6.34%           | 0.476%        |
| 22        | 9.198       | 1205          | 1209        | 1212         | rBV      | 4801044        | 4199384       | 92.04%          | 6.912%        |
| 23        | 9.239       | 1214          | 1216        | 1218         | rBV2     | 85920          | 77042         | 1.69%           | 0.127%        |
| 24        | 9.280       | 1220          | 1223        | 1228         | rBV4     | 539781         | 565364        | 12.39%          | 0.931%        |
| 25        | 9.327       | 1228          | 1231        | 1233         | rBV4     | 172822         | 200612        | 4.40%           | 0.330%        |
| 26        | 9.351       | 1233          | 1235        | 1237         | rVB2     | 190169         | 141270        | 3.10%           | 0.233%        |
| 27        | 9.475       | 1254          | 1256        | 1259         | rBV4     | 177142         | 153943        | 3.37%           | 0.253%        |
| 28        | 9.510       | 1260          | 1262        | 1264         | rBV3     | 249340         | 258434        | 5.66%           | 0.425%        |
| 29        | 9.539       | 1264          | 1267        | 1269         | rVV3     | 311806         | 233239        | 5.11%           | 0.384%        |
| 30        | 9.563       | 1269          | 1271        | 1273         | rVB2     | 391513         | 320605        | 7.03%           | 0.528%        |
| 31        | 9.592       | 1273          | 1276        | 1280         | rBV2     | 2862490        | 3076227       | 67.42%          | 5.063%        |
| 32        | 9.751       | 1299          | 1303        | 1305         | rBV3     | 959059         | 1057162       | 23.17%          | 1.740%        |
| 33        | 9.798       | 1308          | 1311        | 1313         | rVB      | 1938184        | 1744635       | 38.24%          | 2.871%        |
| 34        | 9.833       | 1315          | 1317        | 1319         | rVV2     | 287424         | 194009        | 4.25%           | 0.319%        |
| 35        | 9.880       | 1319          | 1325        | 1328         | rBV5     | 2594172        | 3585578       | 78.58%          | 5.901%        |
| 36        | 9.922       | 1328          | 1332        | 1335         | rVV5     | 470361         | 766154        | 16.79%          | 1.261%        |

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 ClientSampleId :  
 CHRT-20426

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-BF100721.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

|    |        |      |      |      |      |         |         |         |        |
|----|--------|------|------|------|------|---------|---------|---------|--------|
| 37 | 9.992  | 1341 | 1344 | 1347 | rBV4 | 458528  | 499514  | 10.95%  | 0.822% |
| 38 | 10.022 | 1347 | 1349 | 1351 | rBV  | 565103  | 557798  | 12.23%  | 0.918% |
| 39 | 10.080 | 1357 | 1359 | 1361 | rVB2 | 567731  | 414577  | 9.09%   | 0.682% |
| 40 | 10.145 | 1367 | 1370 | 1372 | rBV2 | 1159114 | 1154488 | 25.30%  | 1.900% |
| 41 | 10.198 | 1376 | 1379 | 1381 | rBV4 | 828975  | 1017115 | 22.29%  | 1.674% |
| 42 | 10.222 | 1381 | 1383 | 1385 | rVB3 | 652637  | 465976  | 10.21%  | 0.767% |
| 43 | 10.257 | 1386 | 1389 | 1391 | rBV2 | 772361  | 816319  | 17.89%  | 1.344% |
| 44 | 10.292 | 1391 | 1395 | 1399 | rBV2 | 1883175 | 2106120 | 46.16%  | 3.466% |
| 45 | 10.345 | 1402 | 1404 | 1406 | rBV2 | 525118  | 511600  | 11.21%  | 0.842% |
| 46 | 10.422 | 1415 | 1417 | 1419 | rBV2 | 670220  | 750682  | 16.45%  | 1.236% |
| 47 | 10.539 | 1435 | 1437 | 1441 | rVB  | 1318797 | 1224571 | 26.84%  | 2.016% |
| 48 | 10.669 | 1455 | 1459 | 1462 | rBV3 | 1763251 | 2157370 | 47.28%  | 3.551% |
| 49 | 10.810 | 1478 | 1483 | 1486 | rBV2 | 3382119 | 4562744 | 100.00% | 7.510% |
| 50 | 11.274 | 1560 | 1562 | 1566 | rVB  | 1991738 | 1904199 | 41.73%  | 3.134% |
| 51 | 11.369 | 1576 | 1578 | 1580 | rBV  | 879904  | 810707  | 17.77%  | 1.334% |
| 52 | 12.580 | 1782 | 1784 | 1790 | rVB  | 1331694 | 1238447 | 27.14%  | 2.038% |
| 53 | 12.810 | 1821 | 1823 | 1830 | rVB  | 961572  | 872854  | 19.13%  | 1.437% |
| 54 | 12.951 | 1843 | 1847 | 1850 | rBV  | 2938650 | 2535052 | 55.56%  | 4.172% |
| 55 | 13.998 | 2020 | 2025 | 2028 | rBV2 | 1451711 | 1724173 | 37.79%  | 2.838% |
| 56 | 14.021 | 2028 | 2029 | 2035 | rVB  | 571358  | 482732  | 10.58%  | 0.795% |
| 57 | 14.762 | 2152 | 2155 | 2159 | rVB  | 174598  | 188998  | 4.14%   | 0.311% |
| 58 | 15.027 | 2196 | 2200 | 2210 | rVB  | 377564  | 703351  | 15.42%  | 1.158% |
| 59 | 15.327 | 2247 | 2251 | 2256 | rVB  | 206659  | 270488  | 5.93%   | 0.445% |
| 60 | 15.386 | 2258 | 2261 | 2265 | rVB  | 233342  | 232571  | 5.10%   | 0.383% |
| 61 | 15.451 | 2268 | 2272 | 2276 | rBV  | 849834  | 1005055 | 22.03%  | 1.654% |
| 62 | 16.892 | 2512 | 2517 | 2524 | rVB2 | 82087   | 168181  | 3.69%   | 0.277% |
| 63 | 17.327 | 2587 | 2591 | 2598 | rVB  | 71044   | 126293  | 2.77%   | 0.208% |

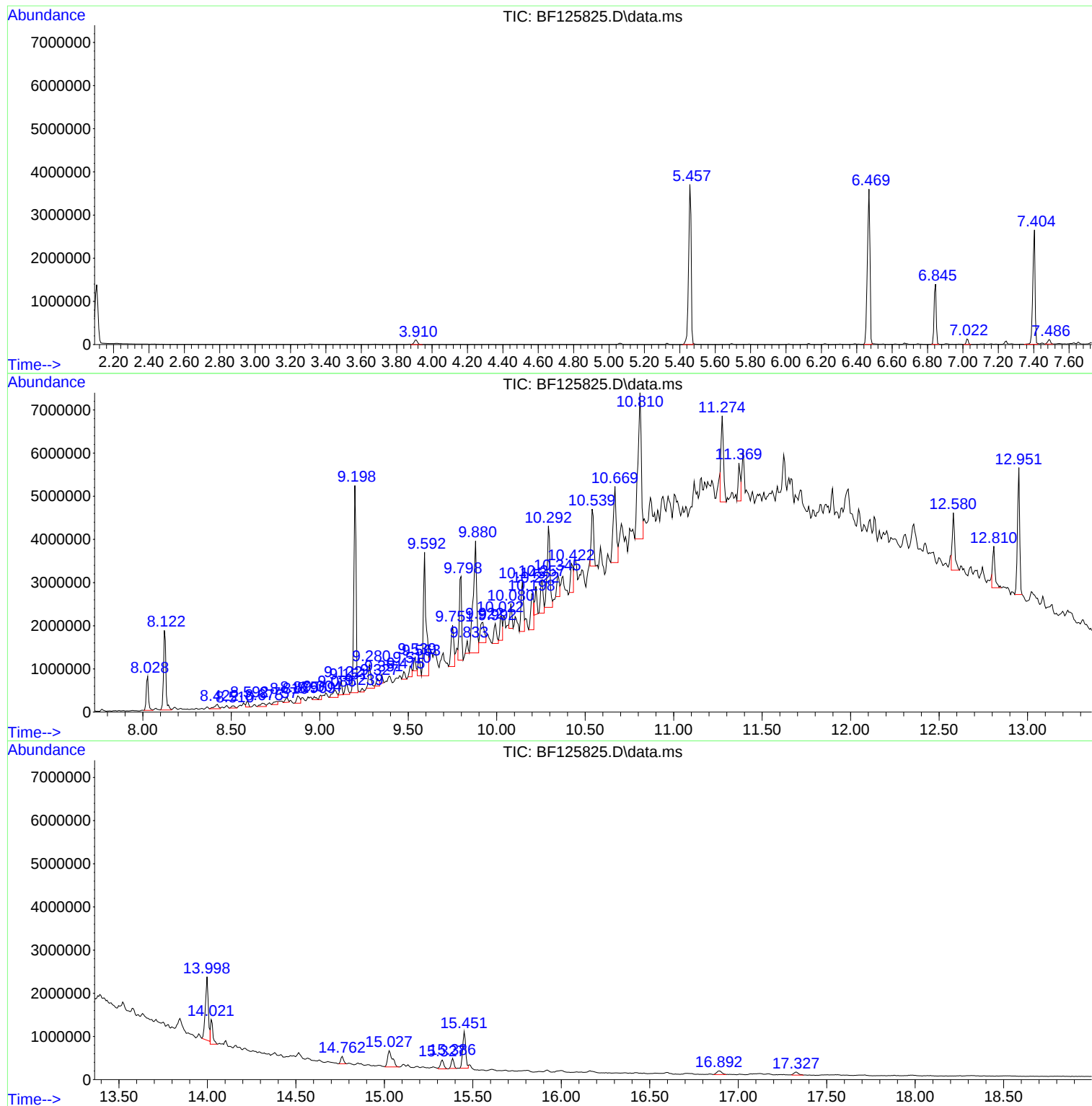
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Sample : M4127-05  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
CHRT-20426

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF100721.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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Misc :  
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Instrument :  
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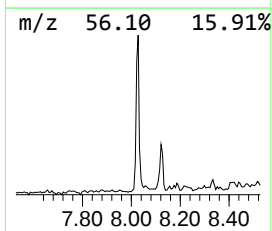
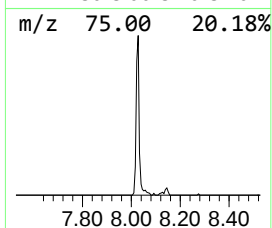
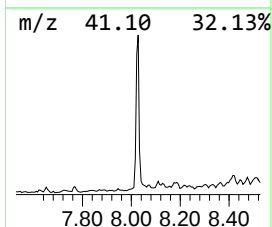
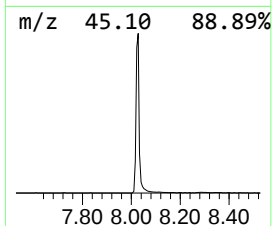
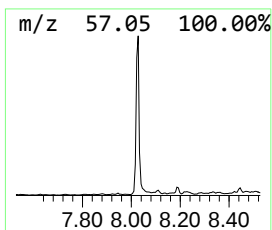
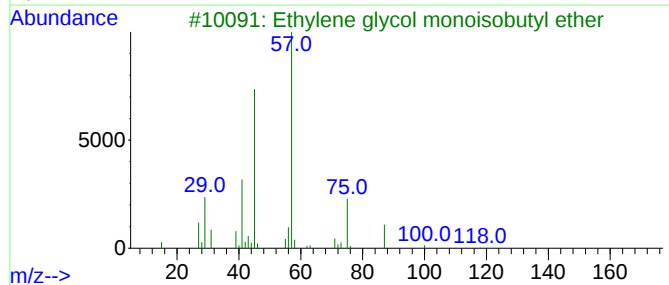
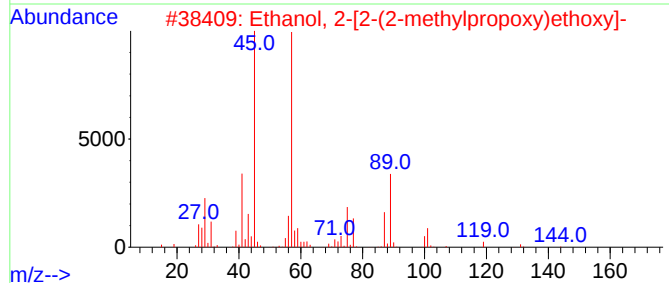
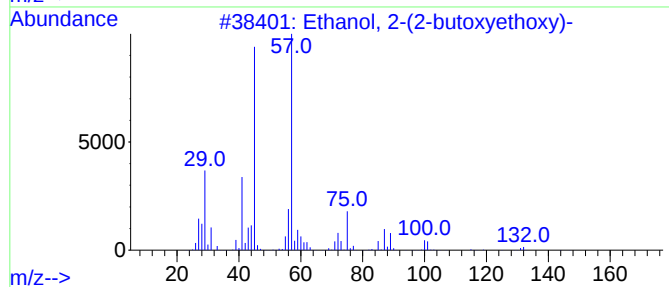
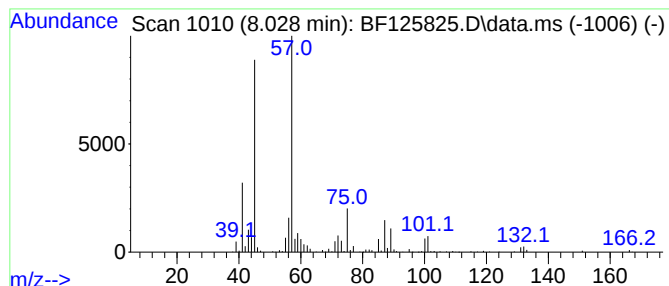
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 2 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 5

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 8.028 | 7.89 ng | 685909 | Naphthalene-d8   | 8.122 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|---------|-------------|------|
| 1    |    |   | Ethanol, 2-(2-butoxyethoxy)-        | 162 | C8H18O3 | 000112-34-5 | 90   |
| 2    |    |   | Ethanol, 2-[2-(2-methylpropoxy)e... | 162 | C8H18O3 | 018912-80-6 | 64   |
| 3    |    |   | Ethylene glycol monoisobutyl ether  | 118 | C6H14O2 | 004439-24-1 | 64   |
| 4    |    |   | Ethanol, 1-(2-butoxyethoxy)-        | 162 | C8H18O3 | 054446-78-5 | 58   |
| 5    |    |   | Methoxyacetic acid, 2-butyl ester   | 146 | C7H14O3 | 070159-90-9 | 50   |



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Instrument :  
BNA\_F  
ClientSampleId :  
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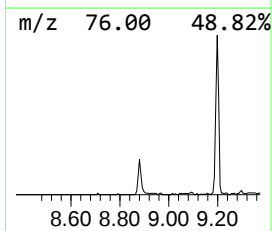
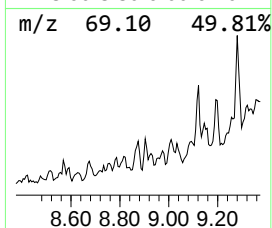
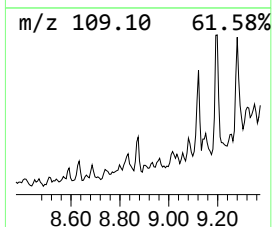
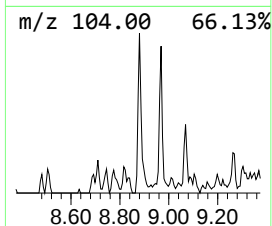
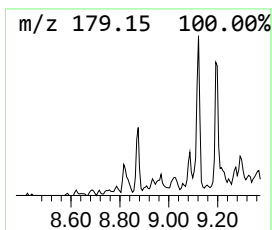
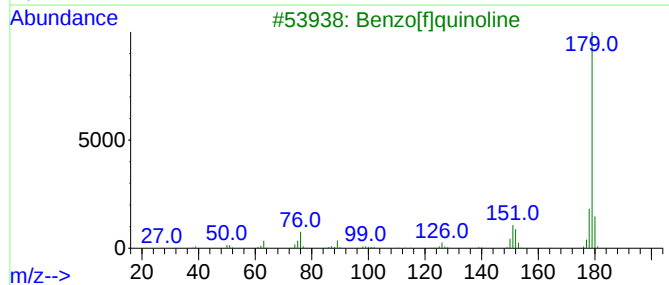
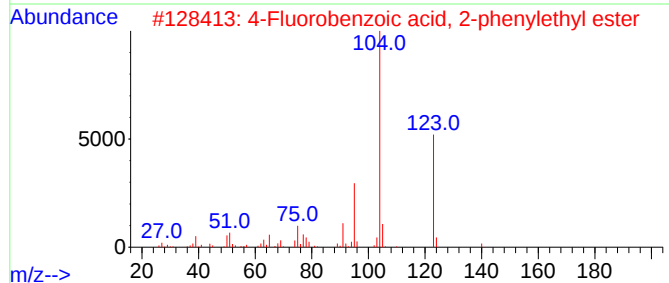
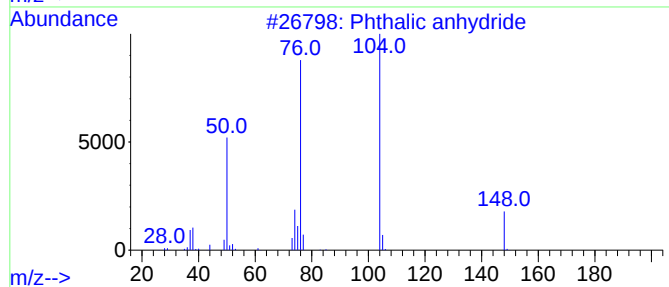
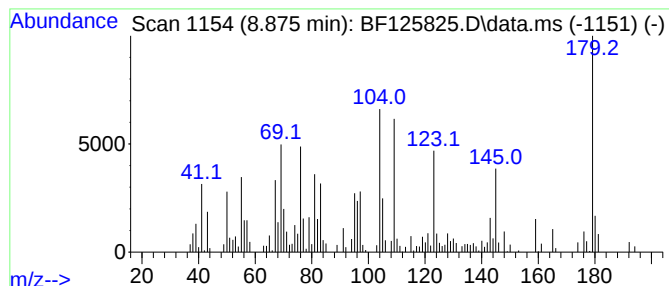
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF100721.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 3 unknown8.875 Concentration Rank 18

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 8.875 | 2.58 ng | 224716 | Naphthalene-d8   | 8.122 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|-----------|-------------|------|
| 1    |    |   | Phthalic anhydride                  | 148 | C8H4O3    | 000085-44-9 | 15   |
| 2    |    |   | 4-Fluorobenzoic acid, 2-phenylet... | 244 | C15H13F02 | 090172-19-3 | 14   |
| 3    |    |   | Benzo[f]quinoline                   | 179 | C13H9N    | 000085-02-9 | 11   |
| 4    |    |   | 3,5-Dimethyl-1-adamantanol          | 180 | C12H20O   | 000707-37-9 | 11   |
| 5    |    |   | Ninhydrin                           | 178 | C9H6O4    | 000485-47-2 | 11   |



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Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF100721.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

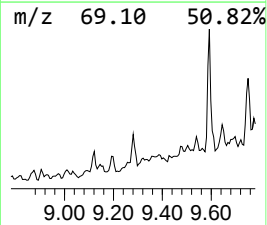
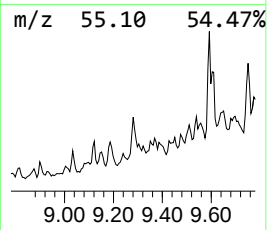
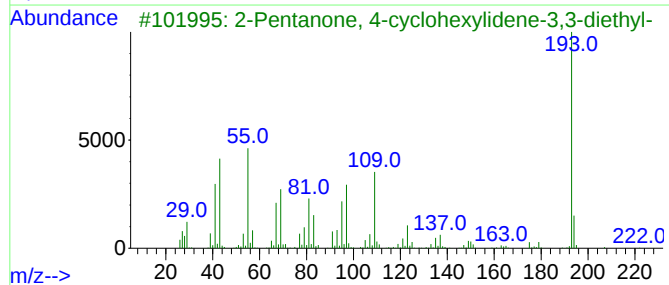
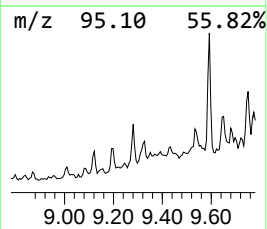
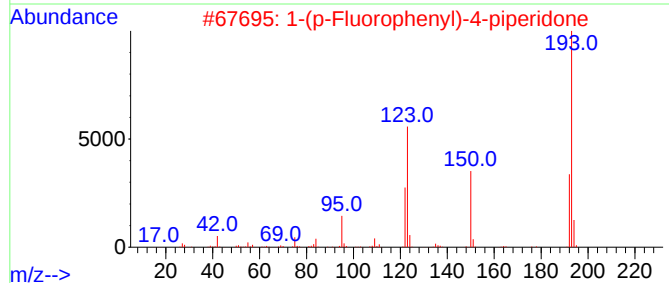
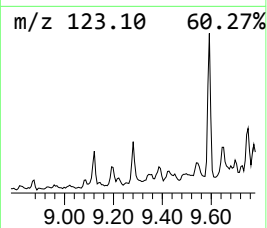
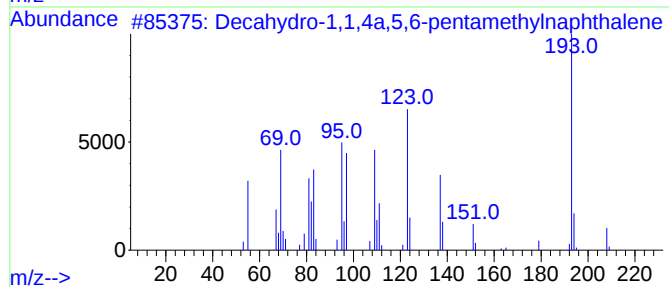
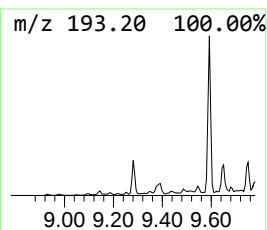
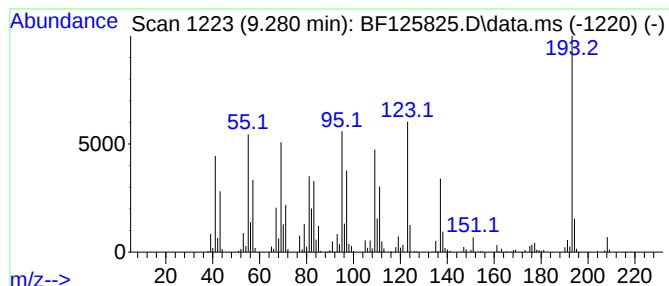
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

\*\*\*\*\*

Peak Number 4 unknown9.280 Concentration Rank 13

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 9.280 | 3.15 ng | 565364 | Acenaphthene-d10 | 9.880 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | Decahydro-1,1,4a,5,6-pentamethyl... | 208 | C15H28    | 080655-44-3  | 49   |
| 2    |    |   | 1-(p-Fluorophenyl)-4-piperidone     | 193 | C11H12FNO | 1000238-56-7 | 43   |
| 3    |    |   | 2-Pentanone, 4-cyclohexylidene-3... | 222 | C15H26O   | 313253-65-5  | 38   |
| 4    |    |   | Toxoflavin                          | 193 | C7H7N5O2  | 000084-82-2  | 30   |
| 5    |    |   | Cyclopentene, 1-isopropyl-2,3-di... | 138 | C10H18    | 007712-73-4  | 25   |



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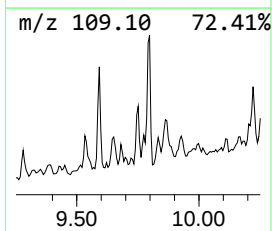
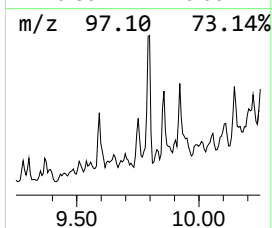
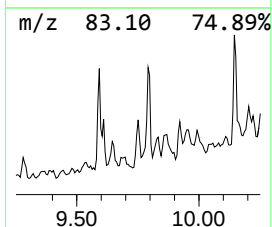
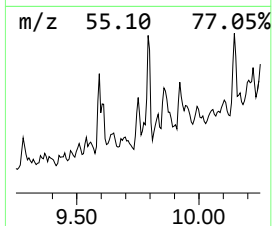
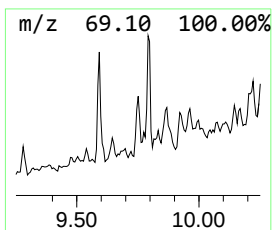
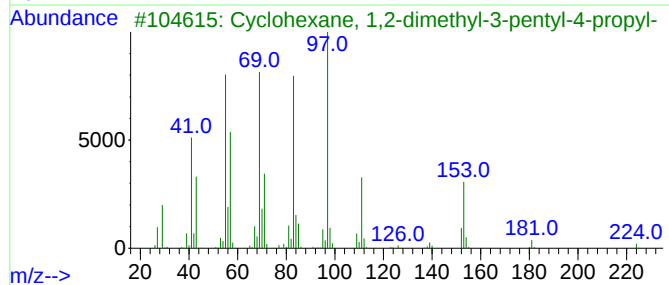
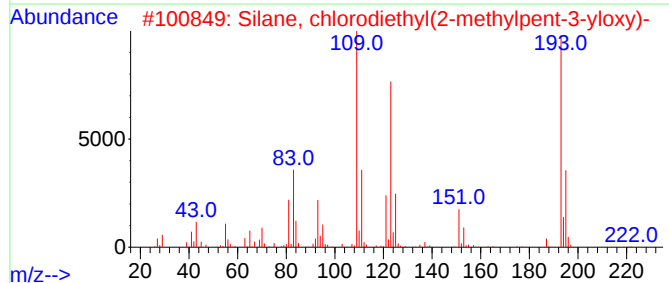
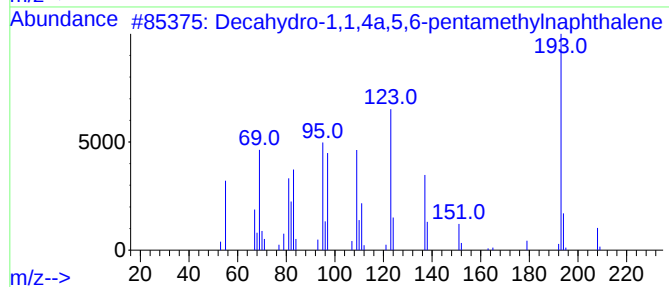
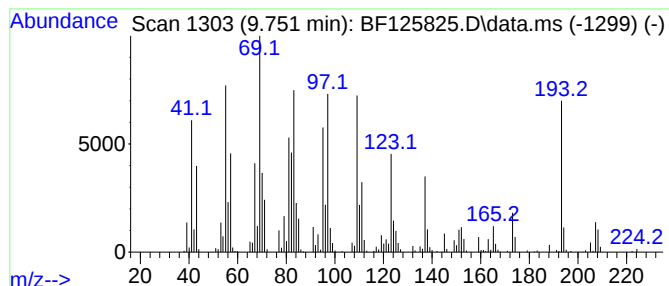
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ClientSampleId :  
CHRT-20426

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF100721.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 5 Decahydro-1,1,4a,5,6-pentam... Concentration Rank 7

| R.T.      | EstConc                             | Area    | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|---------|------------------|--------------|------|
| 9.751     | 5.90 ng                             | 1057160 | Acenaphthene-d10 | 9.880        |      |
| Hit# of 5 | Tentative ID                        | MW      | MolForm          | CAS#         | Qual |
| 1         | Decahydro-1,1,4a,5,6-pentamethyl... | 208     | C15H28           | 080655-44-3  | 93   |
| 2         | Silane, chlorodiethyl(2-methylpe... | 222     | C10H23ClOSi      | 1000363-79-1 | 50   |
| 3         | Cyclohexane, 1,2-dimethyl-3-pent... | 224     | C16H32           | 062376-17-4  | 42   |
| 4         | Neopentylidenecyclohexane           | 152     | C11H20           | 039546-80-0  | 38   |
| 5         | Silane, chlorodiethylhexyloxy-      | 222     | C10H23ClOSi      | 1000363-74-4 | 38   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101321\  
Data File : BF125825.D  
Acq On : 13 Oct 2021 14:14  
Operator : JU/CG  
Sample : M4127-05  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
CHRT-20426

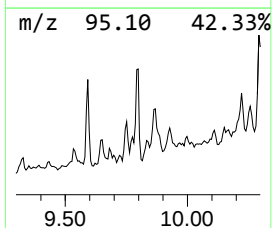
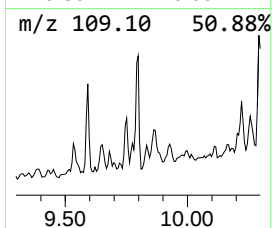
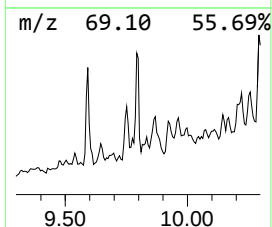
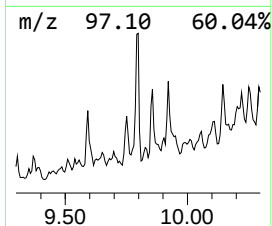
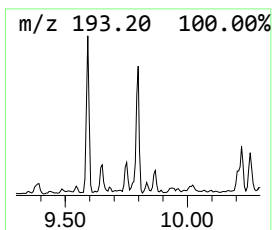
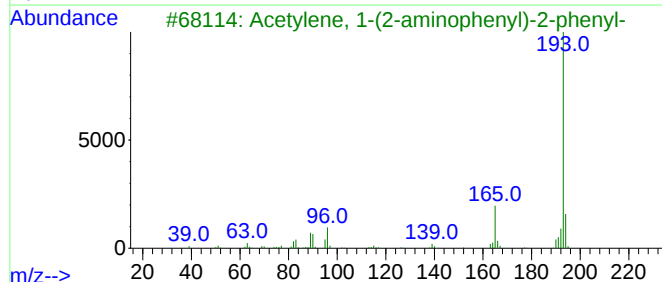
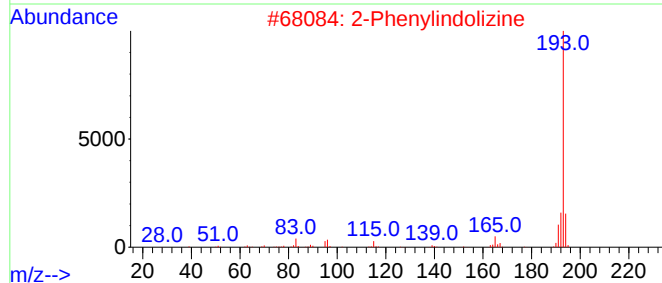
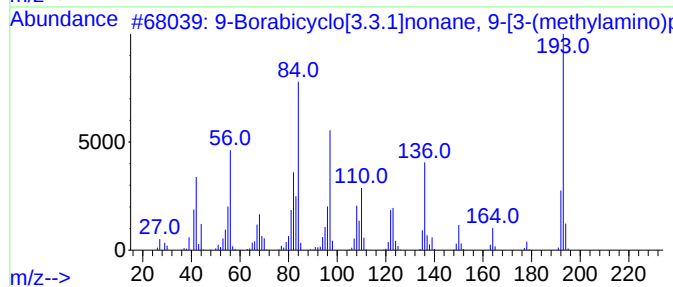
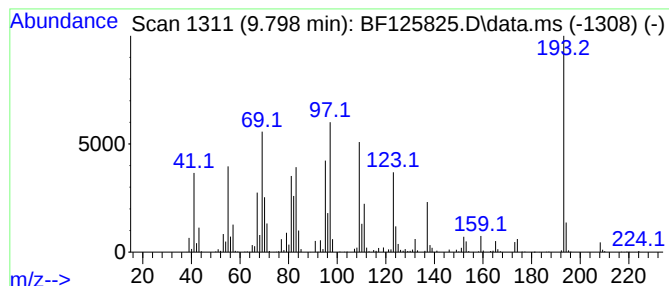
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 6 unknown9.798 Concentration Rank 4

| R.T.  | EstConc | Area    | Relative to ISTD | R.T.  |
|-------|---------|---------|------------------|-------|
| 9.798 | 9.73 ng | 1744640 | Acenaphthene-d10 | 9.880 |

| Hit# | of                                  | 5   | Tentative ID | MW           | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|--------------|---------|------|------|
| 1    | 9-Borabicyclo[3.3.1]nonane, 9-[3... | 193 | C12H24BN     | 1010162-56-0 | 38      |      |      |
| 2    | 2-Phenylindolizine                  | 193 | C14H11N      | 025379-20-8  | 35      |      |      |
| 3    | Acetylene, 1-(2-aminophenyl)-2-p... | 193 | C14H11N      | 1000380-73-4 | 18      |      |      |
| 4    | 2-((Trimethylsilyl)thio)nicotino... | 208 | C9H12N2SSi   | 1000474-02-5 | 14      |      |      |
| 5    | 9-Vinylcarbazole                    | 193 | C14H11N      | 001484-13-5  | 14      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101321\  
Data File : BF125825.D  
Acq On : 13 Oct 2021 14:14  
Operator : JU/CG  
Sample : M4127-05  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
CHRT-20426

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF100721.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

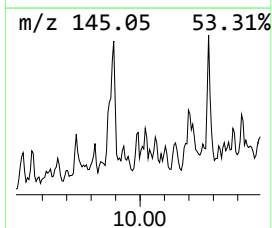
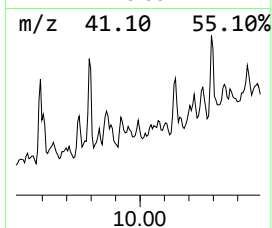
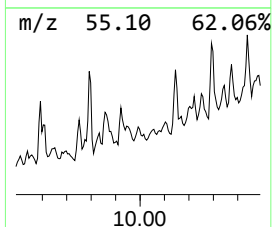
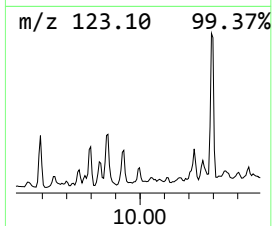
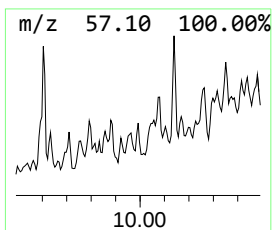
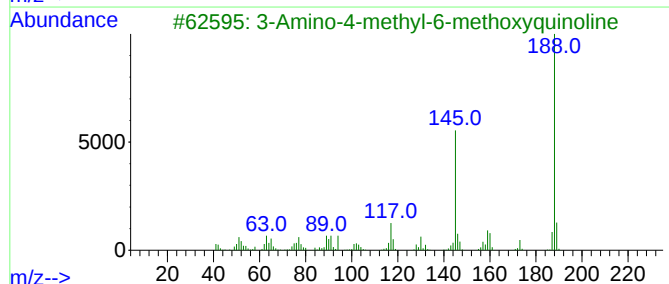
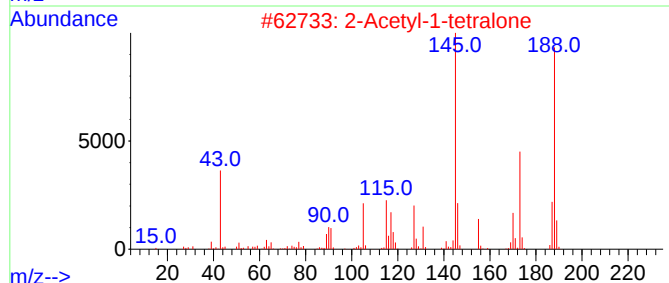
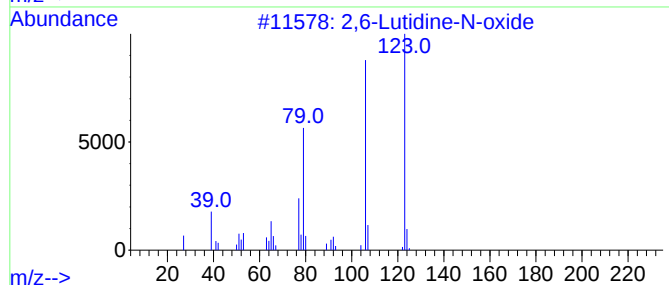
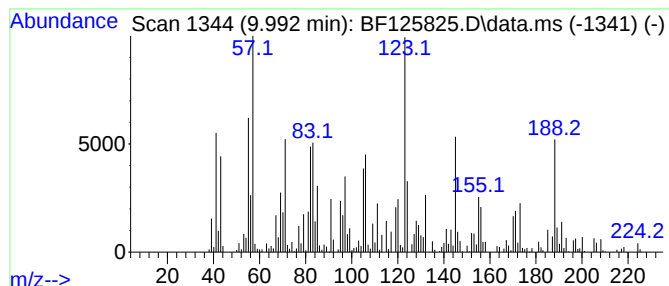
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TIC Integration Parameters: LSCINT.P

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Peak Number 7 unknown9.992 Concentration Rank 16

| R.T.  | EstConc | Area   | Relative to ISTD | R.T.  |
|-------|---------|--------|------------------|-------|
| 9.992 | 2.79 ng | 499514 | Acenaphthene-d10 | 9.880 |

| Hit# | of                                  | 5 | Tentative ID | MW  | MolForm   | CAS#        | Qual |
|------|-------------------------------------|---|--------------|-----|-----------|-------------|------|
| 1    | 2,6-Lutidine-N-oxide                |   |              | 123 | C7H9NO    | 001073-23-0 | 25   |
| 2    | 2-Acetyl-1-tetralone                |   |              | 188 | C12H12O2  | 017216-08-9 | 22   |
| 3    | 3-Amino-4-methyl-6-methoxyquinoline |   |              | 188 | C11H12N2O | 070945-24-3 | 15   |
| 4    | Naphthalene, 2,3-dimethoxy-         |   |              | 188 | C12H12O2  | 010103-06-7 | 14   |
| 5    | 1,4-Dimethyl-2-cyclohexylbenzene    |   |              | 188 | C14H20    | 004501-52-4 | 11   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101321\  
Data File : BF125825.D  
Acq On : 13 Oct 2021 14:14  
Operator : JU/CG  
Sample : M4127-05  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
CHRT-20426

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF100721.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

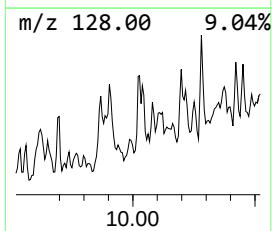
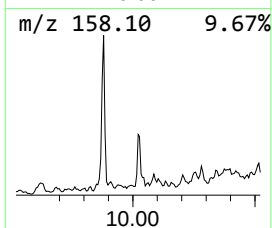
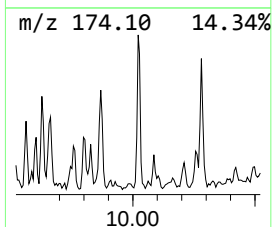
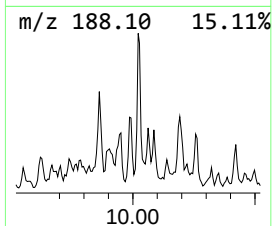
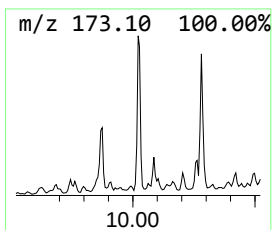
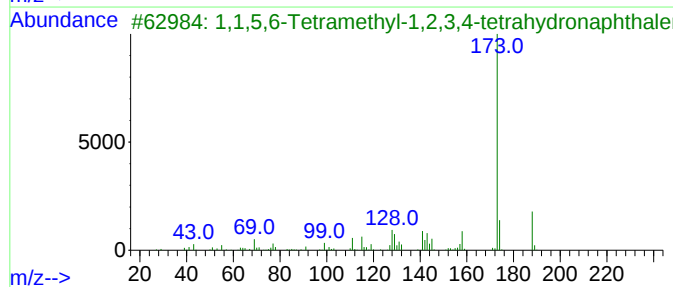
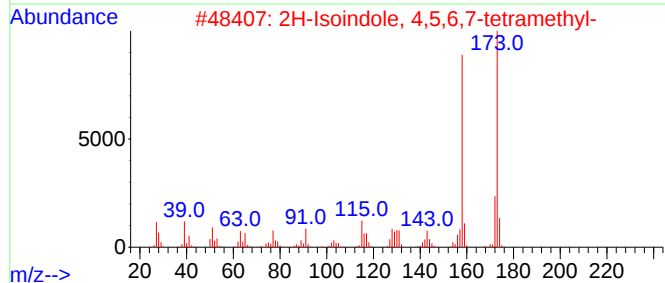
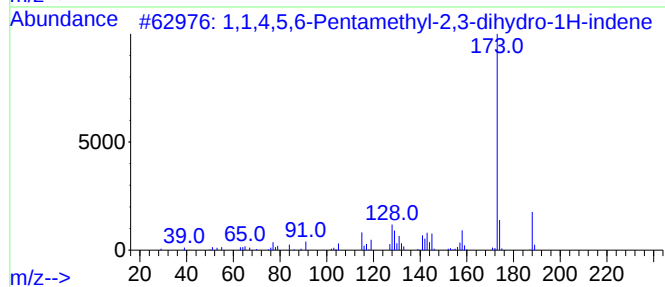
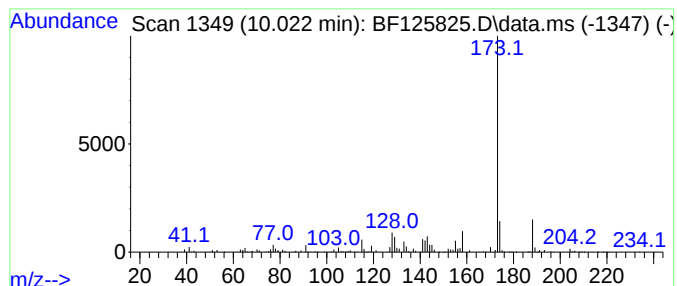
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 1,1,4,5,6-Pentamethyl-2,3-d... Concentration Rank 14

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.  |
|--------|---------|--------|------------------|-------|
| 10.022 | 3.11 ng | 557798 | Acenaphthene-d10 | 9.880 |

| Hit# | of                                  | 5   | Tentative ID | MW           | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|--------------|---------|------|------|
| 1    | 1,1,4,5,6-Pentamethyl-2,3-dihydr... | 188 | C14H20       | 016204-67-4  | 94      |      |      |
| 2    | 2H-Isoindole, 4,5,6,7-tetramethyl-  | 173 | C12H15N      | 070187-61-0  | 72      |      |      |
| 3    | 1,1,5,6-Tetramethyl-1,2,3,4-tetr... | 188 | C14H20       | 031197-54-3  | 64      |      |      |
| 4    | 2-Thioxo-imidazolidin-4-one-5-ac... | 173 | C5H7N3O2S    | 1010146-03-1 | 53      |      |      |
| 5    | 3'-(Trifluoromethyl)acetophenone    | 188 | C9H7F3O      | 000349-76-8  | 53      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101321\  
Data File : BF125825.D  
Acq On : 13 Oct 2021 14:14  
Operator : JU/CG  
Sample : M4127-05  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

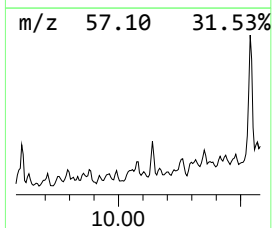
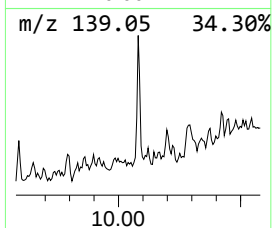
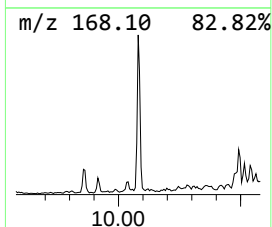
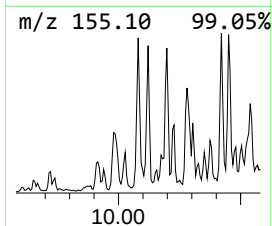
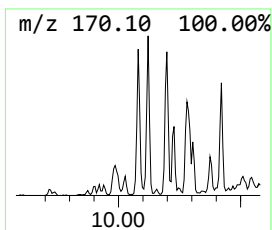
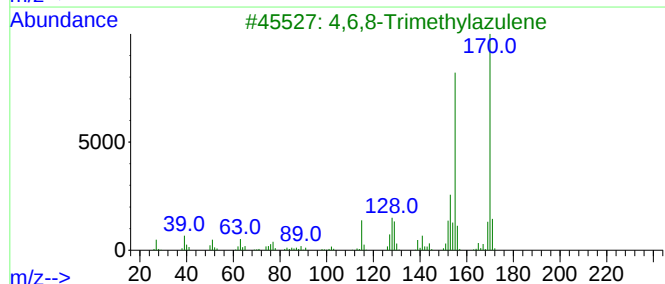
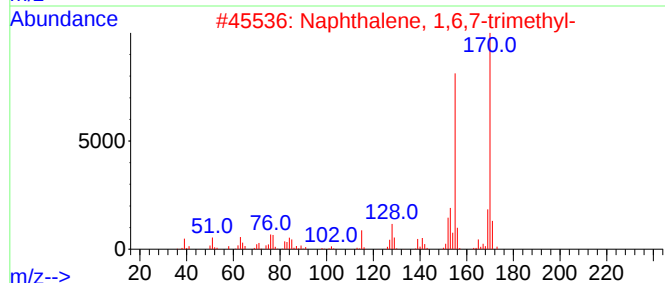
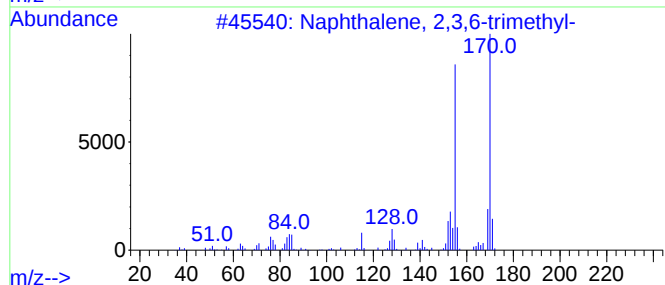
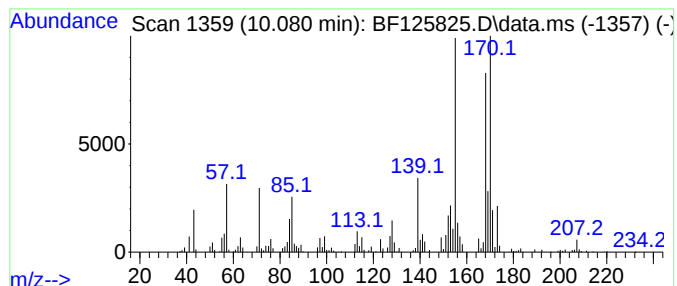
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ClientSampleId :  
CHRT-20426

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF100721.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 9 Naphthalene, 2,3,6-trimethyl- Concentration Rank 20

| R.T.      | EstConc                         | Area   | Relative to ISTD | R.T.         |      |
|-----------|---------------------------------|--------|------------------|--------------|------|
| 10.080    | 2.31 ng                         | 414577 | Acenaphthene-d10 | 9.880        |      |
| Hit# of 5 | Tentative ID                    | MW     | MolForm          | CAS#         | Qual |
| 1         | Naphthalene, 2,3,6-trimethyl-   | 170    | C13H14           | 000829-26-5  | 90   |
| 2         | Naphthalene, 1,6,7-trimethyl-   | 170    | C13H14           | 002245-38-7  | 78   |
| 3         | 4,6,8-Trimethylazulene          | 170    | C13H14           | 000941-81-1  | 70   |
| 4         | 3-(2-Methyl-propenyl)-1H-indene | 170    | C13H14           | 1000187-78-5 | 55   |
| 5         | Naphthalene, 1,4,6-trimethyl-   | 170    | C13H14           | 002131-42-2  | 55   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101321\  
Data File : BF125825.D  
Acq On : 13 Oct 2021 14:14  
Operator : JU/CG  
Sample : M4127-05  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
CHRT-20426

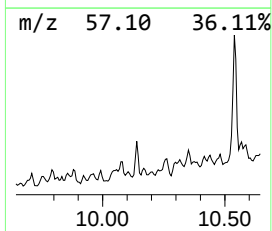
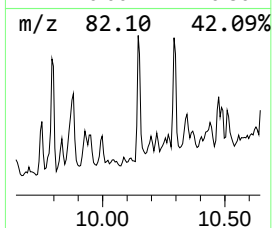
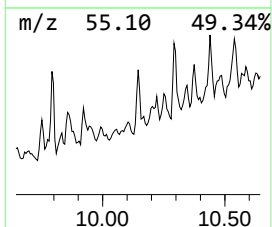
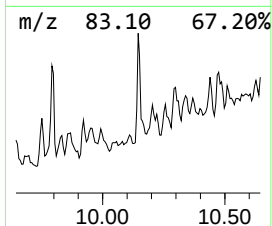
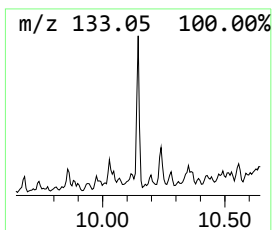
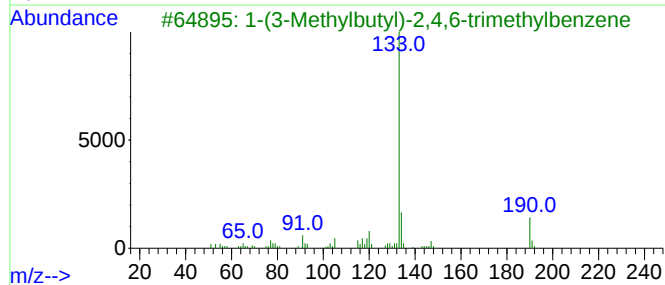
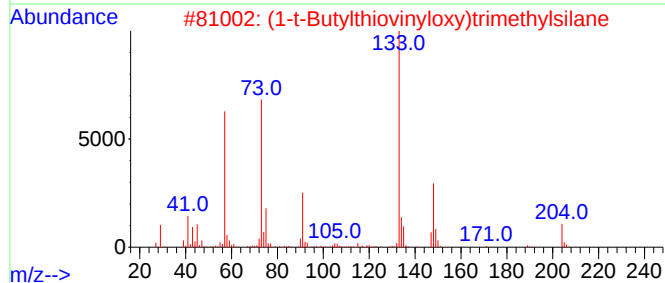
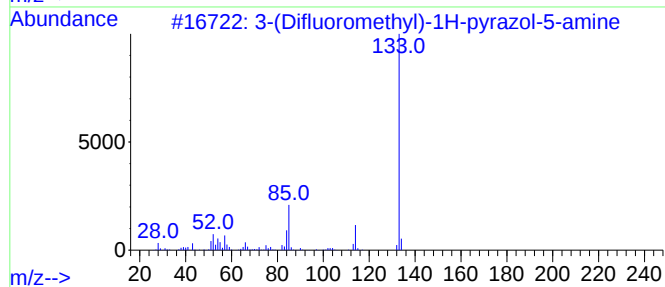
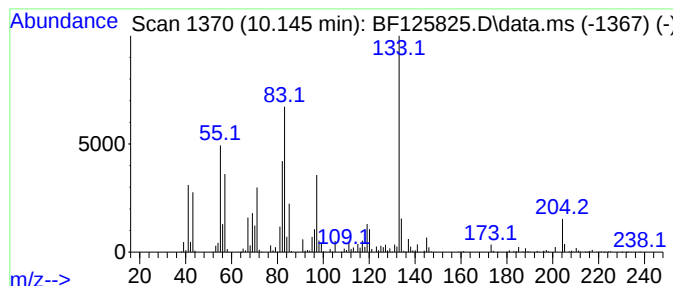
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

\*\*\*\*\*  
Peak Number 10 unknown10.145 Concentration Rank 6

| R.T.   | EstConc | Area    | Relative to ISTD | R.T.  |
|--------|---------|---------|------------------|-------|
| 10.145 | 6.44 ng | 1154490 | Acenaphthene-d10 | 9.880 |

| Hit# | of                                  | 5   | Tentative ID | MW           | MolForm | CAS# | Qual |
|------|-------------------------------------|-----|--------------|--------------|---------|------|------|
| 1    | 3-(Difluoromethyl)-1H-pyrazol-5-... | 133 | C4H5F2N3     | 1284220-49-0 | 30      |      |      |
| 2    | (1-t-Butylthiovinloxy)trimethyl...  | 204 | C9H20OSSi    | 063584-47-4  | 27      |      |      |
| 3    | 1-(3-Methylbutyl)-2,4,6-trimethy... | 190 | C14H22       | 016204-65-2  | 27      |      |      |
| 4    | Succinic acid, 2-methylpent-3-yl... | 284 | C16H28O4     | 1000391-11-2 | 27      |      |      |
| 5    | Cyclohexane, 1,1'-(1,4-butanedi...  | 222 | C16H30       | 006165-44-2  | 25      |      |      |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101321\  
Data File : BF125825.D  
Acq On : 13 Oct 2021 14:14  
Operator : JU/CG  
Sample : M4127-05  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
CHRT-20426

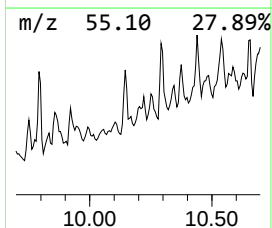
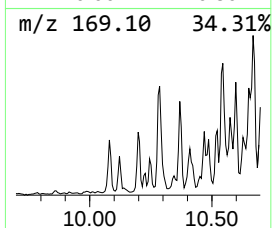
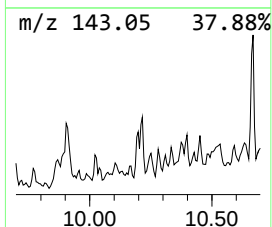
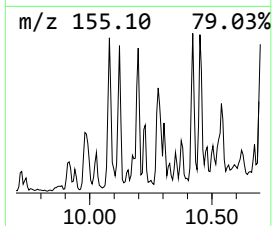
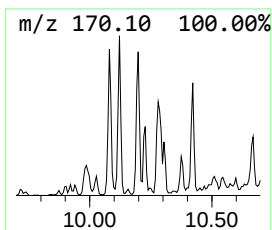
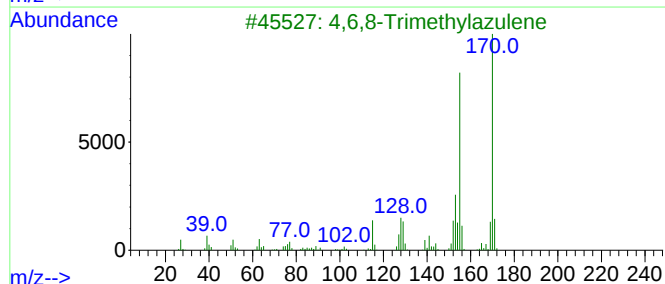
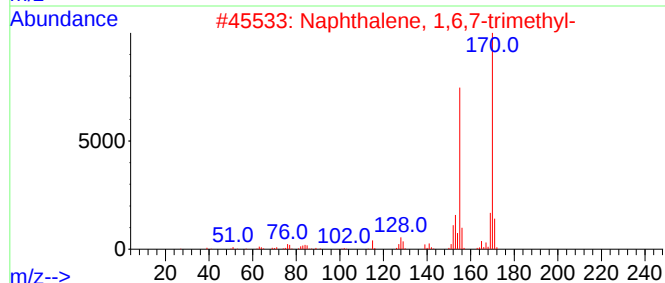
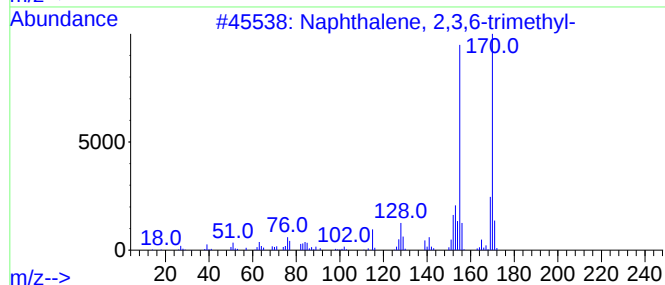
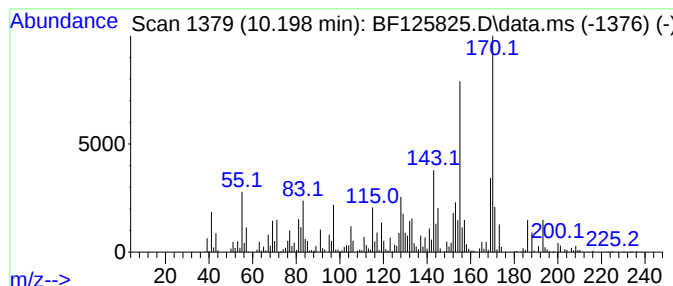
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 11 Naphthalene, 1,6,7-trimethyl- Concentration Rank 8

| R.T.   | EstConc | Area    | Relative to ISTD | R.T.  |
|--------|---------|---------|------------------|-------|
| 10.198 | 5.67 ng | 1017120 | Acenaphthene-d10 | 9.880 |

| Hit# | of | 5 | Tentative ID                  | MW  | MolForm | CAS#        | Qual |
|------|----|---|-------------------------------|-----|---------|-------------|------|
| 1    |    |   | Naphthalene, 2,3,6-trimethyl- | 170 | C13H14  | 000829-26-5 | 96   |
| 2    |    |   | Naphthalene, 1,6,7-trimethyl- | 170 | C13H14  | 002245-38-7 | 96   |
| 3    |    |   | 4,6,8-Trimethylazulene        | 170 | C13H14  | 000941-81-1 | 93   |
| 4    |    |   | Naphthalene, 1,4,6-trimethyl- | 170 | C13H14  | 002131-42-2 | 93   |
| 5    |    |   | Naphthalene, 1,4,5-trimethyl- | 170 | C13H14  | 002131-41-1 | 70   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101321\  
Data File : BF125825.D  
Acq On : 13 Oct 2021 14:14  
Operator : JU/CG  
Sample : M4127-05  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
CHRT-20426

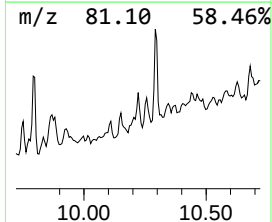
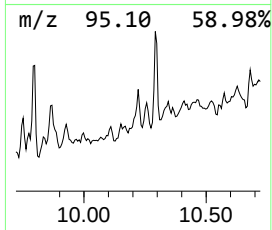
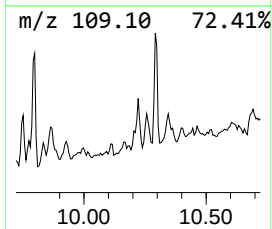
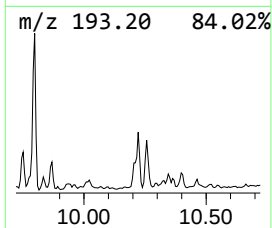
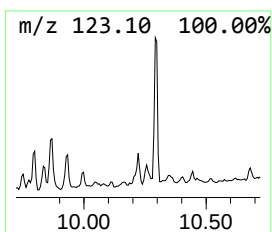
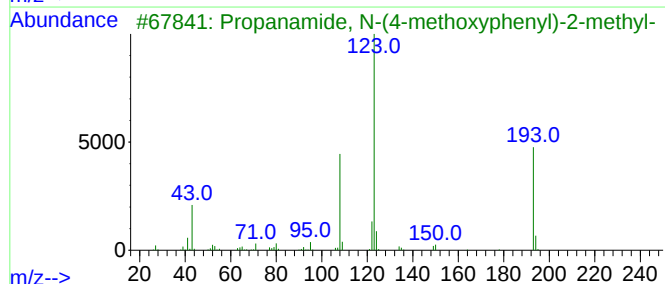
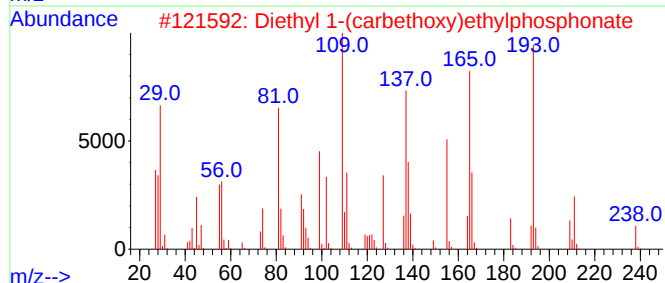
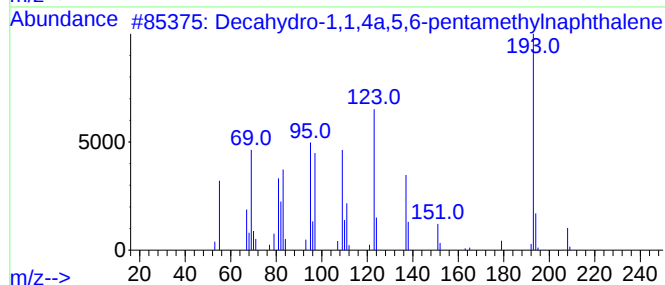
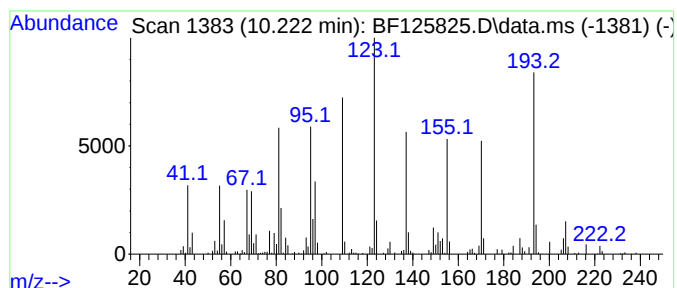
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 12 unknown10.222 Concentration Rank 17

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 10.222    | 2.60 ng                             | 465976 | Acenaphthene-d10 | 9.880        |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | Decahydro-1,1,4a,5,6-pentamethyl... | 208    | C15H28           | 080655-44-3  | 70   |
| 2         | Diethyl 1-(carbethoxy)ethylphosp... | 238    | C9H19O5P         | 1000306-25-7 | 30   |
| 3         | Propanamide, N-(4-methoxyphenyl)... | 193    | C11H15NO2        | 1000307-32-2 | 27   |
| 4         | Diallyldivinylsilane                | 164    | C10H16Si         | 119901-88-1  | 22   |
| 5         | 1,4-Benzenedithiol, S,S'-dimethyl-  | 170    | C8H10S2          | 1000353-07-2 | 15   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101321\  
Data File : BF125825.D  
Acq On : 13 Oct 2021 14:14  
Operator : JU/CG  
Sample : M4127-05  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
CHRT-20426

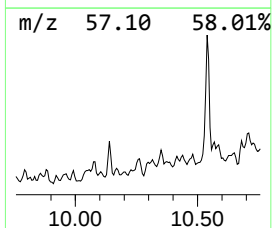
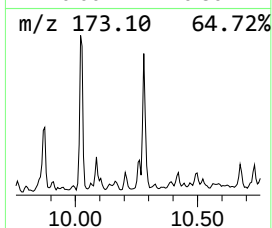
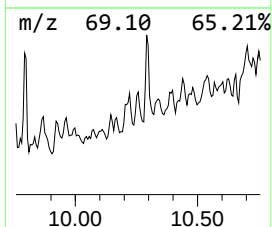
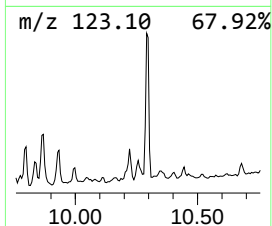
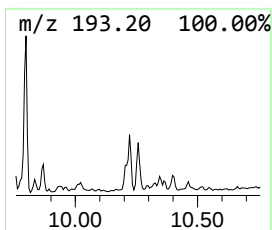
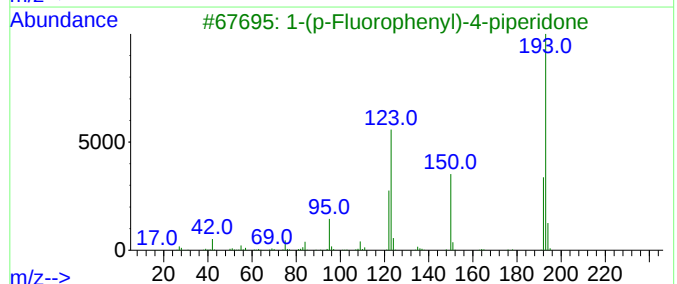
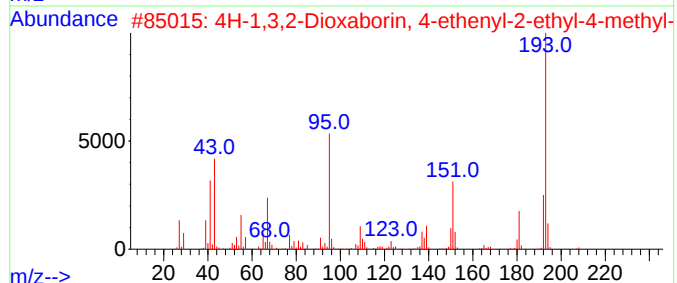
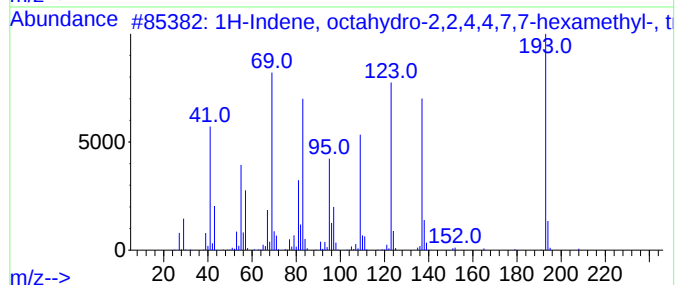
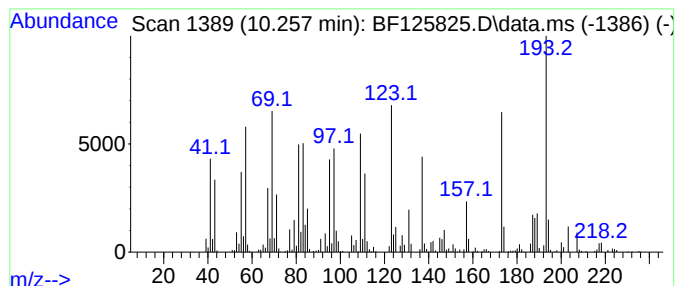
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 13 unknown10.257 Concentration Rank 10

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.  |
|--------|---------|--------|------------------|-------|
| 10.257 | 4.55 ng | 816319 | Acenaphthene-d10 | 9.880 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|-----------|--------------|------|
| 1    |    |   | 1H-Indene, octahydro-2,2,4,7,7...   | 208 | C15H28    | 054832-83-6  | 49   |
| 2    |    |   | 4H-1,3,2-Dioxaborin, 4-ethenyl-2... | 208 | C12H21BO2 | 074630-04-9  | 30   |
| 3    |    |   | 1-(p-Fluorophenyl)-4-piperidone     | 193 | C11H12FNO | 1000238-56-7 | 30   |
| 4    |    |   | 2,6-Heptadienal, 2,4-dimethyl-      | 138 | C9H14O    | 085136-08-9  | 25   |
| 5    |    |   | Benzenamine, 3-ethoxy-              | 137 | C8H11NO   | 000621-33-0  | 15   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101321\  
Data File : BF125825.D  
Acq On : 13 Oct 2021 14:14  
Operator : JU/CG  
Sample : M4127-05  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
CHRT-20426

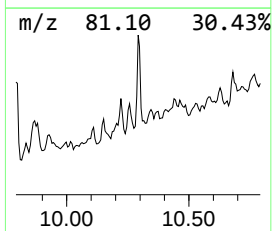
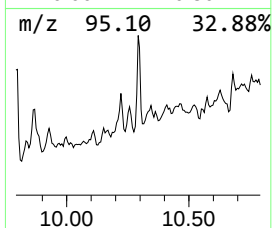
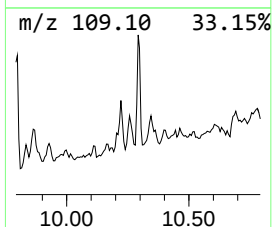
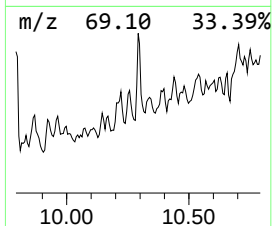
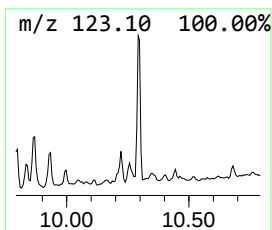
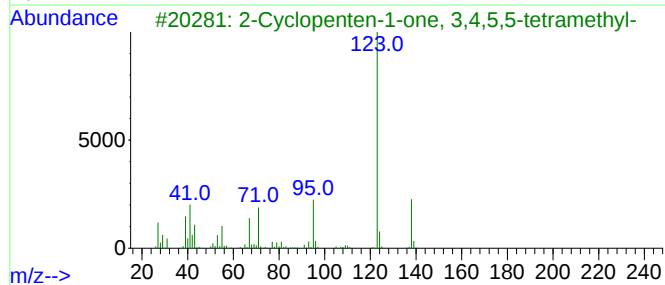
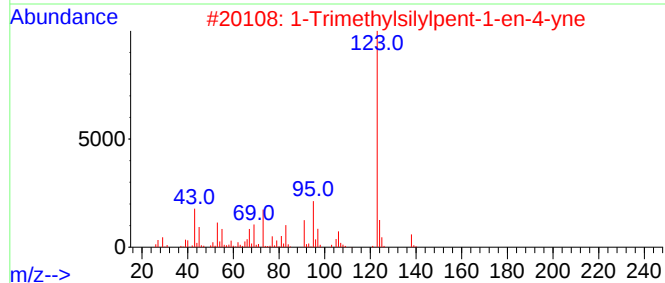
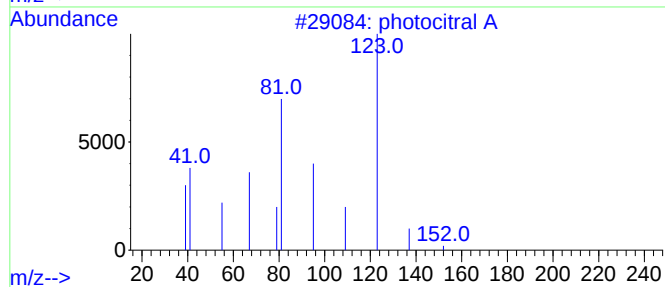
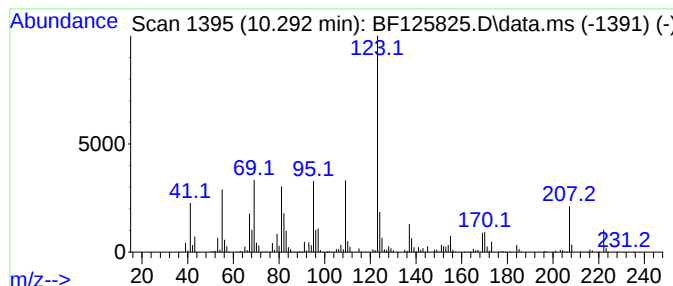
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 14 unknown10.292 Concentration Rank 3

| R.T.   | EstConc  | Area    | Relative to ISTD | R.T.  |
|--------|----------|---------|------------------|-------|
| 10.292 | 11.75 ng | 2106120 | Acenaphthene-d10 | 9.880 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | photocitral A                       | 152 | C10H16O    | 055253-28-6  | 47   |
| 2    |    |   | 1-Trimethylsilylpent-1-en-4-yne     | 138 | C8H14Si    | 079516-25-9  | 47   |
| 3    |    |   | 2-Cyclopenten-1-one, 3,4,5,5-tet... | 138 | C9H14O     | 090611-02-2  | 43   |
| 4    |    |   | 4-Fluorobenzoic acid, 3-methylbu... | 208 | C12H13FO2  | 1000299-15-1 | 43   |
| 5    |    |   | 2-Fluoro-N-(1,3-thiazol-2-yl)ben... | 222 | C10H7FN2OS | 000319-38-0  | 43   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101321\  
Data File : BF125825.D  
Acq On : 13 Oct 2021 14:14  
Operator : JU/CG  
Sample : M4127-05  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
CHRT-20426

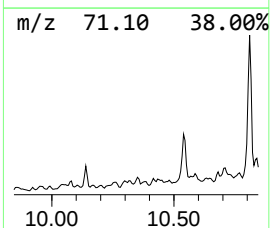
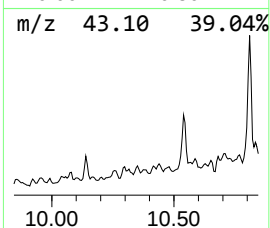
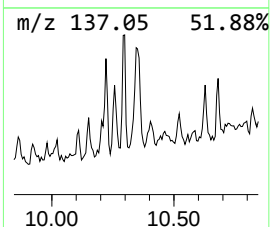
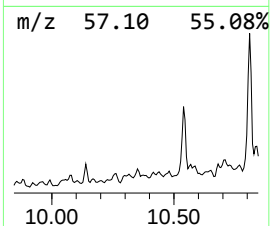
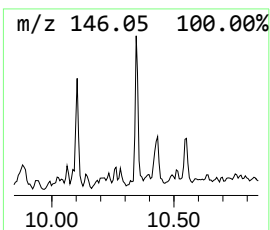
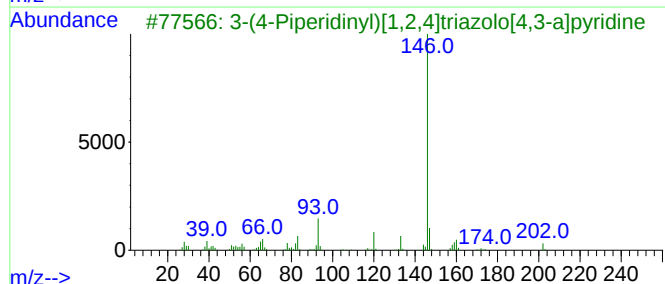
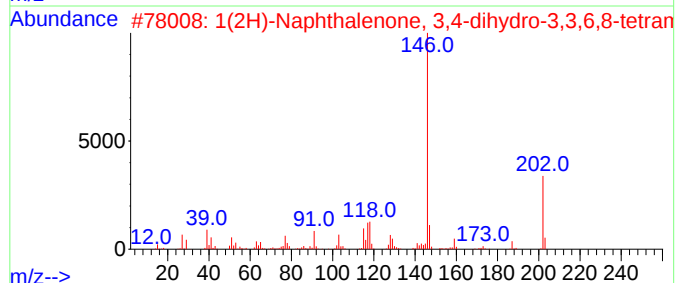
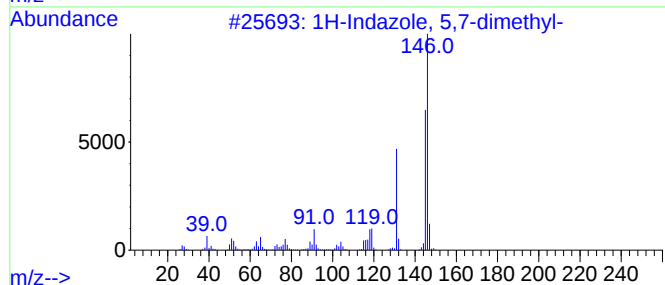
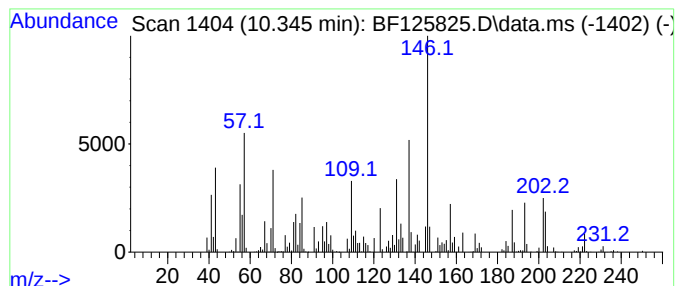
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 15 unknown10.345 Concentration Rank 15

| R.T.      | EstConc                             | Area   | Relative to ISTD | R.T.         |      |
|-----------|-------------------------------------|--------|------------------|--------------|------|
| 10.345    | 2.85 ng                             | 511600 | Acenaphthene-d10 | 9.880        |      |
| Hit# of 5 | Tentative ID                        | MW     | MolForm          | CAS#         | Qual |
| 1         | 1H-Indazole, 5,7-dimethyl-          | 146    | C9H10N2          | 043067-41-0  | 25   |
| 2         | 1(2H)-Naphthalenone, 3,4-dihydro... | 202    | C14H18O          | 005409-55-2  | 22   |
| 3         | 3-(4-Piperidiny)[1,2,4]triazolo...  | 202    | C11H14N4         | 852627-78-2  | 22   |
| 4         | Methyl N-(2-acetylphenyl)carbamate  | 193    | C10H11NO3        | 1000445-05-3 | 22   |
| 5         | Methyl N-(4-acetylphenyl)carbamate  | 193    | C10H11NO3        | 060677-43-2  | 22   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101321\  
Data File : BF125825.D  
Acq On : 13 Oct 2021 14:14  
Operator : JU/CG  
Sample : M4127-05  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
CHRT-20426

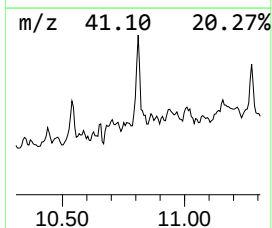
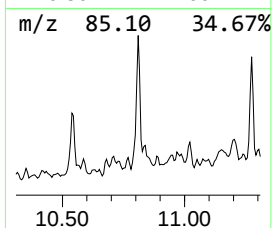
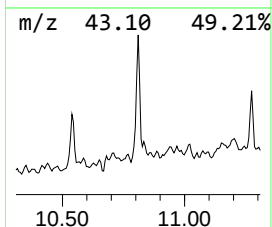
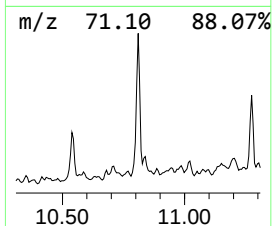
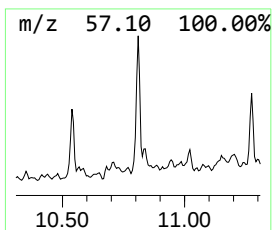
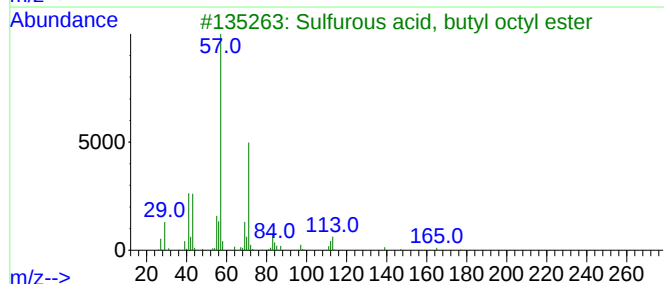
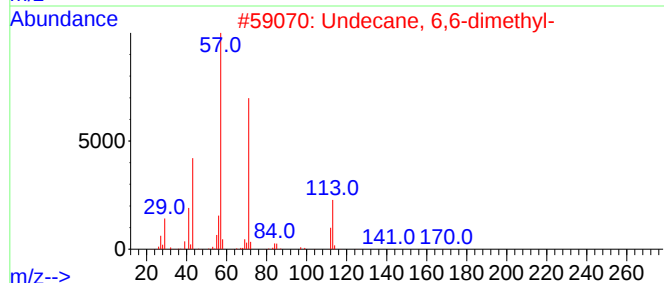
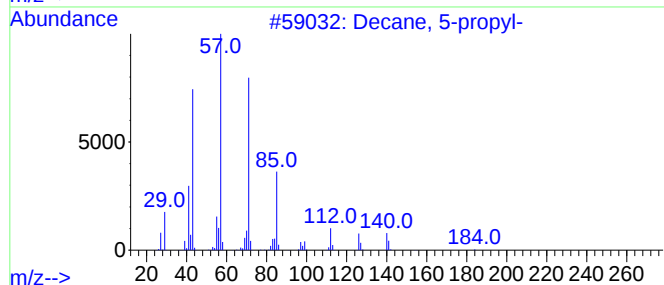
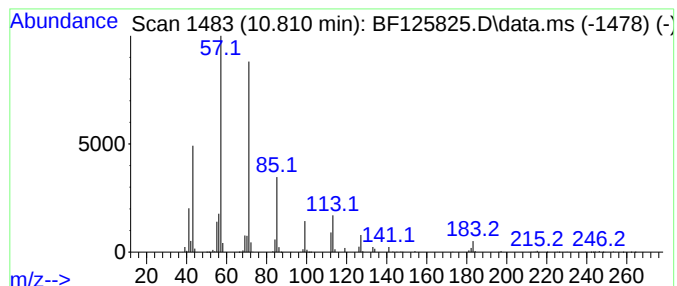
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 16 Decane, 5-propyl- Concentration Rank 1

| R.T.   | EstConc   | Area    | Relative to ISTD | R.T.   |
|--------|-----------|---------|------------------|--------|
| 10.810 | 112.56 ng | 4562740 | Phenanthrene-d10 | 11.368 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm   | CAS#         | Qual |
|------|------|-------------------------------------|-----|-----------|--------------|------|
| 1    |      | Decane, 5-propyl-                   | 184 | C13H28    | 017312-62-8  | 80   |
| 2    |      | Undecane, 6,6-dimethyl-             | 184 | C13H28    | 017312-76-4  | 64   |
| 3    |      | Sulfurous acid, butyl octyl ester   | 250 | C12H26O3S | 1000309-17-5 | 59   |
| 4    |      | Nonane, 5-methyl-5-propyl-          | 184 | C13H28    | 017312-75-3  | 53   |
| 5    |      | Pentadecane, 2,6,10,14-tetramethyl- | 268 | C19H40    | 001921-70-6  | 50   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101321\  
Data File : BF125825.D  
Acq On : 13 Oct 2021 14:14  
Operator : JU/CG  
Sample : M4127-05  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
CHRT-20426

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF100721.M  
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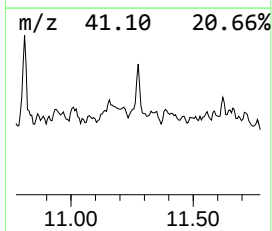
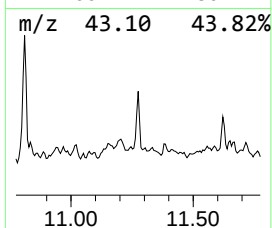
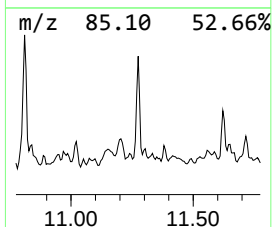
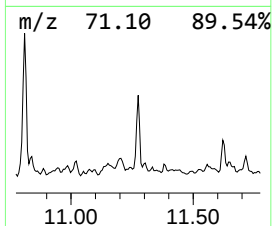
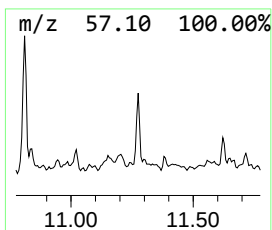
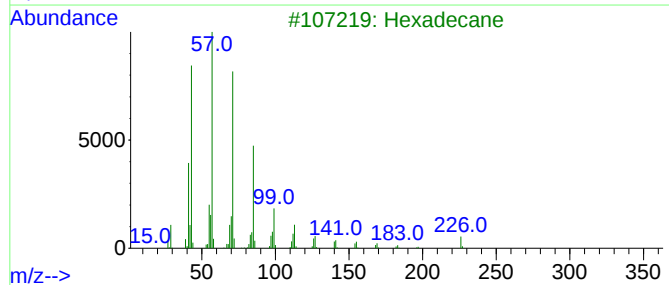
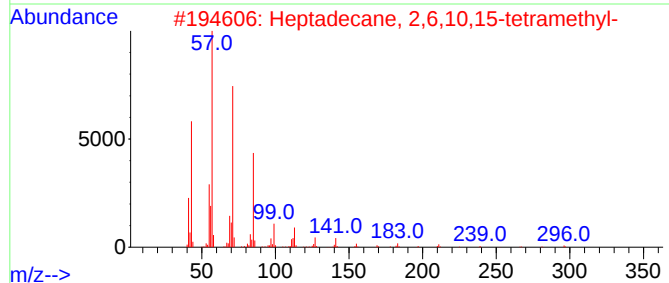
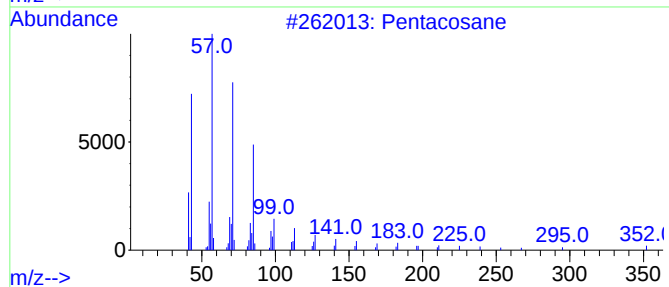
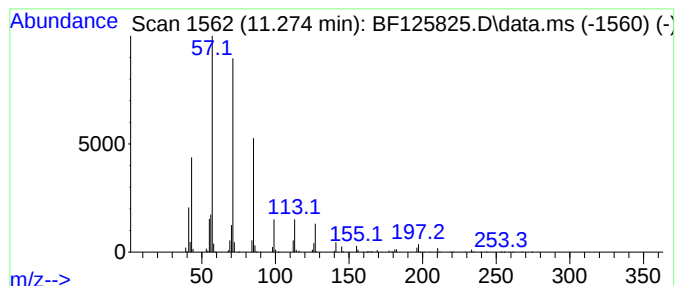
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 17 Pentacosane Concentration Rank 2

| R.T.   | EstConc  | Area    | Relative to ISTD | R.T.   |
|--------|----------|---------|------------------|--------|
| 11.274 | 46.98 ng | 1904200 | Phenanthrene-d10 | 11.368 |

| Hit# | of 5 | Tentative ID                        | MW  | MolForm | CAS#        | Qual |
|------|------|-------------------------------------|-----|---------|-------------|------|
| 1    |      | Pentacosane                         | 352 | C25H52  | 000629-99-2 | 90   |
| 2    |      | Heptadecane, 2,6,10,15-tetramethyl- | 296 | C21H44  | 054833-48-6 | 86   |
| 3    |      | Hexadecane                          | 226 | C16H34  | 000544-76-3 | 86   |
| 4    |      | Tridecane, 7-hexyl-                 | 268 | C19H40  | 007225-66-3 | 86   |
| 5    |      | Hexadecane, 2,6,10,14-tetramethyl-  | 282 | C20H42  | 000638-36-8 | 83   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101321\  
Data File : BF125825.D  
Acq On : 13 Oct 2021 14:14  
Operator : JU/CG  
Sample : M4127-05  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
CHRT-20426

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF100721.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

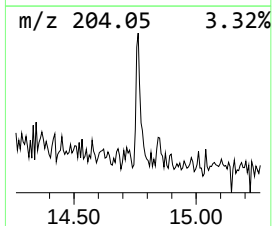
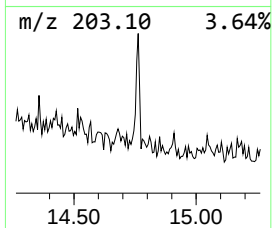
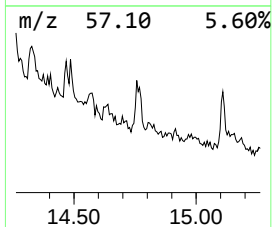
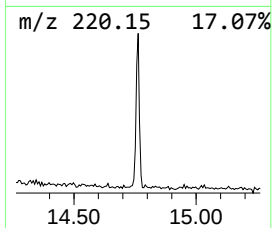
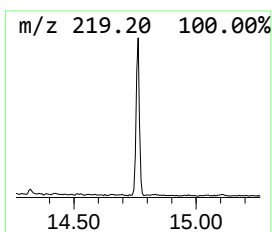
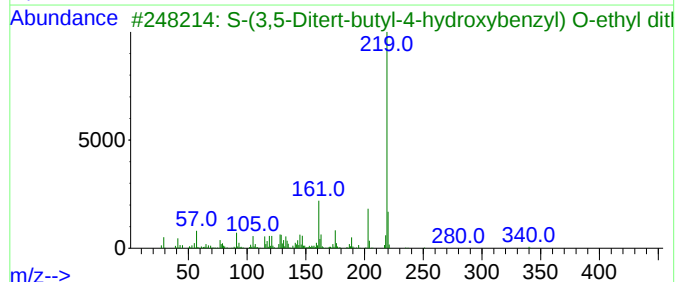
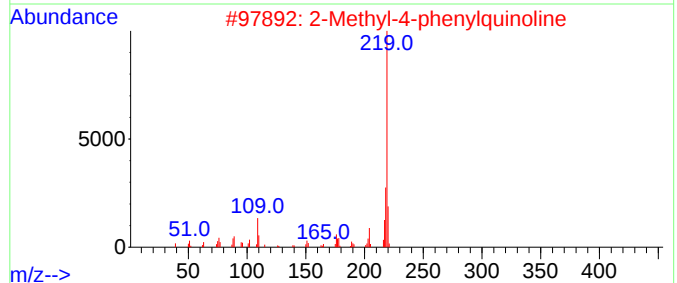
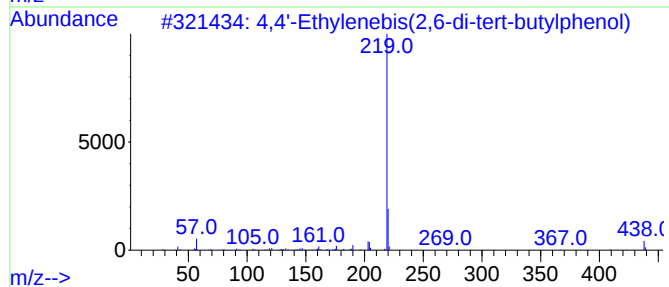
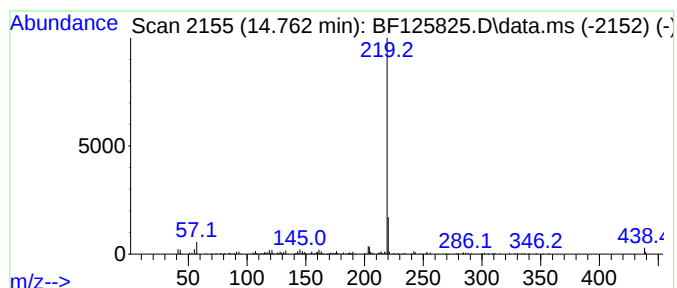
TIC Library : C:\Database\NIST20.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 18 4,4'-Ethylenebis(2,6-di-ter... Concentration Rank 11

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 14.762 | 3.76 ng | 188998 | Perylene-d12     | 15.451 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#         | Qual |
|------|----|---|-------------------------------------|-----|------------|--------------|------|
| 1    |    |   | 4,4'-Ethylenebis(2,6-di-tert-but... | 438 | C30H46O2   | 001516-94-5  | 91   |
| 2    |    |   | 2-Methyl-4-phenylquinoline          | 219 | C16H13N    | 001721-92-2  | 72   |
| 3    |    |   | S-(3,5-Ditert-butyl-4-hydroxyben... | 340 | C18H28O2S2 | 015367-50-7  | 64   |
| 4    |    |   | Terephthalic acid, 3-chloropheny... | 346 | C19H19ClO4 | 1000323-64-4 | 64   |
| 5    |    |   | Phenol, 2,6-bis(1,1-dimethylethy... | 234 | C16H26O    | 004130-42-1  | 56   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101321\  
Data File : BF125825.D  
Acq On : 13 Oct 2021 14:14  
Operator : JU/CG  
Sample : M4127-05  
Misc :  
ALS Vial : 10 Sample Multiplier: 1

Instrument :  
BNA\_F  
ClientSampleId :  
CHRT-20426

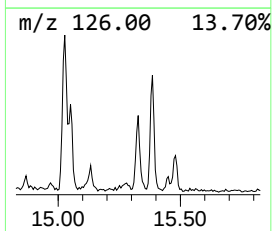
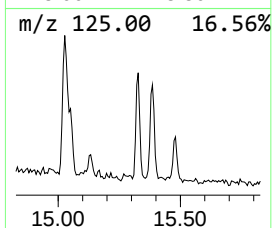
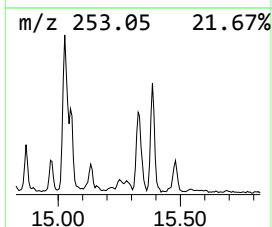
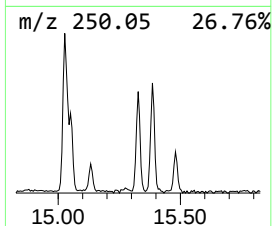
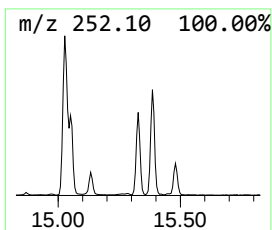
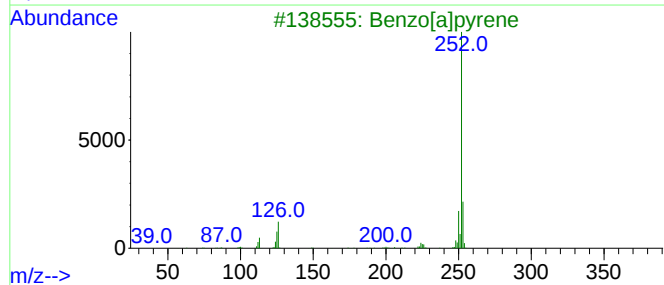
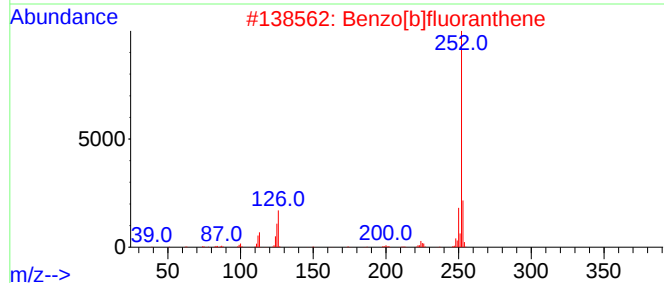
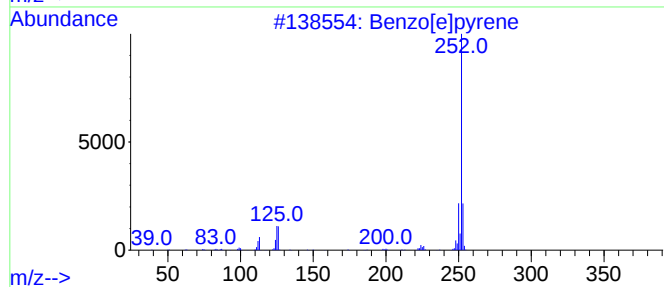
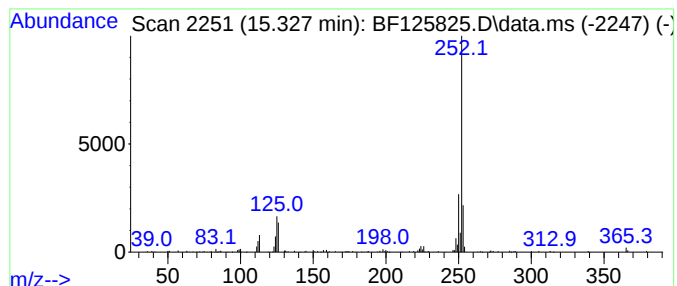
Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF100721.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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Peak Number 19 Benzo[e]pyrene Concentration Rank 9

| R.T.   | EstConc | Area   | Relative to ISTD | R.T.   |
|--------|---------|--------|------------------|--------|
| 15.327 | 5.38 ng | 270488 | Perylene-d12     | 15.451 |

| Hit# | of | 5 | Tentative ID         | MW  | MolForm | CAS#        | Qual |
|------|----|---|----------------------|-----|---------|-------------|------|
| 1    |    |   | Benzo[e]pyrene       | 252 | C20H12  | 000192-97-2 | 98   |
| 2    |    |   | Benzo[b]fluoranthene | 252 | C20H12  | 000205-99-2 | 96   |
| 3    |    |   | Benzo[a]pyrene       | 252 | C20H12  | 000050-32-8 | 95   |
| 4    |    |   | Benzo[k]fluoranthene | 252 | C20H12  | 000207-08-9 | 94   |
| 5    |    |   | Perylene             | 252 | C20H12  | 000198-55-0 | 94   |



Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF101321\  
 Data File : BF125825.D  
 Acq On : 13 Oct 2021 14:14  
 Operator : JU/CG  
 Sample : M4127-05  
 Misc :  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 CHRT-20426

Quant Method : Z:\SVOASRV\HPCHEM1\BNA\_F\METHODS\8270-BF100721.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 |
| Ethanol, 2-(2-b... | 8.028  | 7.9     | ng    | 685909   | 2                     | 8.122  | 1739570 | 20.0 |
| unknown8.875       | 8.875  | 2.6     | ng    | 224716   | 2                     | 8.122  | 1739570 | 20.0 |
| unknown9.280       | 9.280  | 3.1     | ng    | 565364   | 3                     | 9.880  | 3585580 | 20.0 |
| Decahydro-1,1,4... | 9.751  | 5.9     | ng    | 1057160  | 3                     | 9.880  | 3585580 | 20.0 |
| unknown9.798       | 9.798  | 9.7     | ng    | 1744640  | 3                     | 9.880  | 3585580 | 20.0 |
| unknown9.992       | 9.992  | 2.8     | ng    | 499514   | 3                     | 9.880  | 3585580 | 20.0 |
| 1,1,4,5,6-Penta... | 10.022 | 3.1     | ng    | 557798   | 3                     | 9.880  | 3585580 | 20.0 |
| Naphthalene, 2,... | 10.080 | 2.3     | ng    | 414577   | 3                     | 9.880  | 3585580 | 20.0 |
| unknown10.145      | 10.145 | 6.4     | ng    | 1154490  | 3                     | 9.880  | 3585580 | 20.0 |
| Naphthalene, 1,... | 10.198 | 5.7     | ng    | 1017120  | 3                     | 9.880  | 3585580 | 20.0 |
| unknown10.222      | 10.222 | 2.6     | ng    | 465976   | 3                     | 9.880  | 3585580 | 20.0 |
| unknown10.257      | 10.257 | 4.5     | ng    | 816319   | 3                     | 9.880  | 3585580 | 20.0 |
| unknown10.292      | 10.292 | 11.8    | ng    | 2106120  | 3                     | 9.880  | 3585580 | 20.0 |
| unknown10.345      | 10.345 | 2.9     | ng    | 511600   | 3                     | 9.880  | 3585580 | 20.0 |
| Decane, 5-propyl-  | 10.810 | 112.6   | ng    | 4562740  | 4                     | 11.368 | 810707  | 20.0 |
| Pentacosane        | 11.274 | 47.0    | ng    | 1904200  | 4                     | 11.368 | 810707  | 20.0 |
| 4,4'-Ethylenebi... | 14.762 | 3.8     | ng    | 188998   | 6                     | 15.451 | 1005060 | 20.0 |
| Benzo[e]pyrene     | 15.327 | 5.4     | ng    | 270488   | 6                     | 15.451 | 1005060 | 20.0 |