LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110625\
Data File : BF144192.D
Acq On : 06 Nov 2025 17:24
Operator : RC/JU
Sample : Q3534-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** MW-EB-1

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-BF110525.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF144192.D\data.ms

| peak # | R.T. min | first scan | max scan | last scan | PK TY | peak height | corr. area | corr. % max. | % of total |
|---|---|---|---|---|---|---|---|---|---|
| 1 | 2.222 | 3 | 5 | 20 | rVB | 144898 | 209686 | 3.05% | 0.537% |
| 2 | 3.852 | 274 | 282 | 290 | rVB | 266298 | 461324 | 6.71% | 1.182% |
| 3 | 4.516 | 391 | 395 | 403 | rVB | 51544 | 77906 | 1.13% | 0.200% |
| 4 | 5.175 | 502 | 507 | 517 | rVB | 309923 | 425652 | 6.19% | 1.090% |
| 5 | 5.493 | 554 | 561 | 569 | rBV | 726092 | 996173 | 14.50% | 2.551% |
| 6 | 6.046 | 650 | 655 | 659 | rBV | 553751 | 717643 | 10.44% | 1.838% |
| 7 | 6.334 | 695 | 704 | 709 | rBV | 599976 | 785018 | 11.42% | 2.011% |
| 8 | 6.498 | 727 | 732 | 738 | rBV | 500989 | 650826 | 9.47% | 1.667% |
| 9 | 6.822 | 780 | 787 | 791 | rVV2 | 41045 | 75289 | 1.10% | 0.193% |
| 10 | 6.869 | 791 | 795 | 801 | rVV2 | 295870 | 447492 | 6.51% | 1.146% |
| 11 | 6.928 | 801 | 805 | 812 | rVB2 | 70606 | 102258 | 1.49% | 0.262% |
| 12 | 7.051 | 819 | 826 | 832 | rBV | 732095 | 921791 | 13.41% | 2.361% |
| 13 | 7.116 | 832 | 837 | 841 | rBV2 | 78978 | 102663 | 1.49% | 0.263% |
| 14 | 7.193 | 846 | 850 | 857 | rVB | 51668 | 88814 | 1.29% | 0.227% |
| 15 | 7.404 | 881 | 886 | 888 | rBV2 | 301525 | 407084 | 5.92% | 1.043% |
| 16 | 7.434 | 888 | 891 | 901 | rVB | 935760 | 1287981 | 18.74% | 3.299% |
| 17 | 7.593 | 914 | 918 | 921 | rVB2 | 71406 | 85172 | 1.24% | 0.218% |
| 18 | 7.640 | 921 | 926 | 930 | rBV | 163233 | 191208 | 2.78% | 0.490% |
| 19 | 7.687 | 930 | 934 | 938 | rBV3 | 47406 | 77220 | 1.12% | 0.198% |
| 20 | 7.822 | 951 | 957 | 964 | rBV | 86081 | 131489 | 1.91% | 0.337% |
| 21 | 7.893 | 964 | 969 | 972 | rBV3 | 222362 | 354723 | 5.16% | 0.909% |
| 22 | 7.928 | 972 | 975 | 983 | rVB | 513137 | 670743 | 9.76% | 1.718% |
| 23 | 8.151 | 1006 | 1013 | 1018 | rBV | 350286 | 503630 | 7.33% | 1.290% |
| 24 | 8.228 | 1022 | 1026 | 1030 | rVB3 | 62675 | 84289 | 1.23% | 0.216% |
| 25 | 8.328 | 1039 | 1043 | 1051 | rVB2 | 54887 | 94461 | 1.37% | 0.242% |
| 26 | 8.416 | 1052 | 1058 | 1060 | rBV3 | 57137 | 100609 | 1.46% | 0.258% |
| 27 | 8.445 | 1060 | 1063 | 1067 | rVB | 84770 | 107976 | 1.57% | 0.277% |
| 28 | 8.510 | 1067 | 1074 | 1081 | rBV5 | 31599 | 92883 | 1.35% | 0.238% |
| 29 | 8.581 | 1082 | 1086 | 1089 | rBV | 56723 | 75172 | 1.09% | 0.193% |
| 30 | 8.734 | 1104 | 1112 | 1118 | rBV | 3955875 | 6872210 | 100.00% | 17.602% |
| 31 | 8.787 | 1118 | 1121 | 1129 | rVB7 | 38766 | 86852 | 1.26% | 0.222% |
| 32 | 8.934 | 1140 | 1146 | 1150 | rBV3 | 112632 | 172276 | 2.51% | 0.441% |
| 33 | 9.039 | 1160 | 1164 | 1167 | rBV3 | 109407 | 147594 | 2.15% | 0.378% |
| 34 | 9.234 | 1191 | 1197 | 1201 | rBV | 1463248 | 1923497 | 27.99% | 4.927% |
| 35 | 9.339 | 1209 | 1215 | 1220 | rVB | 2789269 | 3906627 | 56.85% | 10.006% |
| 36 | 9.634 | 1261 | 1265 | 1273 | rVB2 | 501232 | 806705 | 11.74% | 2.066% |

LSC Area Percent Report

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Operator : RC/JU
Sample : Q3534-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
MW-EB-1

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-BF110525.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 37 | 9.828 | 1293 | 1298 | 1300 | rBV2 | 80954 | 128609 | 1.87% | 0.329% |
| 38 | 9.910 | 1304 | 1312 | 1318 | rVB3 | 432760 | 667546 | 9.71% | 1.710% |
| 39 | 10.069 | 1334 | 1339 | 1341 | rBV | 76130 | 107264 | 1.56% | 0.275% |
| 40 | 10.104 | 1341 | 1345 | 1351 | rVB | 153956 | 213194 | 3.10% | 0.546% |
| 41 | 10.322 | 1378 | 1382 | 1386 | rVB | 140744 | 167166 | 2.43% | 0.428% |
| 42 | 10.434 | 1396 | 1401 | 1409 | rVB | 146963 | 228689 | 3.33% | 0.586% |
| 43 | 10.545 | 1413 | 1420 | 1422 | rBV2 | 167482 | 260516 | 3.79% | 0.667% |
| 44 | 10.628 | 1429 | 1434 | 1435 | rBV2 | 112862 | 142959 | 2.08% | 0.366% |
| 45 | 10.657 | 1435 | 1439 | 1442 | rVV | 394349 | 530233 | 7.72% | 1.358% |
| 46 | 10.704 | 1442 | 1447 | 1453 | rVB2 | 949557 | 1349638 | 19.64% | 3.457% |
| 47 | 10.804 | 1459 | 1464 | 1465 | rBV2 | 103235 | 138415 | 2.01% | 0.355% |
| 48 | 10.833 | 1465 | 1469 | 1476 | rVV2 | 416672 | 664454 | 9.67% | 1.702% |
| 49 | 10.916 | 1476 | 1483 | 1488 | rVV | 813704 | 1108252 | 16.13% | 2.839% |
| 50 | 10.998 | 1494 | 1497 | 1501 | rBV | 80825 | 115379 | 1.68% | 0.296% |
| 51 | 11.081 | 1507 | 1511 | 1515 | rVV2 | 61955 | 89251 | 1.30% | 0.229% |
| 52 | 11.128 | 1515 | 1519 | 1526 | rVB3 | 323961 | 474777 | 6.91% | 1.216% |
| 53 | 11.233 | 1531 | 1537 | 1541 | rBV | 555476 | 786790 | 11.45% | 2.015% |
| 54 | 11.269 | 1541 | 1543 | 1551 | rVB | 251046 | 296161 | 4.31% | 0.759% |
| 55 | 11.404 | 1561 | 1566 | 1570 | rBV | 404121 | 507550 | 7.39% | 1.300% |
| 56 | 11.704 | 1612 | 1617 | 1623 | rBV | 137717 | 222456 | 3.24% | 0.570% |
| 57 | 11.928 | 1651 | 1655 | 1666 | rBV4 | 69924 | 138436 | 2.01% | 0.355% |
| 58 | 12.045 | 1672 | 1675 | 1681 | rVB | 55357 | 69864 | 1.02% | 0.179% |
| 59 | 12.175 | 1692 | 1697 | 1699 | rBV | 187445 | 275526 | 4.01% | 0.706% |
| 60 | 12.198 | 1699 | 1701 | 1706 | rVB | 201561 | 222957 | 3.24% | 0.571% |
| 61 | 12.363 | 1726 | 1729 | 1738 | rVB5 | 40210 | 90077 | 1.31% | 0.231% |
| 62 | 12.669 | 1778 | 1781 | 1786 | rVB | 109043 | 141820 | 2.06% | 0.363% |
| 63 | 12.880 | 1813 | 1817 | 1826 | rVB | 176541 | 248792 | 3.62% | 0.637% |
| 64 | 12.992 | 1831 | 1836 | 1841 | rBV | 1232175 | 1576654 | 22.94% | 4.038% |
| 65 | 13.551 | 1927 | 1931 | 1934 | rBV2 | 80393 | 104474 | 1.52% | 0.268% |
| 66 | 13.586 | 1934 | 1937 | 1947 | rVB | 98688 | 123519 | 1.80% | 0.316% |
| 67 | 13.751 | 1961 | 1965 | 1968 | rBV2 | 115419 | 175724 | 2.56% | 0.450% |
| 68 | 14.051 | 2012 | 2016 | 2024 | rVB | 343497 | 467083 | 6.80% | 1.196% |
| 69 | 14.222 | 2040 | 2045 | 2052 | rVB | 509939 | 690276 | 10.04% | 1.768% |
| 70 | 14.686 | 2117 | 2124 | 2138 | rVB | 890063 | 1582868 | 23.03% | 4.054% |
| 71 | 15.498 | 2257 | 2262 | 2265 | rBV | 104670 | 174112 | 2.53% | 0.446% |
| 72 | 15.539 | 2265 | 2269 | 2280 | rVB | 329023 | 518627 | 7.55% | 1.328% |

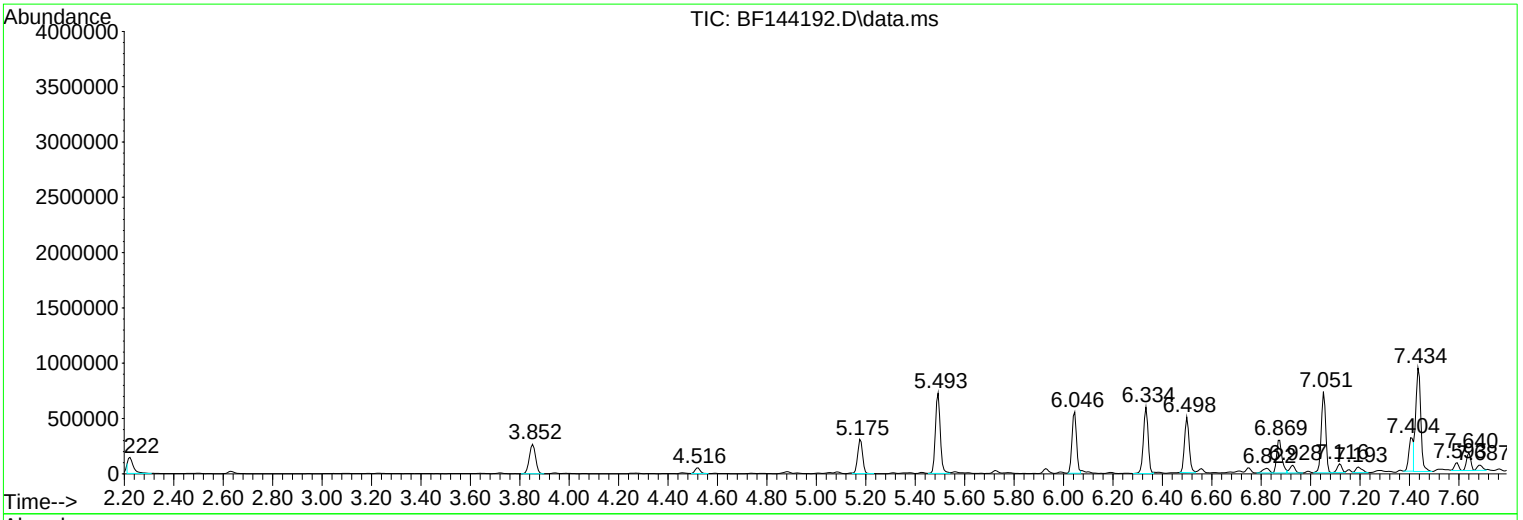
Sum of corrected areas: 39043044

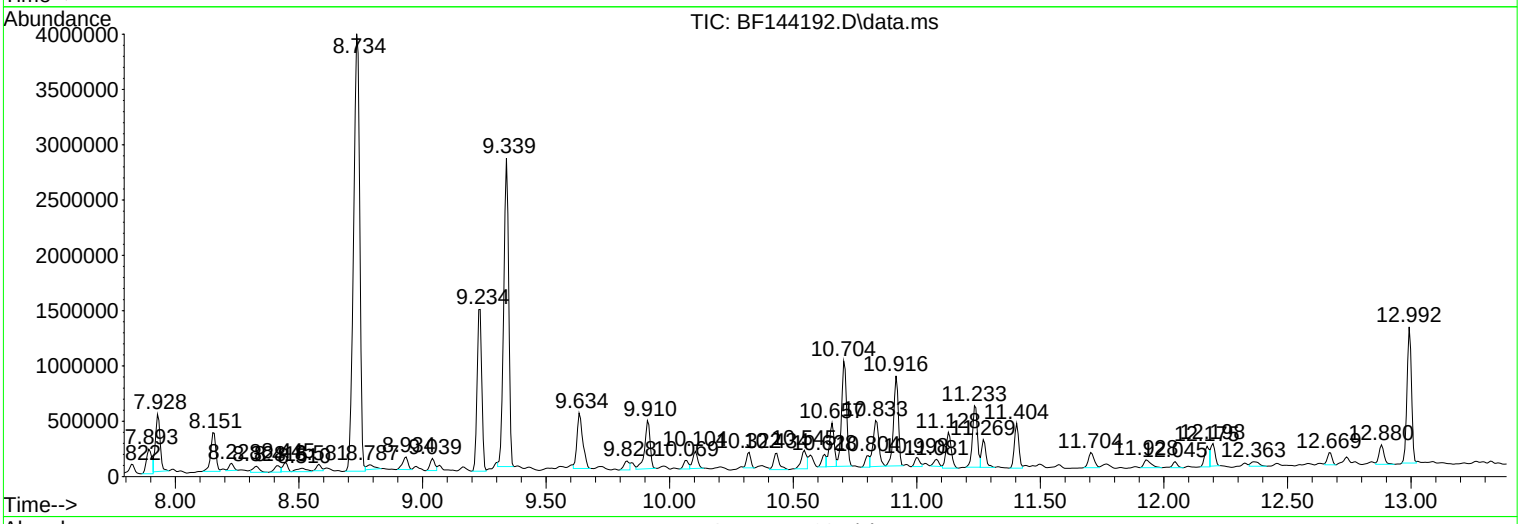
Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110625\
Data File : BF144192.D
Acq On : 06 Nov 2025 17:24
Operator : RC/JU
Sample : Q3534-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

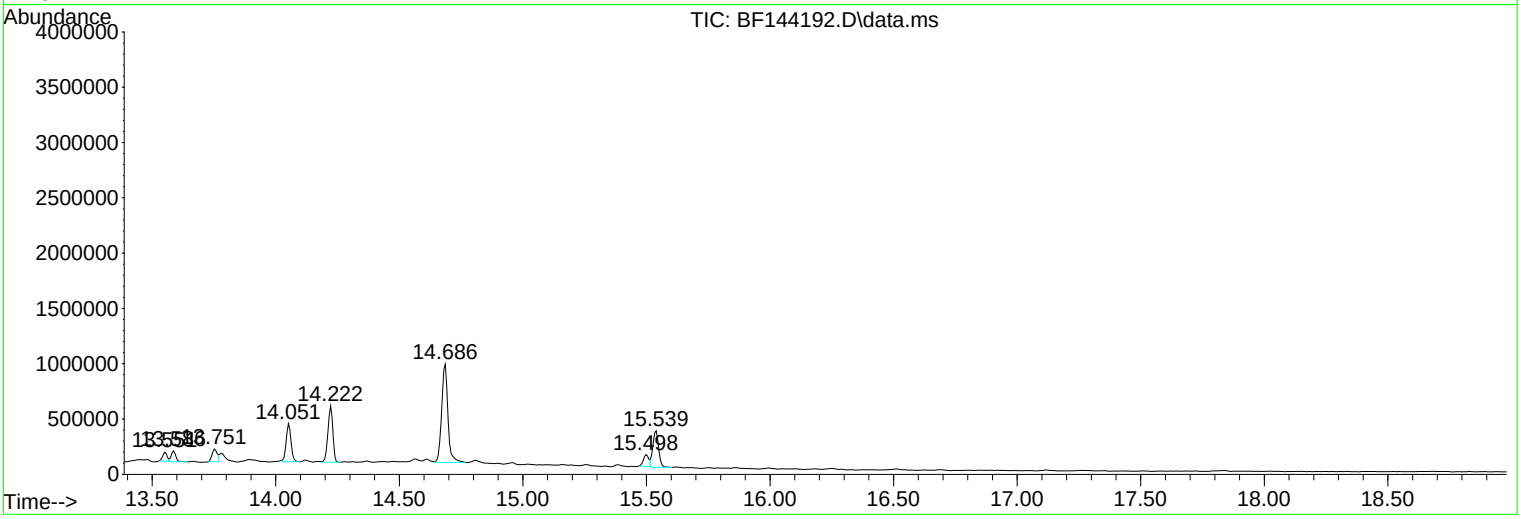
**Instrument :**
BNA_F
**ClientSampleId :**
MW-EB-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110525.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110625\
Data File : BF144192.D
Acq On : 06 Nov 2025 17:24
Operator : RC/JU
Sample : Q3534-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
MW-EB-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110525.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

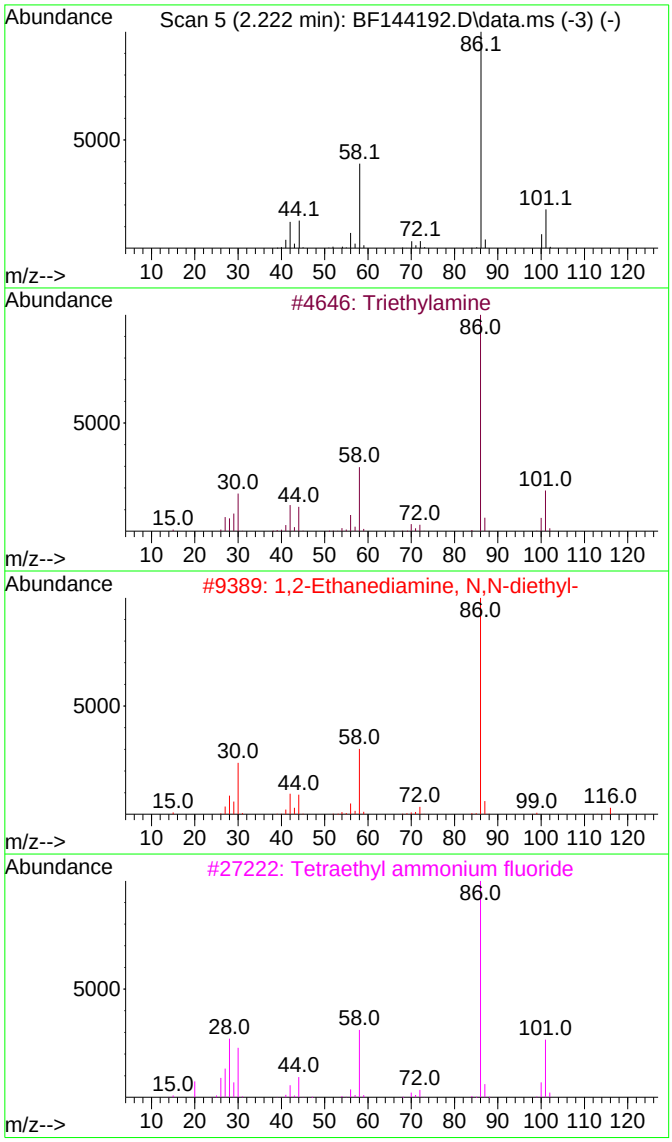
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

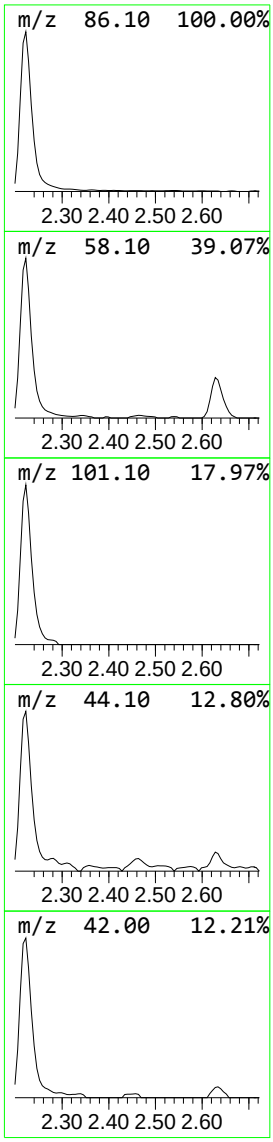
**********

Peak Number 1 Triethylamine Concentration Rank 20

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 2.222 | 9.37 ng | 209686 | 1,4-Dichlorobenzene-d4 | 6.875 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Triethylamine | 101 | C6H15N | 000121-44-8 | 91 |
| 2 | 1,2-Ethanediamine, N,N-diethyl- | 116 | C6H16N2 | 000100-36-7 | 72 |
| 3 | Tetraethyl ammonium fluoride | 149 | C8H20FN | 000665-46-3 | 64 |
| 4 | 1,2-Ethanediamine, N,N-diethyl-N... | 130 | C7H18N2 | 000104-79-0 | 64 |
| 5 | Tetraethylammonium chloride | 165 | C8H20ClN | 000056-34-8 | 64 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110625\
Data File : BF144192.D
Acq On : 06 Nov 2025 17:24
Operator : RC/JU
Sample : Q3534-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
MW-EB-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110525.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

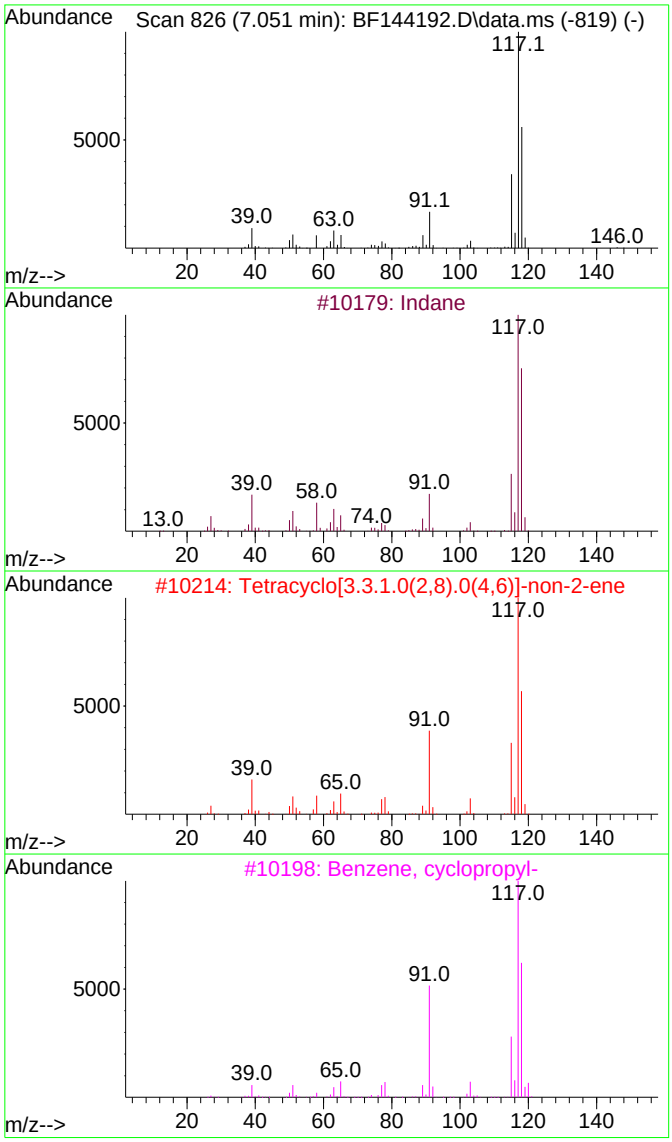
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

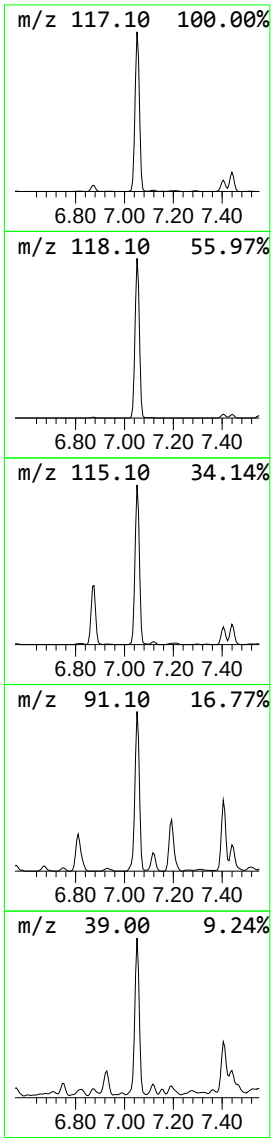
**********

Peak Number 6 Indane Concentration Rank 5

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 7.051 | 41.20 ng | 921791 | 1,4-Dichlorobenzene-d4 | 6.875 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Indane | 118 | C9H10 | 000496-11-7 | 91 |
| 2 | Tetracyclo[3.3.1.0(2,8).0(4,6)]-... | 118 | C9H10 | 1000191-13-7 | 80 |
| 3 | Benzene, cyclopropyl- | 118 | C9H10 | 000873-49-4 | 72 |
| 4 | Benzene, 1-ethenyl-2-methyl- | 118 | C9H10 | 000611-15-4 | 60 |
| 5 | Benzene, 1,1'-(1,5-hexadiene-1,6... | 234 | C18H18 | 004439-45-6 | 59 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110625\
Data File : BF144192.D
Acq On : 06 Nov 2025 17:24
Operator : RC/JU
Sample : Q3534-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
MW-EB-1

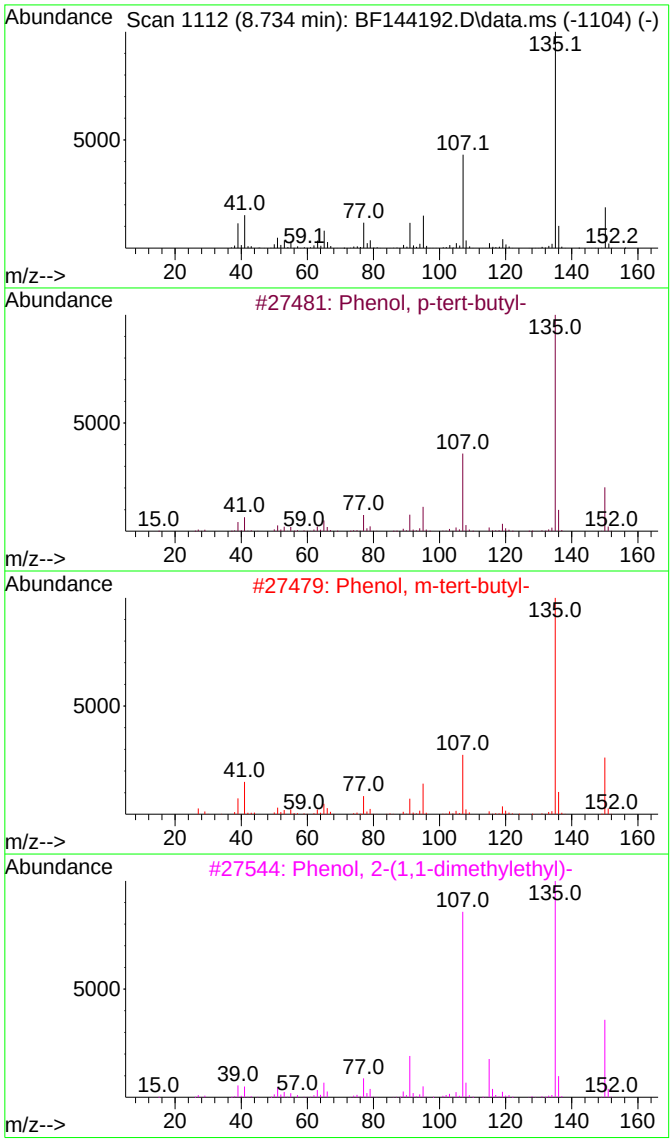
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110525.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

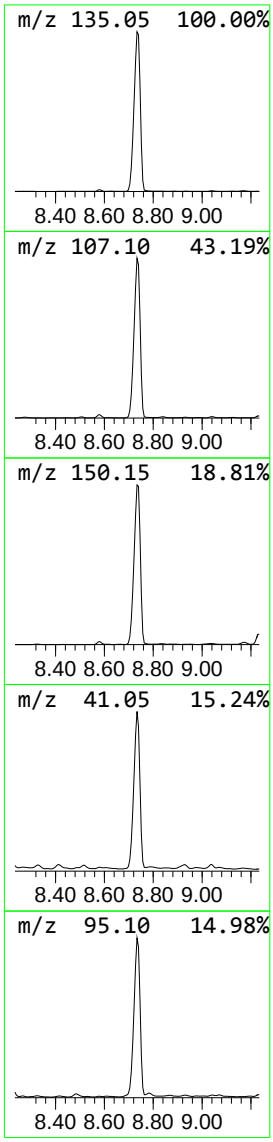
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 9 Phenol, p-tert-butyl- Concentration Rank 1

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 8.734 | 272.91 ng | 6872210 | Naphthalene-d8 | 8.157 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, p-tert-butyl- | 150 | C10H14O | 000098-54-4 | 96 |
| 2 | Phenol, m-tert-butyl- | 150 | C10H14O | 000585-34-2 | 93 |
| 3 | Phenol, 2-(1,1-dimethylethyl)- | 150 | C10H14O | 000088-18-6 | 74 |
| 4 | Imidazo[1,2-c]pyrimidin-5-ol | 135 | C6H5N3O | 055662-66-3 | 58 |
| 5 | 5-Cyano-1,2,3,4-tetrahydro-4,6-d... | 150 | C8H10N2O | 027036-93-7 | 52 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110625\
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Acq On : 06 Nov 2025 17:24
Operator : RC/JU
Sample : Q3534-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
MW-EB-1

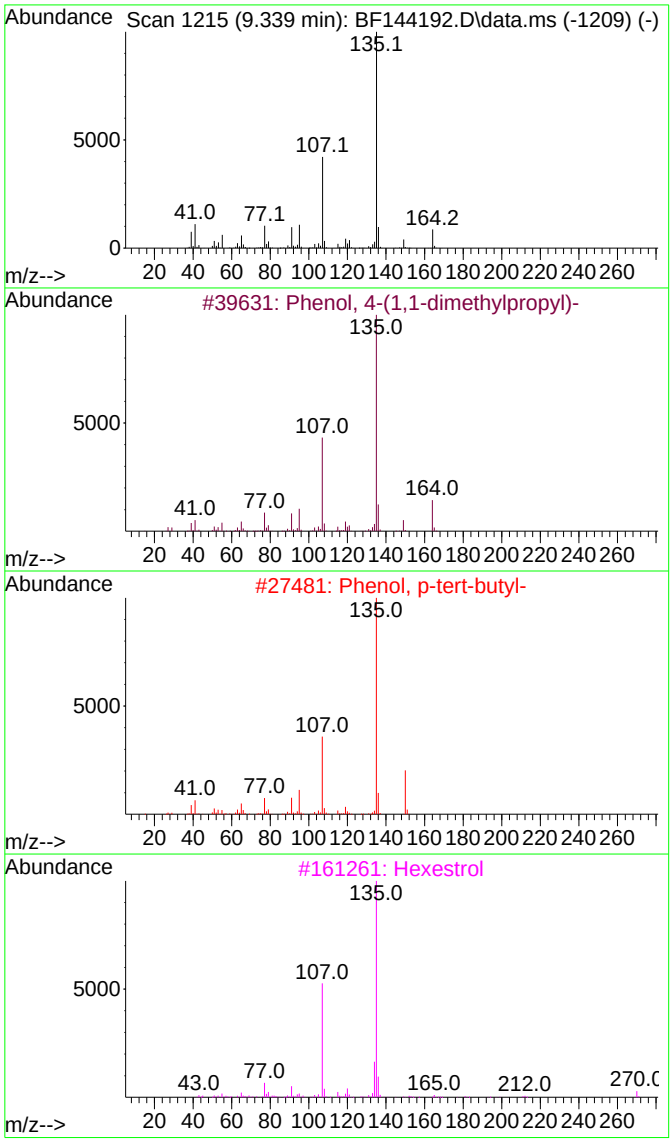
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110525.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

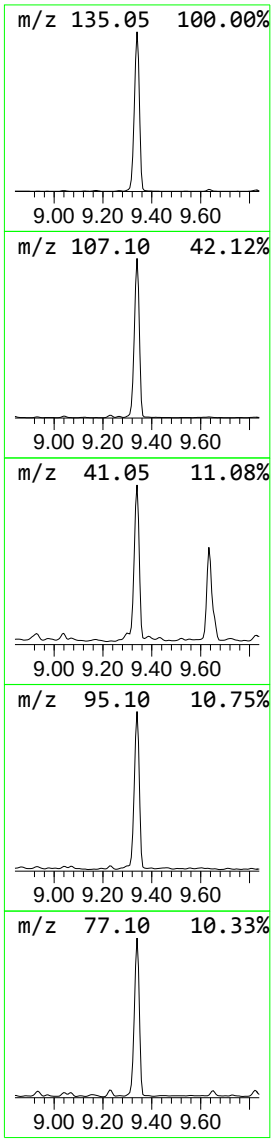
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 10 Phenol, 4-(1,1-dimethylprop... Concentration Rank 2

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 9.339 | 117.04 ng | 3906630 | Acenaphthene-d10 | 9.910 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, 4-(1,1-dimethylpropyl)- | 164 | C11H16O | 000080-46-6 | 95 |
| 2 | Phenol, p-tert-butyl- | 150 | C10H14O | 000098-54-4 | 86 |
| 3 | Hexestrol | 270 | C18H22O2 | 000084-16-2 | 72 |
| 4 | Phenol, 4-(1,1,3,3-tetramethylbu... | 206 | C14H22O | 000140-66-9 | 64 |
| 5 | 4,4'-(1,2-diethylethylene)diphenol | 270 | C18H22O2 | 005635-50-7 | 64 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110625\
Data File : BF144192.D
Acq On : 06 Nov 2025 17:24
Operator : RC/JU
Sample : Q3534-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
MW-EB-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110525.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

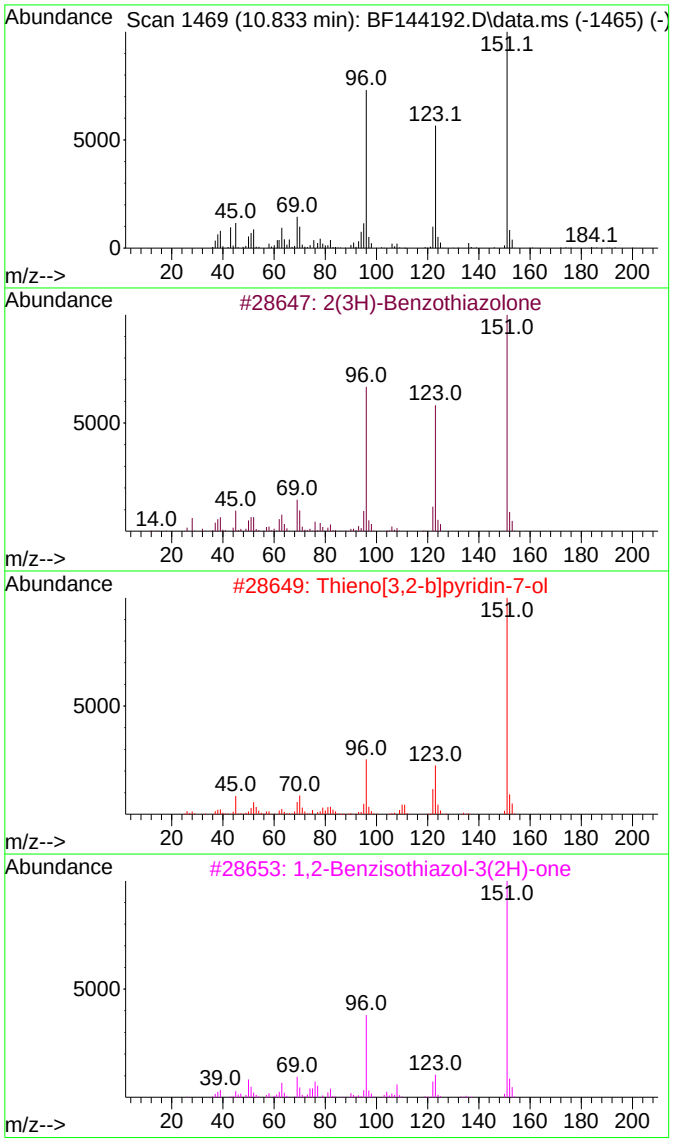
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

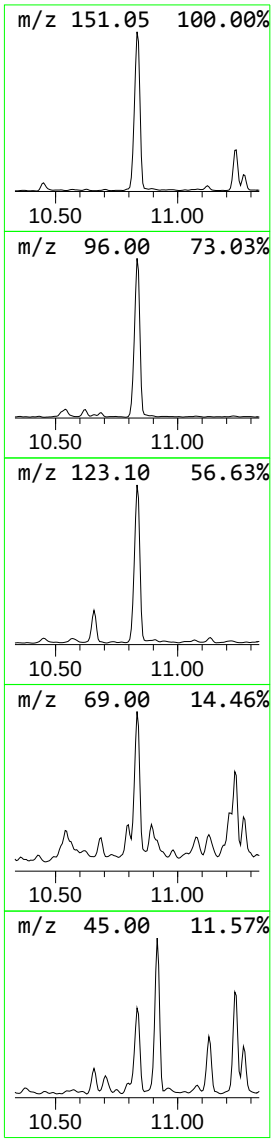
**********

Peak Number 12 2(3H)-Benzothiazolone Concentration Rank 11

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.834 | 26.18 ng | 664454 | Phenanthrene-d10 | 11.404 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2(3H)-Benzothiazolone | 151 | C7H5NOS | 000934-34-9 | 94 |
| 2 | Thieno[3,2-b]pyridin-7-ol | 151 | C7H5NOS | 107818-20-2 | 58 |
| 3 | 1,2-Benzisothiazol-3(2H)-one | 151 | C7H5NOS | 002634-33-5 | 49 |
| 4 | 6-Methyl-7-oxo-4,7-dihydro-triaz... | 151 | C5H5N5O | 057250-39-2 | 47 |
| 5 | 2-Cyano-3-fluorophenylhydrazine | 151 | C7H6FN3 | 1000131-91-9 | 46 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110625\
Data File : BF144192.D
Acq On : 06 Nov 2025 17:24
Operator : RC/JU
Sample : Q3534-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
MW-EB-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110525.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

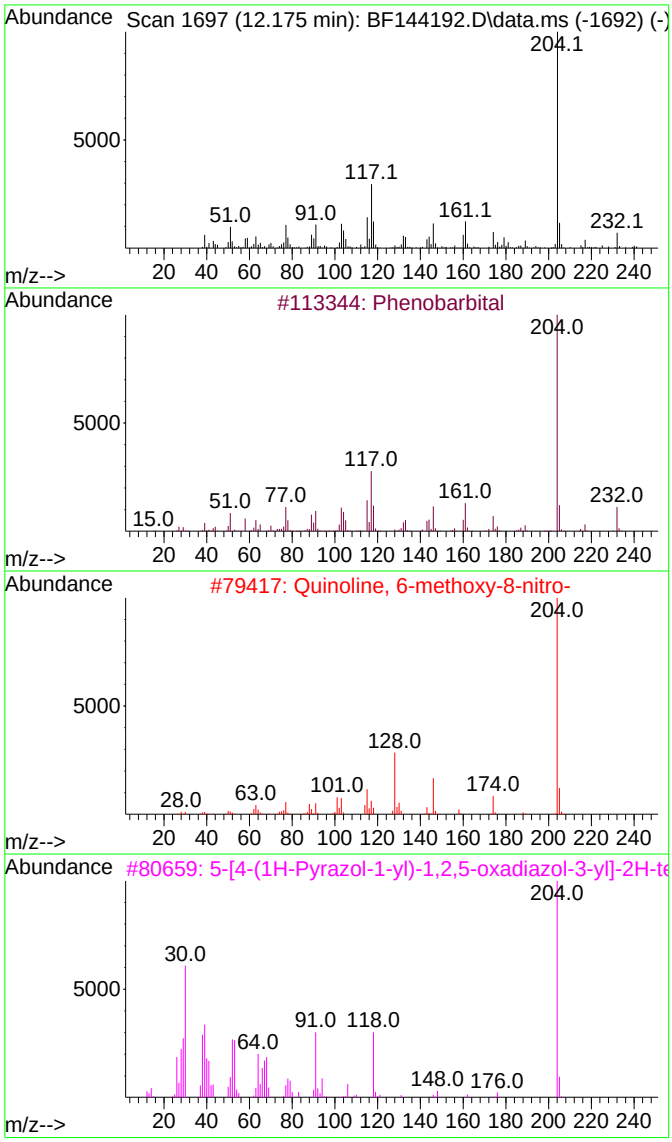
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

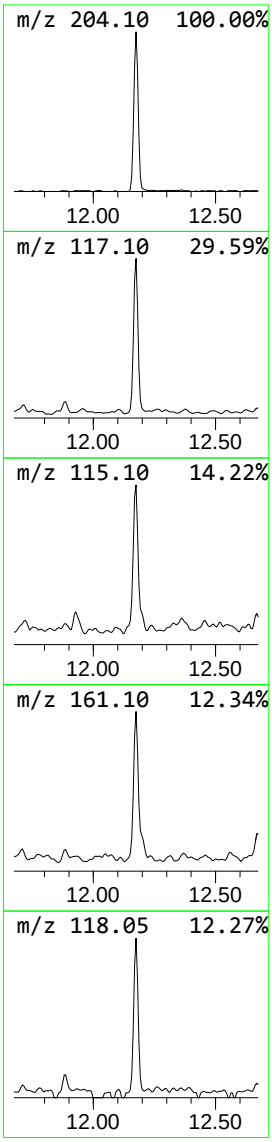
**********

Peak Number 17 Phenobarbital Concentration Rank 18

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.175 | 10.86 ng | 275526 | Phenanthrene-d10 | 11.404 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenobarbital | 232 | C12H12N2O3 | 000050-06-6 | 96 |
| 2 | Quinoline, 6-methoxy-8-nitro- | 204 | C10H8N2O3 | 000085-81-4 | 52 |
| 3 | 5-[4-(1H-Pyrazol-1-yl)-1,2,5-oxa... | 204 | C6H4N8O | 1000460-43-6 | 47 |
| 4 | 2-Methyl-6-nitro-4-quinolinol | 204 | C10H8N2O3 | 001207-82-5 | 47 |
| 5 | 6-Hydroxy-8-nitrolepidine | 204 | C10H8N2O3 | 1000213-59-0 | 47 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110625\
Data File : BF144192.D
Acq On : 06 Nov 2025 17:24
Operator : RC/JU
Sample : Q3534-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
MW-EB-1

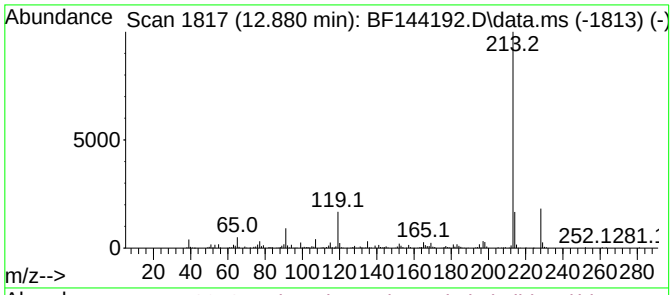
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110525.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

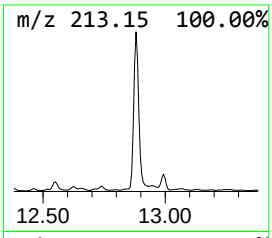
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

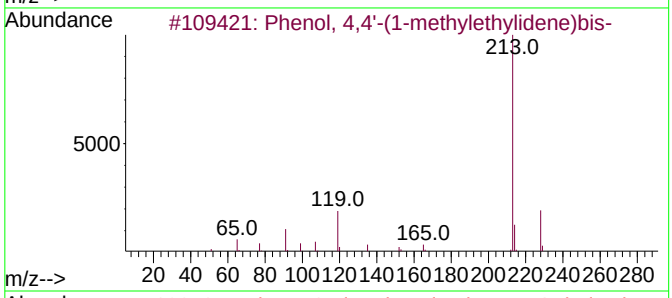
**********
Peak Number 18 Phenol, 4,4'-(1-methylethyl... Concentration Rank 19

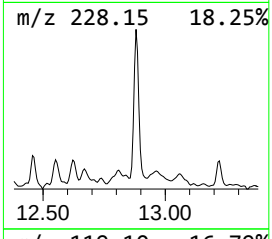
| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 12.880 | 10.65 ng | 248792 | Chrysene-d12 | 14.051 |

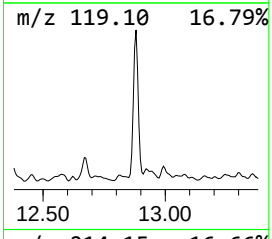
| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, 4,4'-(1-methylethylidene... | 228 | C15H16O2 | 000080-05-7 | 96 |
| 2 | 5-Fluoro-2-nitrophenylamine, TMS... | 228 | C9H13FN2O2Si | 1000484-54-5 | 59 |
| 3 | 2H,8H-Benzo[1,2-b:5,4-b']dipyran... | 228 | C14H12O3 | 000553-19-5 | 59 |
| 4 | 4H-Pyrazolo[5,1-c]-1,4-benzodiaz... | 213 | C11H7N3O2 | 1000277-20-1 | 59 |
| 5 | 2H,8H-Benzo[1,2-b:3,4-b']dipyran... | 228 | C14H12O3 | 000523-59-1 | 59 |

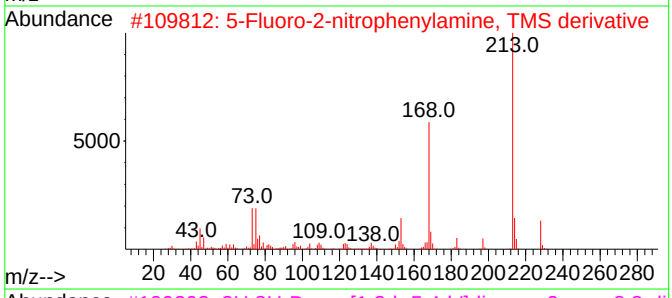


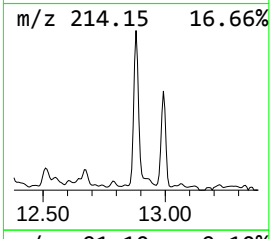



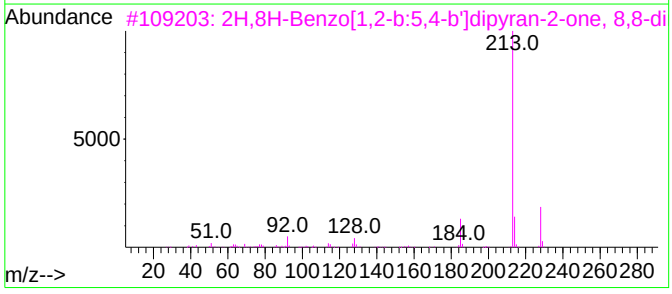



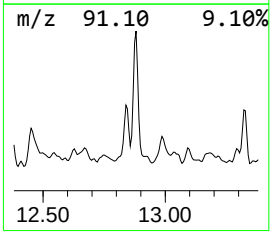



Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110625\
Data File : BF144192.D
Acq On : 06 Nov 2025 17:24
Operator : RC/JU
Sample : Q3534-05
Misc :
ALS Vial : 14 Sample Multiplier: 1

**Instrument :**
BNA_F
**ClientSampleId :**
MW-EB-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110525.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 |
| Triethylamine | 2.222 | 9.4 | ng | 209686 | 1 | 6.875 | 447492 | 20.0 |
| Indane | 7.051 | 41.2 | ng | 921791 | 1 | 6.875 | 447492 | 20.0 |
| Phenol, p-tert-... | 8.734 | 272.9 | ng | 6872210 | 2 | 8.157 | 503630 | 20.0 |
| Phenol, 4-(1,1-... | 9.339 | 117.0 | ng | 3906630 | 3 | 9.910 | 667546 | 20.0 |
| 2(3H)-Benzothia... | 10.834 | 26.2 | ng | 664454 | 4 | 11.404 | 507550 | 20.0 |
| Phenobarbital | 12.175 | 10.9 | ng | 275526 | 4 | 11.404 | 507550 | 20.0 |
| Phenol, 4,4'-(1... | 12.880 | 10.7 | ng | 248792 | 5 | 14.051 | 467083 | 20.0 |