LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070218\
Data File : BF107063.D
Acq On : 3 Jul 2018 1:10
Operator : JU/SJ
Sample : J3799-07
Misc :
ALS Vial : 31 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** MW-6

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-BF061918.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

| peak # | R.T. min | first scan | max scan | last scan | PK TY | peak height | corr. area | corr. % max. | % of total |
|---|---|---|---|---|---|---|---|---|---|
| 1 | 5.189 | 481 | 485 | 491 | rBV | 72207 | 81380 | 4.58% | 0.280% |
| 2 | 5.601 | 551 | 555 | 567 | rBV | 842176 | 813451 | 45.81% | 2.801% |
| 3 | 6.001 | 621 | 623 | 633 | rVB3 | 18407 | 24395 | 1.37% | 0.084% |
| 4 | 6.125 | 639 | 644 | 647 | rBV2 | 37722 | 39081 | 2.20% | 0.135% |
| 5 | 6.413 | 690 | 693 | 697 | rVB | 33509 | 31479 | 1.77% | 0.108% |
| 6 | 6.607 | 720 | 726 | 733 | rBV | 502361 | 615962 | 34.68% | 2.121% |
| 7 | 6.736 | 744 | 748 | 753 | rBV | 1484594 | 1377249 | 77.55% | 4.742% |
| 8 | 6.778 | 753 | 755 | 757 | rVV2 | 36348 | 39322 | 2.21% | 0.135% |
| 9 | 6.795 | 757 | 758 | 762 | rVB2 | 36559 | 27950 | 1.57% | 0.096% |
| 10 | 6.883 | 769 | 773 | 774 | rBV3 | 31004 | 39869 | 2.25% | 0.137% |
| 11 | 6.901 | 774 | 776 | 779 | rVB | 57826 | 47340 | 2.67% | 0.163% |
| 12 | 6.954 | 781 | 785 | 790 | rBV | 434706 | 386513 | 21.76% | 1.331% |
| 13 | 7.113 | 808 | 812 | 815 | rBV | 1421533 | 1317155 | 74.17% | 4.535% |
| 14 | 7.142 | 815 | 817 | 819 | rVB | 29993 | 22011 | 1.24% | 0.076% |
| 15 | 7.195 | 823 | 826 | 828 | rVB2 | 69937 | 53869 | 3.03% | 0.185% |
| 16 | 7.283 | 836 | 841 | 844 | rBV | 111209 | 122448 | 6.89% | 0.422% |
| 17 | 7.342 | 847 | 851 | 854 | rBV2 | 67888 | 67896 | 3.82% | 0.234% |
| 18 | 7.389 | 856 | 859 | 861 | rBV2 | 23043 | 19812 | 1.12% | 0.068% |
| 19 | 7.478 | 865 | 874 | 877 | rBV3 | 119878 | 214301 | 12.07% | 0.738% |
| 20 | 7.525 | 877 | 882 | 885 | rBV | 1072037 | 1087630 | 61.24% | 3.745% |
| 21 | 7.589 | 891 | 893 | 895 | rVB2 | 36677 | 29614 | 1.67% | 0.102% |
| 22 | 7.636 | 899 | 901 | 903 | rBV | 80792 | 62397 | 3.51% | 0.215% |
| 23 | 7.719 | 912 | 915 | 917 | rVB | 155090 | 121144 | 6.82% | 0.417% |
| 24 | 7.742 | 917 | 919 | 921 | rBV2 | 31909 | 27215 | 1.53% | 0.094% |
| 25 | 7.766 | 921 | 923 | 926 | rVB3 | 45834 | 47846 | 2.69% | 0.165% |
| 26 | 7.813 | 926 | 931 | 936 | rBV4 | 72921 | 107999 | 6.08% | 0.372% |
| 27 | 7.878 | 938 | 942 | 945 | rBV3 | 133278 | 175527 | 9.88% | 0.604% |
| 28 | 7.972 | 954 | 958 | 960 | rBV2 | 119072 | 155546 | 8.76% | 0.536% |
| 29 | 8.042 | 966 | 970 | 972 | rBV2 | 74770 | 81597 | 4.59% | 0.281% |
| 30 | 8.095 | 976 | 979 | 981 | rBV | 43758 | 37834 | 2.13% | 0.130% |
| 31 | 8.148 | 985 | 988 | 989 | rVV2 | 55299 | 47280 | 2.66% | 0.163% |
| 32 | 8.178 | 989 | 993 | 1000 | rVV4 | 181079 | 368713 | 20.76% | 1.270% |
| 33 | 8.242 | 1000 | 1004 | 1008 | rVV | 558501 | 680315 | 38.31% | 2.342% |
| 34 | 8.278 | 1008 | 1010 | 1013 | rVB2 | 80540 | 73116 | 4.12% | 0.252% |

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Sample : J3799-07
Misc :
ALS Vial : 31 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** MW-6

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-BF061918.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 35 | 8.354 | 1019 | 1023 | 1026 | rBV2 | 132347 | 166840 | 9.39% | 0.574% |
| 36 | 8.389 | 1026 | 1029 | 1030 | rVV3 | 41247 | 46774 | 2.63% | 0.161% |
| 37 | 8.407 | 1030 | 1032 | 1034 | rVV2 | 71019 | 78392 | 4.41% | 0.270% |
| 38 | 8.430 | 1034 | 1036 | 1038 | rVV | 161984 | 155908 | 8.78% | 0.537% |
| 39 | 8.454 | 1038 | 1040 | 1041 | rVV | 119207 | 111569 | 6.28% | 0.384% |
| 40 | 8.477 | 1041 | 1044 | 1048 | rVV4 | 217282 | 305390 | 17.20% | 1.051% |
| 41 | 8.519 | 1048 | 1051 | 1055 | rVV5 | 93587 | 140748 | 7.93% | 0.485% |
| 42 | 8.554 | 1055 | 1057 | 1059 | rVB3 | 29511 | 23005 | 1.30% | 0.079% |
| 43 | 8.583 | 1059 | 1062 | 1063 | rBV3 | 55547 | 62291 | 3.51% | 0.214% |
| 44 | 8.619 | 1063 | 1068 | 1072 | rVV5 | 100906 | 206736 | 11.64% | 0.712% |
| 45 | 8.660 | 1072 | 1075 | 1077 | rVV3 | 88327 | 134364 | 7.57% | 0.463% |
| 46 | 8.683 | 1077 | 1079 | 1081 | rVV2 | 96963 | 99356 | 5.59% | 0.342% |
| 47 | 8.713 | 1081 | 1084 | 1087 | rVV3 | 98154 | 157895 | 8.89% | 0.544% |
| 48 | 8.748 | 1087 | 1090 | 1093 | rVB4 | 108719 | 130376 | 7.34% | 0.449% |
| 49 | 8.777 | 1093 | 1095 | 1096 | rBV | 46678 | 32853 | 1.85% | 0.113% |
| 50 | 8.813 | 1096 | 1101 | 1105 | rVV5 | 113392 | 187852 | 10.58% | 0.647% |
| 51 | 8.901 | 1114 | 1116 | 1118 | rVB2 | 72067 | 54962 | 3.09% | 0.189% |
| 52 | 8.930 | 1118 | 1121 | 1123 | rBV2 | 137681 | 130487 | 7.35% | 0.449% |
| 53 | 8.954 | 1123 | 1125 | 1127 | rVV2 | 87142 | 82253 | 4.63% | 0.283% |
| 54 | 8.995 | 1130 | 1132 | 1135 | rVV3 | 152217 | 156130 | 8.79% | 0.538% |
| 55 | 9.025 | 1135 | 1137 | 1140 | rVV2 | 170599 | 184516 | 10.39% | 0.635% |
| 56 | 9.060 | 1140 | 1143 | 1147 | rVB4 | 172138 | 200877 | 11.31% | 0.692% |
| 57 | 9.130 | 1153 | 1155 | 1161 | rVB5 | 103845 | 168816 | 9.51% | 0.581% |
| 58 | 9.183 | 1161 | 1164 | 1166 | rVB3 | 64641 | 65953 | 3.71% | 0.227% |
| 59 | 9.219 | 1166 | 1170 | 1172 | rBV4 | 84886 | 123905 | 6.98% | 0.427% |
| 60 | 9.242 | 1172 | 1174 | 1175 | rBV | 92583 | 82280 | 4.63% | 0.283% |
| 61 | 9.313 | 1183 | 1186 | 1190 | rVB | 1719612 | 1775896 | 100.00% | 6.115% |
| 62 | 9.354 | 1190 | 1193 | 1194 | rBV | 153595 | 132891 | 7.48% | 0.458% |
| 63 | 9.419 | 1199 | 1204 | 1207 | rBV | 171667 | 262174 | 14.76% | 0.903% |
| 64 | 9.460 | 1209 | 1211 | 1214 | rVB2 | 120584 | 101818 | 5.73% | 0.351% |
| 65 | 9.489 | 1214 | 1216 | 1217 | rBV | 81678 | 69383 | 3.91% | 0.239% |
| 66 | 9.601 | 1233 | 1235 | 1237 | rBV3 | 59572 | 63423 | 3.57% | 0.218% |
| 67 | 9.666 | 1241 | 1246 | 1250 | rVV6 | 215380 | 295279 | 16.63% | 1.017% |
| 68 | 9.707 | 1251 | 1253 | 1257 | rVB | 131569 | 145360 | 8.19% | 0.500% |
| 69 | 9.842 | 1273 | 1276 | 1278 | rBV2 | 107021 | 116494 | 6.56% | 0.401% |
| 70 | 9.901 | 1284 | 1286 | 1290 | rVB5 | 81922 | 84893 | 4.78% | 0.292% |
| 71 | 10.001 | 1300 | 1303 | 1307 | rVB | 507345 | 453336 | 25.53% | 1.561% |

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Sample : J3799-07
Misc :
ALS Vial : 31 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** MW-6

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-BF061918.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 72 | 10.236 | 1340 | 1343 | 1344 | rBV3 | 54294 | 63107 | 3.55% | 0.217% |
| 73 | 10.283 | 1349 | 1351 | 1353 | rBV3 | 54440 | 61385 | 3.46% | 0.211% |
| 74 | 10.336 | 1358 | 1360 | 1363 | rVV2 | 142051 | 167647 | 9.44% | 0.577% |
| 75 | 10.383 | 1363 | 1368 | 1374 | rVV6 | 147536 | 320351 | 18.04% | 1.103% |
| 76 | 10.442 | 1376 | 1378 | 1385 | rVB6 | 114750 | 164824 | 9.28% | 0.568% |
| 77 | 10.748 | 1427 | 1430 | 1432 | rBV | 209534 | 193974 | 10.92% | 0.668% |
| 78 | 10.771 | 1432 | 1434 | 1435 | rVV | 229831 | 199862 | 11.25% | 0.688% |
| 79 | 10.801 | 1435 | 1439 | 1445 | rVB2 | 957874 | 1117447 | 62.92% | 3.847% |
| 80 | 10.919 | 1456 | 1459 | 1464 | rBV | 457163 | 423437 | 23.84% | 1.458% |
| 81 | 10.971 | 1464 | 1468 | 1470 | rVV | 962289 | 907361 | 51.09% | 3.124% |
| 82 | 11.007 | 1470 | 1474 | 1477 | rVV2 | 1074841 | 1491139 | 83.97% | 5.134% |
| 83 | 11.036 | 1477 | 1479 | 1482 | rVV2 | 1134675 | 1134909 | 63.91% | 3.908% |
| 84 | 11.066 | 1482 | 1484 | 1486 | rVV | 690241 | 556576 | 31.34% | 1.916% |
| 85 | 11.107 | 1486 | 1491 | 1495 | rVV2 | 570094 | 937143 | 52.77% | 3.227% |
| 86 | 11.154 | 1495 | 1499 | 1502 | rVV3 | 926657 | 1088965 | 61.32% | 3.749% |
| 87 | 11.189 | 1502 | 1505 | 1507 | rVV | 1004570 | 951466 | 53.58% | 3.276% |
| 88 | 11.207 | 1507 | 1508 | 1510 | rVV2 | 451567 | 375652 | 21.15% | 1.293% |
| 89 | 11.230 | 1510 | 1512 | 1516 | rVB | 654004 | 543162 | 30.59% | 1.870% |
| 90 | 11.501 | 1555 | 1558 | 1561 | rVB | 445425 | 357964 | 20.16% | 1.233% |
| 91 | 11.913 | 1625 | 1628 | 1632 | rBV | 184510 | 152115 | 8.57% | 0.524% |
| 92 | 12.018 | 1643 | 1646 | 1652 | rVB2 | 77287 | 90275 | 5.08% | 0.311% |
| 93 | 12.101 | 1658 | 1660 | 1662 | rBV | 43197 | 39116 | 2.20% | 0.135% |
| 94 | 12.207 | 1675 | 1678 | 1679 | rBV2 | 71909 | 69931 | 3.94% | 0.241% |
| 95 | 12.224 | 1679 | 1681 | 1683 | rVV2 | 74165 | 59679 | 3.36% | 0.205% |
| 96 | 12.248 | 1683 | 1685 | 1688 | rVB | 80774 | 64272 | 3.62% | 0.221% |
| 97 | 13.083 | 1822 | 1827 | 1830 | rBV | 1709448 | 1612657 | 90.81% | 5.553% |
| 98 | 13.954 | 1973 | 1975 | 1979 | rVB2 | 48871 | 41039 | 2.31% | 0.141% |
| 99 | 14.154 | 2005 | 2009 | 2012 | rBV | 361970 | 348086 | 19.60% | 1.198% |
| 100 | 15.689 | 2266 | 2270 | 2276 | rVB2 | 202960 | 273223 | 15.39% | 0.941% |

Sum of corrected areas: 29043495

LSC Report - Integrated Chromatogram

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070218\
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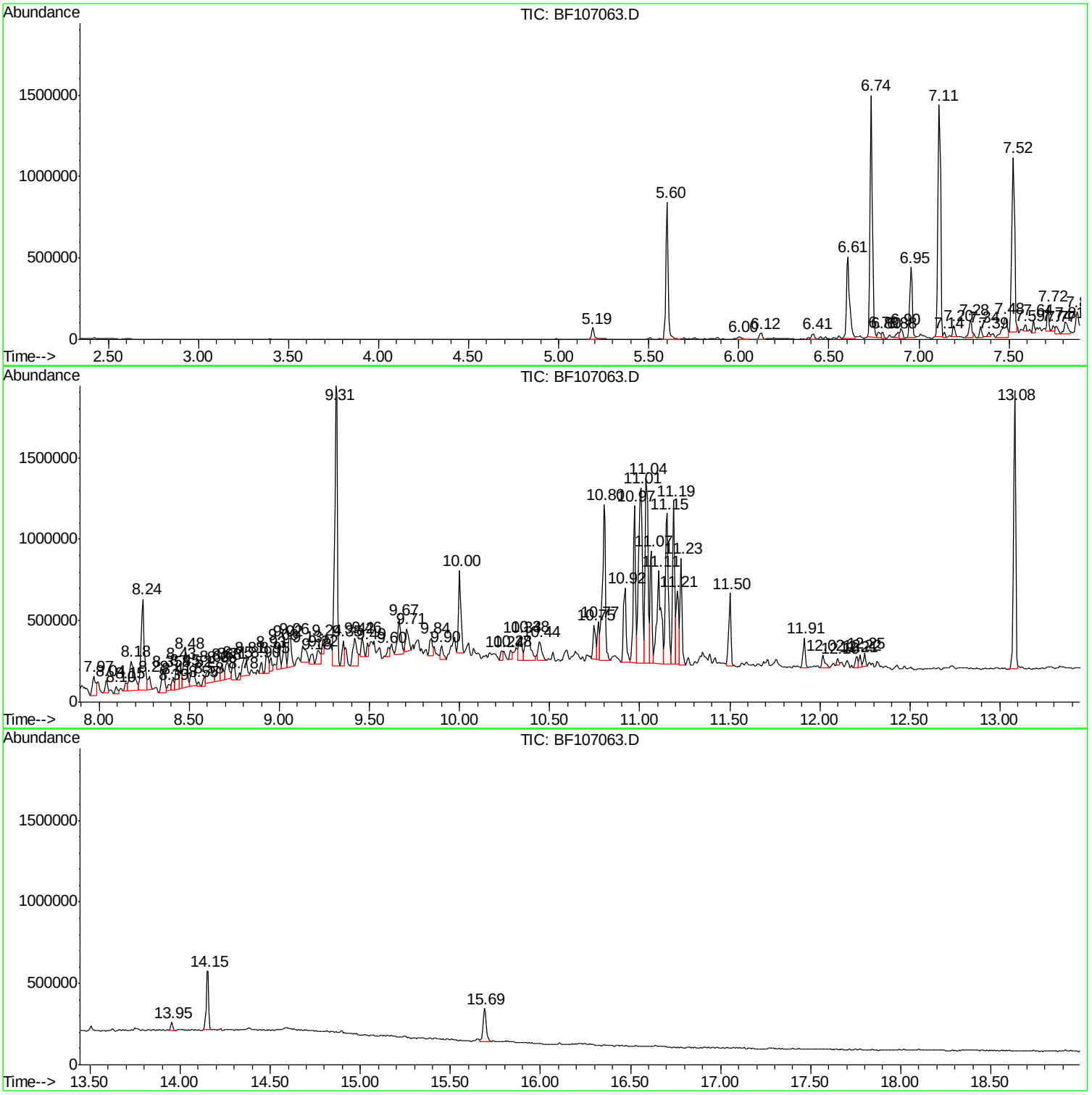
Sample : J3799-07
Misc :
ALS Vial : 31 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** MW-6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



# Library Search Compound Report

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

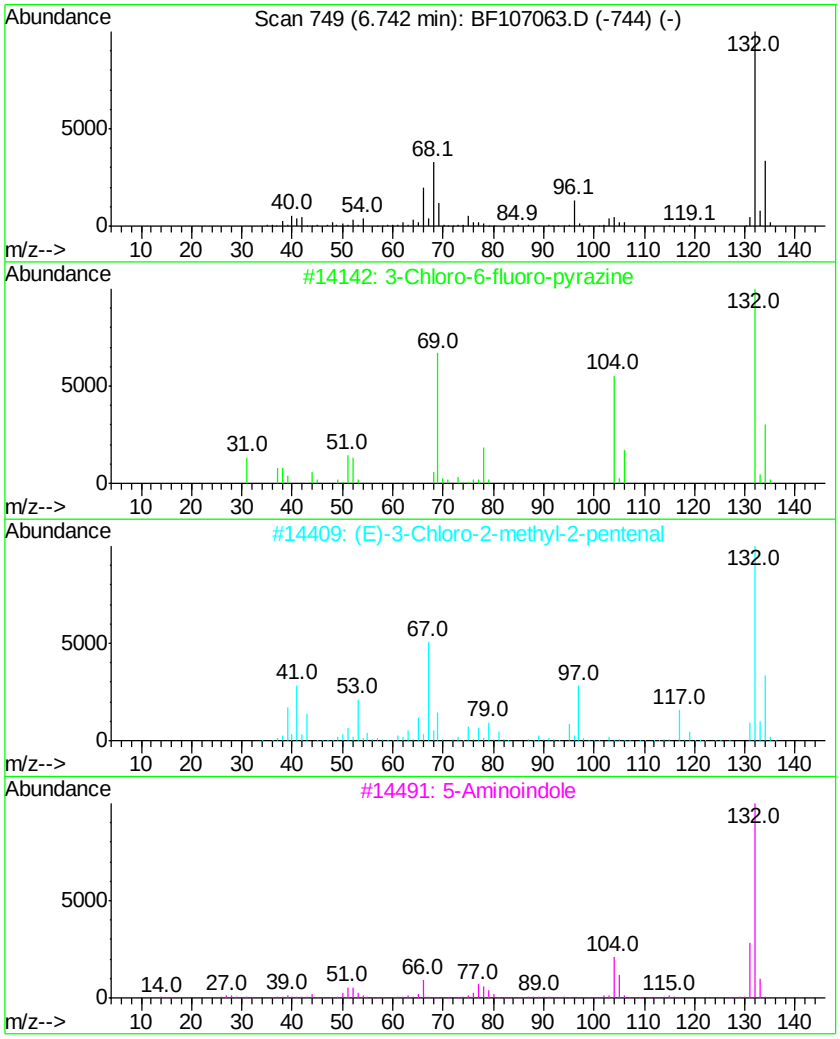
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

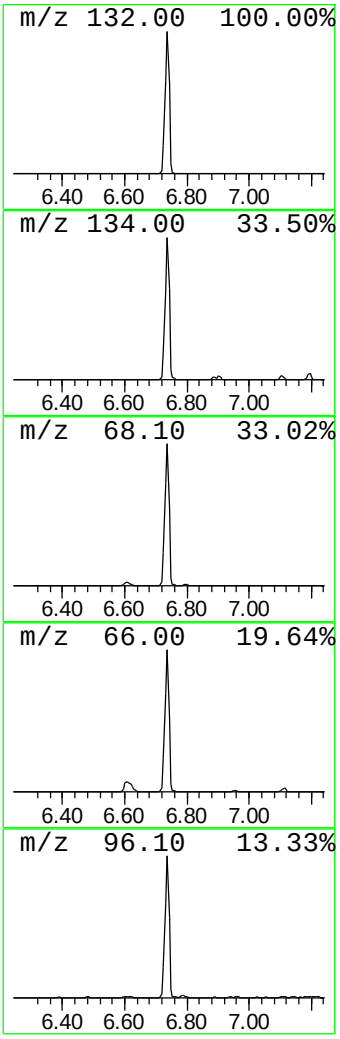
**********

Peak Number 1 unknown6.74 Concentration Rank 2

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 6.74 | 71.27 ng | 1377250 | 1,4-Dichlorobenzene-d4 | 6.95 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 3-Chloro-6-fluoro-pyrazine | 132 | C4H2ClFN2 | 1000146-10-7 | 27 |
| 2 | (E)-3-Chloro-2-methyl-2-pentenal | 132 | C6H9ClO | 031357-76-3 | 16 |
| 3 | 5-Aminoindole | 132 | C8H8N2 | 005192-03-0 | 12 |
| 4 | 1H-Benzimidazole, 2-methyl- | 132 | C8H8N2 | 000615-15-6 | 10 |
| 5 | Tranylcypromine-propionyl | 189 | C12H15NO | 1000123-86-3 | 10 |



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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

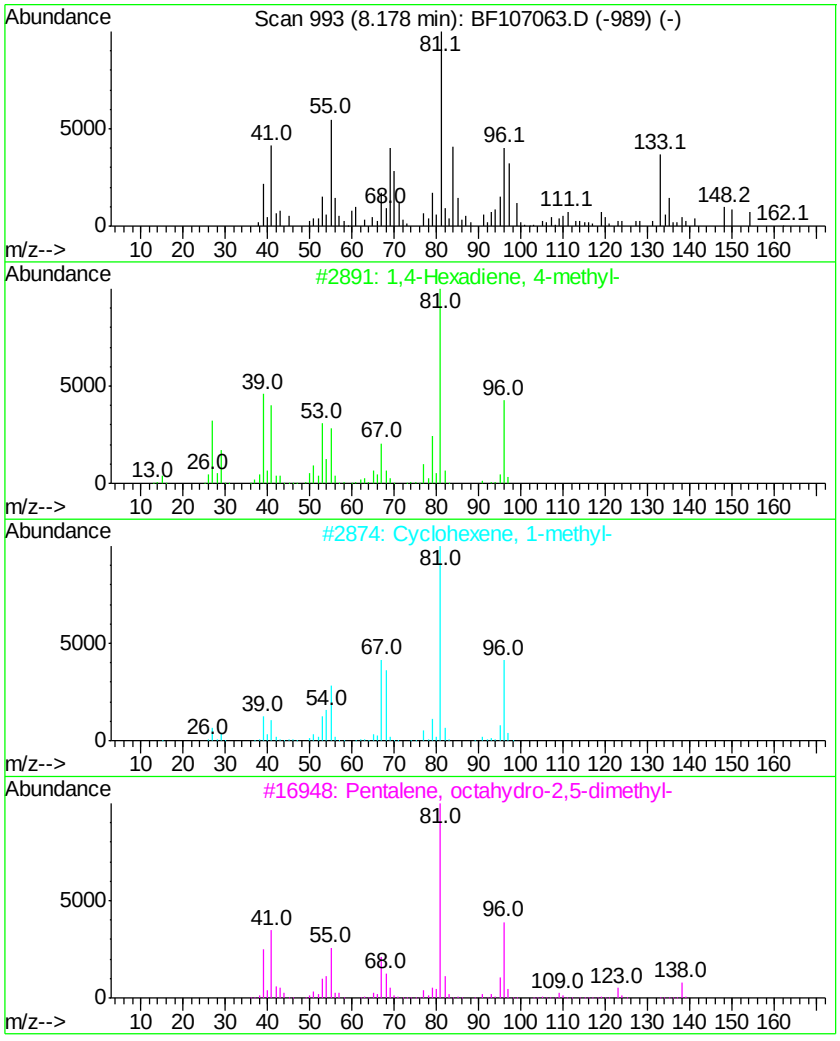
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

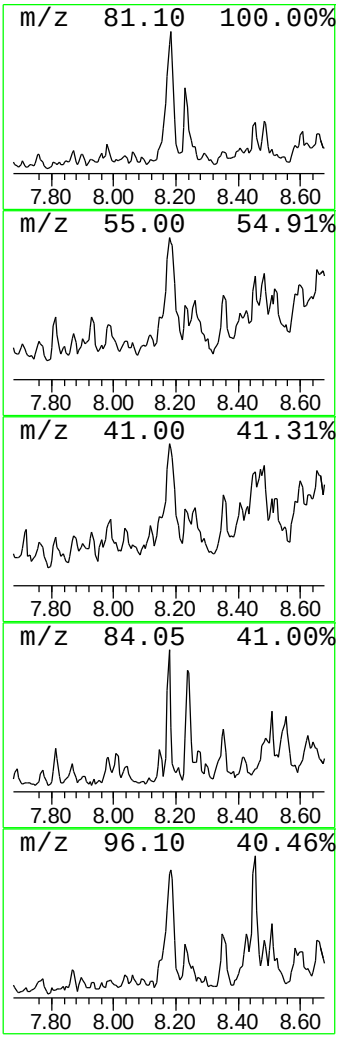
**********

Peak Number 3 1,4-Hexadiene, 4-methyl- Concentration Rank 15

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 8.18 | 10.84 ng | 368713 | Naphthalene-d8 | 8.24 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1,4-Hexadiene, 4-methyl- | 96 | C7H12 | 001116-90-1 | 50 |
| 2 | Cyclohexene, 1-methyl- | 96 | C7H12 | 000591-49-1 | 43 |
| 3 | Pentalene, octahydro-2,5-dimethyl- | 138 | C10H18 | 028588-55-8 | 43 |
| 4 | Cyclohexene, 3-methyl- | 96 | C7H12 | 000591-48-0 | 43 |
| 5 | Cyclohexene, 4-methyl- | 96 | C7H12 | 000591-47-9 | 38 |



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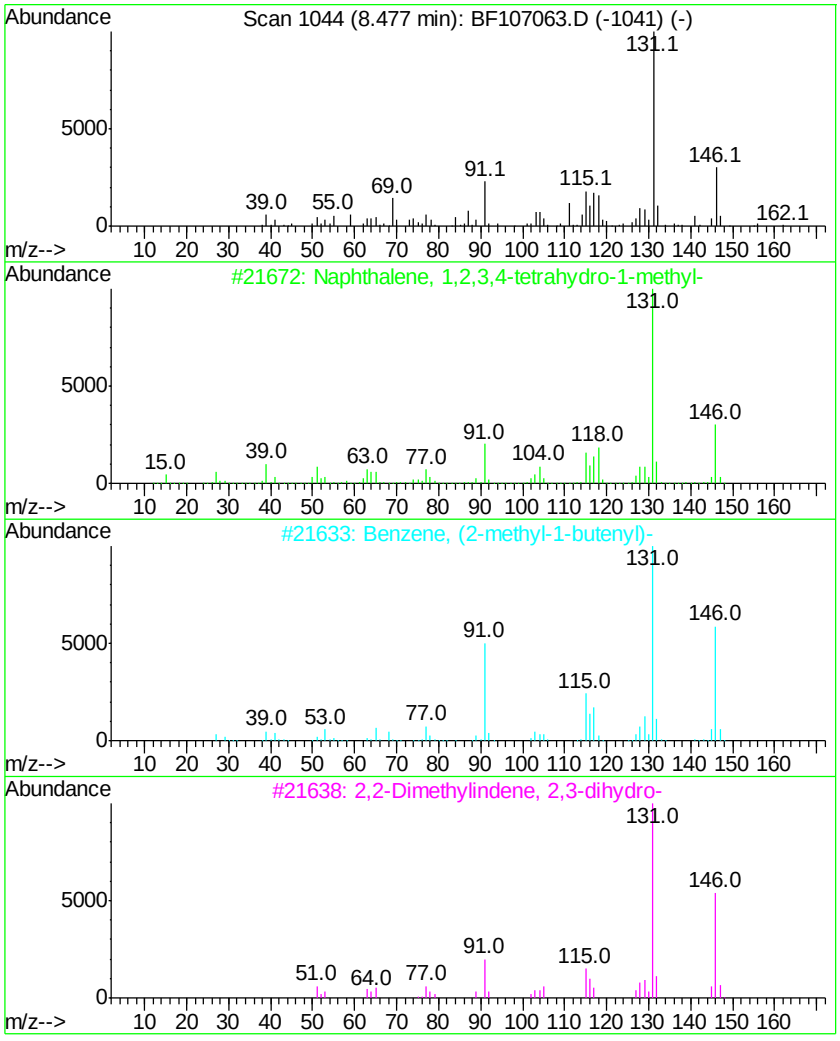
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

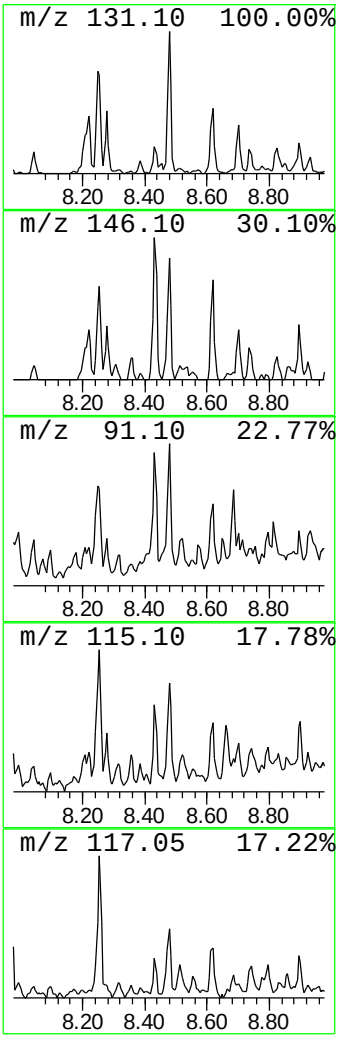
**********

Peak Number 4 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 16

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 8.48 | 8.98 ng | 305390 | Naphthalene-d8 | 8.24 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Naphthalene, 1,2,3,4-tetrahydro-... | 146 | C11H14 | 001559-81-5 | 94 |
| 2 | Benzene, (2-methyl-1-butenyl)- | 146 | C11H14 | 056253-64-6 | 94 |
| 3 | 2,2-Dimethylindene, 2,3-dihydro- | 146 | C11H14 | 020836-11-7 | 90 |
| 4 | Benzene, (1,1-dimethyl-2-propenyl)- | 146 | C11H14 | 018321-36-3 | 81 |
| 5 | Benzene, (3-methyl-2-butenyl)- | 146 | C11H14 | 004489-84-3 | 81 |



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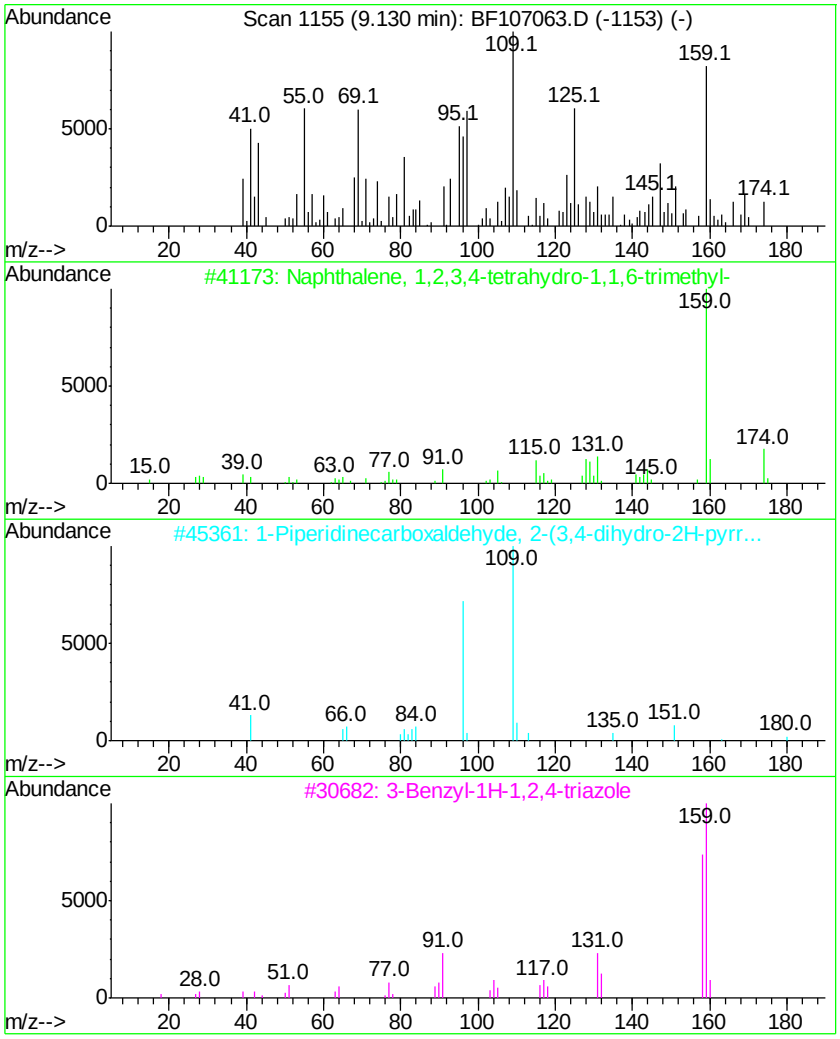
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

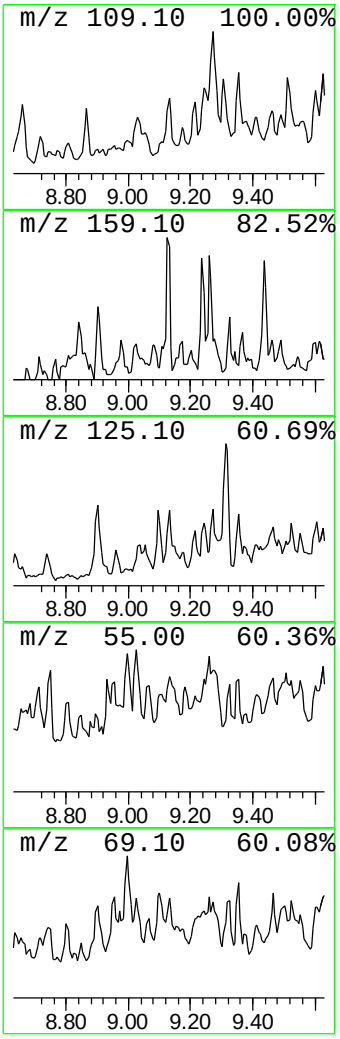
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 5 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 19

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 9.13 | 7.45 ng | 168816 | Acenaphthene-d10 | 10.00 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Naphthalene, 1,2,3,4-tetrahydro-... | 174 | C13H18 | 000475-03-6 | 58 |
| 2 | 1-Piperidinecarboxaldehyde, 2-(3... | 180 | C10H16N2O | 052196-11-9 | 22 |
| 3 | 3-Benzyl-1H-1,2,4-triazole | 159 | C9H9N3 | 021117-34-0 | 11 |
| 4 | 1H-Indene, 2,3-dihydro-1,1,4,6-t... | 174 | C13H18 | 000941-60-6 | 11 |
| 5 | Indan, 1,1,6,7-tetramethyl- | 174 | C13H18 | 016204-58-3 | 11 |



# Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070218\
Data File : BF107063.D
Acq On : 3 Jul 2018 1:10
Operator : JU/SJ
Sample : J3799-07
Misc :
ALS Vial : 31 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** MW-6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

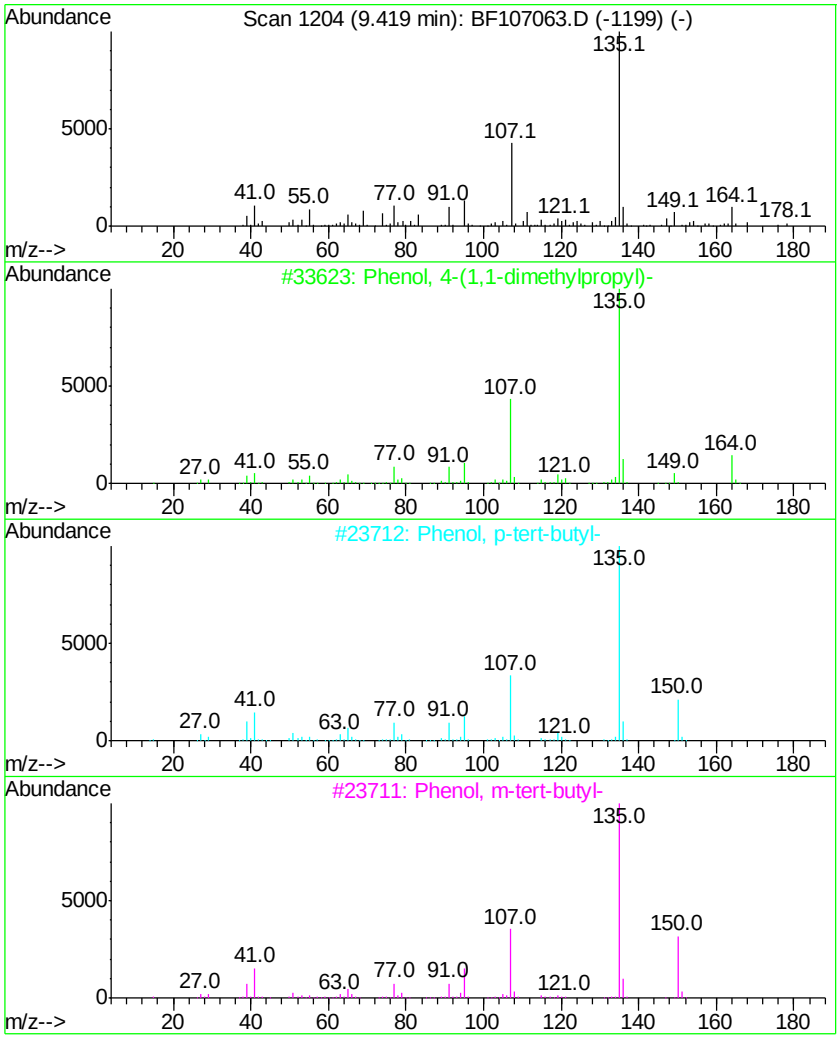
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

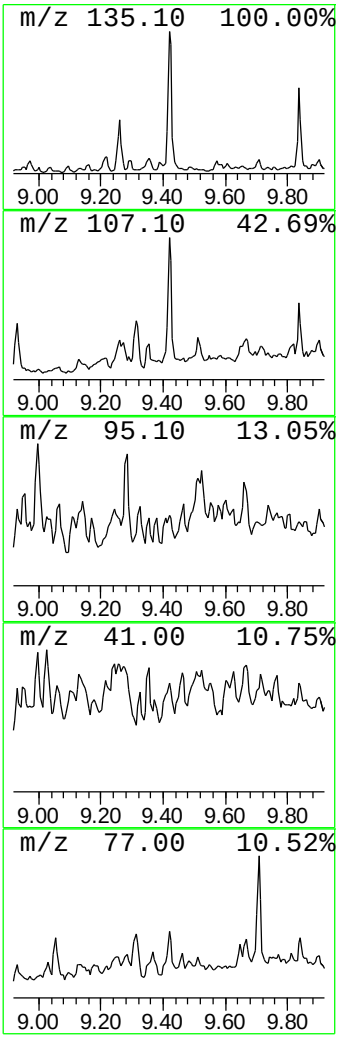
**********

Peak Number 6 Phenol, 4-(1,1-dimethylprop... Concentration Rank 14

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 9.42 | 11.57 ng | 262174 | Acenaphthene-d10 | 10.00 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, 4-(1,1-dimethylpropyl)- | 164 | C11H16O | 000080-46-6 | 94 |
| 2 | Phenol, p-tert-butyl- | 150 | C10H14O | 000098-54-4 | 72 |
| 3 | Phenol, m-tert-butyl- | 150 | C10H14O | 000585-34-2 | 64 |
| 4 | Hexestrol | 270 | C18H22O2 | 005635-50-7 | 59 |
| 5 | Hexestrol, O-trifluoroacetyl- | 366 | C20H21F3O3 | 1000365-31-7 | 59 |



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Data File : BF107063.D
Acq On : 3 Jul 2018 1:10
Operator : JU/SJ
Sample : J3799-07
Misc :
ALS Vial : 31 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** MW-6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

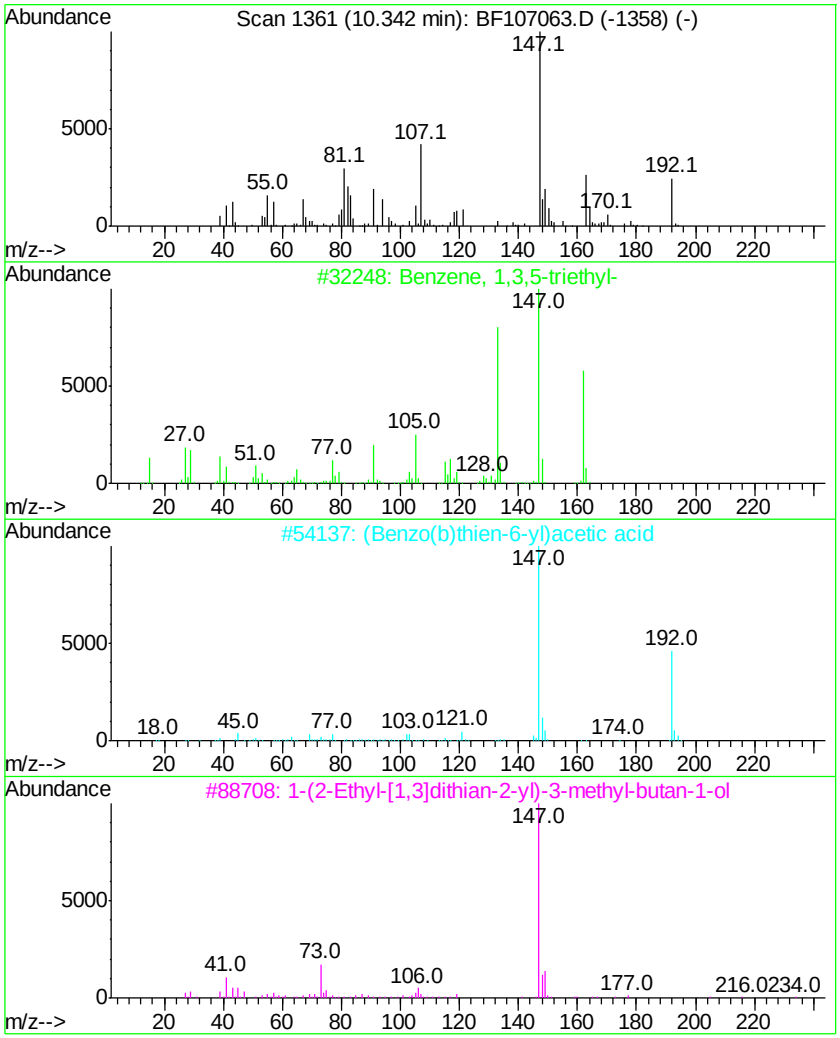
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

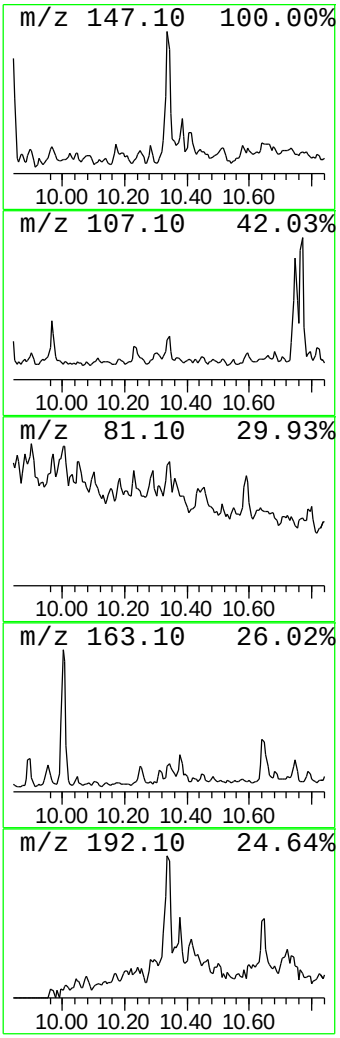
**********

Peak Number 7 Benzene, 1,3,5-triethyl- Concentration Rank 20

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.34 | 7.40 ng | 167647 | Acenaphthene-d10 | 10.00 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzene, 1,3,5-triethyl- | 162 | C12H18 | 000102-25-0 | 50 |
| 2 | (Benzo(b)thien-6-yl)acetic acid | 192 | C10H8O2S | 006177-87-3 | 50 |
| 3 | 1-(2-Ethyl-[1,3]dithian-2-yl)-3-... | 234 | C11H22OS2 | 332048-49-4 | 43 |
| 4 | 1,3-Benzenediol, o-butoxycarbony... | 292 | C15H16O6 | 1000330-80-3 | 38 |
| 5 | (3-Methoxyphenyl)acetonitrile | 147 | C9H9NO | 019924-43-7 | 38 |



# Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070218\
Data File : BF107063.D
Acq On : 3 Jul 2018 1:10
Operator : JU/SJ
Sample : J3799-07
Misc :
ALS Vial : 31 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** MW-6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

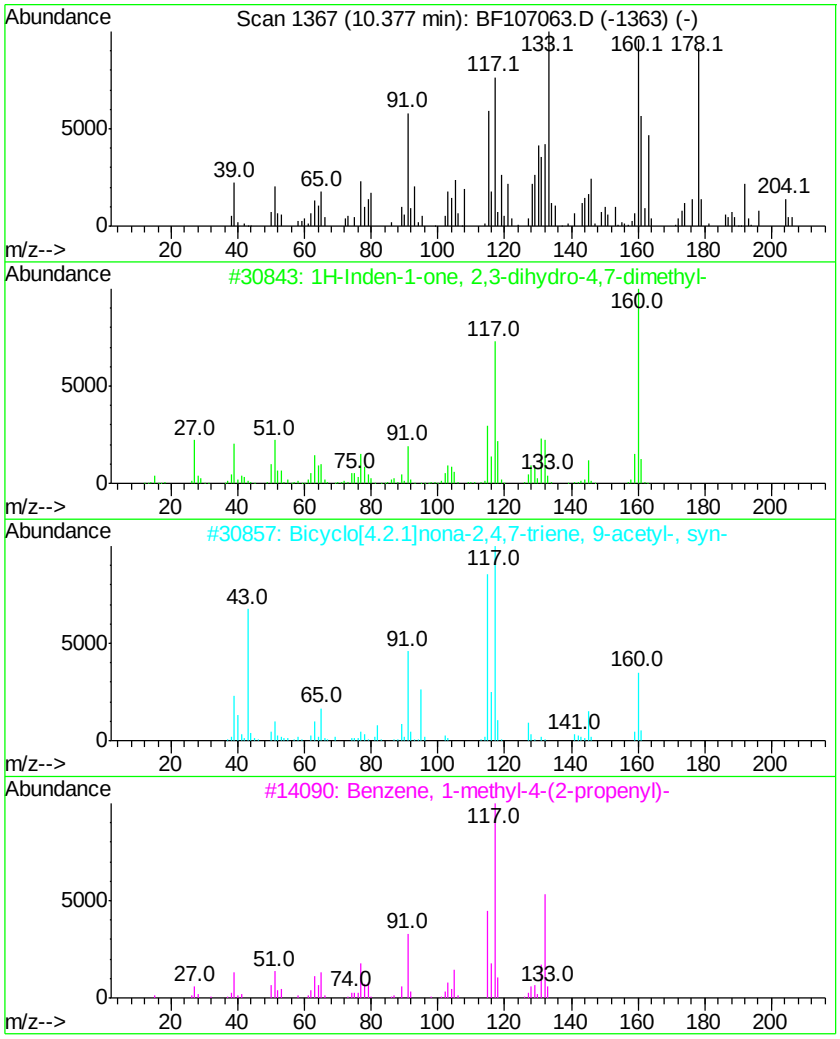
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

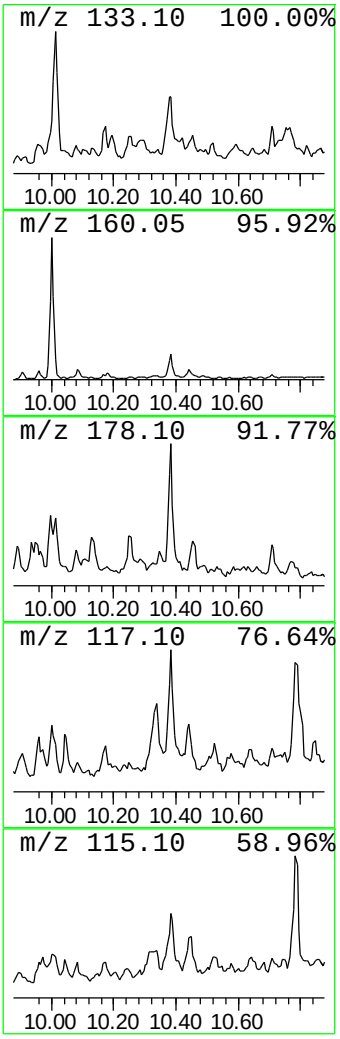
**********

Peak Number 8 unknown10.38 Concentration Rank 13

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.38 | 14.13 ng | 320351 | Acenaphthene-d10 | 10.00 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 1H-Inden-1-one, 2,3-dihydro-4,7-... | 160 | C11H12O | 005037-60-5 | 38 |
| 2 | Bicyclo[4.2.1]nona-2,4,7-triene,... | 160 | C11H12O | 1000141-51-6 | 30 |
| 3 | Benzene, 1-methyl-4-(2-propenyl)- | 132 | C10H12 | 003333-13-9 | 25 |
| 4 | 1H-Indole-2,3-dione, 6-methyl- | 161 | C9H7NO2 | 001128-47-8 | 25 |
| 5 | Benzene, 2-ethenyl-1,4-dimethyl- | 132 | C10H12 | 002039-89-6 | 25 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070218\
Data File : BF107063.D
Acq On : 3 Jul 2018 1:10
Operator : JU/SJ
Sample : J3799-07
Misc :
ALS Vial : 31 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** MW-6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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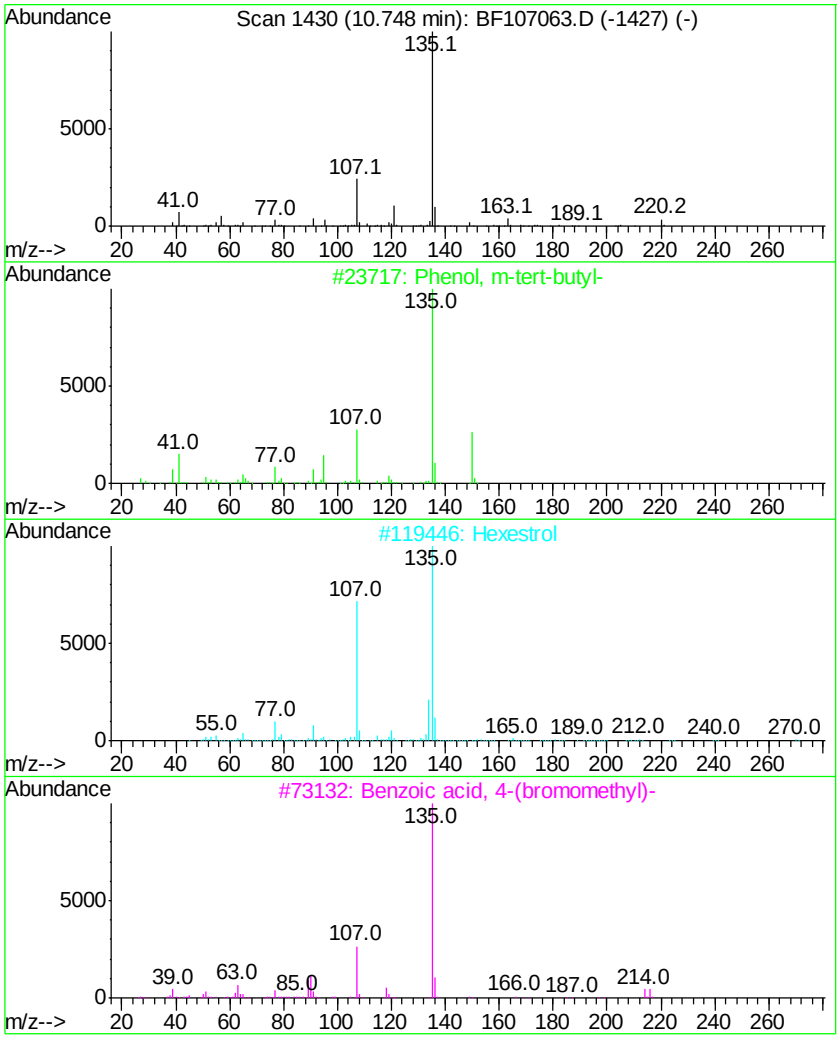
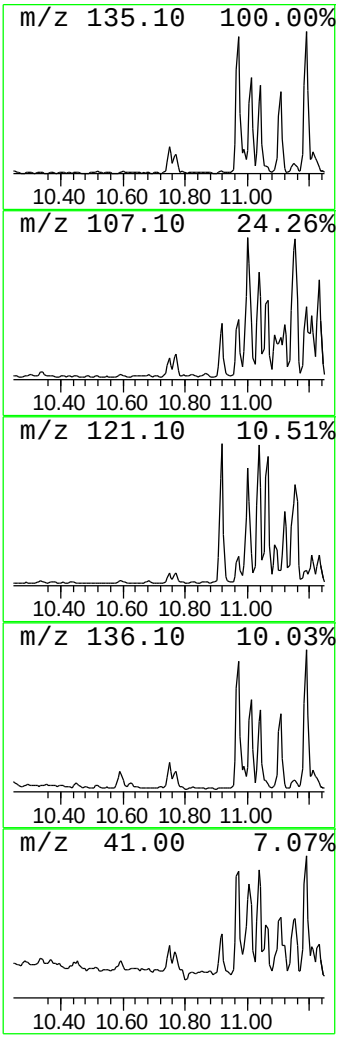
Peak Number 9 Phenol, m-tert-butyl- Concentration Rank 17

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.75 | 8.56 ng | 193974 | Acenaphthene-d10 | 10.00 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, m-tert-butyl- | 150 | C10H14O | 000585-34-2 | 64 |
| 2 | Hexestrol | 270 | C18H22O2 | 000084-16-2 | 59 |
| 3 | Benzoic acid, 4-(bromomethyl)- | 214 | C8H7BrO2 | 006232-88-8 | 59 |
| 4 | Phenol, 4-(1,1,3,3-tetramethylbu... | 206 | C14H22O | 000140-66-9 | 53 |
| 5 | 1-Adamantyl bromomethyl ketone | 256 | C12H17BrO | 005122-82-7 | 53 |



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070218\
Data File : BF107063.D
Acq On : 3 Jul 2018 1:10
Operator : JU/SJ
Sample : J3799-07
Misc :
ALS Vial : 31 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** MW-6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

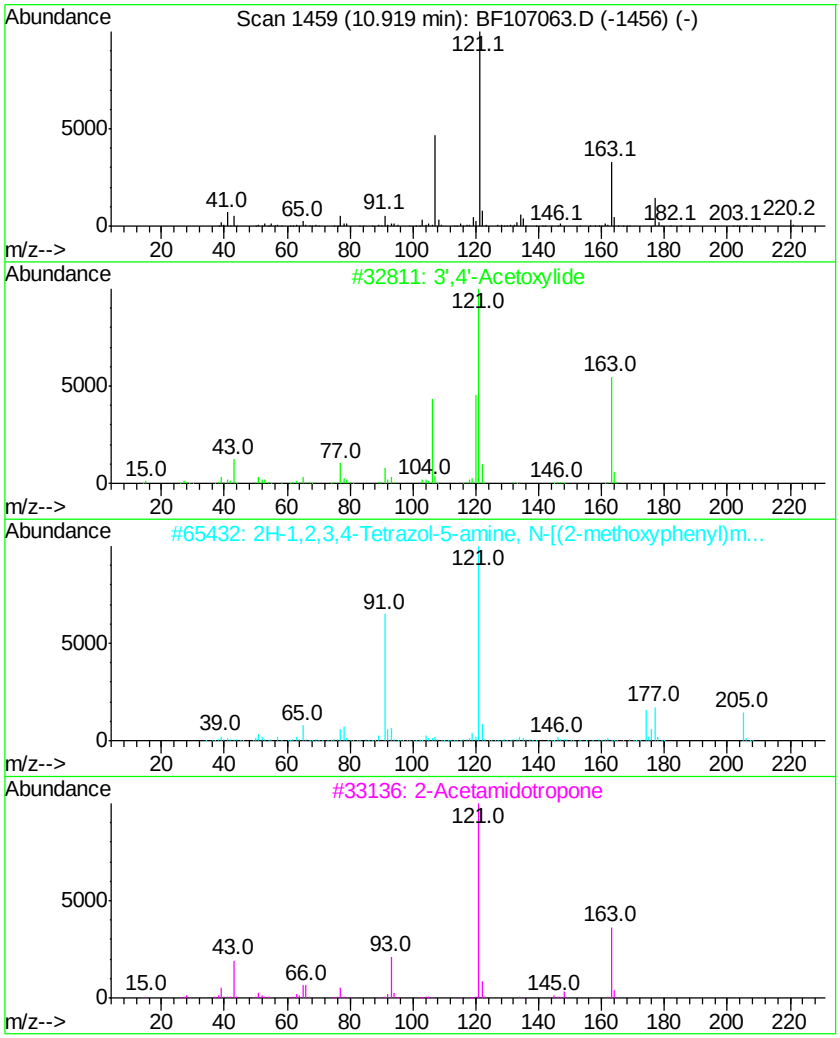
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

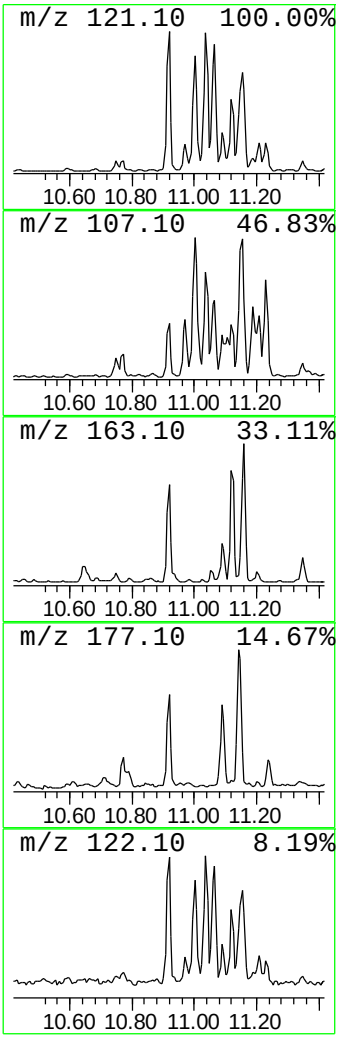
**********

Peak Number 10 3',4'-Acetoxylide Concentration Rank 11

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.92 | 23.66 ng | 423437 | Phenanthrene-d10 | 11.50 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 3',4'-Acetoxylide | 163 | C10H13NO | 002198-54-1 | 52 |
| 2 | 2H-1,2,3,4-Tetrazol-5-amine, N-[... | 205 | C9H11N5O | 1000337-56-1 | 47 |
| 3 | 2-Acetamidotropone | 163 | C9H9NO2 | 006422-12-4 | 47 |
| 4 | Carbonic acid, isobutyl 4-isopro... | 236 | C14H20O3 | 1000331-40-6 | 38 |
| 5 | N-Allyl-p-hydroxybenzamide | 177 | C10H11NO2 | 090921-72-5 | 38 |



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Data File : BF107063.D
Acq On : 3 Jul 2018 1:10
Operator : JU/SJ
Sample : J3799-07
Misc :
ALS Vial : 31 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** MW-6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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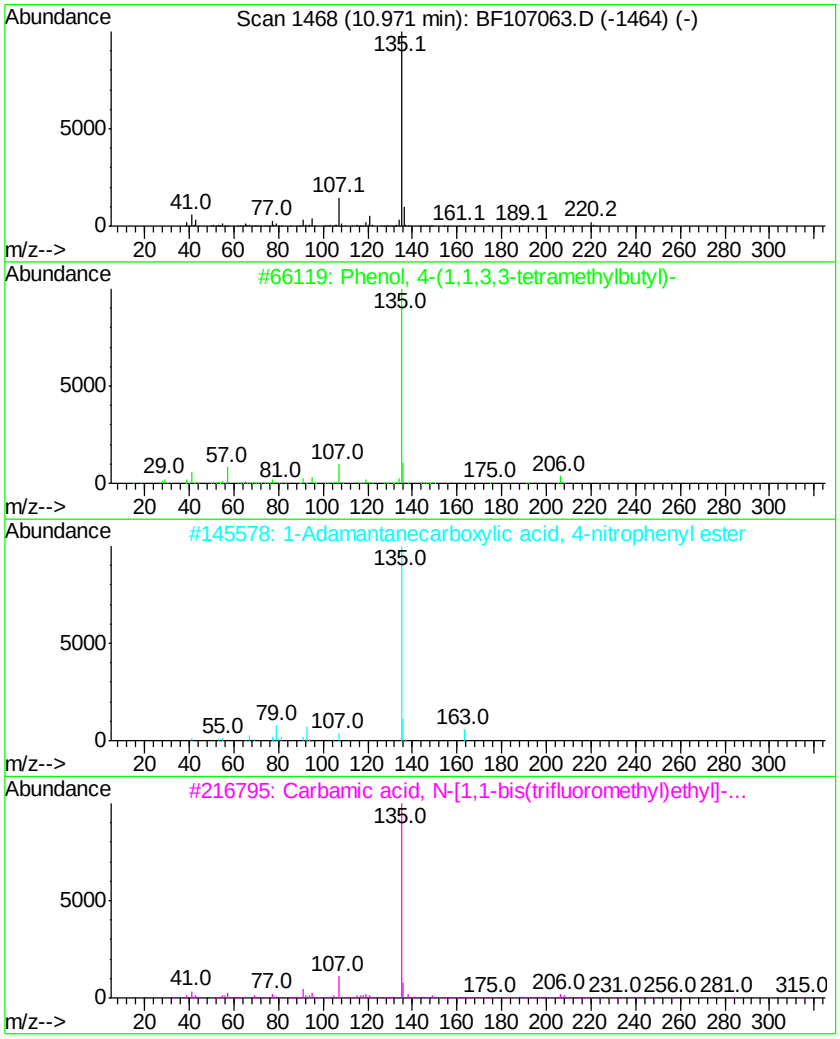
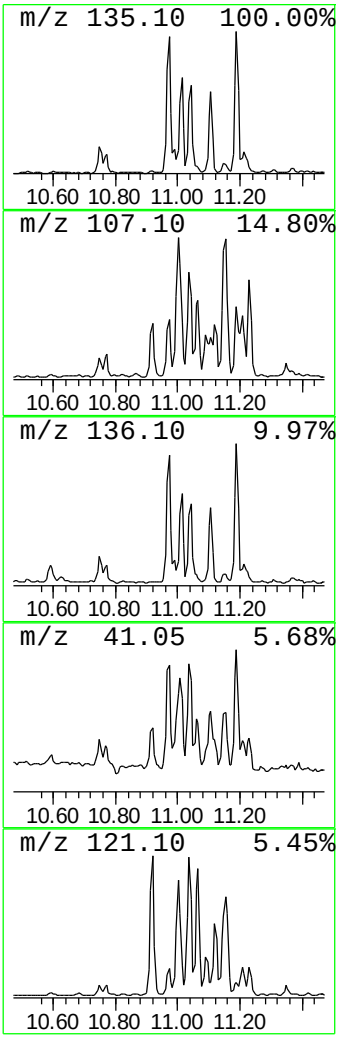
Peak Number 11 Phenol, 4-(1,1,3,3-tetramet... Concentration Rank 8

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.97 | 50.70 ng | 907361 | Phenanthrene-d10 | 11.50 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, 4-(1,1,3,3-tetramethylbu... | 206 | C14H22O | 000140-66-9 | 78 |
| 2 | 1-Adamantanecarboxylic acid, 4-n... | 301 | C17H19NO4 | 1000307-66-5 | 64 |
| 3 | Carbamic acid, N-[1,1-bis(triflu... | 413 | C19H25F6NO2 | 296242-69-8 | 64 |
| 4 | 1-n-Hexyladamantane | 220 | C16H28 | 022458-75-9 | 64 |
| 5 | 4-tert-Butylphenyl acetate | 192 | C12H16O2 | 003056-64-2 | 64 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070218\
Data File : BF107063.D
Acq On : 3 Jul 2018 1:10
Operator : JU/SJ
Sample : J3799-07
Misc :
ALS Vial : 31 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** MW-6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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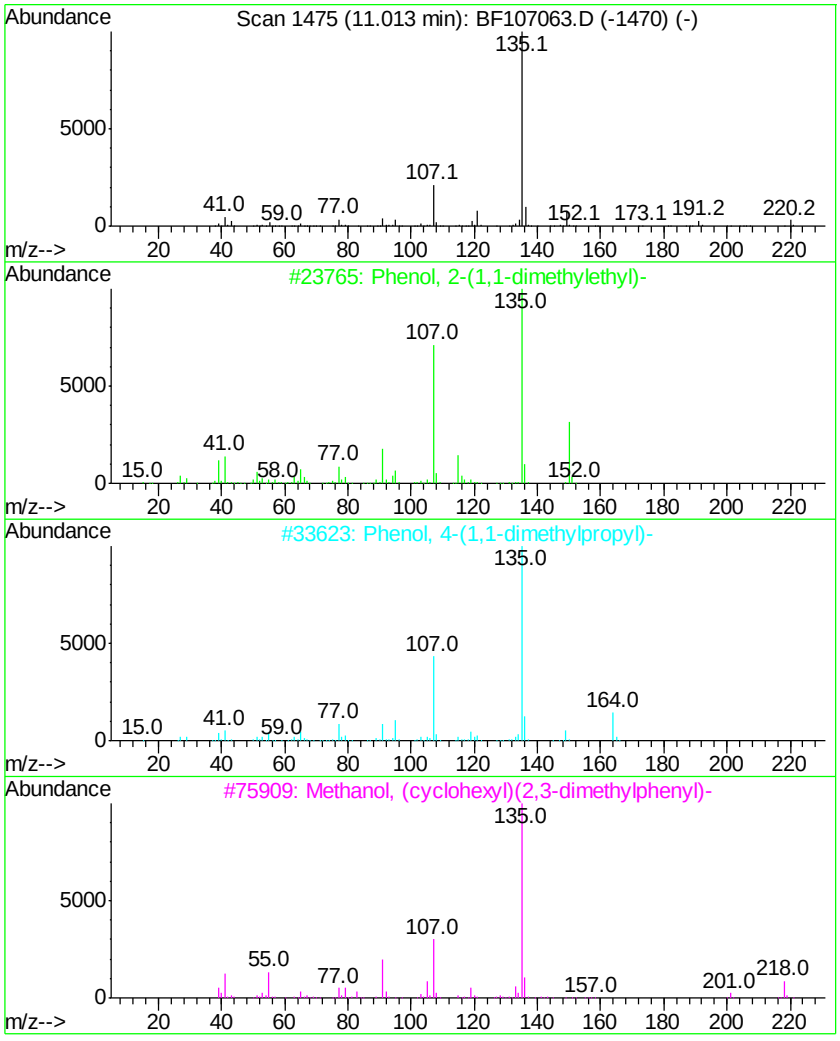
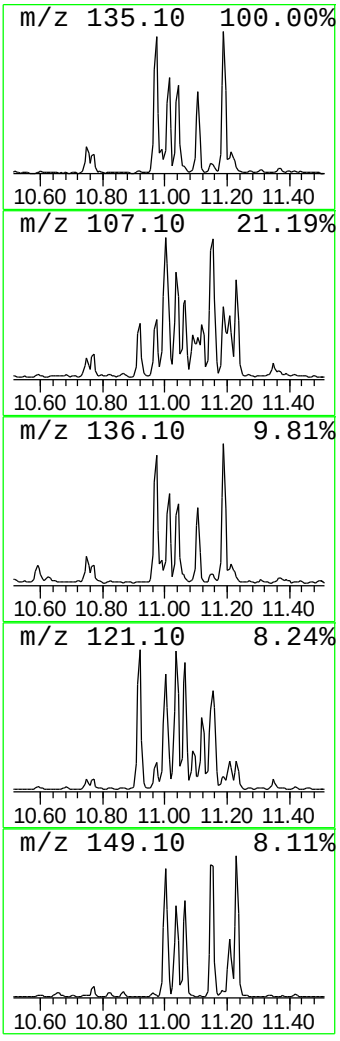
Peak Number 12 Phenol, 2-(1,1-dimethylethyl)- Concentration Rank 1

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.01 | 83.31 ng | 1491140 | Phenanthrene-d10 | 11.50 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, 2-(1,1-dimethylethyl)- | 150 | C10H14O | 000088-18-6 | 52 |
| 2 | Phenol, 4-(1,1-dimethylpropyl)- | 164 | C11H16O | 000080-46-6 | 50 |
| 3 | Methanol, (cyclohexyl)(2,3-dimet... | 218 | C15H22O | 349498-47-1 | 27 |
| 4 | Phenol, 4-propyl- | 136 | C9H12O | 000645-56-7 | 22 |
| 5 | 1-Adamantan-1-yl-6-chloro-7-fluo... | 389 | C21H21ClFNO3 | 1000210-85-2 | 22 |



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070218\
Data File : BF107063.D
Acq On : 3 Jul 2018 1:10
Operator : JU/SJ
Sample : J3799-07
Misc :
ALS Vial : 31 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** MW-6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

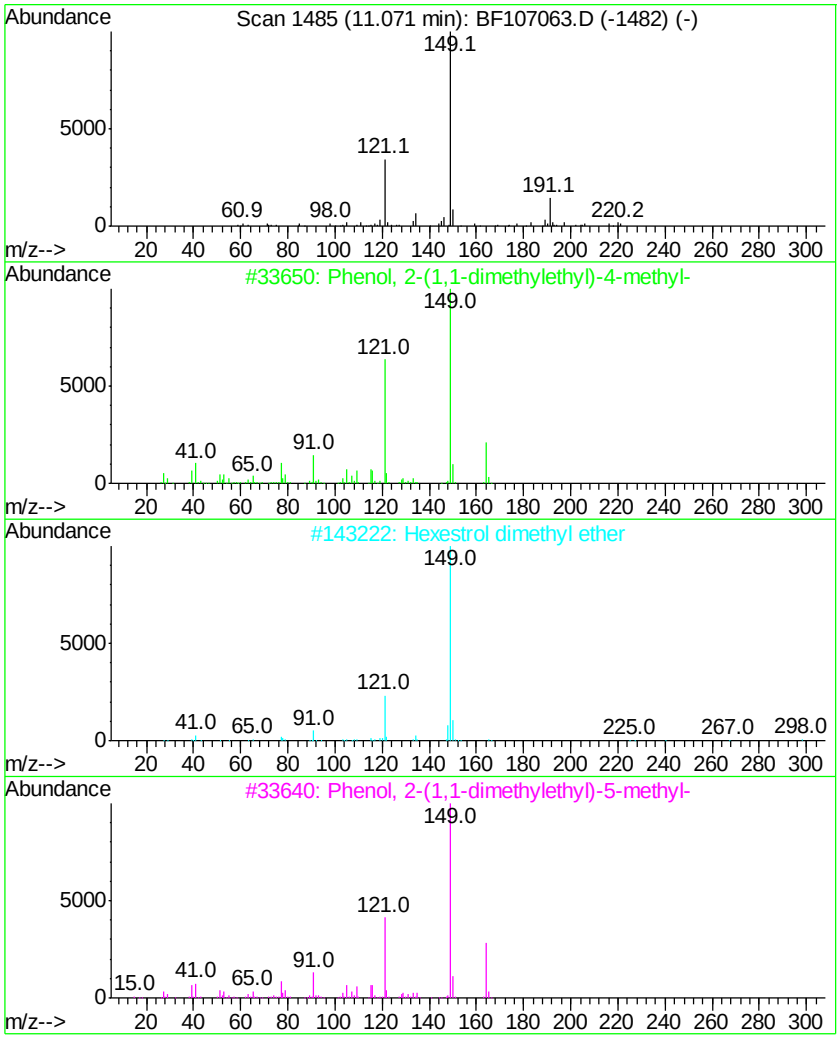
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

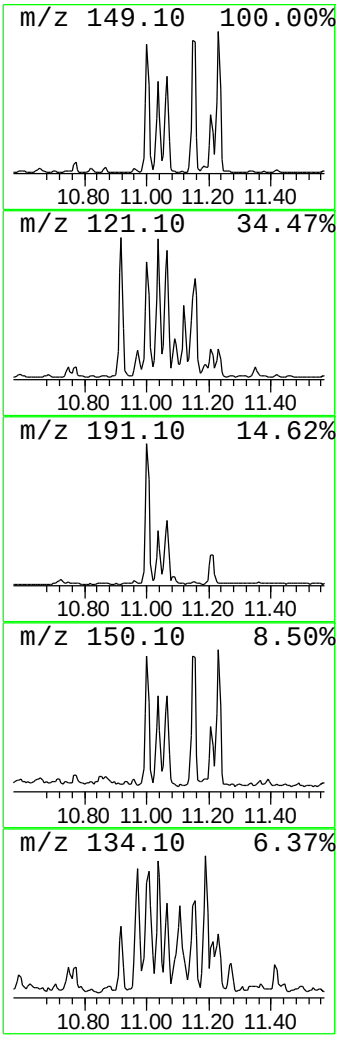
**********

Peak Number 14 unknown11.07 Concentration Rank 9

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.07 | 31.10 ng | 556576 | Phenanthrene-d10 | 11.50 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, 2-(1,1-dimethylethyl)-4-... | 164 | C11H16O | 002409-55-4 | 45 |
| 2 | Hexestrol dimethyl ether | 298 | C20H26O2 | 000130-78-9 | 45 |
| 3 | Phenol, 2-(1,1-dimethylethyl)-5-... | 164 | C11H16O | 000088-60-8 | 45 |
| 4 | 4-Ethoxybenzoyl chloride | 184 | C9H9ClO2 | 016331-46-7 | 42 |
| 5 | 4-Isopropoxy-4-phenylbutyronitrile | 203 | C13H17NO | 1000370-35-3 | 37 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070218\
Data File : BF107063.D
Acq On : 3 Jul 2018 1:10
Operator : JU/SJ
Sample : J3799-07
Misc :
ALS Vial : 31 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** MW-6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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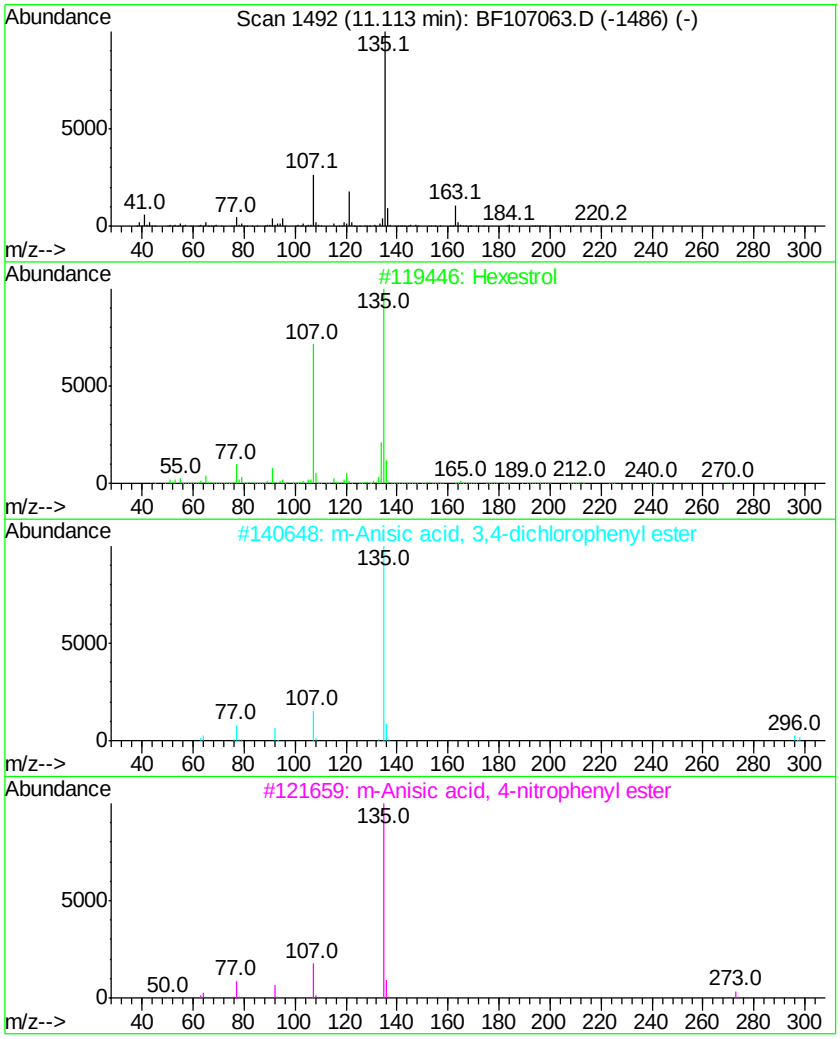
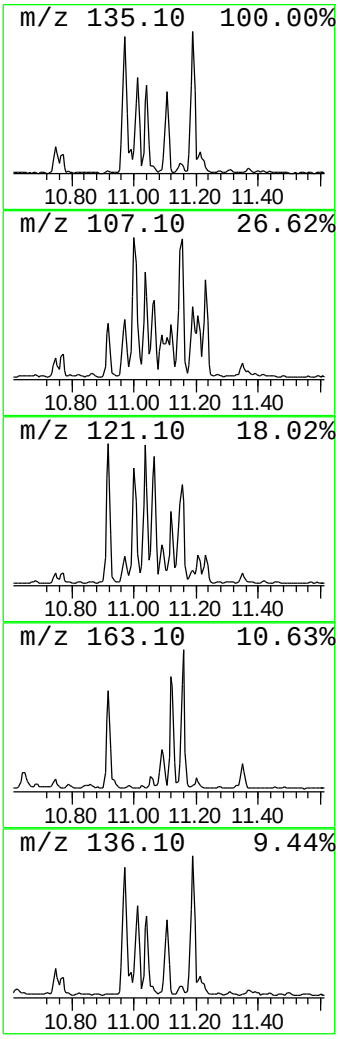
Peak Number 15 Hexestrol Concentration Rank 7

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.11 | 52.36 ng | 937143 | Phenanthrene-d10 | 11.50 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Hexestrol | 270 | C18H22O2 | 000084-16-2 | 56 |
| 2 | m-Anisic acid, 3,4-dichlorophenyl... | 296 | C14H10Cl2O3 | 1000307-80-3 | 39 |
| 3 | m-Anisic acid, 4-nitrophenyl ester | 273 | C14H11NO5 | 1000307-80-4 | 39 |
| 4 | 1-Adamantanecarboxylic acid, 2-a... | 314 | C21H30O2 | 1000282-94-3 | 39 |
| 5 | 1-Adamantanethiol, S-methacryloyl- | 236 | C14H20OS | 1000358-48-9 | 38 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070218\
Data File : BF107063.D
Acq On : 3 Jul 2018 1:10
Operator : JU/SJ
Sample : J3799-07
Misc :
ALS Vial : 31 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** MW-6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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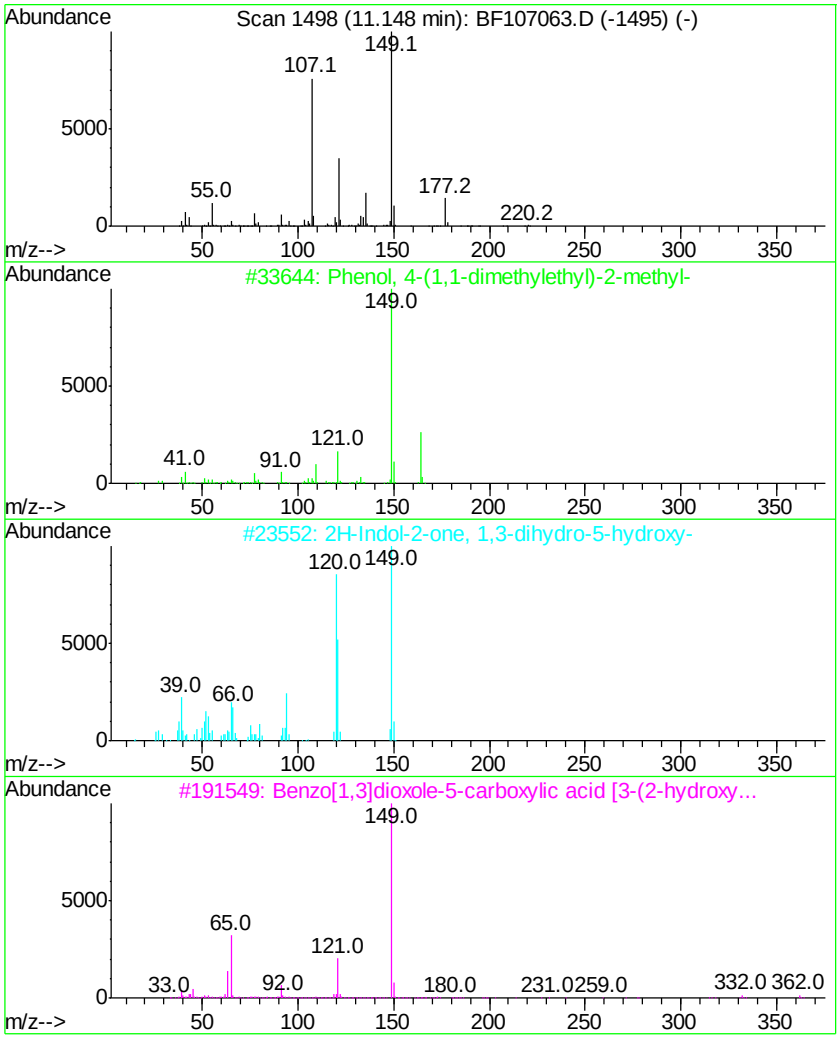
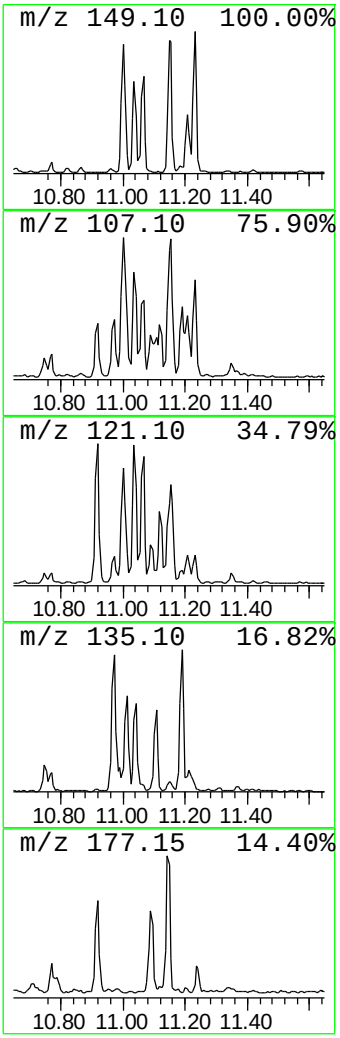
Peak Number 16 unknown11.15 Concentration Rank 5

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.15 | 60.84 ng | 1088970 | Phenanthrene-d10 | 11.50 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, 4-(1,1-dimethylethyl)-2-... | 164 | C11H16O | 000098-27-1 | 38 |
| 2 | 2H-Indol-2-one, 1,3-dihydro-5-hy... | 149 | C8H7NO2 | 003416-18-0 | 35 |
| 3 | Benzo[1,3]dioxole-5-carboxylic a... | 362 | C16H14N2O6S | 1000296-89-7 | 35 |
| 4 | Phenol, 2-(1,1-dimethylethyl)-4-... | 164 | C11H16O | 002409-55-4 | 35 |
| 5 | 1,2-propanedione,1-(3,4-methylen... | 192 | C10H8O4 | 1000378-93-0 | 35 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070218\
Data File : BF107063.D
Acq On : 3 Jul 2018 1:10
Operator : JU/SJ
Sample : J3799-07
Misc :
ALS Vial : 31 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** MW-6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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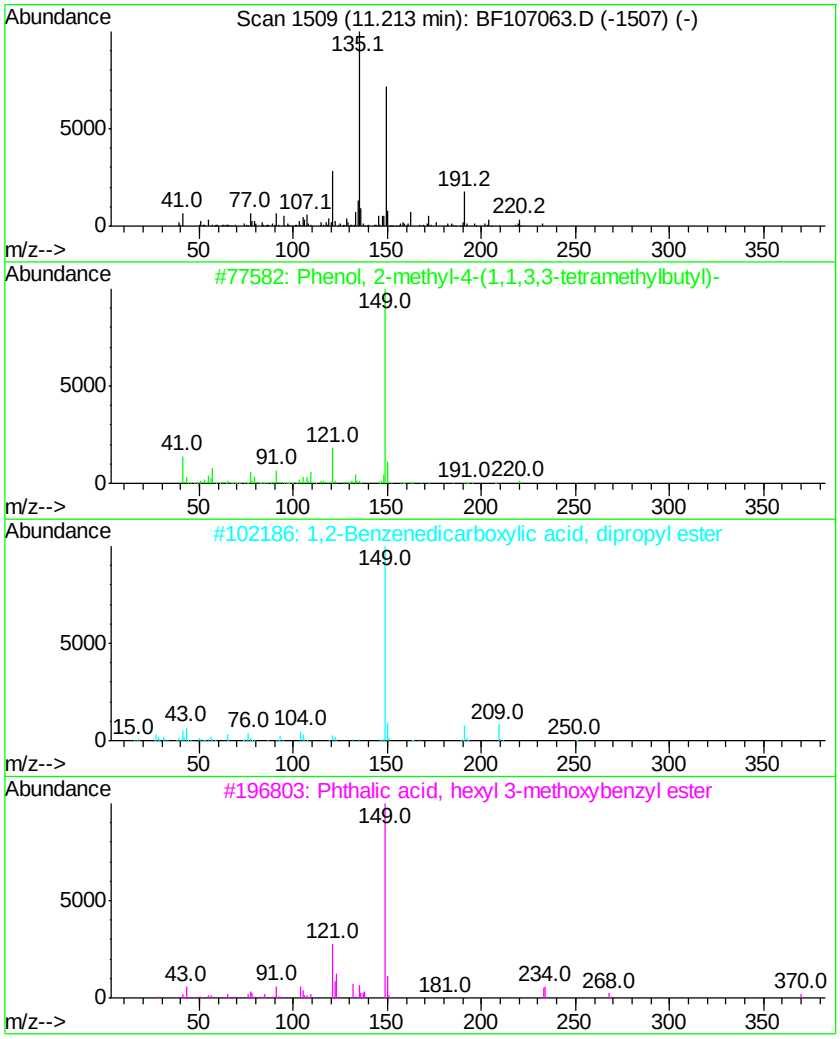
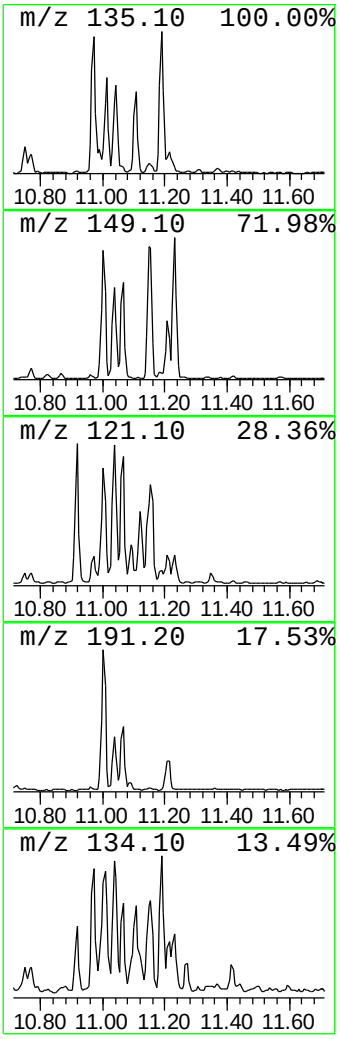
Peak Number 18 unknown11.21 Concentration Rank 12

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.21 | 20.99 ng | 375652 | Phenanthrene-d10 | 11.50 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, 2-methyl-4-(1,1,3,3-tetr... | 220 | C15H24O | 002219-84-3 | 49 |
| 2 | 1,2-Benzenedicarboxylic acid, di... | 250 | C14H18O4 | 000131-16-8 | 47 |
| 3 | Phthalic acid, hexyl 3-methoxybe... | 370 | C22H26O5 | 1000377-97-9 | 47 |
| 4 | Phthalic acid, cyclohexylmethyl ... | 304 | C18H24O4 | 1000309-08-7 | 47 |
| 5 | Phthalic acid, propyl tridec-2-y... | 386 | C24H34O4 | 1000315-43-6 | 47 |



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Data File : BF107063.D
Acq On : 3 Jul 2018 1:10
Operator : JU/SJ
Sample : J3799-07
Misc :
ALS Vial : 31 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** MW-6

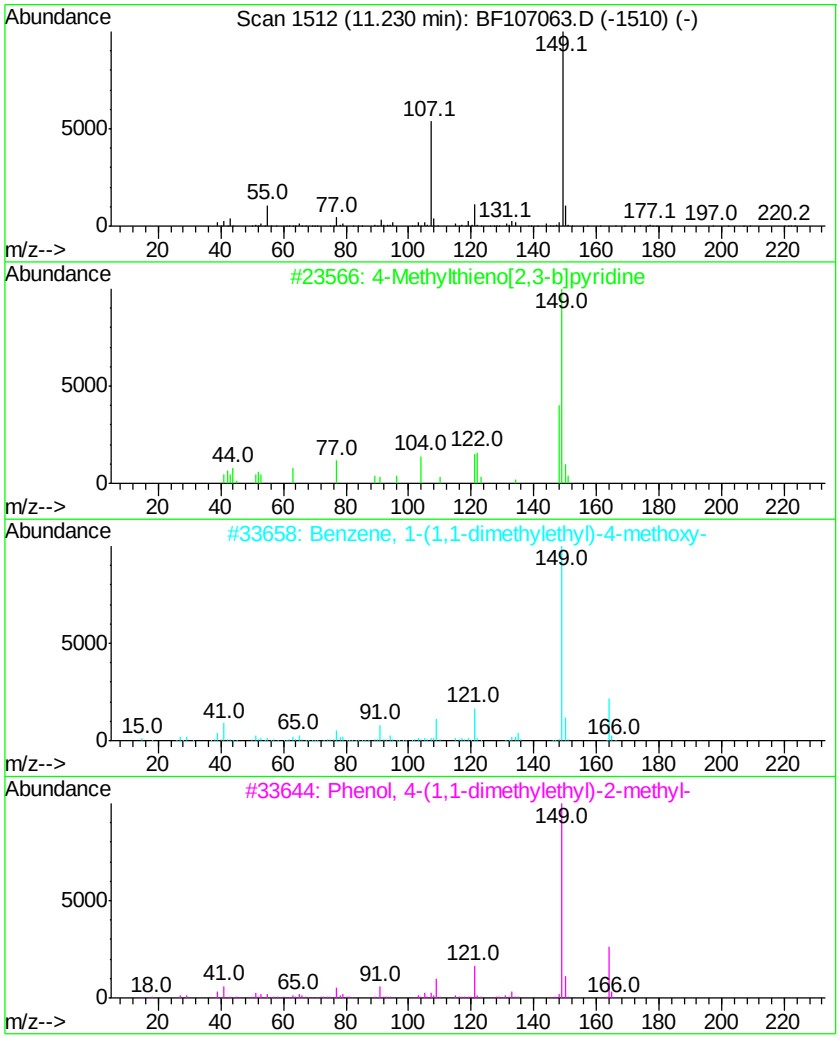
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

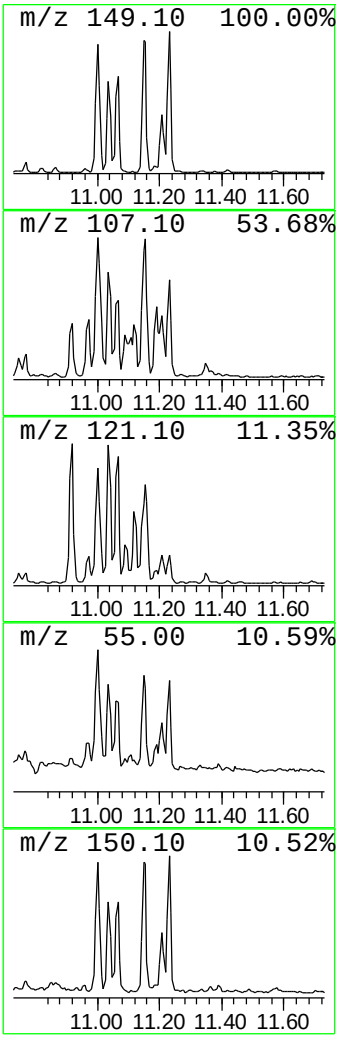
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 19 4-Methylthieno[2,3-b]pyridine Concentration Rank 10

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.23 | 30.35 ng | 543162 | Phenanthrene-d10 | 11.50 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 4-Methylthieno[2,3-b]pyridine | 149 | C8H7NS | 013362-81-7 | 50 |
| 2 | Benzene, 1-(1,1-dimethylethyl)-4... | 164 | C11H16O | 005396-38-3 | 40 |
| 3 | Phenol, 4-(1,1-dimethylethyl)-2-... | 164 | C11H16O | 000098-27-1 | 40 |
| 4 | Phthalic acid, butyl 2-propylphe... | 340 | C21H24O4 | 1000357-01-9 | 38 |
| 5 | Phthalic acid, propyl tridec-2-y... | 386 | C24H34O4 | 1000315-43-6 | 37 |



Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070218\
Data File : BF107063.D
Acq On : 3 Jul 2018 1:10
Operator : JU/SJ
Sample : J3799-07
Misc :
ALS Vial : 31 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** MW-6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

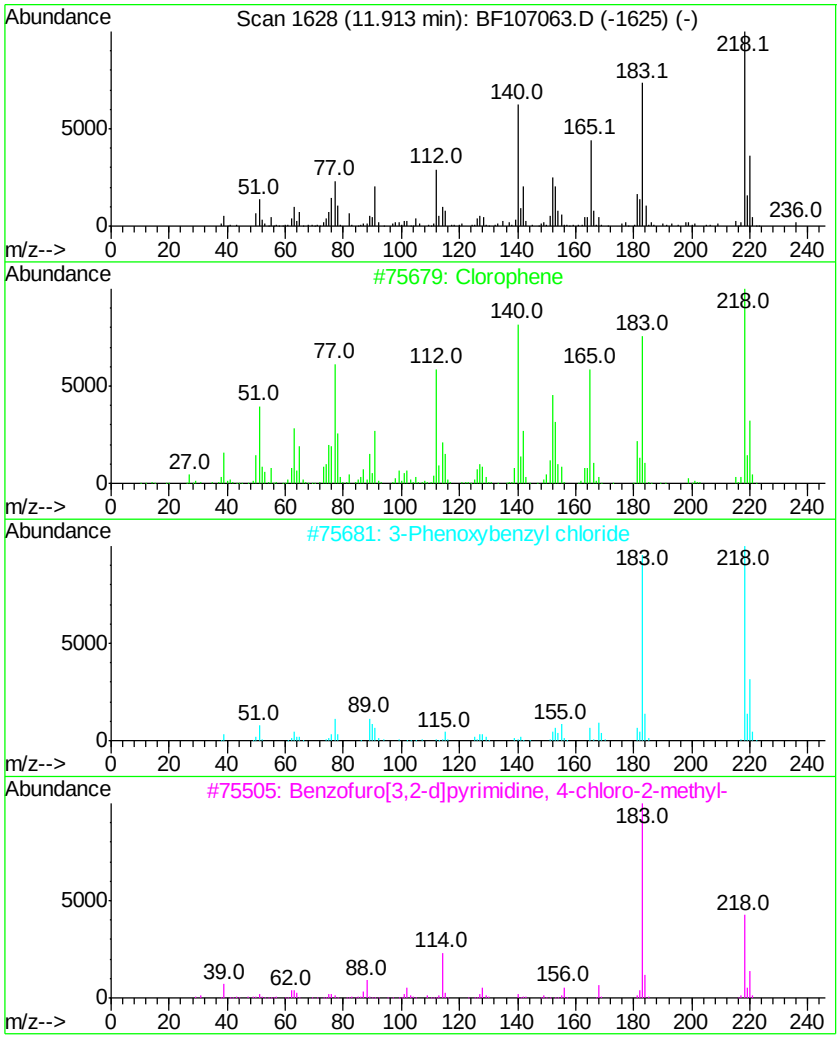
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

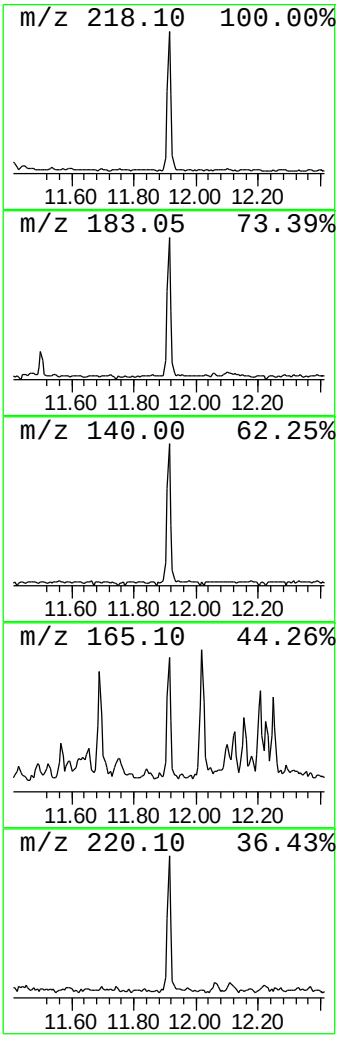
**********

Peak Number 20 Clorophene Concentration Rank 18

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.91 | 8.50 ng | 152115 | Phenanthrene-d10 | 11.50 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Clorophene | 218 | C13H11ClO | 000120-32-1 | 99 |
| 2 | 3-Phenoxybenzyl chloride | 218 | C13H11ClO | 1000353-82-2 | 60 |
| 3 | Benzofuro[3,2-d]pyrimidine, 4-ch... | 218 | C11H7ClN2O | 039786-40-8 | 49 |
| 4 | 6-Methoxy-2-naphthonitrile | 183 | C12H9NO | 067886-70-8 | 46 |
| 5 | Pyrimidine, tetrachloro- | 216 | C4Cl4N2 | 001780-40-1 | 43 |



Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF070218\
Data File : BF107063.D
Acq On : 3 Jul 2018 1:10
Operator : JU/SJ
Sample : J3799-07
Misc :
ALS Vial : 31 Sample Multiplier: 1

**Instrument :** BNA_F
**ClientSampleId :** MW-6

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF061918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

| | | | | | Internal Standard | | | |
|---|---|---|---|---|---|---|---|---|
| TIC Top Hit name | RT | EstConc | Units | Response | # | RT | Resp | Conc |
| unknown6.74 | 6.74 | 71.3 | ng | 1377250 | 1 | 6.95 | 386513 | 20.0 |
| 1,4-Hexadiene, 4-... | 8.18 | 10.8 | ng | 368713 | 2 | 8.24 | 680315 | 20.0 |
| Naphthalene, 1,2,... | 8.48 | 9.0 | ng | 305390 | 2 | 8.24 | 680315 | 20.0 |
| Naphthalene, 1,2,... | 9.13 | 7.5 | ng | 168816 | 3 | 10.00 | 453336 | 20.0 |
| Phenol, 4-(1,1-di... | 9.42 | 11.6 | ng | 262174 | 3 | 10.00 | 453336 | 20.0 |
| Benzene, 1,3,5-tr... | 10.34 | 7.4 | ng | 167647 | 3 | 10.00 | 453336 | 20.0 |
| unknown10.38 | 10.38 | 14.1 | ng | 320351 | 3 | 10.00 | 453336 | 20.0 |
| Phenol, m-tert-bu... | 10.75 | 8.6 | ng | 193974 | 3 | 10.00 | 453336 | 20.0 |
| 3',4'-Acetoxylide | 10.92 | 23.7 | ng | 423437 | 4 | 11.50 | 357964 | 20.0 |
| Phenol, 4-(1,1,3,... | 10.97 | 50.7 | ng | 907361 | 4 | 11.50 | 357964 | 20.0 |
| Phenol, 2-(1,1-di... | 11.01 | 83.3 | ng | 1491140 | 4 | 11.50 | 357964 | 20.0 |
| unknown11.07 | 11.07 | 31.1 | ng | 556576 | 4 | 11.50 | 357964 | 20.0 |
| Hexestrol | 11.11 | 52.4 | ng | 937143 | 4 | 11.50 | 357964 | 20.0 |
| unknown11.15 | 11.15 | 60.8 | ng | 1088970 | 4 | 11.50 | 357964 | 20.0 |
| unknown11.21 | 11.21 | 21.0 | ng | 375652 | 4 | 11.50 | 357964 | 20.0 |
| 4-Methylthieno[2,... | 11.23 | 30.4 | ng | 543162 | 4 | 11.50 | 357964 | 20.0 |
| Clorophene | 11.91 | 8.5 | ng | 152115 | 4 | 11.50 | 357964 | 20.0 |