LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042123\
Data File : BM039565.D
Acq On : 21 Apr 2023 23:29
Operator : CG/JU
Sample : 02386-11
Misc :
ALS Vial : 13 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** SPA-3WC-23-04

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_M\Methods\8270-BM041723.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BM039565.D\data.ms

| peak # | R.T. min | first scan | max scan | last scan | PK TY | peak height | corr. area | corr. % max. | % of total |
|---|---|---|---|---|---|---|---|---|---|
| 1 | 3.966 | 149 | 158 | 170 | rVB | 2196762 | 3500061 | 9.71% | 0.374% |
| 2 | 5.107 | 345 | 352 | 362 | rBV | 4276032 | 7400345 | 20.54% | 0.790% |
| 3 | 5.307 | 381 | 386 | 393 | rVV | 11276127 | 16974591 | 47.12% | 1.813% |
| 4 | 5.390 | 393 | 400 | 403 | rVV2 | 2104103 | 4186470 | 11.62% | 0.447% |
| 5 | 5.443 | 403 | 409 | 425 | rVB | 17624518 | 36027671 | 100.00% | 3.848% |
| 6 | 5.813 | 465 | 472 | 480 | rVV | 11313525 | 17467327 | 48.48% | 1.866% |
| 7 | 6.066 | 507 | 515 | 523 | rBV4 | 2029850 | 4749483 | 13.18% | 0.507% |
| 8 | 6.290 | 546 | 553 | 559 | rBV4 | 2697223 | 5661047 | 15.71% | 0.605% |
| 9 | 6.354 | 559 | 564 | 569 | rBV | 3073958 | 4343434 | 12.06% | 0.464% |
| 10 | 6.437 | 574 | 578 | 581 | rVV2 | 2960586 | 5027273 | 13.95% | 0.537% |
| 11 | 6.472 | 581 | 584 | 591 | rVB2 | 2103195 | 4252893 | 11.80% | 0.454% |
| 12 | 6.701 | 614 | 623 | 628 | rBV2 | 1788545 | 4429590 | 12.29% | 0.473% |
| 13 | 6.795 | 628 | 639 | 644 | rBV3 | 8213322 | 16288195 | 45.21% | 1.740% |
| 14 | 6.854 | 644 | 649 | 652 | rVV | 3918330 | 5228296 | 14.51% | 0.558% |
| 15 | 6.901 | 652 | 657 | 661 | rVV | 8279119 | 12239562 | 33.97% | 1.307% |
| 16 | 6.960 | 661 | 667 | 672 | rVV3 | 7234629 | 13909072 | 38.61% | 1.486% |
| 17 | 7.037 | 672 | 680 | 686 | rVB3 | 5552111 | 11016021 | 30.58% | 1.177% |
| 18 | 7.201 | 701 | 708 | 714 | rBV3 | 3299337 | 6873483 | 19.08% | 0.734% |
| 19 | 7.442 | 742 | 749 | 751 | rBV | 12774279 | 22195435 | 61.61% | 2.371% |
| 20 | 7.466 | 751 | 753 | 760 | rVB | 12717368 | 17122950 | 47.53% | 1.829% |
| 21 | 7.713 | 787 | 795 | 802 | rBV6 | 2981663 | 8536974 | 23.70% | 0.912% |
| 22 | 7.795 | 802 | 809 | 814 | rVB | 8251902 | 12622867 | 35.04% | 1.348% |
| 23 | 7.860 | 814 | 820 | 828 | rVB3 | 7824476 | 14629959 | 40.61% | 1.563% |
| 24 | 7.937 | 828 | 833 | 840 | rBV3 | 5307846 | 10999375 | 30.53% | 1.175% |
| 25 | 8.001 | 840 | 844 | 847 | rBV2 | 2333579 | 3062619 | 8.50% | 0.327% |
| 26 | 8.042 | 847 | 851 | 857 | rBV | 5321677 | 8527210 | 23.67% | 0.911% |
| 27 | 8.101 | 857 | 861 | 868 | rVB2 | 3554877 | 6049532 | 16.79% | 0.646% |
| 28 | 8.178 | 868 | 874 | 881 | rVB3 | 5915384 | 10909430 | 30.28% | 1.165% |
| 29 | 8.313 | 891 | 897 | 899 | rBV3 | 1767882 | 3359106 | 9.32% | 0.359% |
| 30 | 8.348 | 899 | 903 | 910 | rBV2 | 5007620 | 10831538 | 30.06% | 1.157% |
| 31 | 8.401 | 910 | 912 | 917 | rVB | 3346554 | 4443284 | 12.33% | 0.475% |
| 32 | 8.472 | 917 | 924 | 935 | rBV4 | 8061865 | 20468894 | 56.81% | 2.186% |
| 33 | 8.584 | 935 | 943 | 946 | rBV | 3943167 | 6423926 | 17.83% | 0.686% |
| 34 | 8.778 | 971 | 976 | 980 | rBV | 2285125 | 3646787 | 10.12% | 0.389% |
| 35 | 8.825 | 980 | 984 | 991 | rVB | 2749915 | 4528159 | 12.57% | 0.484% |
| 36 | 8.931 | 996 | 1002 | 1008 | rVB | 6476319 | 11021959 | 30.59% | 1.177% |

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042123\
Data File : BM039565.D
Acq On : 21 Apr 2023 23:29
Operator : CG/JU
Sample : 02386-11
Misc :
ALS Vial : 13 Sample Multiplier: 1

**Instrument :**
BNA_M
**ClientSampleId :**
SPA-3WC-23-04

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_M\Methods\8270-BM041723.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 37 | 9.013 | 1009 | 1016 | 1018 | rBV3 | 2288157 | 4684148 | 13.00% | 0.500% |
| 38 | 9.054 | 1018 | 1023 | 1032 | rVB2 | 14150847 | 27958639 | 77.60% | 2.986% |
| 39 | 9.178 | 1039 | 1044 | 1050 | rVB4 | 2652874 | 5527996 | 15.34% | 0.590% |
| 40 | 9.266 | 1052 | 1059 | 1063 | rBV3 | 1821245 | 4595149 | 12.75% | 0.491% |
| 41 | 9.389 | 1074 | 1080 | 1086 | rBV4 | 1298317 | 3014481 | 8.37% | 0.322% |
| 42 | 9.454 | 1086 | 1091 | 1096 | rBV2 | 4290759 | 7116322 | 19.75% | 0.760% |
| 43 | 9.513 | 1096 | 1101 | 1109 | rVB | 5142866 | 9043726 | 25.10% | 0.966% |
| 44 | 9.972 | 1173 | 1179 | 1183 | rBV2 | 1910812 | 3993971 | 11.09% | 0.427% |
| 45 | 10.025 | 1183 | 1188 | 1196 | rVV3 | 4829201 | 12307023 | 34.16% | 1.314% |
| 46 | 10.089 | 1196 | 1199 | 1204 | rVV2 | 2166562 | 3523983 | 9.78% | 0.376% |
| 47 | 10.207 | 1212 | 1219 | 1223 | rVV4 | 1241929 | 2927538 | 8.13% | 0.313% |
| 48 | 10.260 | 1223 | 1228 | 1234 | rVB | 3102806 | 4919113 | 13.65% | 0.525% |
| 49 | 10.325 | 1234 | 1239 | 1246 | rVB | 1521277 | 2834488 | 7.87% | 0.303% |
| 50 | 10.495 | 1259 | 1268 | 1277 | rBV5 | 2257369 | 5667338 | 15.73% | 0.605% |
| 51 | 10.595 | 1279 | 1285 | 1293 | rVV2 | 7421223 | 12868295 | 35.72% | 1.374% |
| 52 | 10.683 | 1293 | 1300 | 1306 | rVB2 | 5851087 | 11517872 | 31.97% | 1.230% |
| 53 | 10.778 | 1312 | 1316 | 1322 | rVB | 2283798 | 3367299 | 9.35% | 0.360% |
| 54 | 11.054 | 1357 | 1363 | 1369 | rBV2 | 2449187 | 4315931 | 11.98% | 0.461% |
| 55 | 11.536 | 1440 | 1445 | 1452 | rVB3 | 1968080 | 3620133 | 10.05% | 0.387% |
| 56 | 11.648 | 1459 | 1464 | 1468 | rBV | 2850485 | 4746213 | 13.17% | 0.507% |
| 57 | 12.048 | 1527 | 1532 | 1543 | rVB | 6334380 | 9793951 | 27.18% | 1.046% |
| 58 | 12.301 | 1570 | 1575 | 1581 | rVB | 6615156 | 10616021 | 29.47% | 1.134% |
| 59 | 12.519 | 1606 | 1612 | 1620 | rVB | 4240174 | 7508581 | 20.84% | 0.802% |
| 60 | 12.683 | 1635 | 1640 | 1644 | rBV2 | 1790347 | 3031274 | 8.41% | 0.324% |
| 61 | 12.742 | 1644 | 1650 | 1655 | rVB3 | 1388088 | 3218568 | 8.93% | 0.344% |
| 62 | 12.954 | 1682 | 1686 | 1690 | rBV | 1901402 | 3008726 | 8.35% | 0.321% |
| 63 | 13.007 | 1690 | 1695 | 1700 | rVB | 6193568 | 8942660 | 24.82% | 0.955% |
| 64 | 13.101 | 1707 | 1711 | 1715 | rVB | 2184533 | 2986261 | 8.29% | 0.319% |
| 65 | 13.307 | 1738 | 1746 | 1756 | rBV2 | 11307182 | 23709329 | 65.81% | 2.532% |
| 66 | 13.389 | 1757 | 1760 | 1769 | rVB5 | 1603151 | 3820966 | 10.61% | 0.408% |
| 67 | 13.542 | 1777 | 1786 | 1791 | rBV | 11509540 | 20280432 | 56.29% | 2.166% |
| 68 | 13.595 | 1791 | 1795 | 1799 | rBV | 1712259 | 2934158 | 8.14% | 0.313% |
| 69 | 13.642 | 1799 | 1803 | 1806 | rBV2 | 2765687 | 4746859 | 13.18% | 0.507% |
| 70 | 13.807 | 1825 | 1831 | 1837 | rBV4 | 7244597 | 19312909 | 53.61% | 2.063% |
| 71 | 13.866 | 1837 | 1841 | 1849 | rVV4 | 4903705 | 12107529 | 33.61% | 1.293% |
| 72 | 13.960 | 1849 | 1857 | 1861 | rVV2 | 10500577 | 17922080 | 49.75% | 1.914% |
| 73 | 14.001 | 1861 | 1864 | 1868 | rVV2 | 4057281 | 5341837 | 14.83% | 0.571% |
| 74 | 14.077 | 1874 | 1877 | 1881 | rVB2 | 3077013 | 3811278 | 10.58% | 0.407% |
| 75 | 14.313 | 1912 | 1917 | 1922 | rVB4 | 4423474 | 8406136 | 23.33% | 0.898% |
| 76 | 14.377 | 1923 | 1928 | 1932 | rBV | 10514059 | 16657809 | 46.24% | 1.779% |

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042123\
Data File : BM039565.D
Acq On : 21 Apr 2023 23:29
Operator : CG/JU
Sample : 02386-11
Misc :
ALS Vial : 13 Sample Multiplier: 1

**Instrument :** BNA_M
**ClientSampleId :** SPA-3WC-23-04

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_M\Methods\8270-BM041723.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 77 | 14.448 | 1937 | 1940 | 1943 | rVV4 | 2958943 | 4654142 | 12.92% | 0.497% |
| 78 | 14.489 | 1944 | 1947 | 1953 | rVB3 | 3872768 | 5807919 | 16.12% | 0.620% |
| 79 | 14.883 | 2009 | 2014 | 2019 | rVB2 | 4487319 | 7235357 | 20.08% | 0.773% |
| 80 | 14.995 | 2030 | 2033 | 2042 | rVB4 | 2935259 | 5290772 | 14.69% | 0.565% |
| 81 | 15.271 | 2075 | 2080 | 2082 | rBV3 | 1743914 | 3231872 | 8.97% | 0.345% |
| 82 | 15.330 | 2086 | 2090 | 2095 | rVB | 7376174 | 10089597 | 28.01% | 1.078% |
| 83 | 15.495 | 2113 | 2118 | 2127 | rBV | 5655414 | 10528440 | 29.22% | 1.124% |
| 84 | 15.742 | 2154 | 2160 | 2164 | rBV2 | 6013114 | 10709680 | 29.73% | 1.144% |
| 85 | 15.783 | 2164 | 2167 | 2171 | rVB2 | 2815597 | 3877096 | 10.76% | 0.414% |
| 86 | 15.948 | 2191 | 2195 | 2198 | rVV3 | 2769669 | 4885798 | 13.56% | 0.522% |
| 87 | 16.001 | 2198 | 2204 | 2208 | rVB | 13917157 | 24696120 | 68.55% | 2.638% |
| 88 | 16.107 | 2218 | 2222 | 2228 | rBV | 7230373 | 9979884 | 27.70% | 1.066% |
| 89 | 16.207 | 2234 | 2239 | 2240 | rBV | 9985964 | 14140302 | 39.25% | 1.510% |
| 90 | 16.336 | 2255 | 2261 | 2266 | rBV | 8566983 | 16778660 | 46.57% | 1.792% |
| 91 | 16.395 | 2266 | 2271 | 2277 | rVB2 | 9381745 | 18241695 | 50.63% | 1.948% |
| 92 | 16.454 | 2277 | 2281 | 2285 | rBV2 | 9456557 | 14856191 | 41.24% | 1.587% |
| 93 | 16.501 | 2285 | 2289 | 2291 | rBV | 3638203 | 4103853 | 11.39% | 0.438% |
| 94 | 16.542 | 2291 | 2296 | 2298 | rBV3 | 12005978 | 21778259 | 60.45% | 2.326% |
| 95 | 16.654 | 2310 | 2315 | 2321 | rVB3 | 5835248 | 10890664 | 30.23% | 1.163% |
| 96 | 16.718 | 2321 | 2326 | 2329 | rBV | 7094527 | 11792241 | 32.73% | 1.259% |
| 97 | 16.995 | 2369 | 2373 | 2377 | rBV | 5888116 | 8878711 | 24.64% | 0.948% |
| 98 | 17.042 | 2377 | 2381 | 2389 | rVB | 6274735 | 11130807 | 30.90% | 1.189% |
| 99 | 17.736 | 2495 | 2499 | 2506 | rVB | 5178088 | 8230238 | 22.84% | 0.879% |
| 100 | 18.689 | 2658 | 2661 | 2670 | rVB4 | 4739278 | 8826183 | 24.50% | 0.943% |

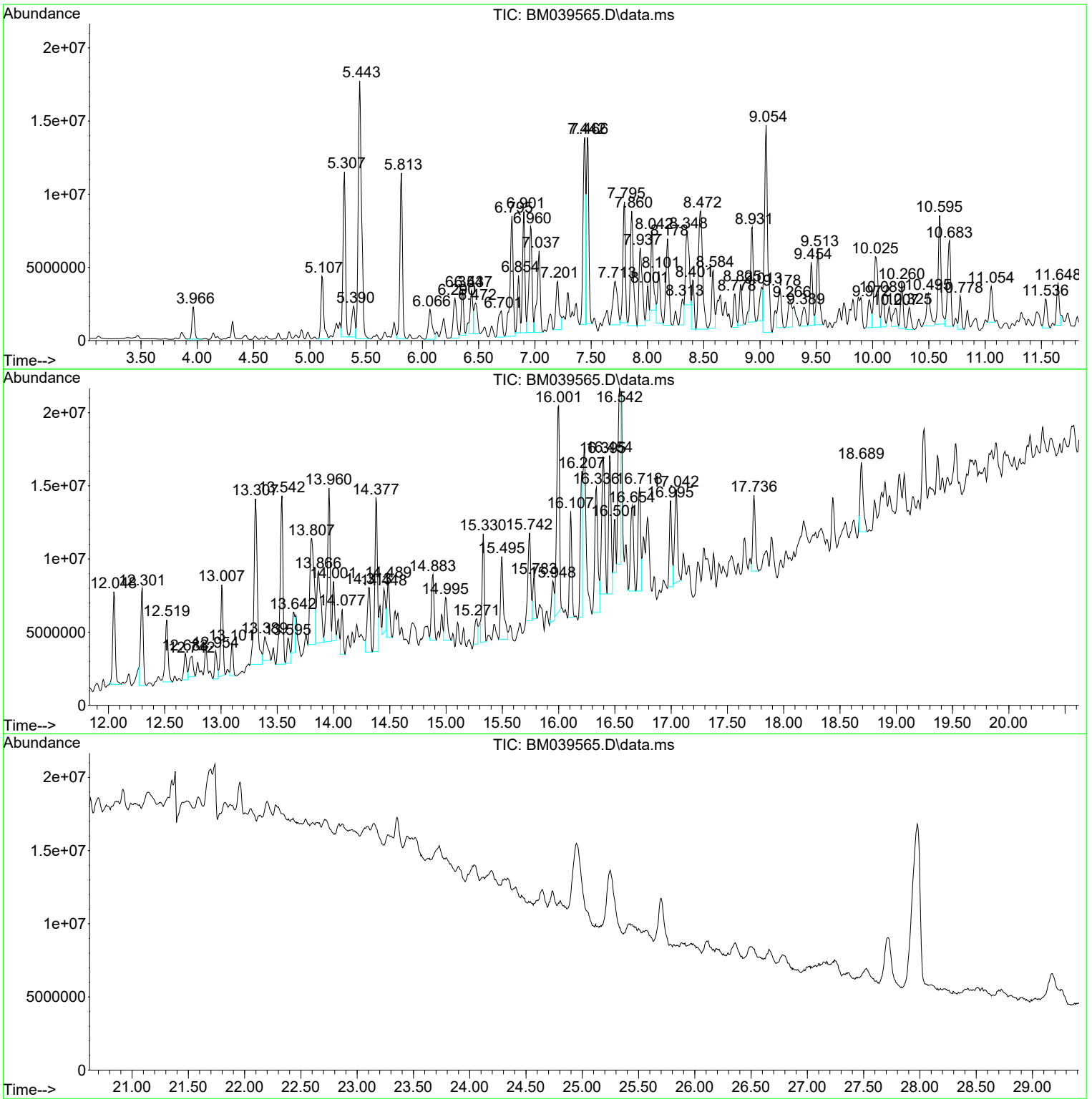
Sum of corrected areas: 936296321

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042123\
Data File : BM039565.D
Acq On : 21 Apr 2023 23:29
Operator : CG/JU
Sample : 02386-11
Misc :
ALS Vial : 13 Sample Multiplier: 1

**Instrument :**
BNA_M
**ClientSampleId :**
SPA-3WC-23-04

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM041723.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_M\Data\BM042123\
Data File : BM039565.D
Acq On : 21 Apr 2023 23:29
Operator : CG/JU
Sample : 02386-11
Misc :
ALS Vial : 13 Sample Multiplier: 1

**Instrument :**
BNA_M
**ClientSampleId :**
SPA-3WC-23-04

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM041723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

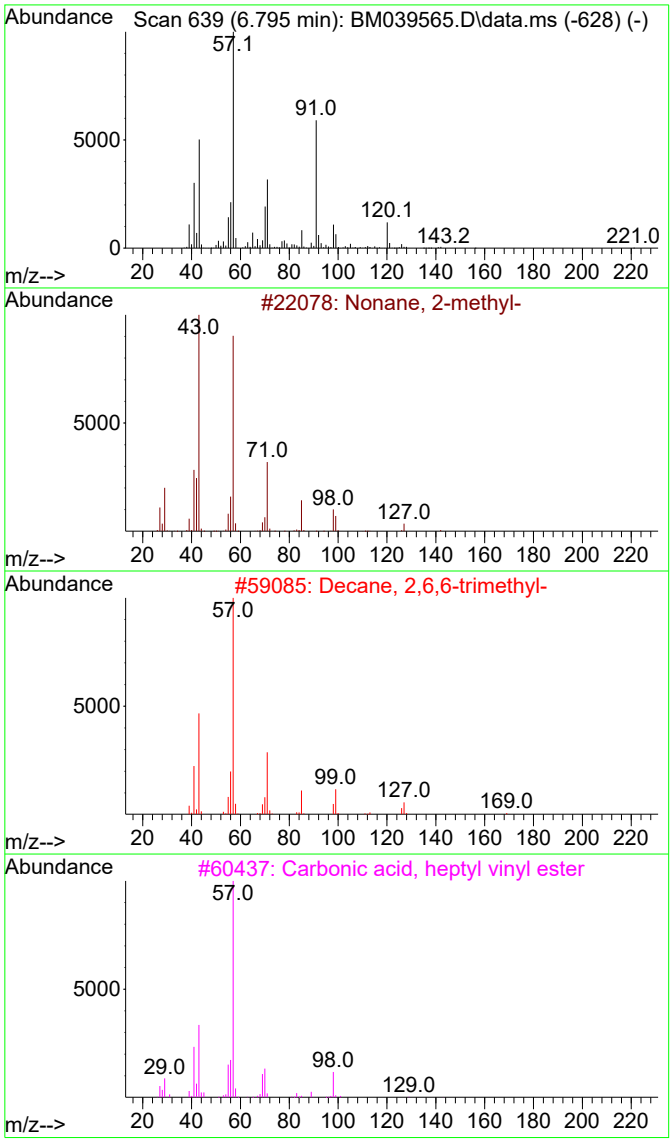
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

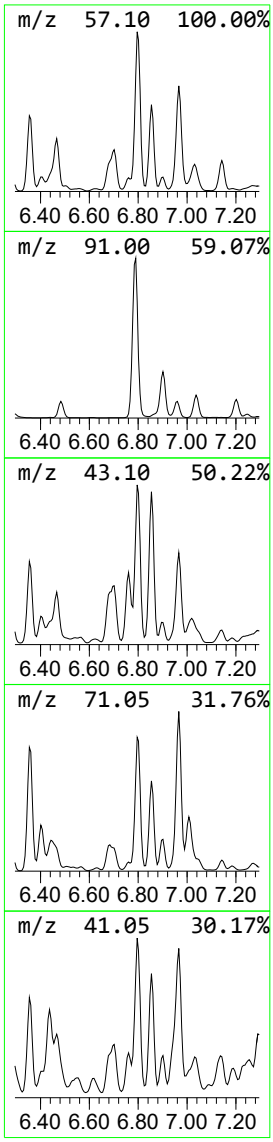
**********

Peak Number 3 unknown6.795 Concentration Rank 19

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 6.795 | 25.81 ng | 16288200 | 1,4-Dichlorobenzene-d4 | 7.825 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Nonane, 2-methyl- | 142 | C10H22 | 000871-83-0 | 49 |
| 2 | Decane, 2,6,6-trimethyl- | 184 | C13H28 | 062108-24-1 | 47 |
| 3 | Carbonic acid, heptyl vinyl ester | 186 | C10H18O3 | 1010383-25-4 | 43 |
| 4 | Nonane, 4-methyl- | 142 | C10H22 | 017301-94-9 | 43 |
| 5 | Undecane, 3-methyl- | 170 | C12H26 | 001002-43-3 | 43 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042123\
Data File : BM039565.D
Acq On : 21 Apr 2023 23:29
Operator : CG/JU
Sample : 02386-11
Misc :
ALS Vial : 13 Sample Multiplier: 1

**Instrument :**
BNA_M
**ClientSampleId :**
SPA-3WC-23-04

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM041723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

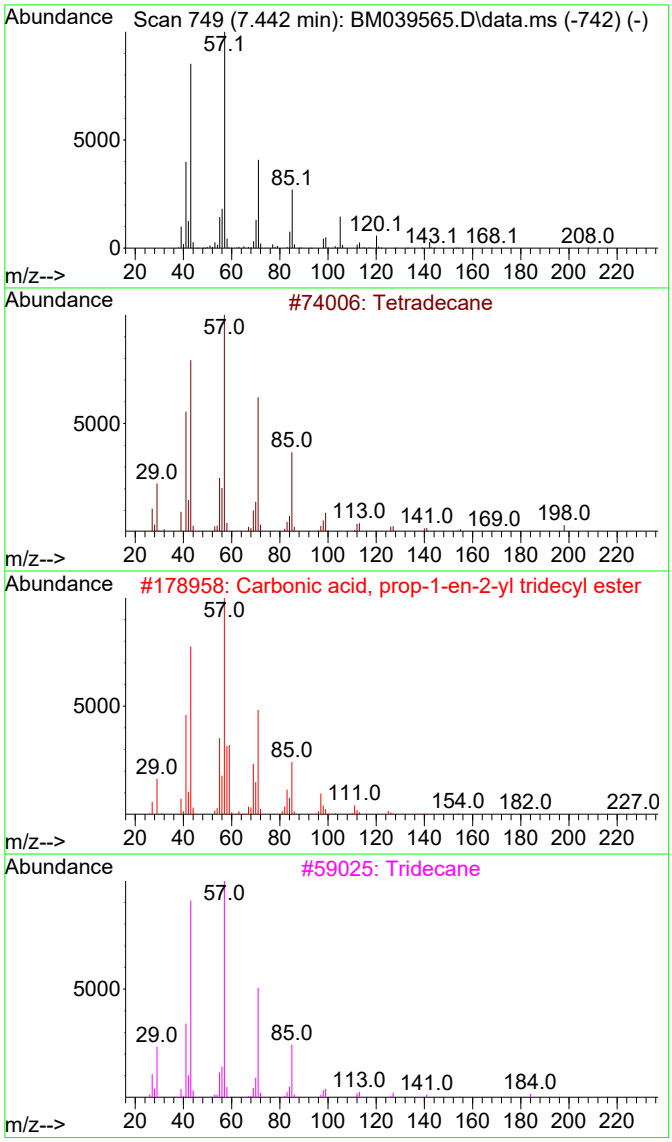
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

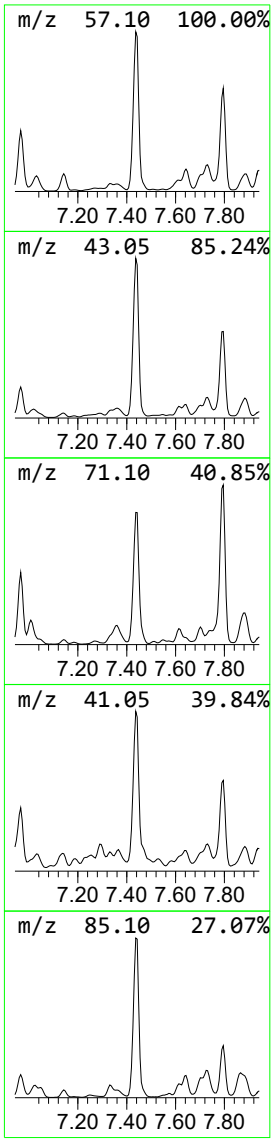
**********

Peak Number 4 Tetradecane Concentration Rank 8

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 7.442 | 35.17 ng | 22195400 | 1,4-Dichlorobenzene-d4 | 7.825 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Tetradecane | 198 | C14H30 | 000629-59-4 | 80 |
| 2 | Carbonic acid, prop-1-en-2-yl tr... | 284 | C17H32O3 | 1000382-90-8 | 72 |
| 3 | Tridecane | 184 | C13H28 | 000629-50-5 | 64 |
| 4 | Decane | 142 | C10H22 | 000124-18-5 | 64 |
| 5 | Decane, 2,3,6-trimethyl- | 184 | C13H28 | 062238-12-4 | 59 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042123\
Data File : BM039565.D
Acq On : 21 Apr 2023 23:29
Operator : CG/JU
Sample : 02386-11
Misc :
ALS Vial : 13 Sample Multiplier: 1

**Instrument :**
BNA_M
**ClientSampleId :**
SPA-3WC-23-04

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM041723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

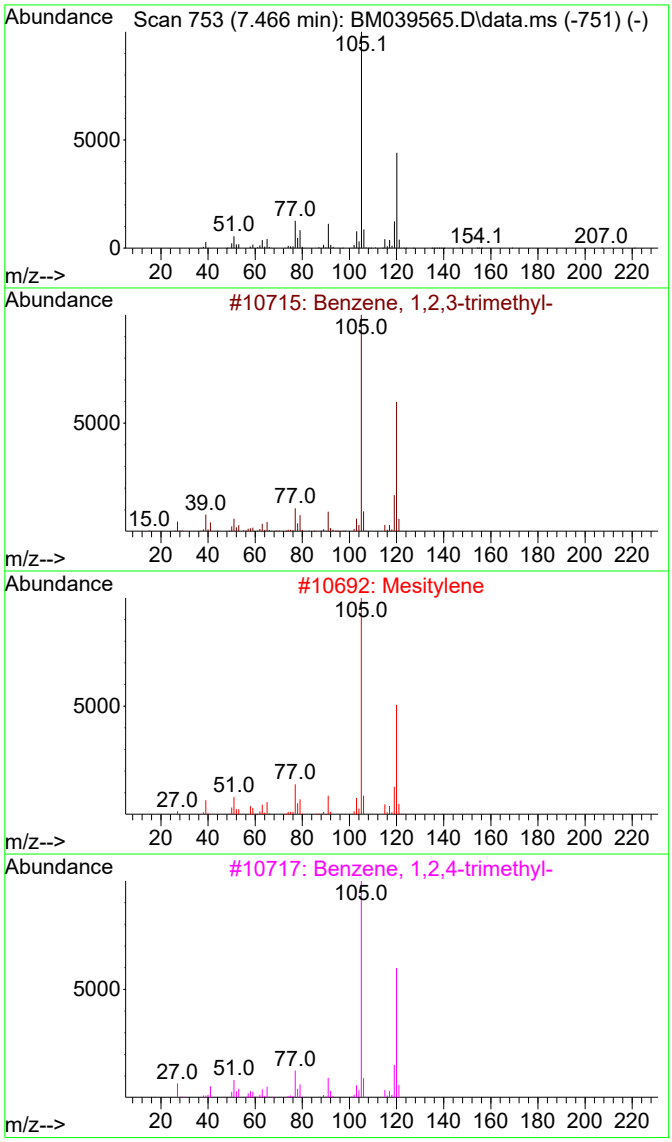
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

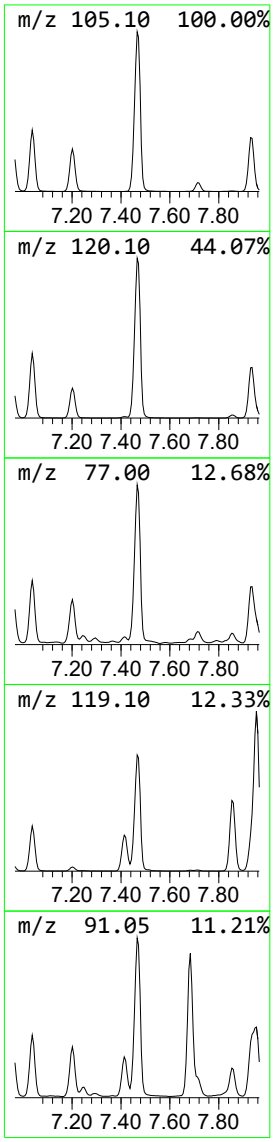
**********

Peak Number 5 Benzene, 1,2,3-trimethyl- Concentration Rank 16

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 7.466 | 27.13 ng | 17123000 | 1,4-Dichlorobenzene-d4 | 7.825 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, 1,2,3-trimethyl- | 120 | C9H12 | 000526-73-8 | 97 |
| 2 | Mesitylene | 120 | C9H12 | 000108-67-8 | 97 |
| 3 | Benzene, 1,2,4-trimethyl- | 120 | C9H12 | 000095-63-6 | 94 |
| 4 | Benzene, 1-ethyl-2-methyl- | 120 | C9H12 | 000611-14-3 | 90 |
| 5 | Benzene, 1-ethyl-4-methyl- | 120 | C9H12 | 000622-96-8 | 90 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042123\
Data File : BM039565.D
Acq On : 21 Apr 2023 23:29
Operator : CG/JU
Sample : 02386-11
Misc :
ALS Vial : 13 Sample Multiplier: 1

**Instrument :**
BNA_M
**ClientSampleId :**
SPA-3WC-23-04

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM041723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

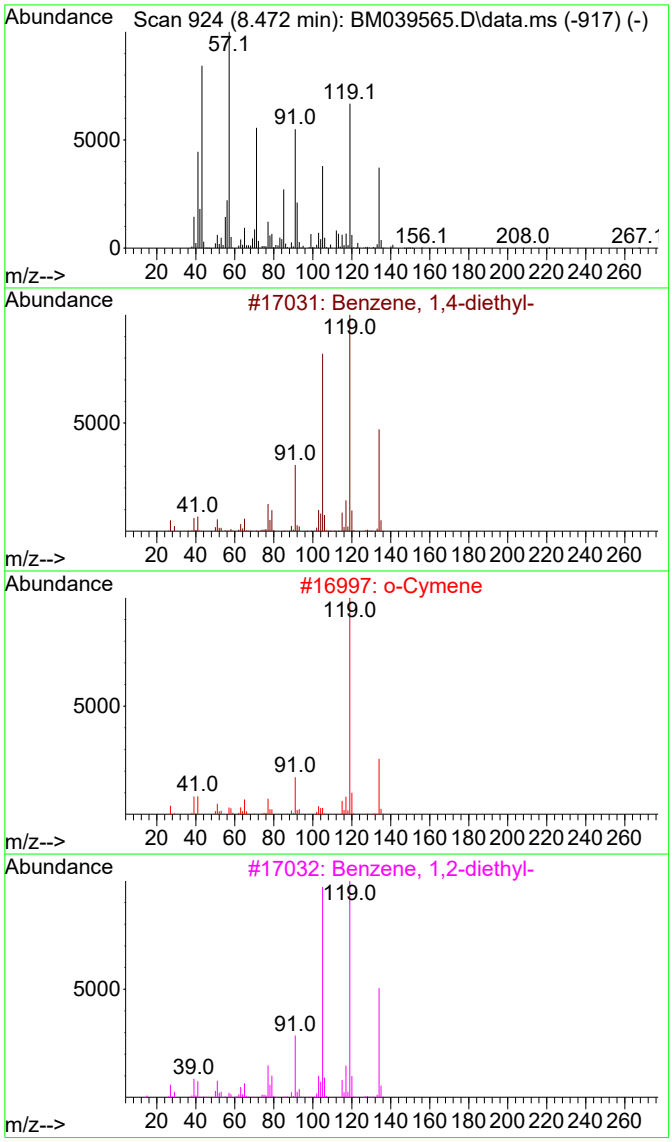
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

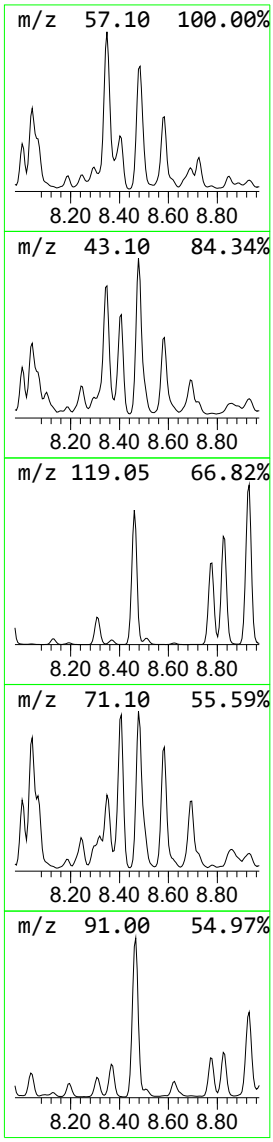
**********

Peak Number 6 unknown8.472 Concentration Rank 11

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 8.472 | 32.43 ng | 20468900 | 1,4-Dichlorobenzene-d4 | 7.825 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, 1,4-diethyl- | 134 | C10H14 | 000105-05-5 | 42 |
| 2 | o-Cymene | 134 | C10H14 | 000527-84-4 | 38 |
| 3 | Benzene, 1,2-diethyl- | 134 | C10H14 | 000135-01-3 | 38 |
| 4 | Benzene, 1-ethyl-3,5-dimethyl- | 134 | C10H14 | 000934-74-7 | 38 |
| 5 | Benzene, 1-methyl-3-(1-methyleth... | 134 | C10H14 | 000535-77-3 | 38 |



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Data File : BM039565.D
Acq On : 21 Apr 2023 23:29
Operator : CG/JU
Sample : 02386-11
Misc :
ALS Vial : 13 Sample Multiplier: 1

**Instrument :**
BNA_M
**ClientSampleId :**
SPA-3WC-23-04

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM041723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

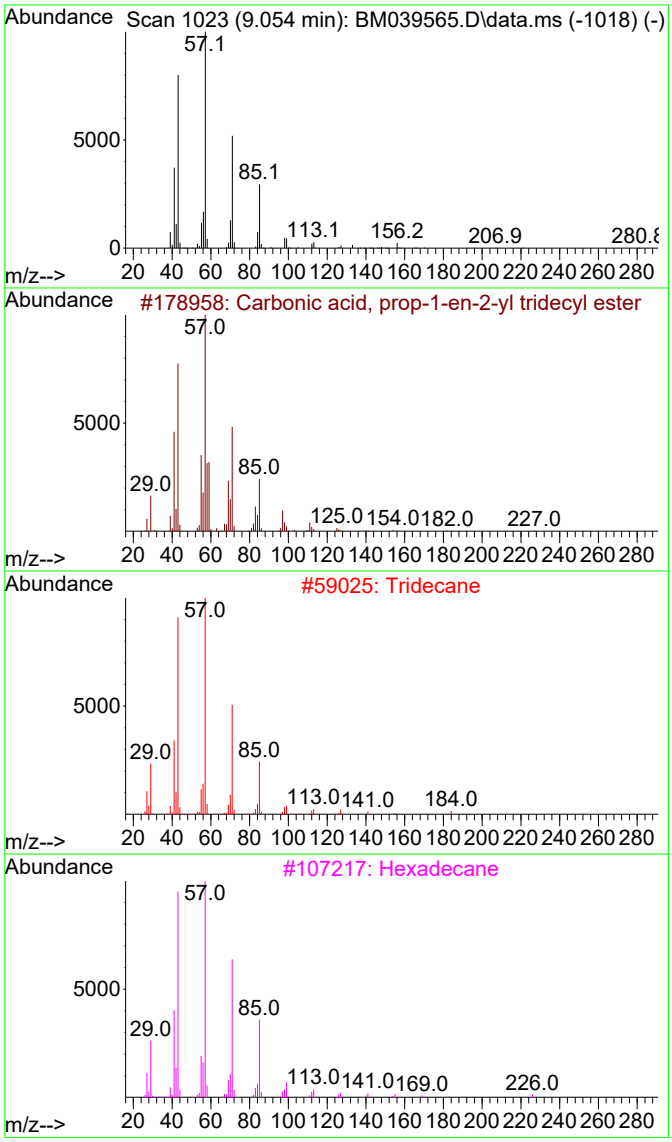
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

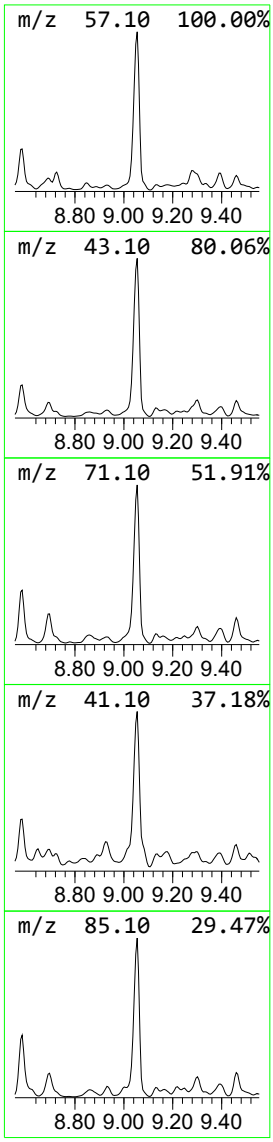
**********

Peak Number 7 Carbonic acid, prop-1-en-2-... Concentration Rank 5

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 9.054 | 44.30 ng | 27958600 | 1,4-Dichlorobenzene-d4 | 7.825 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Carbonic acid, prop-1-en-2-yl tr... | 284 | C17H32O3 | 1000382-90-8 | 90 |
| 2 | Tridecane | 184 | C13H28 | 000629-50-5 | 86 |
| 3 | Hexadecane | 226 | C16H34 | 000544-76-3 | 83 |
| 4 | Carbonic acid, prop-1-en-2-yl te... | 298 | C18H34O3 | 1000382-90-9 | 83 |
| 5 | Tetradecane | 198 | C14H30 | 000629-59-4 | 83 |



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Data File : BM039565.D
Acq On : 21 Apr 2023 23:29
Operator : CG/JU
Sample : 02386-11
Misc :
ALS Vial : 13 Sample Multiplier: 1

**Instrument :**
BNA_M
**ClientSampleId :**
SPA-3WC-23-04

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM041723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

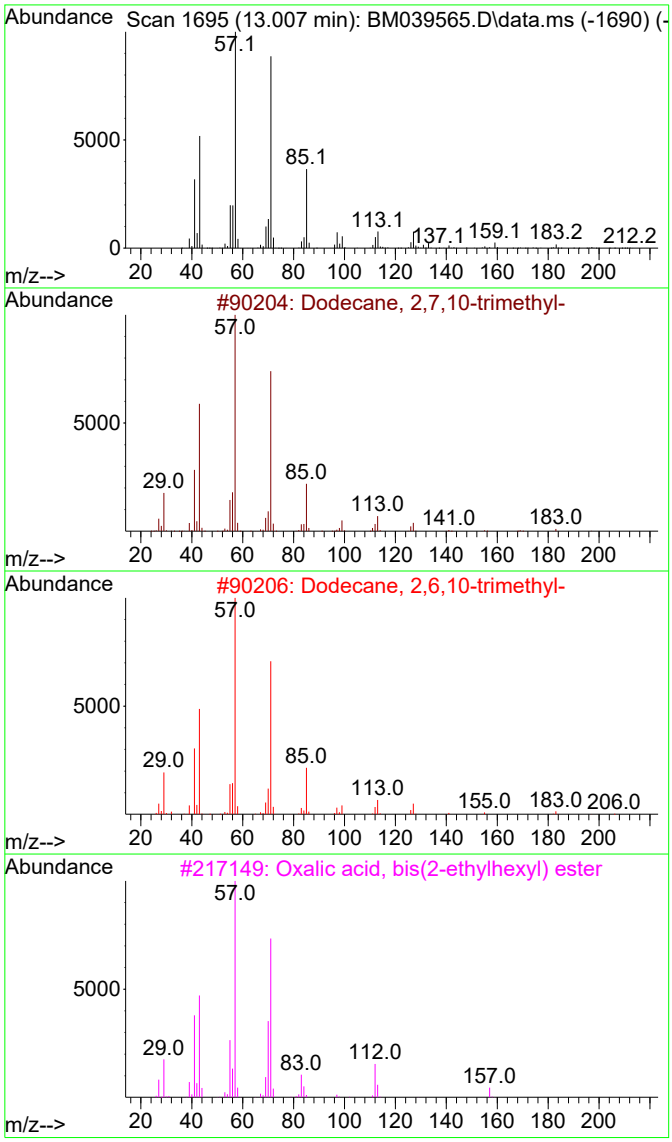
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

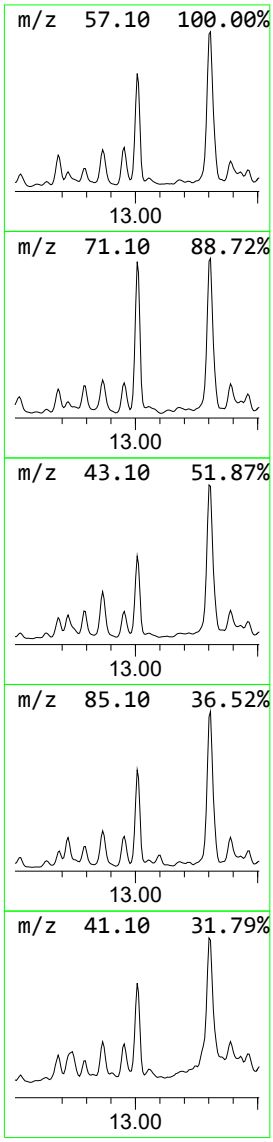
**********

Peak Number 8 Dodecane, 2,7,10-trimethyl- Concentration Rank 12

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.007 | 30.79 ng | 8942660 | Acenaphthene-d10 | 14.483 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Dodecane, 2,7,10-trimethyl- | 212 | C15H32 | 074645-98-0 | 91 |
| 2 | Dodecane, 2,6,10-trimethyl- | 212 | C15H32 | 003891-98-3 | 90 |
| 3 | Oxalic acid, bis(2-ethylhexyl) e... | 314 | C18H34O4 | 1000309-39-0 | 53 |
| 4 | Sulfurous acid, 2-ethylhexyl pen... | 264 | C13H28O3S | 1000309-18-9 | 53 |
| 5 | Nonane, 5-methyl-5-propyl- | 184 | C13H28 | 017312-75-3 | 53 |



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Acq On : 21 Apr 2023 23:29
Operator : CG/JU
Sample : 02386-11
Misc :
ALS Vial : 13 Sample Multiplier: 1

**Instrument :**
BNA_M
**ClientSampleId :**
SPA-3WC-23-04

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM041723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

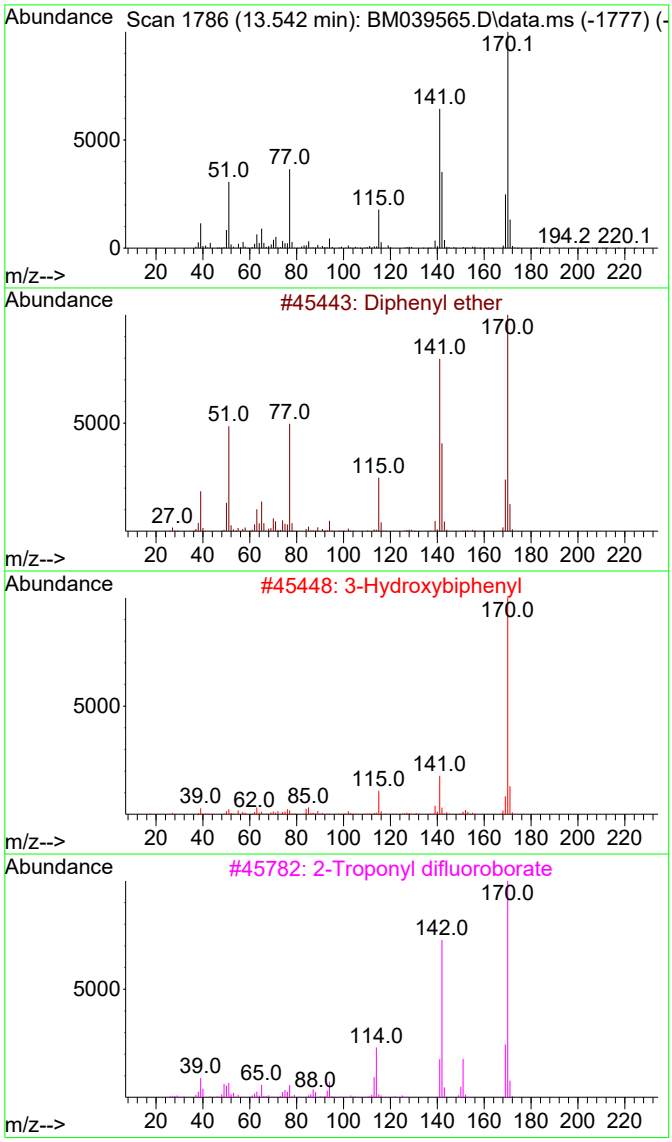
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

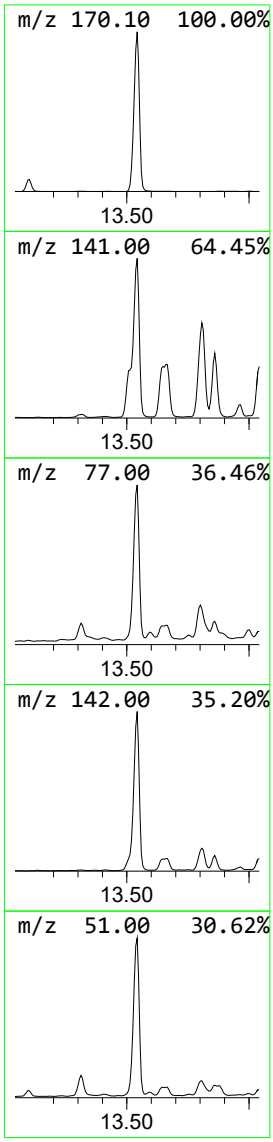
**********

Peak Number 9 Diphenyl ether Concentration Rank 1

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.542 | 69.84 ng | 20280400 | Acenaphthene-d10 | 14.483 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Diphenyl ether | 170 | C12H10O | 000101-84-8 | 96 |
| 2 | 3-Hydroxybiphenyl | 170 | C12H10O | 000580-51-8 | 50 |
| 3 | 2-Troponyl difluoroborate | 170 | C7H5BF2O2 | 022502-10-9 | 47 |
| 4 | 8-Hydroxyquinoline-2-carbonitrile | 170 | C10H6N2O | 006759-78-0 | 43 |
| 5 | Phenol, O-(ethylsulfinyl)- | 170 | C8H10O2S | 029634-40-0 | 40 |



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Data File : BM039565.D
Acq On : 21 Apr 2023 23:29
Operator : CG/JU
Sample : 02386-11
Misc :
ALS Vial : 13 Sample Multiplier: 1

**Instrument :**
BNA_M
**ClientSampleId :**
SPA-3WC-23-04

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM041723.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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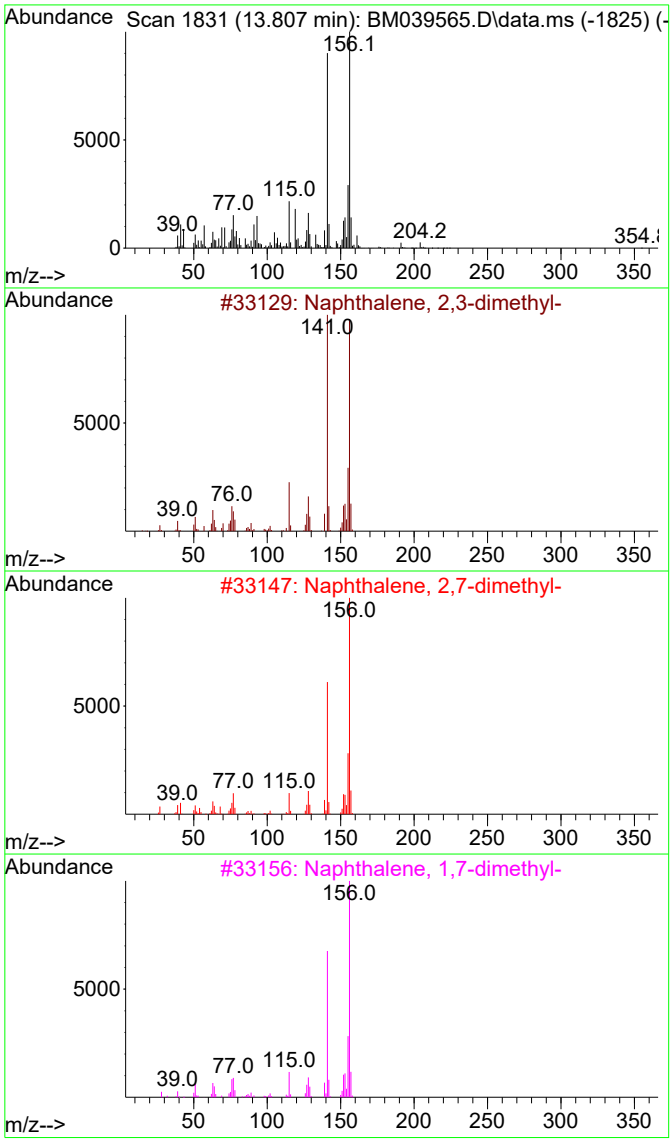
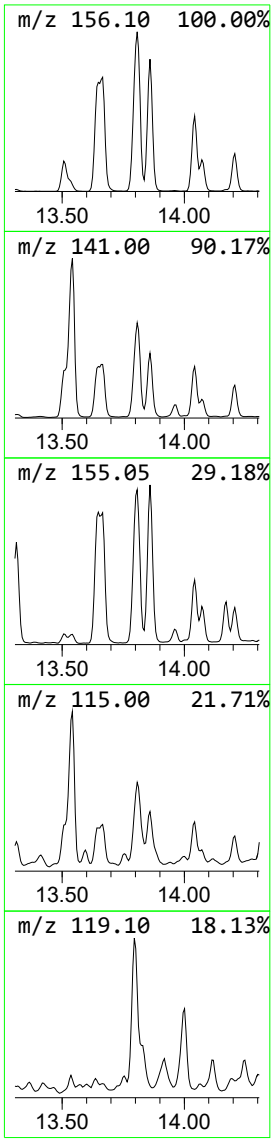
Peak Number 10 Naphthalene, 2,3-dimethyl- Concentration Rank 2

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.807 | 66.51 ng | 19312900 | Acenaphthene-d10 | 14.483 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Naphthalene, 2,3-dimethyl- | 156 | C12H12 | 000581-40-8 | 97 |
| 2 | Naphthalene, 2,7-dimethyl- | 156 | C12H12 | 000582-16-1 | 97 |
| 3 | Naphthalene, 1,7-dimethyl- | 156 | C12H12 | 000575-37-1 | 97 |
| 4 | Naphthalene, 1,3-dimethyl- | 156 | C12H12 | 000575-41-7 | 96 |
| 5 | Naphthalene, 1,4-dimethyl- | 156 | C12H12 | 000571-58-4 | 96 |



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Acq On : 21 Apr 2023 23:29
Operator : CG/JU
Sample : 02386-11
Misc :
ALS Vial : 13 Sample Multiplier: 1

**Instrument :**
BNA_M
**ClientSampleId :**
SPA-3WC-23-04

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM041723.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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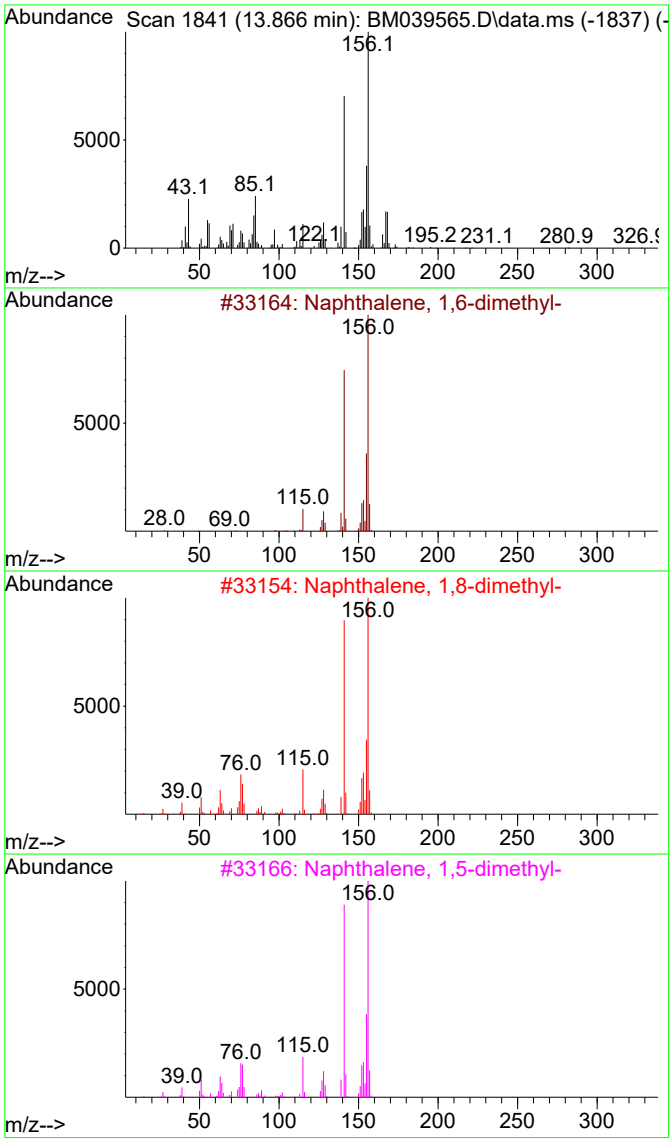
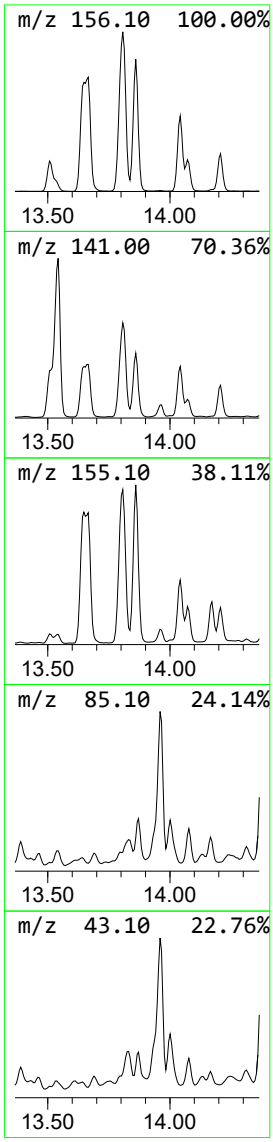
Peak Number 11 Naphthalene, 1,6-dimethyl- Concentration Rank 6

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.866 | 41.69 ng | 12107500 | Acenaphthene-d10 | 14.483 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Naphthalene, 1,6-dimethyl- | 156 | C12H12 | 000575-43-9 | 95 |
| 2 | Naphthalene, 1,8-dimethyl- | 156 | C12H12 | 000569-41-5 | 94 |
| 3 | Naphthalene, 1,5-dimethyl- | 156 | C12H12 | 000571-61-9 | 94 |
| 4 | Naphthalene, 1,7-dimethyl- | 156 | C12H12 | 000575-37-1 | 93 |
| 5 | Naphthalene, 2,7-dimethyl- | 156 | C12H12 | 000582-16-1 | 89 |



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Data File : BM039565.D
Acq On : 21 Apr 2023 23:29
Operator : CG/JU
Sample : 02386-11
Misc :
ALS Vial : 13 Sample Multiplier: 1

**Instrument :**
BNA_M
**ClientSampleId :**
SPA-3WC-23-04

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM041723.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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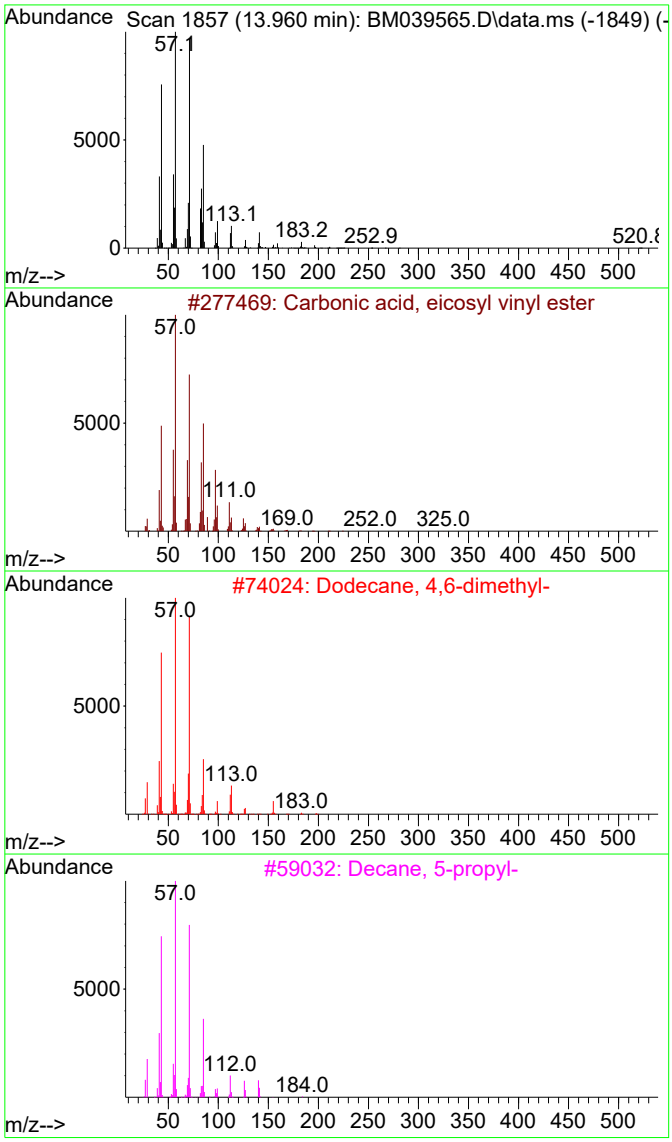
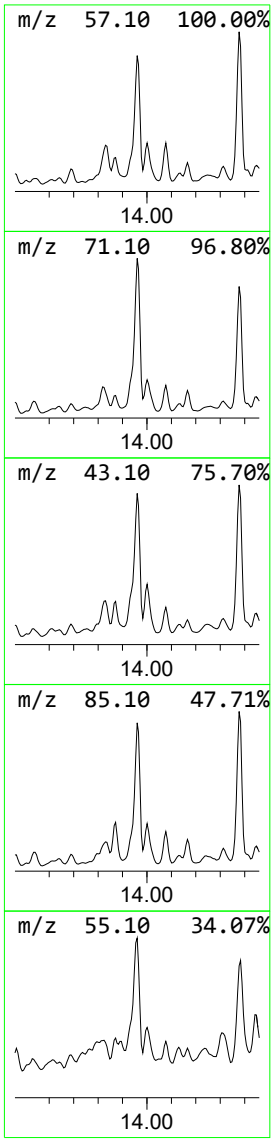
Peak Number 12 Carbonic acid, eicosyl viny... Concentration Rank 3

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.960 | 61.72 ng | 17922100 | Acenaphthene-d10 | 14.483 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Carbonic acid, eicosyl vinyl ester | 368 | C23H44O3 | 1000382-54-3 | 64 |
| 2 | Dodecane, 4,6-dimethyl- | 198 | C14H30 | 061141-72-8 | 64 |
| 3 | Decane, 5-propyl- | 184 | C13H28 | 017312-62-8 | 62 |
| 4 | 1,3-Dimethylcyclopentanol | 114 | C7H14O | 019550-46-0 | 52 |
| 5 | Oxalic acid, isohexyl neopentyl ... | 244 | C13H24O4 | 1000309-73-0 | 50 |



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Data File : BM039565.D
Acq On : 21 Apr 2023 23:29
Operator : CG/JU
Sample : 02386-11
Misc :
ALS Vial : 13 Sample Multiplier: 1

**Instrument :**
BNA_M
**ClientSampleId :**
SPA-3WC-23-04

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM041723.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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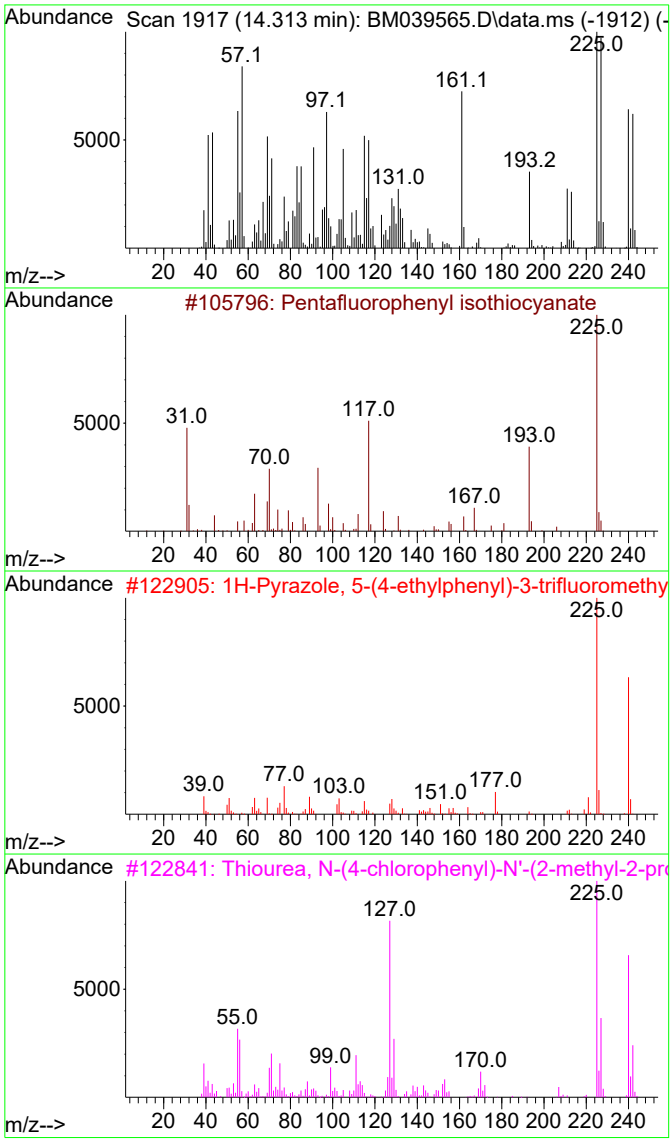
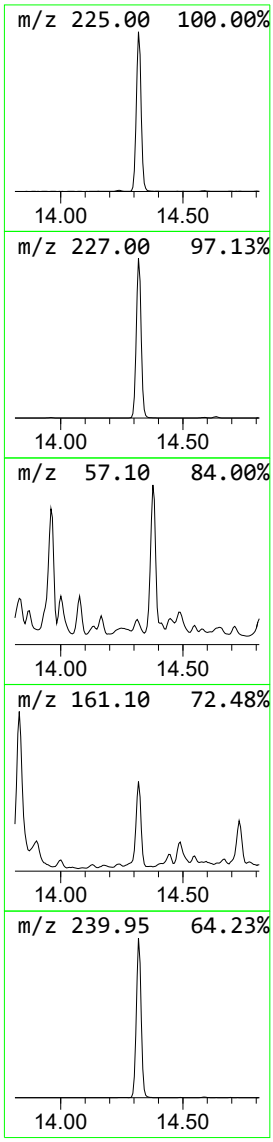
Peak Number 13 unknown14.313 Concentration Rank 14

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.313 | 28.95 ng | 8406140 | Acenaphthene-d10 | 14.483 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Pentafluorophenyl isothiocyanate | 225 | C7F5NS | 035923-79-6 | 38 |
| 2 | 1H-Pyrazole, 5-(4-ethylphenyl)-3... | 240 | C12H11F3N2 | 1000310-25-4 | 35 |
| 3 | Thiourea, N-(4-chlorophenyl)-N'-... | 240 | C11H13ClN2S | 1000351-06-5 | 30 |
| 4 | Benzothiazole, 2-(m-tolyl)- | 225 | C14H11NS | 1000302-53-5 | 25 |
| 5 | 7-Bromo-4-methoxy-indan-1-one | 240 | C10H9BrO2 | 1010202-16-3 | 15 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042123\
Data File : BM039565.D
Acq On : 21 Apr 2023 23:29
Operator : CG/JU
Sample : 02386-11
Misc :
ALS Vial : 13 Sample Multiplier: 1

**Instrument :**
BNA_M
**ClientSampleId :**
SPA-3WC-23-04

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM041723.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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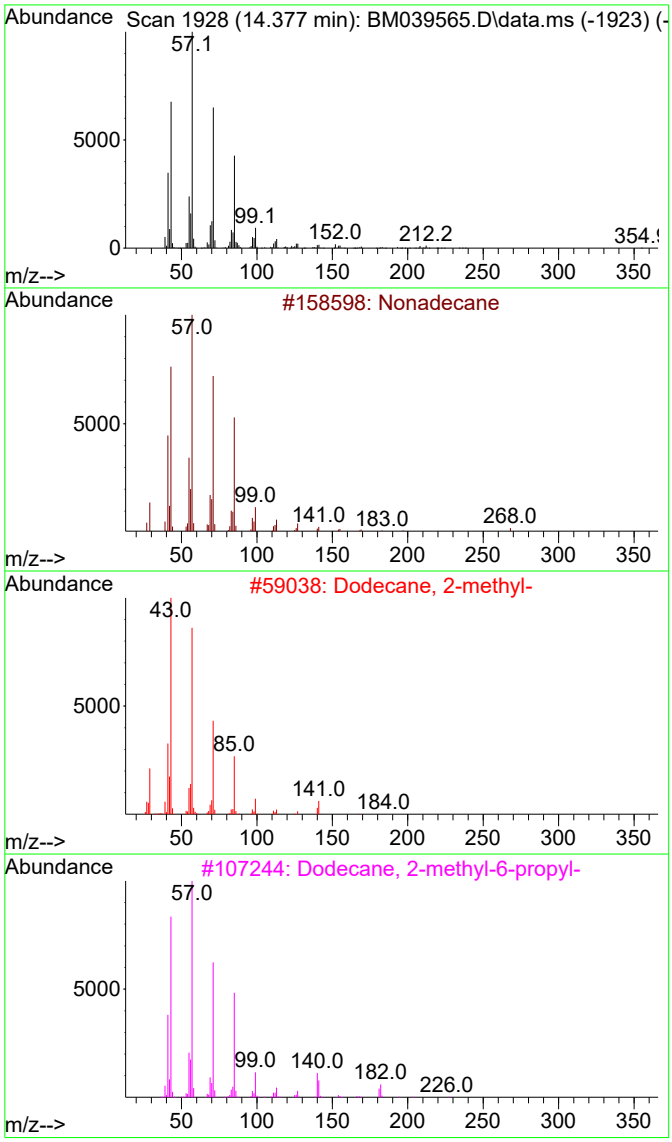
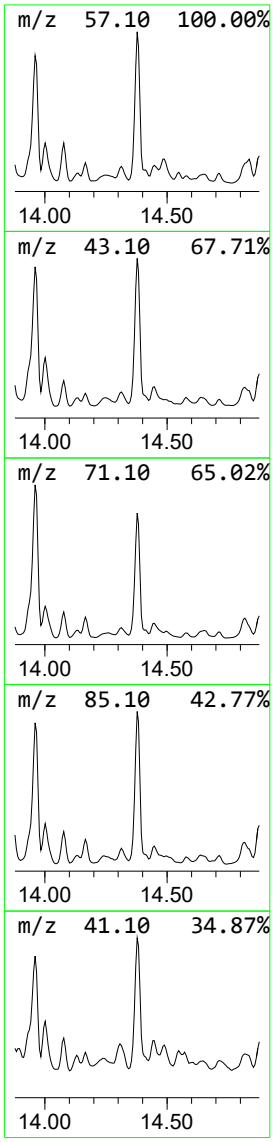
Peak Number 14 Nonadecane Concentration Rank 4

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 14.377 | 57.36 ng | 16657800 | Acenaphthene-d10 | 14.483 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Nonadecane | 268 | C19H40 | 000629-92-5 | 87 |
| 2 | Dodecane, 2-methyl- | 184 | C13H28 | 001560-97-0 | 86 |
| 3 | Dodecane, 2-methyl-6-propyl- | 226 | C16H34 | 055045-08-4 | 86 |
| 4 | Tetradecane | 198 | C14H30 | 000629-59-4 | 86 |
| 5 | Heptadecane | 240 | C17H36 | 000629-78-7 | 83 |



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Data File : BM039565.D
Acq On : 21 Apr 2023 23:29
Operator : CG/JU
Sample : 02386-11
Misc :
ALS Vial : 13 Sample Multiplier: 1

**Instrument :**
BNA_M
**ClientSampleId :**
SPA-3WC-23-04

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM041723.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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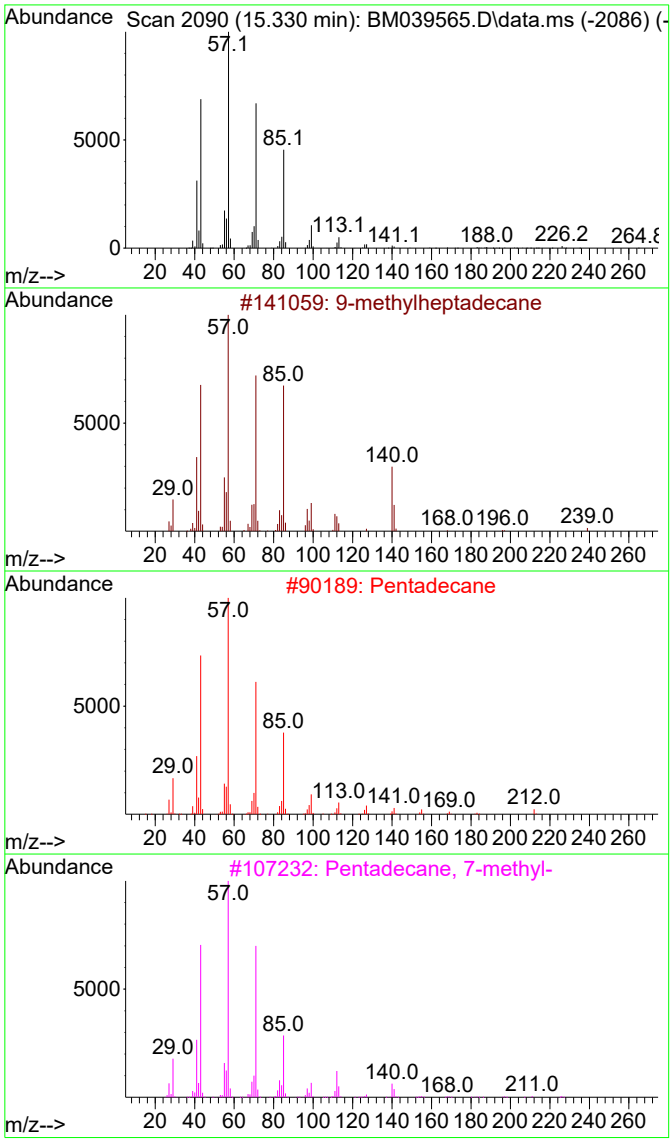
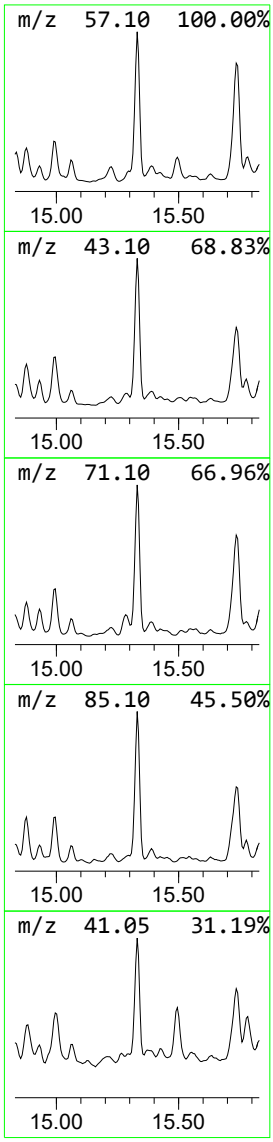
Peak Number 15 9-methylheptadecane Concentration Rank 9

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.330 | 34.74 ng | 10089600 | Acenaphthene-d10 | 14.483 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 9-methylheptadecane | 254 | C18H38 | 026741-18-4 | 90 |
| 2 | Pentadecane | 212 | C15H32 | 000629-62-9 | 90 |
| 3 | Pentadecane, 7-methyl- | 226 | C16H34 | 006165-40-8 | 87 |
| 4 | 3-Ethyl-2,6,10-trimethylundecane | 226 | C16H34 | 1000432-25-9 | 87 |
| 5 | Decane, 2,3,5-trimethyl- | 184 | C13H28 | 062238-11-3 | 83 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042123\
Data File : BM039565.D
Acq On : 21 Apr 2023 23:29
Operator : CG/JU
Sample : 02386-11
Misc :
ALS Vial : 13 Sample Multiplier: 1

**Instrument :**
BNA_M
**ClientSampleId :**
SPA-3WC-23-04

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM041723.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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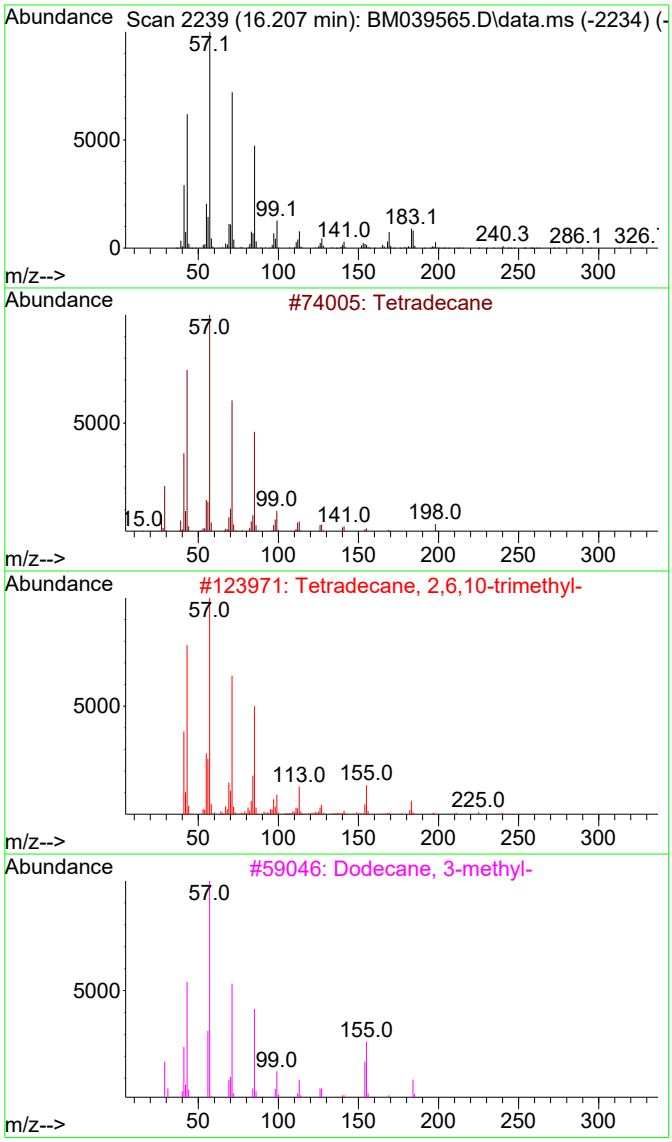
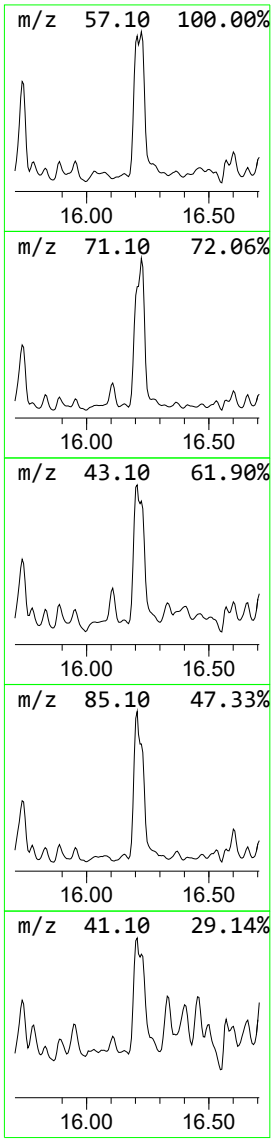
Peak Number 16 Tetradecane, 2,6,10-trimethyl- Concentration Rank 20

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.207 | 25.41 ng | 14140300 | Phenanthrene-d10 | 17.248 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Tetradecane | 198 | C14H30 | 000629-59-4 | 95 |
| 2 | Tetradecane, 2,6,10-trimethyl- | 240 | C17H36 | 014905-56-7 | 87 |
| 3 | Dodecane, 3-methyl- | 184 | C13H28 | 017312-57-1 | 87 |
| 4 | Tridecane, 1-iodo- | 310 | C13H27I | 035599-77-0 | 87 |
| 5 | Heptacosane | 380 | C27H56 | 000593-49-7 | 83 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042123\
Data File : BM039565.D
Acq On : 21 Apr 2023 23:29
Operator : CG/JU
Sample : 02386-11
Misc :
ALS Vial : 13 Sample Multiplier: 1

**Instrument :**
BNA_M
**ClientSampleId :**
SPA-3WC-23-04

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM041723.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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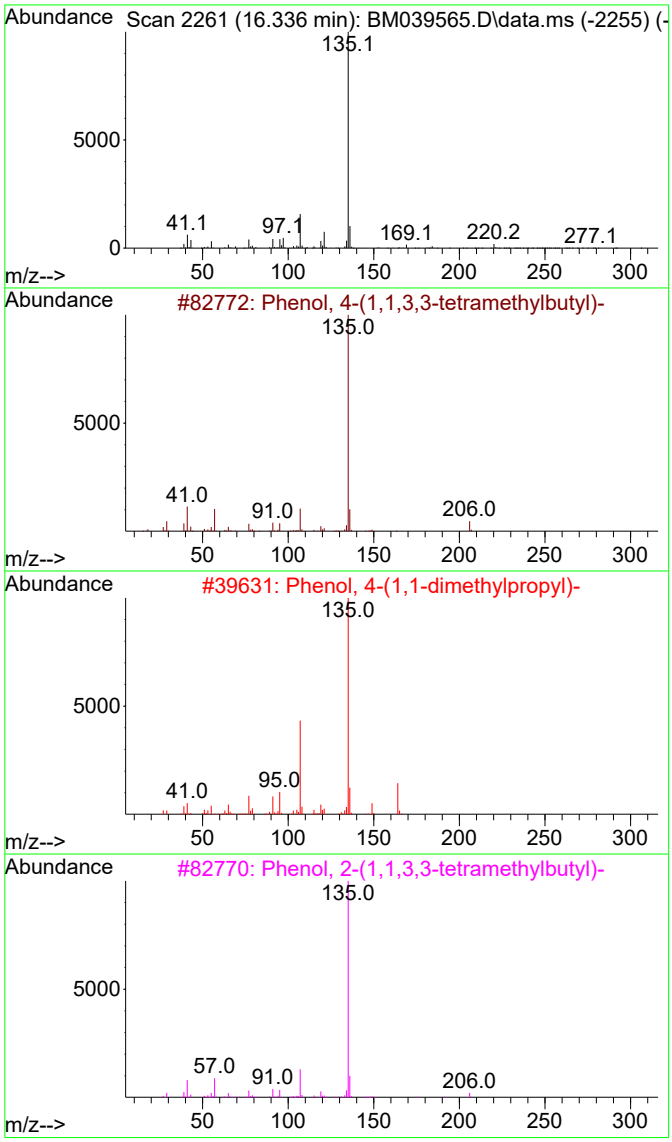
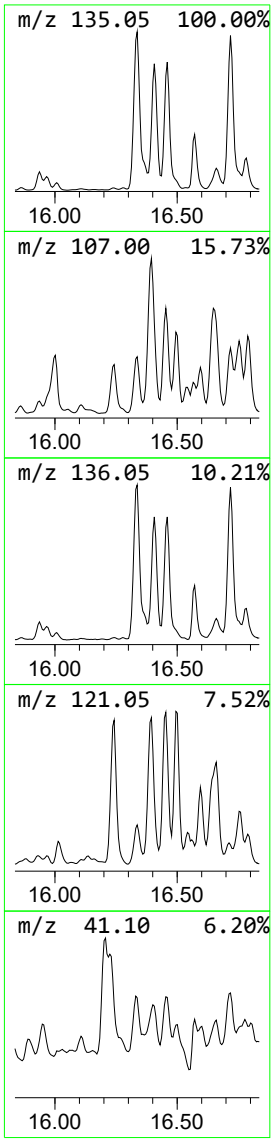
Peak Number 17 Phenol, 4-(1,1,3,3-tetramet... Concentration Rank 13

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.336 | 30.15 ng | 16778700 | Phenanthrene-d10 | 17.248 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, 4-(1,1,3,3-tetramethylbu... | 206 | C14H22O | 000140-66-9 | 72 |
| 2 | Phenol, 4-(1,1-dimethylpropyl)- | 164 | C11H16O | 000080-46-6 | 72 |
| 3 | Phenol, 2-(1,1,3,3-tetramethylbu... | 206 | C14H22O | 003884-95-5 | 72 |
| 4 | 4-tert-Butylphenyl acetate | 192 | C12H16O2 | 003056-64-2 | 72 |
| 5 | m-Anisoyl amide, N-(2-phenylethy... | 297 | C19H23NO2 | 1000451-73-8 | 64 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042123\
Data File : BM039565.D
Acq On : 21 Apr 2023 23:29
Operator : CG/JU
Sample : 02386-11
Misc :
ALS Vial : 13 Sample Multiplier: 1

**Instrument :**
BNA_M
**ClientSampleId :**
SPA-3WC-23-04

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM041723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

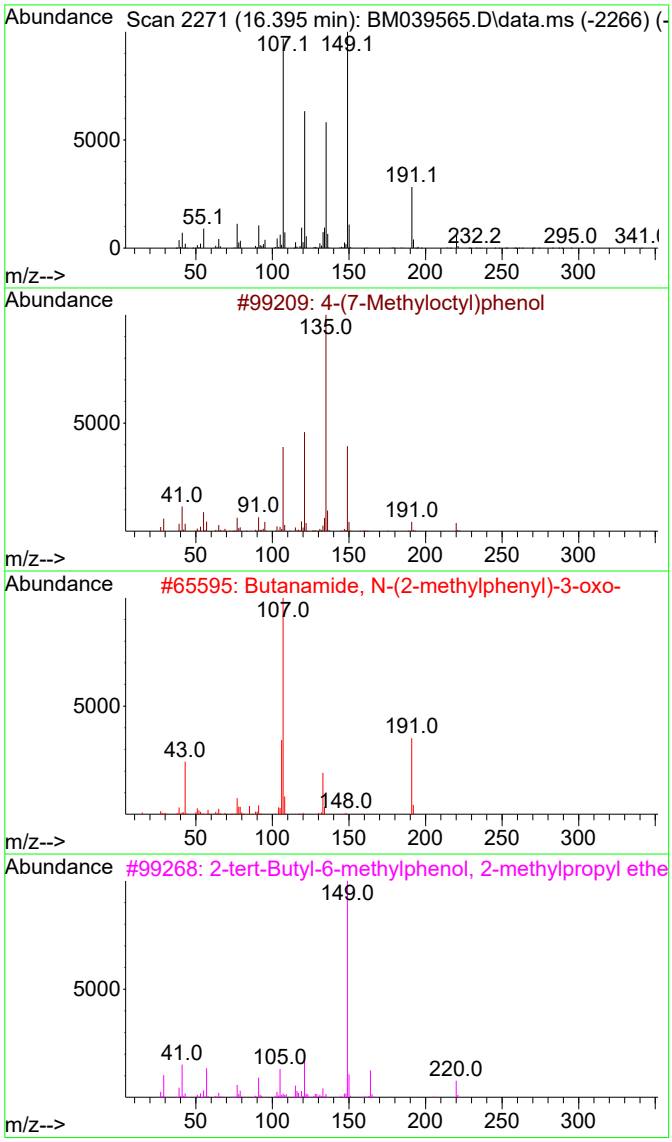
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

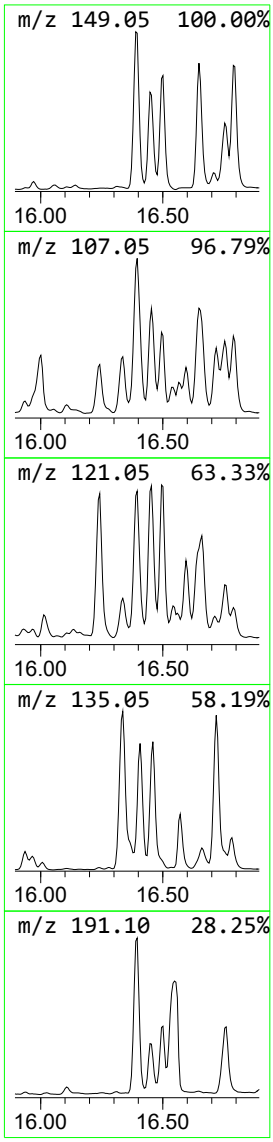
**********

Peak Number 18 unknown16.395 Concentration Rank 10

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.395 | 32.78 ng | 18241700 | Phenanthrene-d10 | 17.248 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 4-(7-Methyloctyl)phenol | 220 | C15H24O | 024518-48-7 | 46 |
| 2 | Butanamide, N-(2-methylphenyl)-3... | 191 | C11H13NO2 | 000093-68-5 | 30 |
| 3 | 2-tert-Butyl-6-methylphenol, 2-m... | 220 | C15H24O | 1000395-19-4 | 30 |
| 4 | Pyridine, pentamethyl- | 149 | C10H15N | 003748-83-2 | 30 |
| 5 | Methyl (1s,3s,5R,7S)-3-methylada... | 208 | C13H20O2 | 1010504-39-1 | 27 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042123\
Data File : BM039565.D
Acq On : 21 Apr 2023 23:29
Operator : CG/JU
Sample : 02386-11
Misc :
ALS Vial : 13 Sample Multiplier: 1

**Instrument :**
BNA_M
**ClientSampleId :**
SPA-3WC-23-04

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM041723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

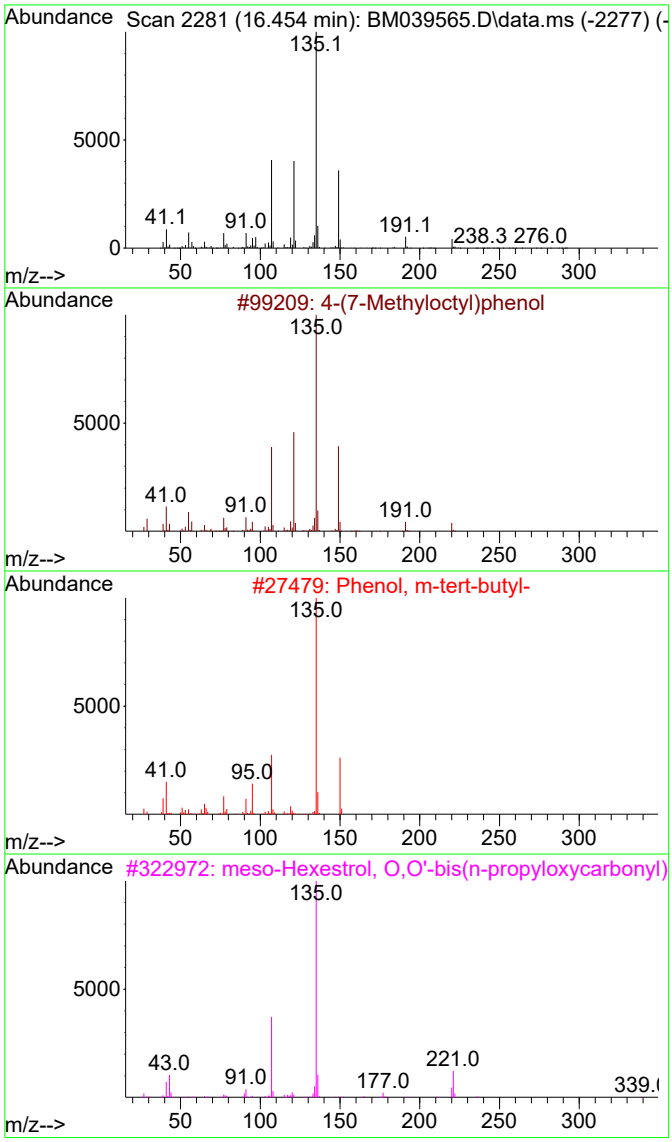
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

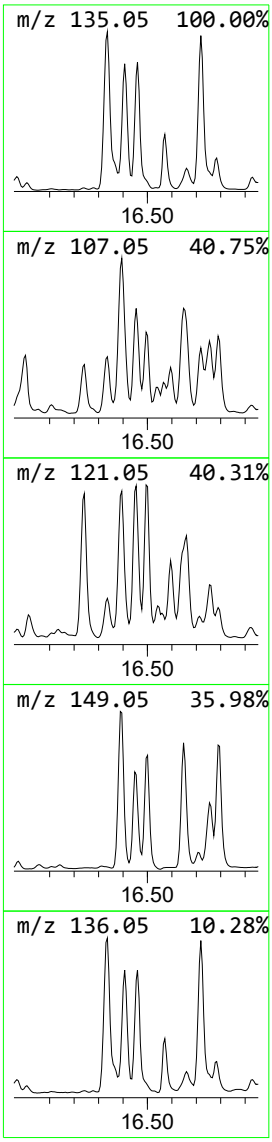
**********

Peak Number 19 4-(7-Methyloctyl)phenol Concentration Rank 18

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.454 | 26.69 ng | 14856200 | Phenanthrene-d10 | 17.248 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 4-(7-Methyloctyl)phenol | 220 | C15H24O | 024518-48-7 | 98 |
| 2 | Phenol, m-tert-butyl- | 150 | C10H14O | 000585-34-2 | 64 |
| 3 | meso-Hexestrol, O,O'-bis(n-propy... | 442 | C26H34O6 | 1000487-65-0 | 59 |
| 4 | Phenol, 4-(1,1-dimethylpropyl)- | 164 | C11H16O | 000080-46-6 | 59 |
| 5 | Hexestrol | 270 | C18H22O2 | 000084-16-2 | 53 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042123\
Data File : BM039565.D
Acq On : 21 Apr 2023 23:29
Operator : CG/JU
Sample : 02386-11
Misc :
ALS Vial : 13 Sample Multiplier: 1

**Instrument :**
BNA_M
**ClientSampleId :**
SPA-3WC-23-04

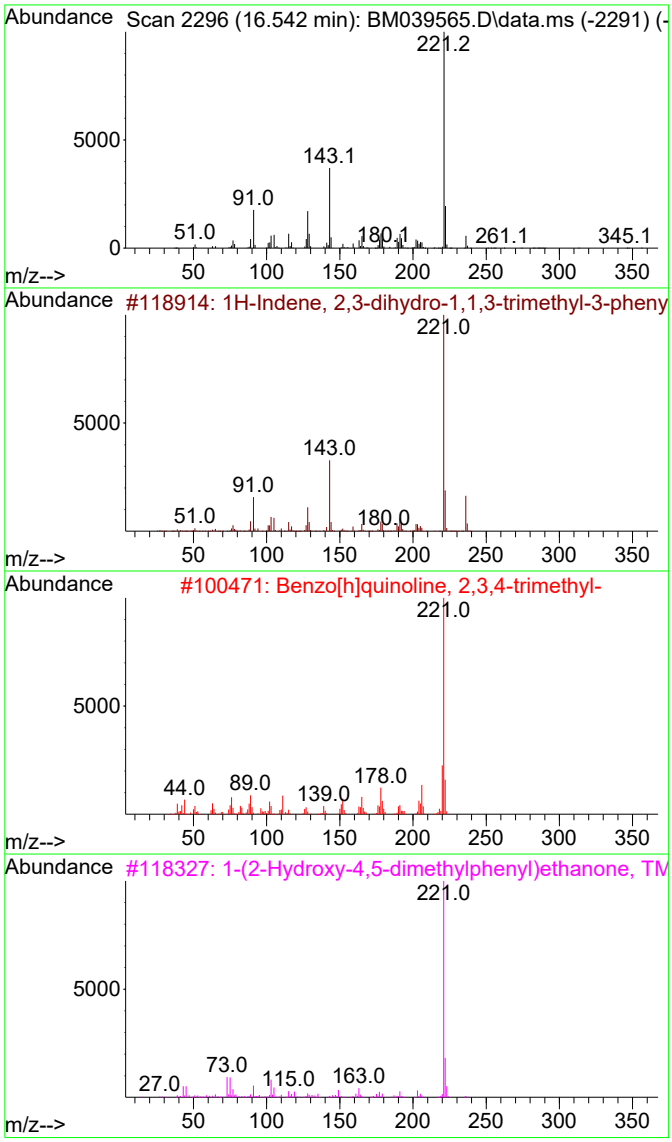
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM041723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

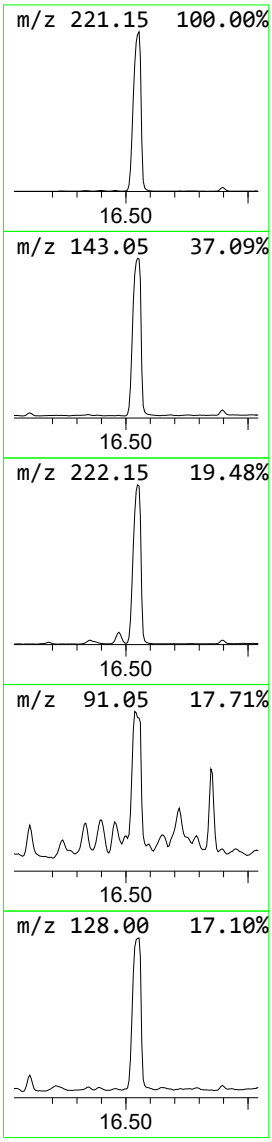
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 20 1H-Indene, 2,3-dihydro-1,1,... Concentration Rank 7

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.542 | 39.13 ng | 21778300 | Phenanthrene-d10 | 17.248 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1H-Indene, 2,3-dihydro-1,1,3-tri... | 236 | C18H20 | 003910-35-8 | 86 |
| 2 | Benzo[h]quinoline, 2,3,4-trimethyl- | 221 | C16H15N | 063042-66-0 | 52 |
| 3 | 1-(2-Hydroxy-4,5-dimethylphenyl)... | 236 | C13H20O2Si | 1000484-73-7 | 50 |
| 4 | Benz[j]isoquinoline, 3,6,8-trime... | 221 | C16H15N | 061171-16-2 | 50 |
| 5 | (E)-2-Hydroxy-4'-cyano-stilbene | 221 | C15H11NO | 1000147-85-5 | 45 |



Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042123\
Data File : BM039565.D
Acq On : 21 Apr 2023 23:29
Operator : CG/JU
Sample : 02386-11
Misc :
ALS Vial : 13 Sample Multiplier: 1

**Instrument :**
BNA_M
**ClientSampleId :**
SPA-3WC-23-04

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM041723.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

| | | | | | --Internal Standard--- | | | |
|---|---|---|---|---|---|---|---|---|
| TIC Top Hit name | RT | EstConc | Units | Response | # | RT | Resp | Conc |
| unknown6.795 | 6.795 | 25.8 | ng | 16288200 | 1 | 7.825 | 12622900 | 20.0 |
| Tetradecane | 7.442 | 35.2 | ng | 22195400 | 1 | 7.825 | 12622900 | 20.0 |
| Benzene, 1,2,3-... | 7.466 | 27.1 | ng | 17123000 | 1 | 7.825 | 12622900 | 20.0 |
| unknown8.472 | 8.472 | 32.4 | ng | 20468900 | 1 | 7.825 | 12622900 | 20.0 |
| Carbonic acid, ... | 9.054 | 44.3 | ng | 27958600 | 1 | 7.825 | 12622900 | 20.0 |
| Dodecane, 2,7,1... | 13.007 | 30.8 | ng | 8942660 | 3 | 14.483 | 5807920 | 20.0 |
| Diphenyl ether | 13.542 | 69.8 | ng | 20280400 | 3 | 14.483 | 5807920 | 20.0 |
| Naphthalene, 2,... | 13.807 | 66.5 | ng | 19312900 | 3 | 14.483 | 5807920 | 20.0 |
| Naphthalene, 1,... | 13.866 | 41.7 | ng | 12107500 | 3 | 14.483 | 5807920 | 20.0 |
| Carbonic acid, ... | 13.960 | 61.7 | ng | 17922100 | 3 | 14.483 | 5807920 | 20.0 |
| unknown14.313 | 14.313 | 28.9 | ng | 8406140 | 3 | 14.483 | 5807920 | 20.0 |
| Nonadecane | 14.377 | 57.4 | ng | 16657800 | 3 | 14.483 | 5807920 | 20.0 |
| 9-methylheptade... | 15.330 | 34.7 | ng | 10089600 | 3 | 14.483 | 5807920 | 20.0 |
| Tetradecane, 2,... | 16.207 | 25.4 | ng | 14140300 | 4 | 17.248 | 11130800 | 20.0 |
| Phenol, 4-(1,1,... | 16.336 | 30.1 | ng | 16778700 | 4 | 17.248 | 11130800 | 20.0 |
| unknown16.395 | 16.395 | 32.8 | ng | 18241700 | 4 | 17.248 | 11130800 | 20.0 |
| 4-(7-Methylocty... | 16.454 | 26.7 | ng | 14856200 | 4 | 17.248 | 11130800 | 20.0 |
| 1H-Indene, 2,3-... | 16.542 | 39.1 | ng | 21778300 | 4 | 17.248 | 11130800 | 20.0 |