LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023955.D
Acq On : 06 Feb 2025 04:32
Operator : RC/JU
Sample : Q1202-18
Misc :
ALS Vial : 28 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
E29B1

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % 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_P\Methods\SFAM-EPA-BP012925.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BP023955.D\data.ms

| peak # | R.T. min | first scan | max scan | last scan | PK TY | peak height | corr. area | corr. % max. | % of total |
|---|---|---|---|---|---|---|---|---|---|
| 1 | 3.293 | 31 | 35 | 38 | rBV | 113735 | 139323 | 0.21% | 0.037% |
| 2 | 3.334 | 38 | 42 | 55 | rVB | 248485 | 323923 | 0.48% | 0.087% |
| 3 | 3.628 | 88 | 92 | 101 | rVB | 214003 | 317153 | 0.47% | 0.085% |
| 4 | 3.711 | 101 | 106 | 117 | rBV | 506945 | 930104 | 1.38% | 0.250% |
| 5 | 3.858 | 126 | 131 | 143 | rBV2 | 70346 | 153775 | 0.23% | 0.041% |
| 6 | 4.022 | 155 | 159 | 171 | rVB | 100062 | 141989 | 0.21% | 0.038% |
| 7 | 5.340 | 376 | 383 | 388 | rVV2 | 79449 | 185362 | 0.28% | 0.050% |
| 8 | 5.452 | 397 | 402 | 412 | rVB | 101918 | 192111 | 0.29% | 0.052% |
| 9 | 6.852 | 632 | 640 | 650 | rVV2 | 131842 | 304562 | 0.45% | 0.082% |
| 10 | 6.952 | 650 | 657 | 671 | rVB | 626851 | 1135261 | 1.68% | 0.305% |
| 11 | 7.111 | 677 | 684 | 692 | rBV | 1743494 | 2745825 | 4.07% | 0.739% |
| 12 | 7.311 | 712 | 718 | 726 | rBV | 2115531 | 3215089 | 4.77% | 0.865% |
| 13 | 7.769 | 785 | 796 | 806 | rBV | 1735261 | 2750786 | 4.08% | 0.740% |
| 14 | 8.210 | 866 | 871 | 876 | rVB | 127580 | 205819 | 0.31% | 0.055% |
| 15 | 8.381 | 891 | 900 | 909 | rBV2 | 181141 | 449168 | 0.67% | 0.121% |
| 16 | 8.475 | 909 | 916 | 921 | rVV | 1794303 | 3005582 | 4.46% | 0.808% |
| 17 | 8.540 | 921 | 927 | 940 | rVB | 1066058 | 2045660 | 3.04% | 0.550% |
| 18 | 8.658 | 940 | 947 | 950 | rBV6 | 86769 | 199902 | 0.30% | 0.054% |
| 19 | 8.863 | 975 | 982 | 985 | rBV | 323225 | 545260 | 0.81% | 0.147% |
| 20 | 8.910 | 985 | 990 | 997 | rVV | 2037174 | 3287741 | 4.88% | 0.884% |
| 21 | 8.993 | 997 | 1004 | 1011 | rVB5 | 243998 | 643013 | 0.95% | 0.173% |
| 22 | 9.099 | 1011 | 1022 | 1027 | rBV | 911436 | 2082234 | 3.09% | 0.560% |
| 23 | 9.181 | 1028 | 1036 | 1050 | rVB2 | 656669 | 1402624 | 2.08% | 0.377% |
| 24 | 9.310 | 1052 | 1058 | 1061 | rBV3 | 144859 | 243066 | 0.36% | 0.065% |
| 25 | 9.346 | 1061 | 1064 | 1076 | rVB2 | 129664 | 292943 | 0.43% | 0.079% |
| 26 | 9.457 | 1076 | 1083 | 1089 | rBV4 | 110564 | 318262 | 0.47% | 0.086% |
| 27 | 9.522 | 1089 | 1094 | 1102 | rVB | 279189 | 458790 | 0.68% | 0.123% |
| 28 | 9.634 | 1104 | 1113 | 1120 | rVB | 1682340 | 2723057 | 4.04% | 0.732% |
| 29 | 9.728 | 1122 | 1129 | 1132 | rBV | 3350875 | 5920559 | 8.79% | 1.593% |
| 30 | 9.757 | 1132 | 1134 | 1141 | rVB | 4164143 | 5739013 | 8.52% | 1.544% |
| 31 | 9.993 | 1165 | 1174 | 1177 | rBV | 1567257 | 3432168 | 5.09% | 0.923% |
| 32 | 10.034 | 1177 | 1181 | 1188 | rVB | 2338780 | 3709149 | 5.50% | 0.998% |
| 33 | 10.110 | 1188 | 1194 | 1197 | rBV6 | 79876 | 164158 | 0.24% | 0.044% |
| 34 | 10.175 | 1197 | 1205 | 1215 | rVB3 | 3818830 | 7438172 | 11.04% | 2.001% |
| 35 | 10.369 | 1233 | 1238 | 1242 | rVV2 | 202999 | 335719 | 0.50% | 0.090% |
| 36 | 10.422 | 1242 | 1247 | 1249 | rVV | 823719 | 1343513 | 1.99% | 0.361% |

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023955.D
Acq On : 06 Feb 2025 04:32
Operator : RC/JU
Sample : Q1202-18
Misc :
ALS Vial : 28 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
E29B1

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % 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_P\Methods\SFAM-EPA-BP012925.MA.M
Title : SVOA CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 37 | 10.475 | 1249 | 1256 | 1262 | rVV | 24075092 | 43074106 | 63.92% | 11.587% |
| 38 | 10.546 | 1262 | 1268 | 1273 | rVV | 2281824 | 3960831 | 5.88% | 1.065% |
| 39 | 10.593 | 1273 | 1276 | 1285 | rVB | 504130 | 857578 | 1.27% | 0.231% |
| 40 | 10.681 | 1285 | 1291 | 1305 | rBV2 | 1040847 | 2231620 | 3.31% | 0.600% |
| 41 | 10.822 | 1310 | 1315 | 1318 | rVV | 187564 | 278365 | 0.41% | 0.075% |
| 42 | 10.899 | 1319 | 1328 | 1335 | rVV | 7632317 | 11746233 | 17.43% | 3.160% |
| 43 | 10.993 | 1335 | 1344 | 1348 | rVB5 | 469604 | 1066422 | 1.58% | 0.287% |
| 44 | 11.046 | 1348 | 1353 | 1356 | rBV3 | 219019 | 386591 | 0.57% | 0.104% |
| 45 | 11.087 | 1356 | 1360 | 1365 | rVV | 496665 | 806508 | 1.20% | 0.217% |
| 46 | 11.163 | 1369 | 1373 | 1382 | rVB2 | 333176 | 753716 | 1.12% | 0.203% |
| 47 | 11.316 | 1393 | 1399 | 1400 | rBV4 | 128518 | 218842 | 0.32% | 0.059% |
| 48 | 11.375 | 1403 | 1409 | 1414 | rVV2 | 1653896 | 3112846 | 4.62% | 0.837% |
| 49 | 11.428 | 1414 | 1418 | 1424 | rVB3 | 453665 | 897197 | 1.33% | 0.241% |
| 50 | 11.557 | 1433 | 1440 | 1446 | rBV4 | 1980076 | 3902239 | 5.79% | 1.050% |
| 51 | 11.622 | 1446 | 1451 | 1456 | rBV2 | 903661 | 1807621 | 2.68% | 0.486% |
| 52 | 11.716 | 1464 | 1467 | 1471 | rVB | 155784 | 170390 | 0.25% | 0.046% |
| 53 | 11.904 | 1489 | 1499 | 1505 | rBV2 | 31471343 | 67382391 | 100.00% | 18.126% |
| 54 | 11.957 | 1505 | 1508 | 1511 | rVB2 | 112051 | 173850 | 0.26% | 0.047% |
| 55 | 12.122 | 1531 | 1536 | 1544 | rVB4 | 419562 | 1060725 | 1.57% | 0.285% |
| 56 | 12.198 | 1545 | 1549 | 1552 | rBV5 | 94319 | 187301 | 0.28% | 0.050% |
| 57 | 12.304 | 1564 | 1567 | 1570 | rBV2 | 156525 | 216396 | 0.32% | 0.058% |
| 58 | 12.540 | 1599 | 1607 | 1612 | rBV3 | 394135 | 815933 | 1.21% | 0.219% |
| 59 | 12.681 | 1628 | 1631 | 1634 | rVB3 | 187094 | 223334 | 0.33% | 0.060% |
| 60 | 12.769 | 1642 | 1646 | 1648 | rBV2 | 169725 | 222494 | 0.33% | 0.060% |
| 61 | 12.798 | 1649 | 1651 | 1658 | rVB4 | 207790 | 330268 | 0.49% | 0.089% |
| 62 | 12.893 | 1664 | 1667 | 1671 | rBV3 | 152293 | 243583 | 0.36% | 0.066% |
| 63 | 12.963 | 1672 | 1679 | 1688 | rVB7 | 379246 | 1005921 | 1.49% | 0.271% |
| 64 | 13.234 | 1718 | 1725 | 1733 | rBV | 24489906 | 34418222 | 51.08% | 9.258% |
| 65 | 13.298 | 1733 | 1736 | 1742 | rVB3 | 284565 | 379955 | 0.56% | 0.102% |
| 66 | 13.369 | 1744 | 1748 | 1750 | rVV | 250532 | 342707 | 0.51% | 0.092% |
| 67 | 13.398 | 1751 | 1753 | 1761 | rVB6 | 204383 | 342127 | 0.51% | 0.092% |
| 68 | 13.504 | 1767 | 1771 | 1776 | rBV6 | 180587 | 406734 | 0.60% | 0.109% |
| 69 | 13.551 | 1776 | 1779 | 1781 | rVV2 | 220182 | 314755 | 0.47% | 0.085% |
| 70 | 13.593 | 1781 | 1786 | 1790 | rVV | 8326209 | 11564752 | 17.16% | 3.111% |
| 71 | 13.634 | 1790 | 1793 | 1800 | rVV | 2560573 | 3586325 | 5.32% | 0.965% |
| 72 | 13.704 | 1800 | 1805 | 1811 | rVV | 836858 | 1351183 | 2.01% | 0.363% |
| 73 | 13.793 | 1815 | 1820 | 1827 | rVB | 4740334 | 6389123 | 9.48% | 1.719% |
| 74 | 13.869 | 1829 | 1833 | 1837 | rVV | 819623 | 1165547 | 1.73% | 0.314% |
| 75 | 13.922 | 1837 | 1842 | 1848 | rVB5 | 622791 | 1132091 | 1.68% | 0.305% |
| 76 | 14.075 | 1855 | 1868 | 1876 | rBV | 5537497 | 9132353 | 13.55% | 2.457% |

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023955.D
Acq On : 06 Feb 2025 04:32
Operator : RC/JU
Sample : Q1202-18
Misc :
ALS Vial : 28 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
E29B1

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % 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_P\Methods\SFAM-EPA-BP012925.MA.M
Title : SVOA CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 77 | 14.145 | 1876 | 1880 | 1884 | rBV2 | 505021 | 671098 | 1.00% | 0.181% |
| 78 | 14.198 | 1884 | 1889 | 1891 | rBV5 | 211876 | 355264 | 0.53% | 0.096% |
| 79 | 14.387 | 1915 | 1921 | 1935 | rBV2 | 4055322 | 6742747 | 10.01% | 1.814% |
| 80 | 14.540 | 1944 | 1947 | 1951 | rVB | 361836 | 523793 | 0.78% | 0.141% |
| 81 | 14.604 | 1952 | 1958 | 1961 | rBV3 | 1180826 | 2211118 | 3.28% | 0.595% |
| 82 | 14.740 | 1977 | 1981 | 1987 | rVB | 883308 | 1405538 | 2.09% | 0.378% |
| 83 | 14.869 | 1997 | 2003 | 2009 | rBV7 | 320257 | 851406 | 1.26% | 0.229% |
| 84 | 14.928 | 2010 | 2013 | 2018 | rVV | 284854 | 371809 | 0.55% | 0.100% |
| 85 | 15.075 | 2033 | 2038 | 2045 | rBV2 | 771659 | 1814495 | 2.69% | 0.488% |
| 86 | 15.192 | 2055 | 2058 | 2061 | rBV | 404410 | 537720 | 0.80% | 0.145% |
| 87 | 15.392 | 2086 | 2092 | 2099 | rBV2 | 6119950 | 10406855 | 15.44% | 2.799% |
| 88 | 15.445 | 2099 | 2101 | 2106 | rVV4 | 346136 | 564569 | 0.84% | 0.152% |
| 89 | 15.504 | 2106 | 2111 | 2117 | rVV | 2287721 | 3382040 | 5.02% | 0.910% |
| 90 | 15.563 | 2118 | 2121 | 2130 | rVB | 707489 | 1065497 | 1.58% | 0.287% |
| 91 | 15.934 | 2180 | 2184 | 2189 | rBV3 | 677109 | 1022495 | 1.52% | 0.275% |
| 92 | 17.175 | 2390 | 2395 | 2405 | rVB | 4443810 | 6847494 | 10.16% | 1.842% |
| 93 | 17.275 | 2406 | 2412 | 2419 | rBV2 | 7859527 | 12909863 | 19.16% | 3.473% |
| 94 | 17.563 | 2457 | 2461 | 2465 | rBV | 1196063 | 1590386 | 2.36% | 0.428% |
| 95 | 18.122 | 2554 | 2556 | 2559 | rBV2 | 709109 | 714798 | 1.06% | 0.192% |
| 96 | 18.157 | 2559 | 2562 | 2567 | rVB | 1281694 | 1741730 | 2.58% | 0.469% |
| 97 | 19.639 | 2809 | 2814 | 2820 | rBV | 10031549 | 12840314 | 19.06% | 3.454% |
| 98 | 21.610 | 3145 | 3149 | 3157 | rVB | 4318657 | 7048790 | 10.46% | 1.896% |
| 99 | 24.751 | 3675 | 3683 | 3694 | rVB | 4811099 | 12621127 | 18.73% | 3.395% |
| 100 | 24.986 | 3715 | 3723 | 3735 | rVB | 2853439 | 7327570 | 10.87% | 1.971% |

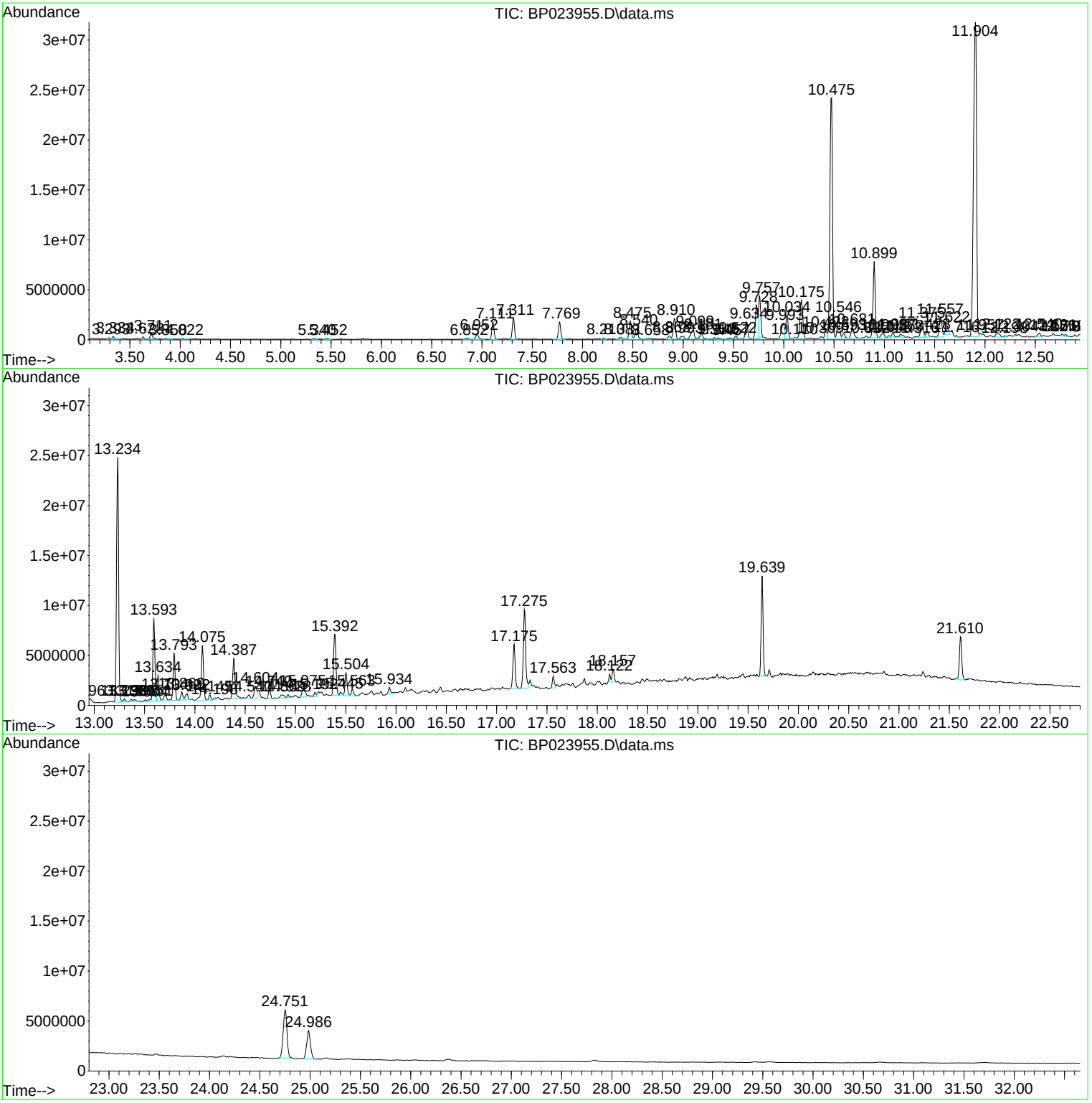
Sum of corrected areas: 371748526

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023955.D
Acq On : 06 Feb 2025 04:32
Operator : RC/JU
Sample : Q1202-18
Misc :
ALS Vial : 28 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
E29B1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023955.D
Acq On : 06 Feb 2025 04:32
Operator : RC/JU
Sample : Q1202-18
Misc :
ALS Vial : 28 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
E29B1

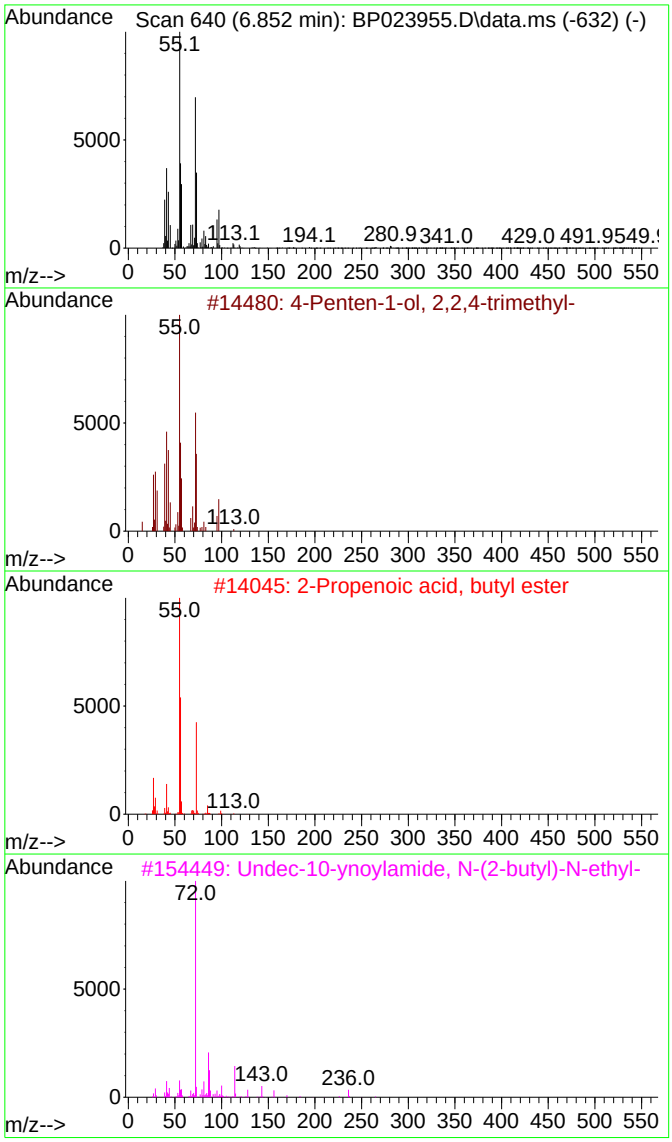
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

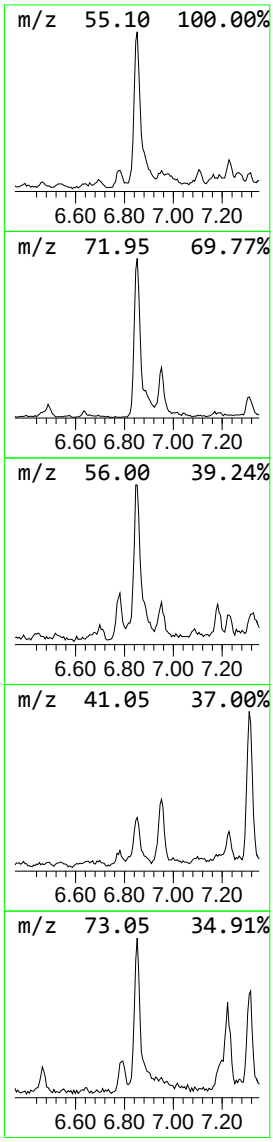
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 2 4-Penten-1-ol, 2,2,4-trimet... Concentration Rank 36

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 6.852 | 2.21 ng/ul | 304562 | 1,4-Dichlorobenzene-d4 | 7.769 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 4-Penten-1-ol, 2,2,4-trimethyl- | 128 | C8H16O | 053907-70-3 | 90 |
| 2 | 2-Propenoic acid, butyl ester | 128 | C7H12O2 | 000141-32-2 | 17 |
| 3 | Undec-10-ynoylamide, N-(2-butyl)... | 265 | C17H31NO | 1000491-72-1 | 10 |
| 4 | 5-Chlorovaleryl amide, N-(2-phen... | 281 | C16H24ClNO | 1000451-75-4 | 9 |
| 5 | L-Valyl-L-leucine, TMS derivative | 302 | C14H30N2O3Si | 1000333-69-7 | 9 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023955.D
Acq On : 06 Feb 2025 04:32
Operator : RC/JU
Sample : Q1202-18
Misc :
ALS Vial : 28 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
E29B1

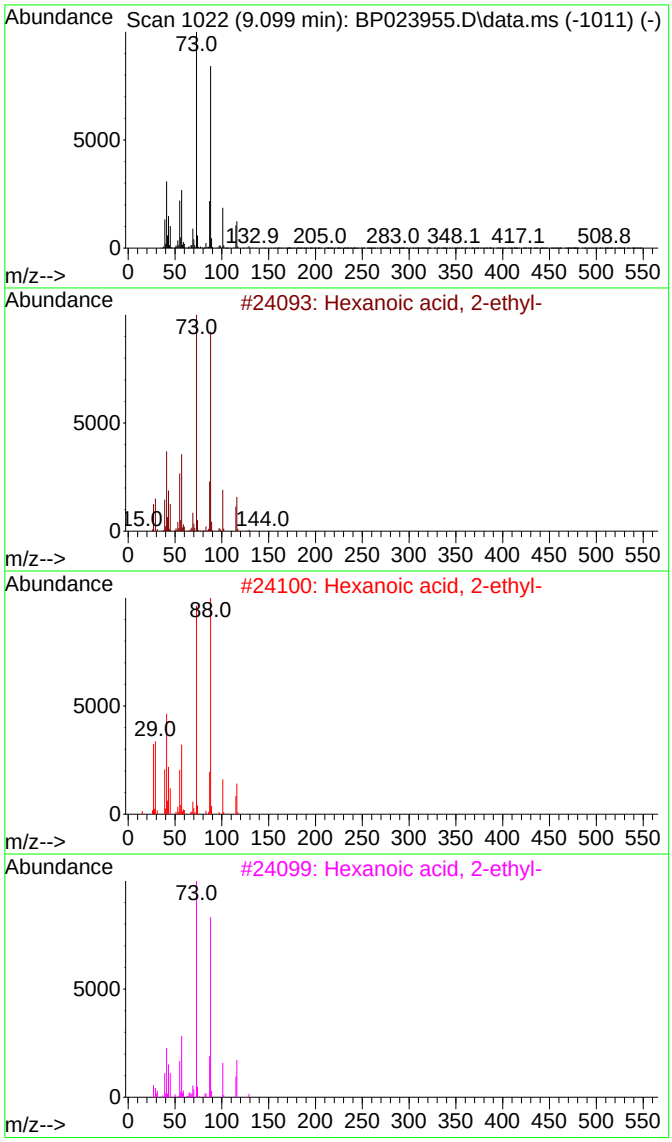
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

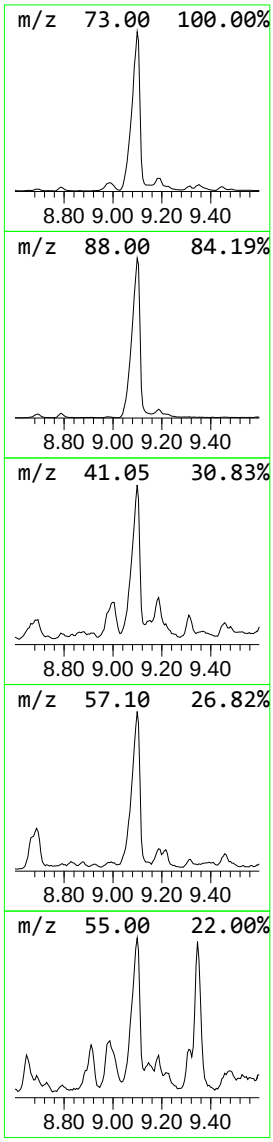
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 5 Hexanoic acid, 2-ethyl- Concentration Rank 10

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 9.099 | 15.14 ng/ul | 2082230 | 1,4-Dichlorobenzene-d4 | 7.769 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Hexanoic acid, 2-ethyl- | 144 | C8H16O2 | 000149-57-5 | 90 |
| 2 | Hexanoic acid, 2-ethyl- | 144 | C8H16O2 | 000149-57-5 | 90 |
| 3 | Hexanoic acid, 2-ethyl- | 144 | C8H16O2 | 000149-57-5 | 83 |
| 4 | Hexanoic acid, 2-ethyl- | 144 | C8H16O2 | 000149-57-5 | 72 |
| 5 | Hexanoic acid, 2-ethyl- | 144 | C8H16O2 | 000149-57-5 | 72 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023955.D
Acq On : 06 Feb 2025 04:32
Operator : RC/JU
Sample : Q1202-18
Misc :
ALS Vial : 28 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
E29B1

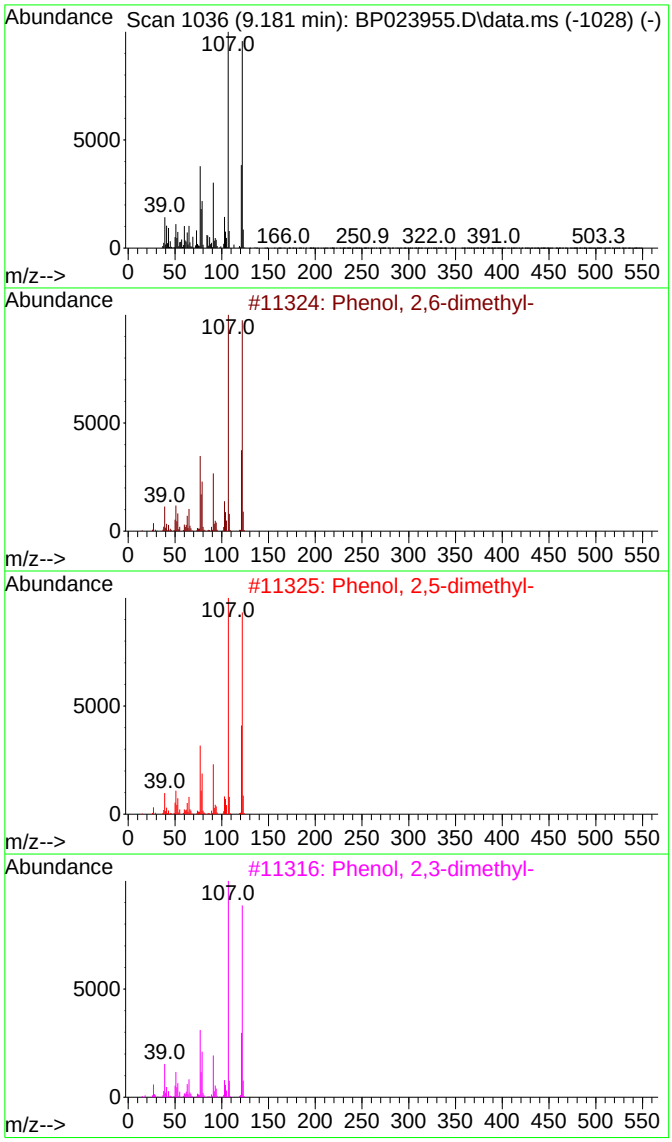
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

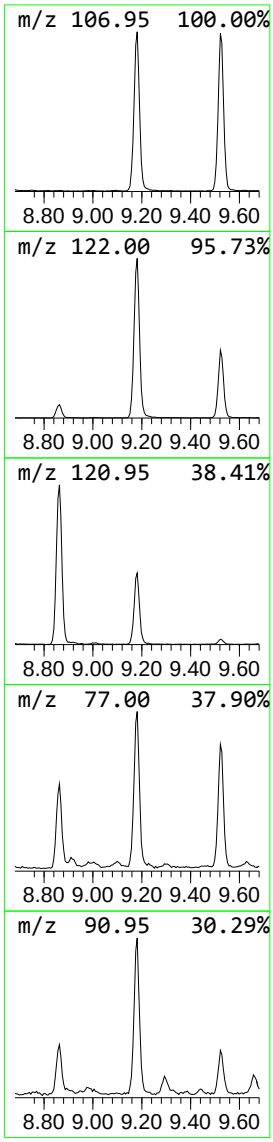
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 6 Phenol, 2,6-dimethyl- Concentration Rank 13

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 9.181 | 7.08 ng/ul | 1402620 | Naphthalene-d8 | 10.546 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, 2,6-dimethyl- | 122 | C8H10O | 000576-26-1 | 97 |
| 2 | Phenol, 2,5-dimethyl- | 122 | C8H10O | 000095-87-4 | 96 |
| 3 | Phenol, 2,3-dimethyl- | 122 | C8H10O | 000526-75-0 | 96 |
| 4 | Phenol, 2,5-dimethyl- | 122 | C8H10O | 000095-87-4 | 95 |
| 5 | Phenol, 2,6-dimethyl- | 122 | C8H10O | 000576-26-1 | 95 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023955.D
Acq On : 06 Feb 2025 04:32
Operator : RC/JU
Sample : Q1202-18
Misc :
ALS Vial : 28 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
E29B1

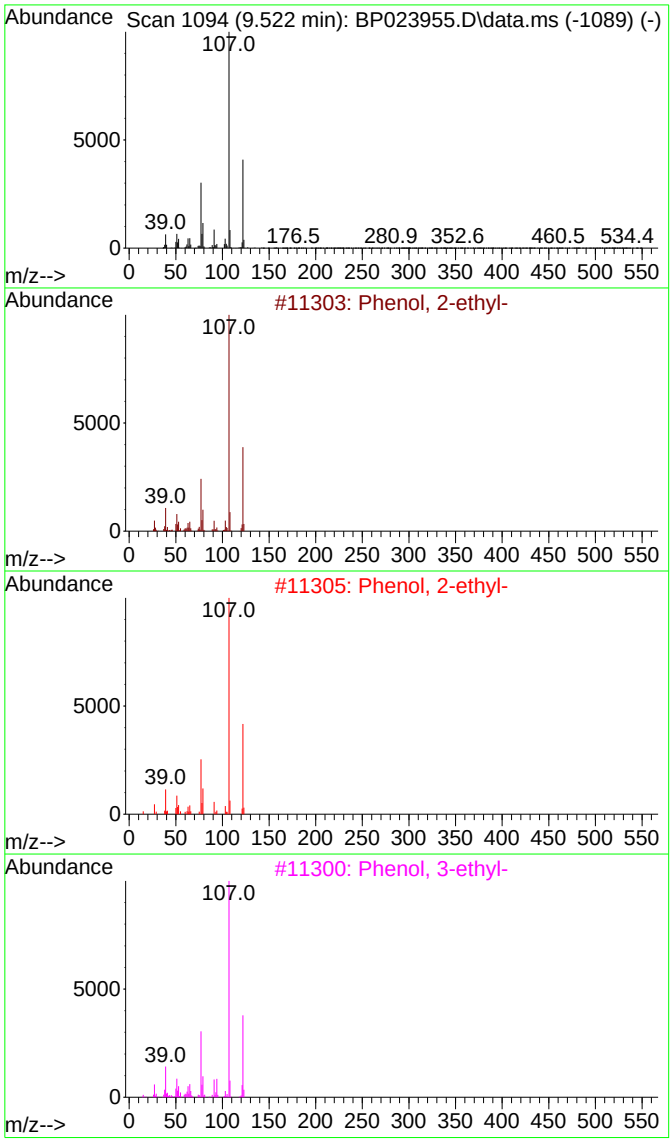
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

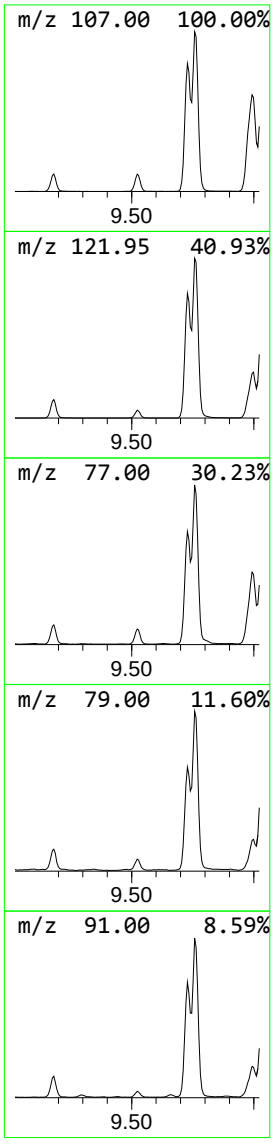
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 7 Phenol, 2-ethyl- Concentration Rank 34

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 9.522 | 2.32 ng/ul | 458790 | Naphthalene-d8 | 10.546 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, 2-ethyl- | 122 | C8H10O | 000090-00-6 | 95 |
| 2 | Phenol, 2-ethyl- | 122 | C8H10O | 000090-00-6 | 95 |
| 3 | Phenol, 3-ethyl- | 122 | C8H10O | 000620-17-7 | 91 |
| 4 | Phenol, 3,4-dimethyl- | 122 | C8H10O | 000095-65-8 | 91 |
| 5 | Phenol, 3,4-dimethyl- | 122 | C8H10O | 000095-65-8 | 91 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023955.D
Acq On : 06 Feb 2025 04:32
Operator : RC/JU
Sample : Q1202-18
Misc :
ALS Vial : 28 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
E29B1

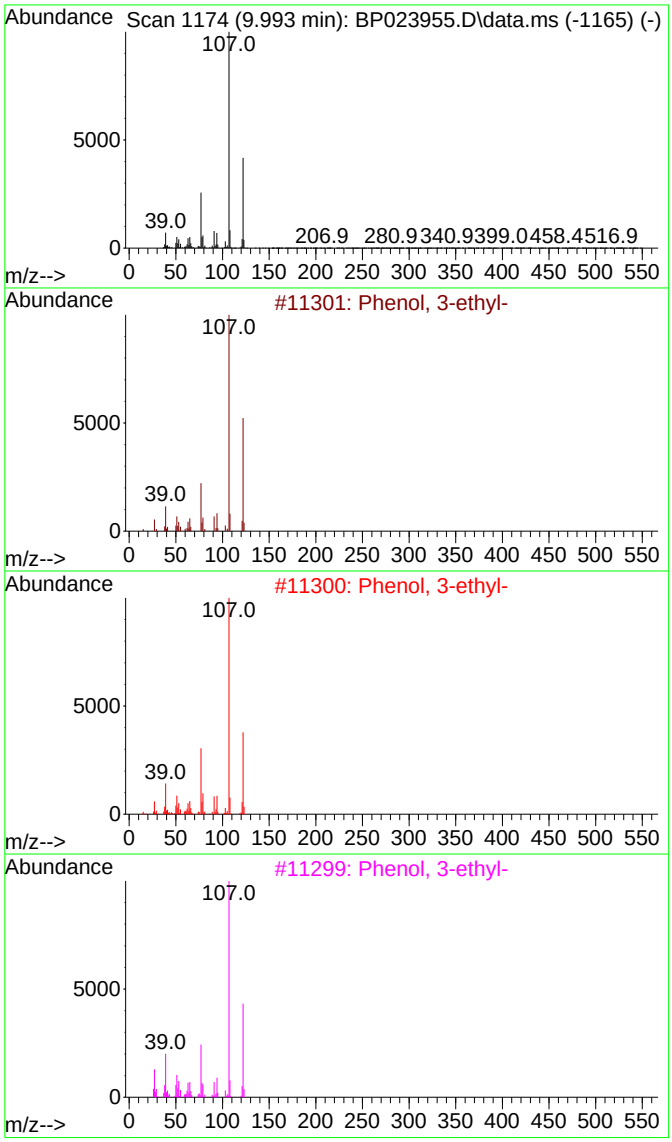
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

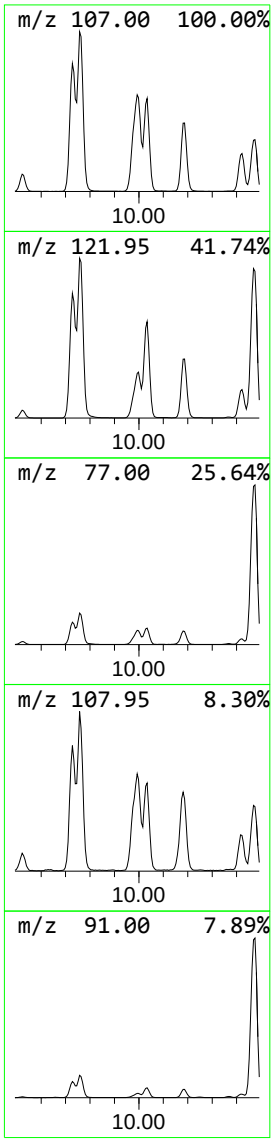
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 8 Phenol, 3-ethyl- Concentration Rank 8

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 9.993 | 17.33 ng/ul | 3432170 | Naphthalene-d8 | 10.546 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, 3-ethyl- | 122 | C8H10O | 000620-17-7 | 97 |
| 2 | Phenol, 3-ethyl- | 122 | C8H10O | 000620-17-7 | 94 |
| 3 | Phenol, 3-ethyl- | 122 | C8H10O | 000620-17-7 | 94 |
| 4 | Phenol, 2-ethyl- | 122 | C8H10O | 000090-00-6 | 91 |
| 5 | Phenol, 2-ethyl- | 122 | C8H10O | 000090-00-6 | 91 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023955.D
Acq On : 06 Feb 2025 04:32
Operator : RC/JU
Sample : Q1202-18
Misc :
ALS Vial : 28 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
E29B1

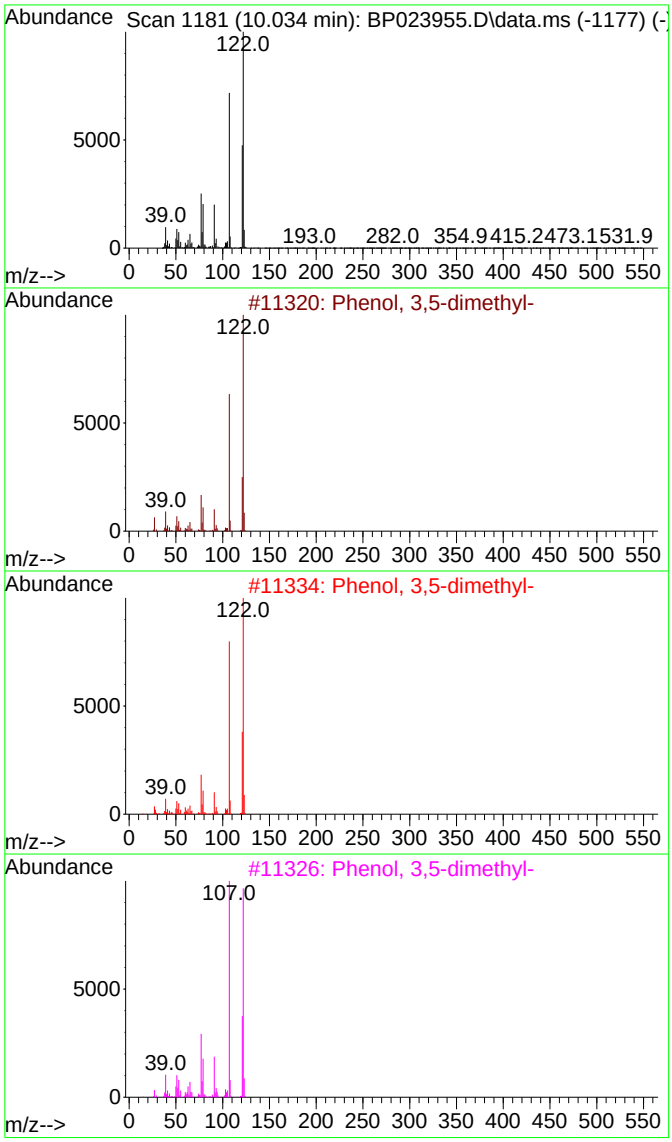
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

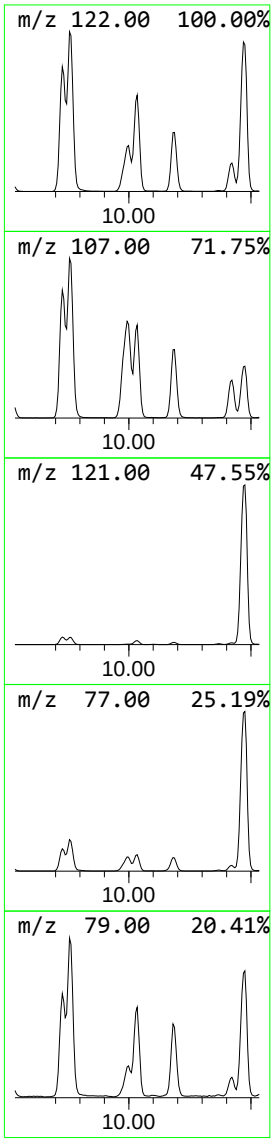
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 9 Phenol, 3,5-dimethyl- Concentration Rank 7

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.034 | 18.73 ng/ul | 3709150 | Naphthalene-d8 | 10.546 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, 3,5-dimethyl- | 122 | C8H10O | 000108-68-9 | 91 |
| 2 | Phenol, 3,5-dimethyl- | 122 | C8H10O | 000108-68-9 | 91 |
| 3 | Phenol, 3,5-dimethyl- | 122 | C8H10O | 000108-68-9 | 90 |
| 4 | Phenol, 2,4-dimethyl- | 122 | C8H10O | 000105-67-9 | 87 |
| 5 | 1,3-Cyclopentadiene, 1,2,5,5-tet... | 122 | C9H14 | 004249-12-1 | 86 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023955.D
Acq On : 06 Feb 2025 04:32
Operator : RC/JU
Sample : Q1202-18
Misc :
ALS Vial : 28 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
E29B1

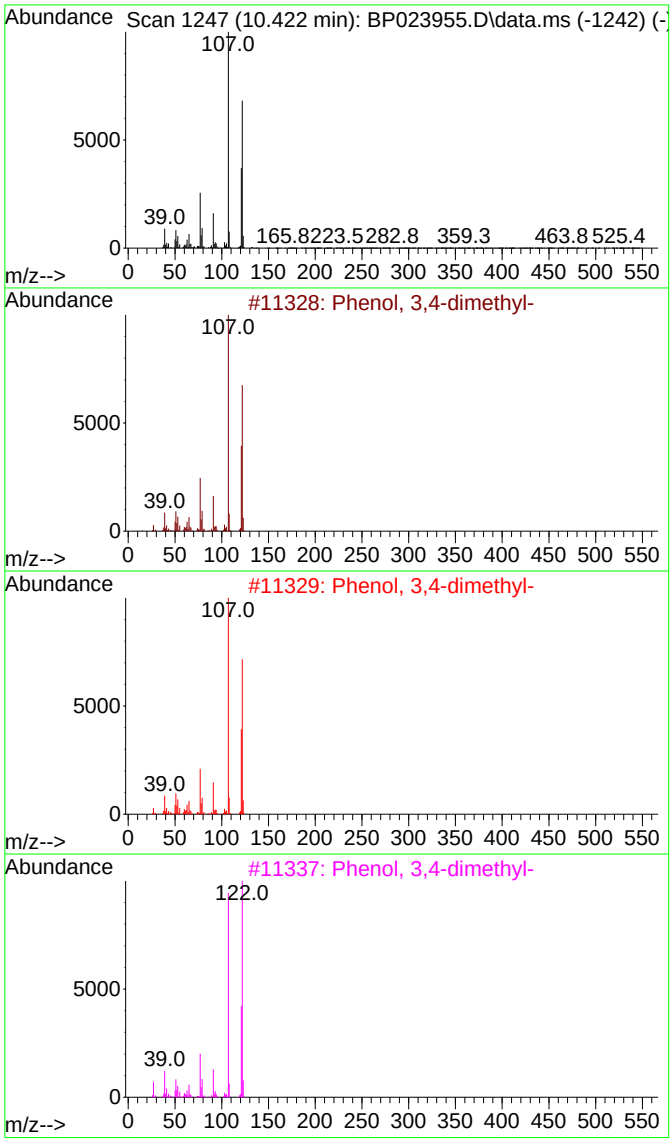
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

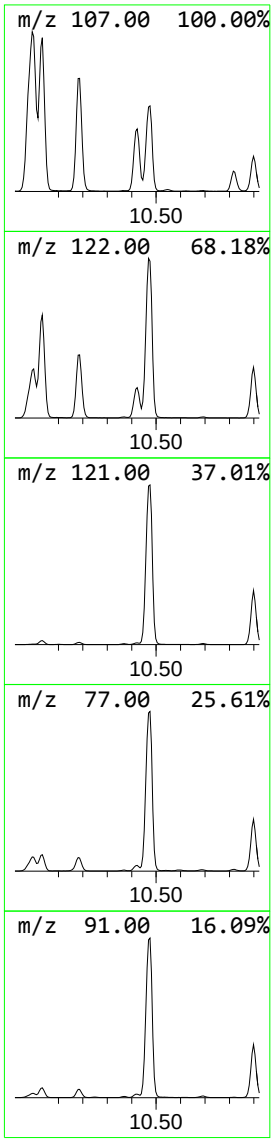
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 10 Phenol, 3,4-dimethyl- Concentration Rank 14

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.422 | 6.78 ng/ul | 1343510 | Naphthalene-d8 | 10.546 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, 3,4-dimethyl- | 122 | C8H10O | 000095-65-8 | 96 |
| 2 | Phenol, 3,4-dimethyl- | 122 | C8H10O | 000095-65-8 | 96 |
| 3 | Phenol, 3,4-dimethyl- | 122 | C8H10O | 000095-65-8 | 95 |
| 4 | Phenol, 3,4-dimethyl- | 122 | C8H10O | 000095-65-8 | 95 |
| 5 | Phenol, 2-ethyl- | 122 | C8H10O | 000090-00-6 | 91 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023955.D
Acq On : 06 Feb 2025 04:32
Operator : RC/JU
Sample : Q1202-18
Misc :
ALS Vial : 28 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
E29B1

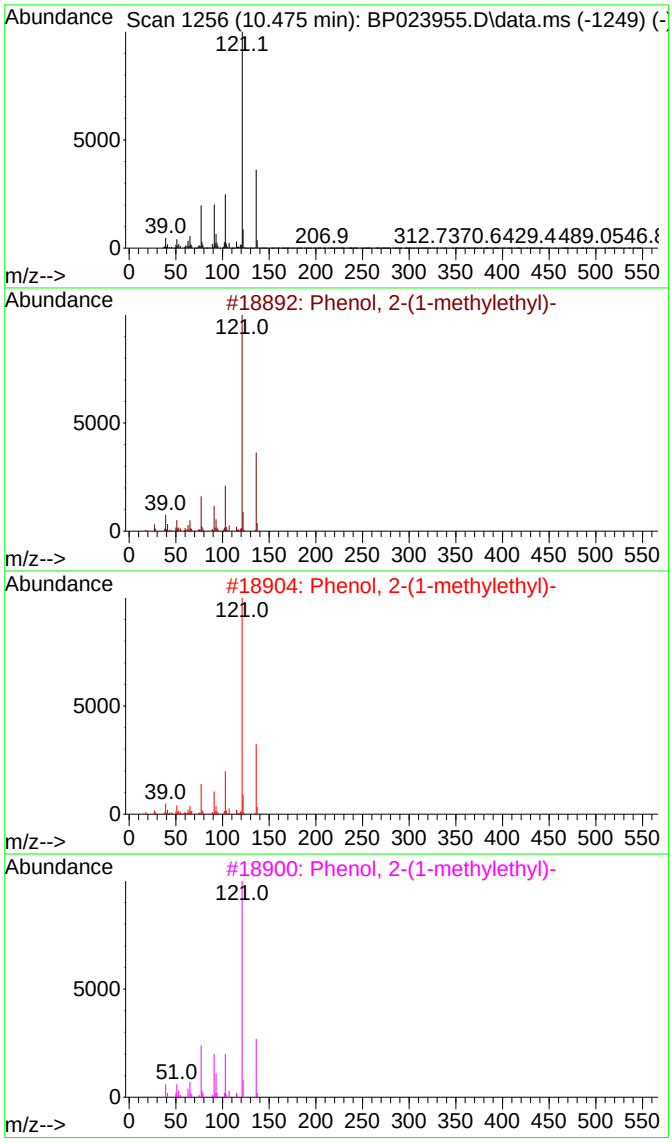
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

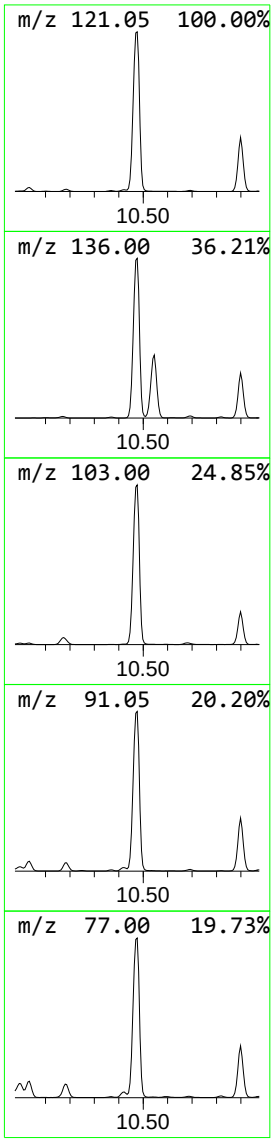
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 11 Phenol, 2-(1-methylethyl)- Concentration Rank 2

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.475 | 217.50 ng/ul | 43074100 | Naphthalene-d8 | 10.546 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, 2-(1-methylethyl)- | 136 | C9H12O | 000088-69-7 | 95 |
| 2 | Phenol, 2-(1-methylethyl)- | 136 | C9H12O | 000088-69-7 | 91 |
| 3 | Phenol, 2-(1-methylethyl)- | 136 | C9H12O | 000088-69-7 | 90 |
| 4 | Phenol, 3-(1-methylethyl)- | 136 | C9H12O | 000618-45-1 | 87 |
| 5 | Phenol, 3-(1-methylethyl)- | 136 | C9H12O | 000618-45-1 | 86 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023955.D
Acq On : 06 Feb 2025 04:32
Operator : RC/JU
Sample : Q1202-18
Misc :
ALS Vial : 28 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
E29B1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

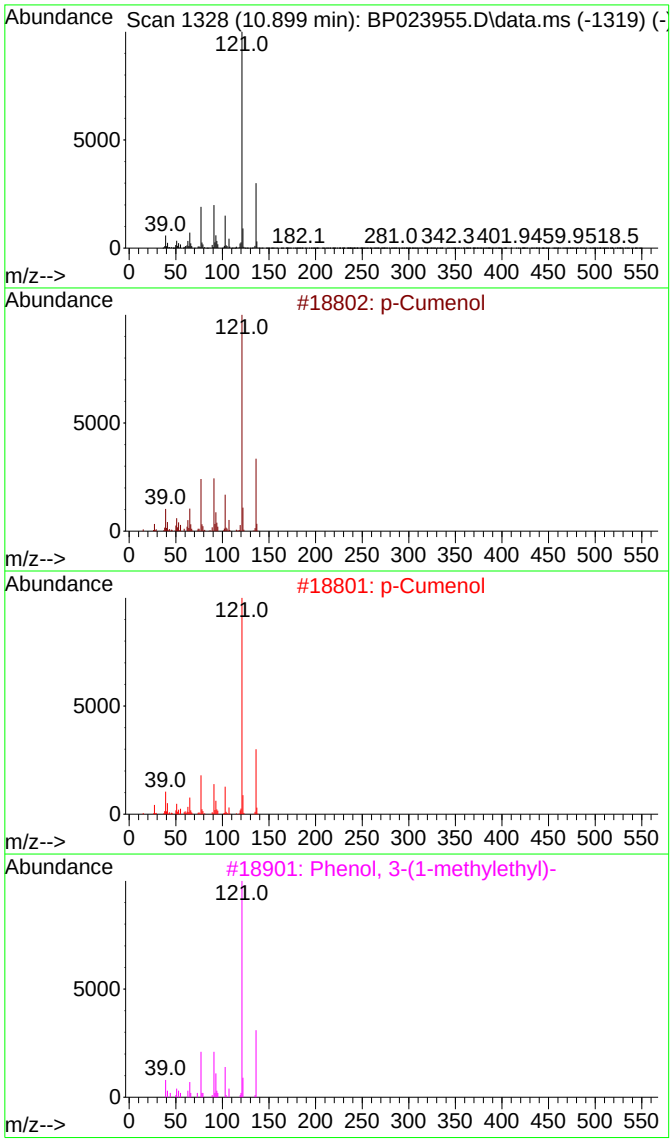
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

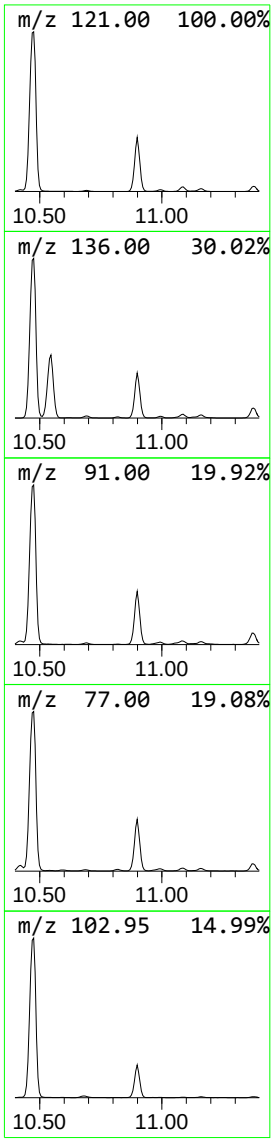
**********

Peak Number 12 p-Cumenol Concentration Rank 4

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.899 | 59.31 ng/ul | 11746200 | Naphthalene-d8 | 10.546 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | p-Cumenol | 136 | C9H12O | 000099-89-8 | 96 |
| 2 | p-Cumenol | 136 | C9H12O | 000099-89-8 | 95 |
| 3 | Phenol, 3-(1-methylethyl)- | 136 | C9H12O | 000618-45-1 | 94 |
| 4 | p-Cumenol | 136 | C9H12O | 000099-89-8 | 91 |
| 5 | Phenol, 3-(1-methylethyl)- | 136 | C9H12O | 000618-45-1 | 91 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023955.D
Acq On : 06 Feb 2025 04:32
Operator : RC/JU
Sample : Q1202-18
Misc :
ALS Vial : 28 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
E29B1

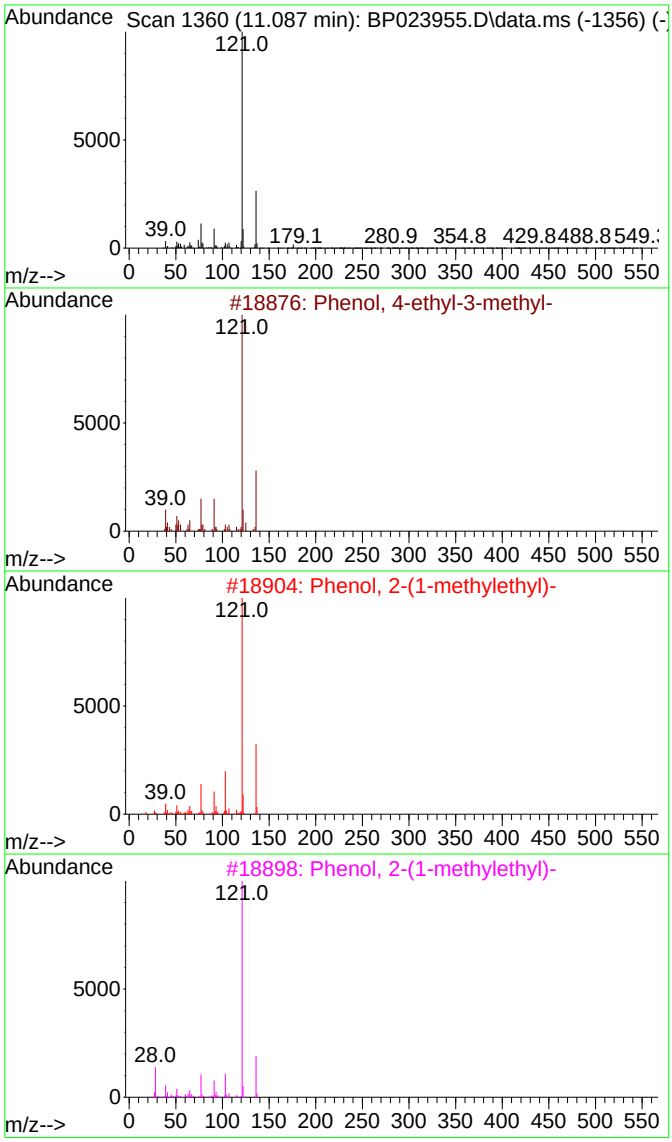
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

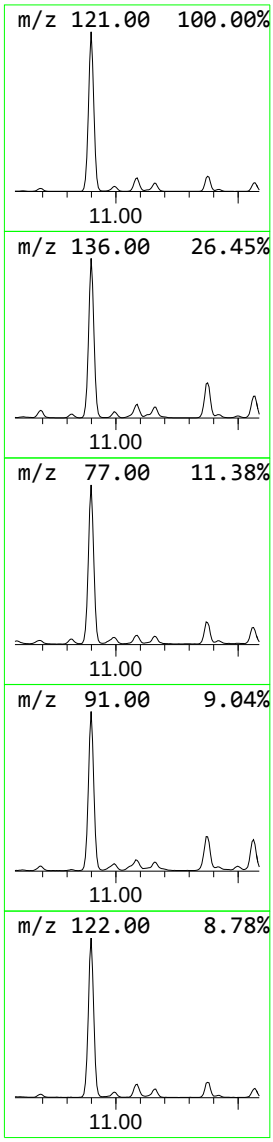
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 14 Phenol, 4-ethyl-3-methyl- Concentration Rank 23

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.087 | 4.07 ng/ul | 806508 | Naphthalene-d8 | 10.546 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, 4-ethyl-3-methyl- | 136 | C9H12O | 001123-94-0 | 91 |
| 2 | Phenol, 2-(1-methylethyl)- | 136 | C9H12O | 000088-69-7 | 91 |
| 3 | Phenol, 2-(1-methylethyl)- | 136 | C9H12O | 000088-69-7 | 90 |
| 4 | p-Cumenol | 136 | C9H12O | 000099-89-8 | 90 |
| 5 | 1,3-Cyclohexadiene, 1,5,5,6-tetr... | 136 | C10H16 | 000514-94-3 | 86 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023955.D
Acq On : 06 Feb 2025 04:32
Operator : RC/JU
Sample : Q1202-18
Misc :
ALS Vial : 28 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
E29B1

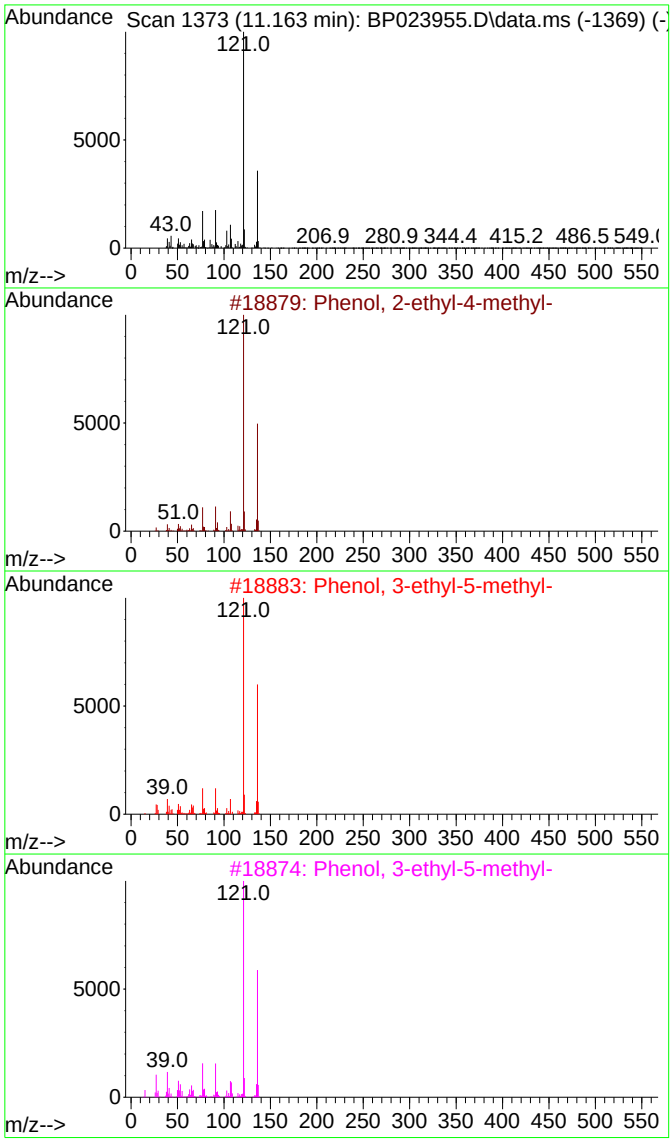
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

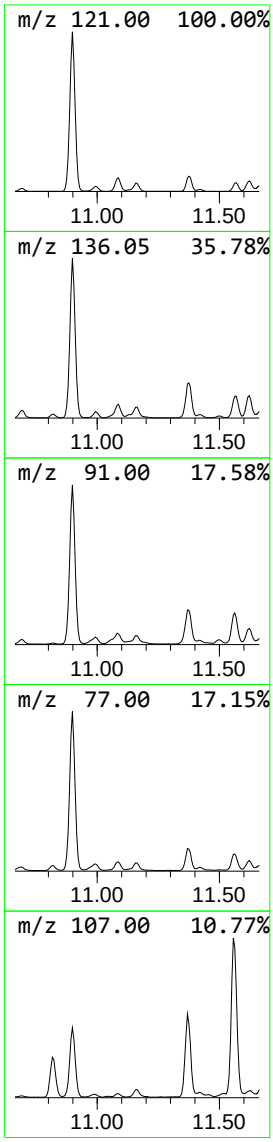
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 15 Phenol, 2-ethyl-4-methyl- Concentration Rank 25

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.163 | 3.81 ng/ul | 753716 | Naphthalene-d8 | 10.546 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, 2-ethyl-4-methyl- | 136 | C9H12O | 003855-26-3 | 91 |
| 2 | Phenol, 3-ethyl-5-methyl- | 136 | C9H12O | 000698-71-5 | 87 |
| 3 | Phenol, 3-ethyl-5-methyl- | 136 | C9H12O | 000698-71-5 | 86 |
| 4 | Phenol, 2-(1-methylethyl)- | 136 | C9H12O | 000088-69-7 | 83 |
| 5 | Phenol, 2-ethyl-5-methyl- | 136 | C9H12O | 001687-61-2 | 83 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023955.D
Acq On : 06 Feb 2025 04:32
Operator : RC/JU
Sample : Q1202-18
Misc :
ALS Vial : 28 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
E29B1

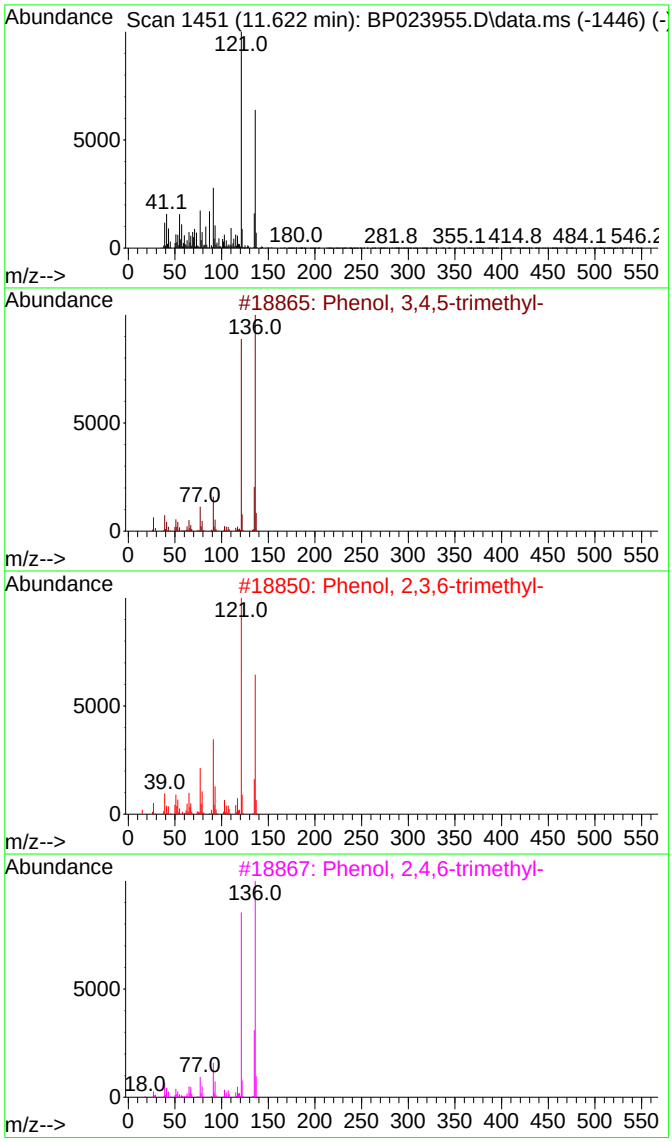
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

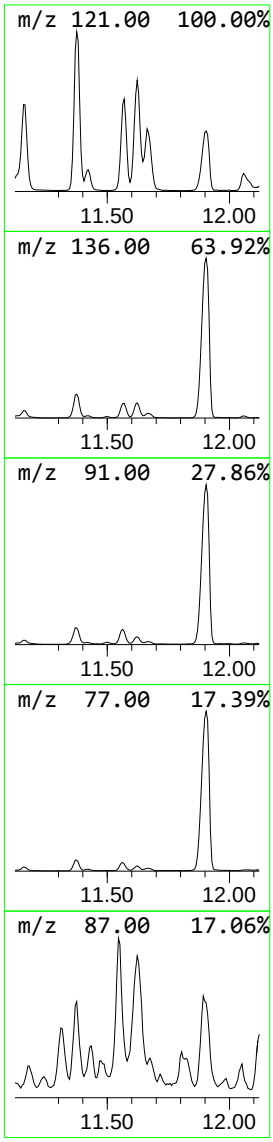
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 19 Phenol, 3,4,5-trimethyl- Concentration Rank 12

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.622 | 9.13 ng/ul | 1807620 | Naphthalene-d8 | 10.546 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, 3,4,5-trimethyl- | 136 | C9H12O | 000527-54-8 | 93 |
| 2 | Phenol, 2,3,6-trimethyl- | 136 | C9H12O | 002416-94-6 | 93 |
| 3 | Phenol, 2,4,6-trimethyl- | 136 | C9H12O | 000527-60-6 | 91 |
| 4 | Phenol, 2,3,5-trimethyl- | 136 | C9H12O | 000697-82-5 | 91 |
| 5 | Phenol, 2,3,5-trimethyl- | 136 | C9H12O | 000697-82-5 | 87 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023955.D
Acq On : 06 Feb 2025 04:32
Operator : RC/JU
Sample : Q1202-18
Misc :
ALS Vial : 28 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
E29B1

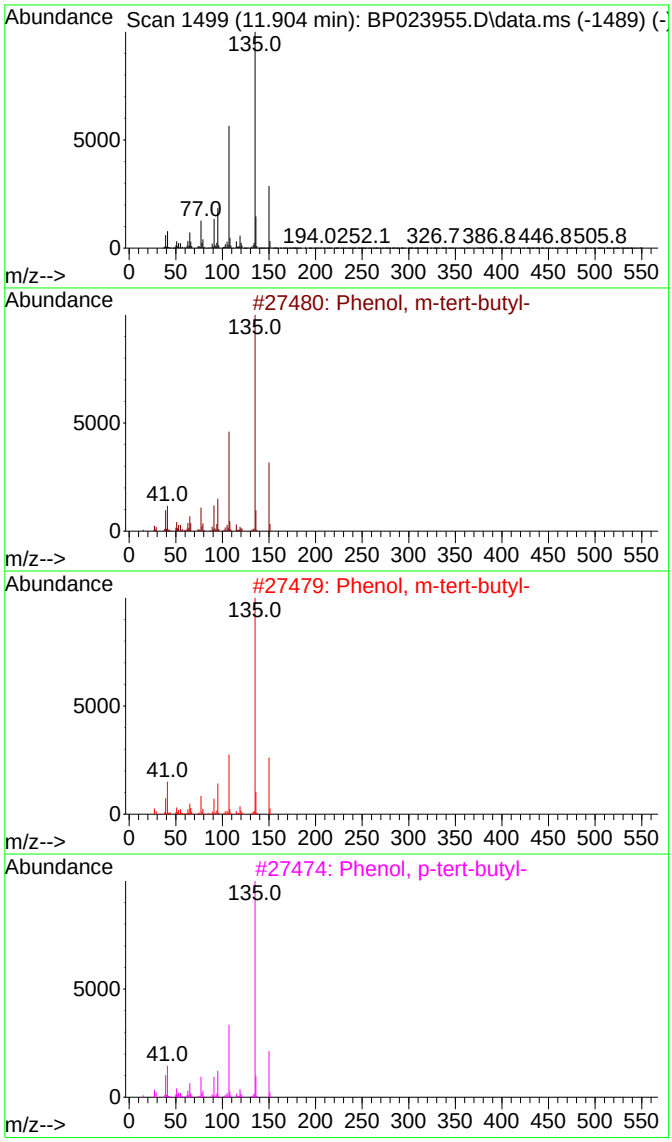
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

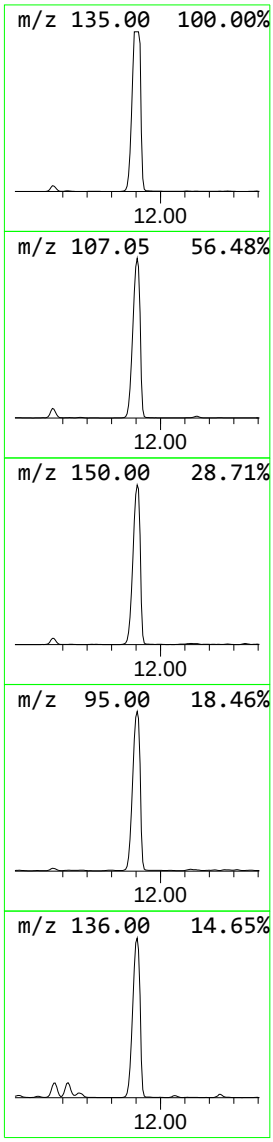
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 20 Phenol, m-tert-butyl- Concentration Rank 1

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 11.904 | 340.24 ng/ul | 67382400 | Naphthalene-d8 | 10.546 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, m-tert-butyl- | 150 | C10H14O | 000585-34-2 | 97 |
| 2 | Phenol, m-tert-butyl- | 150 | C10H14O | 000585-34-2 | 94 |
| 3 | Phenol, p-tert-butyl- | 150 | C10H14O | 000098-54-4 | 93 |
| 4 | Phenol, m-tert-butyl- | 150 | C10H14O | 000585-34-2 | 91 |
| 5 | Phenol, 2-(1,1-dimethylethyl)- | 150 | C10H14O | 000088-18-6 | 90 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023955.D
Acq On : 06 Feb 2025 04:32
Operator : RC/JU
Sample : Q1202-18
Misc :
ALS Vial : 28 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
E29B1

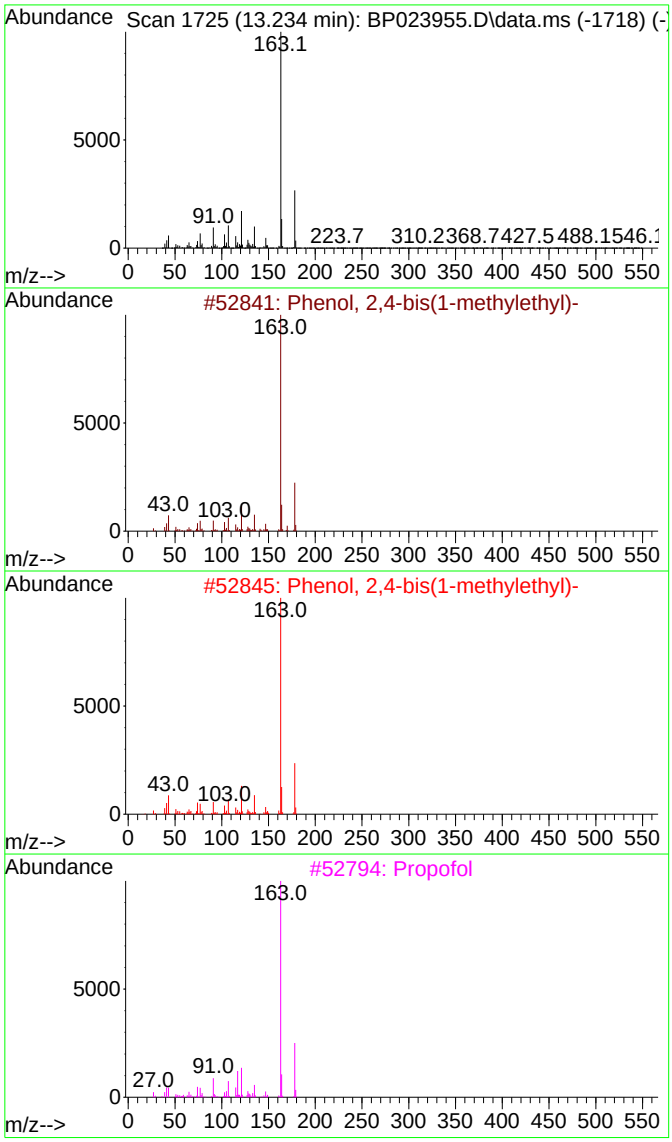
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

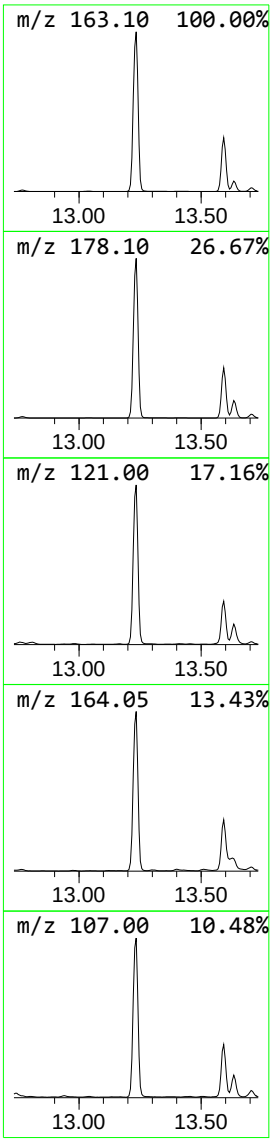
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 24 Phenol, 2,4-bis(1-methyleth... Concentration Rank 3

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.234 | 102.09 ng/ul | 34418200 | Acenaphthene-d10 | 14.387 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, 2,4-bis(1-methylethyl)- | 178 | C12H18O | 002934-05-6 | 97 |
| 2 | Phenol, 2,4-bis(1-methylethyl)- | 178 | C12H18O | 002934-05-6 | 97 |
| 3 | Propofol | 178 | C12H18O | 002078-54-8 | 94 |
| 4 | Propofol | 178 | C12H18O | 002078-54-8 | 91 |
| 5 | Propofol | 178 | C12H18O | 002078-54-8 | 90 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023955.D
Acq On : 06 Feb 2025 04:32
Operator : RC/JU
Sample : Q1202-18
Misc :
ALS Vial : 28 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
E29B1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

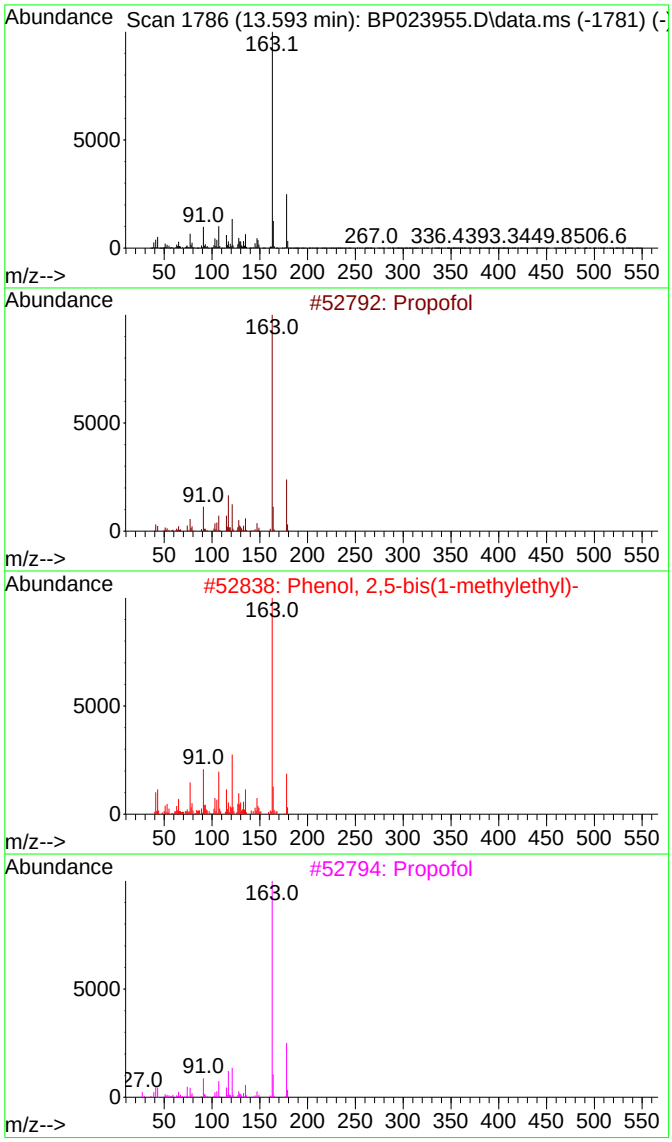
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

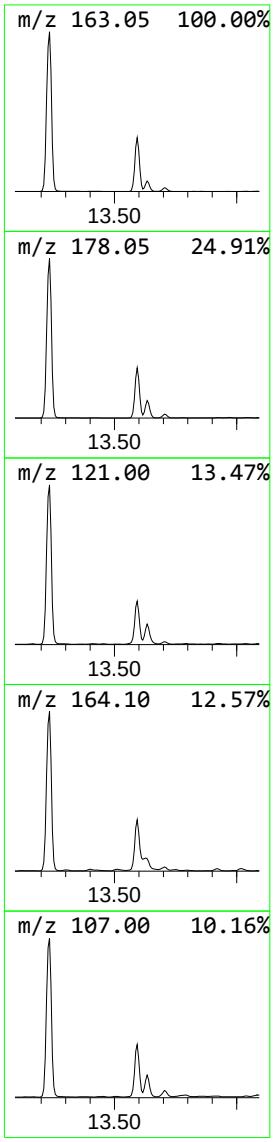
**********

Peak Number 25 Propofol Concentration Rank 5

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.592 | 34.30 ng/ul | 11564800 | Acenaphthene-d10 | 14.387 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Propofol | 178 | C12H18O | 002078-54-8 | 97 |
| 2 | Phenol, 2,5-bis(1-methylethyl)- | 178 | C12H18O | 035946-91-9 | 95 |
| 3 | Propofol | 178 | C12H18O | 002078-54-8 | 93 |
| 4 | Phenol, 2,4-bis(1-methylethyl)- | 178 | C12H18O | 002934-05-6 | 93 |
| 5 | Phenol, 2,4-bis(1-methylethyl)- | 178 | C12H18O | 002934-05-6 | 91 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023955.D
Acq On : 06 Feb 2025 04:32
Operator : RC/JU
Sample : Q1202-18
Misc :
ALS Vial : 28 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
E29B1

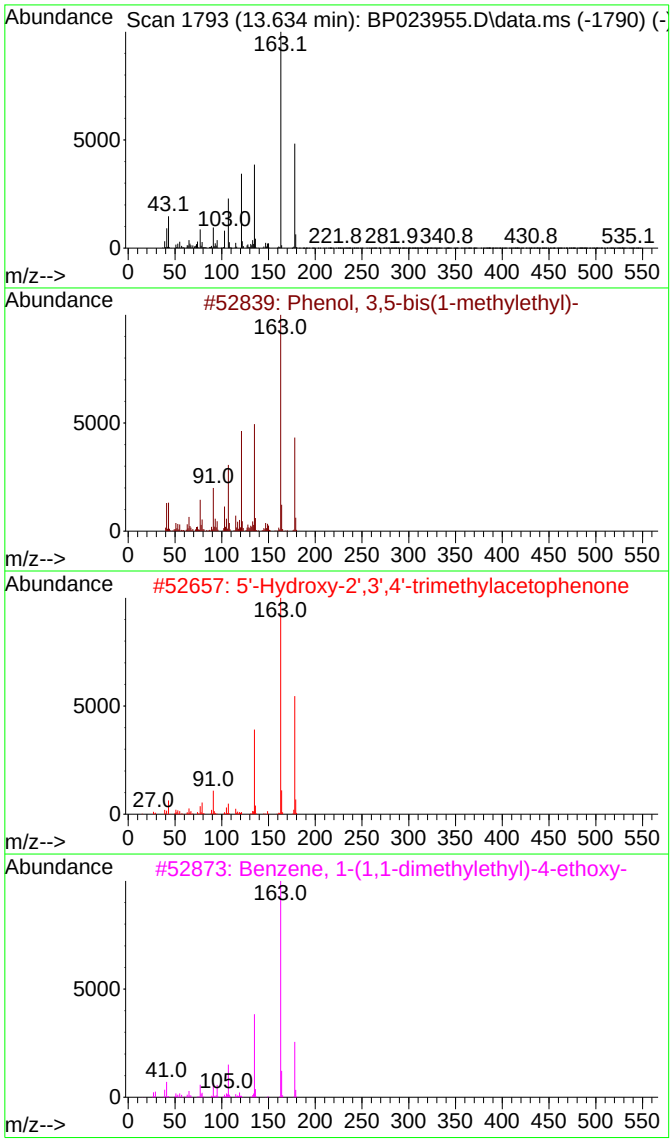
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

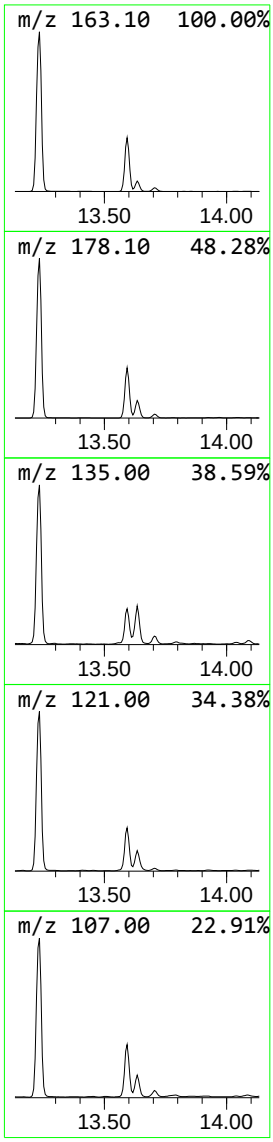
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 26 Phenol, 3,5-bis(1-methyleth... Concentration Rank 11

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.634 | 10.64 ng/ul | 3586330 | Acenaphthene-d10 | 14.387 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, 3,5-bis(1-methylethyl)- | 178 | C12H18O | 026886-05-5 | 94 |
| 2 | 5'-Hydroxy-2',3',4'-trimethylace... | 178 | C11H14O2 | 013667-28-2 | 80 |
| 3 | Benzene, 1-(1,1-dimethylethyl)-4... | 178 | C12H18O | 017269-94-2 | 78 |
| 4 | 2,6-Lutidine 4-diethylamino- | 178 | C11H18N2 | 1000252-95-3 | 64 |
| 5 | Ethanone, 1-(2,3-dihydro-1,4-ben... | 178 | C10H10O3 | 002879-20-1 | 59 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023955.D
Acq On : 06 Feb 2025 04:32
Operator : RC/JU
Sample : Q1202-18
Misc :
ALS Vial : 28 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
E29B1

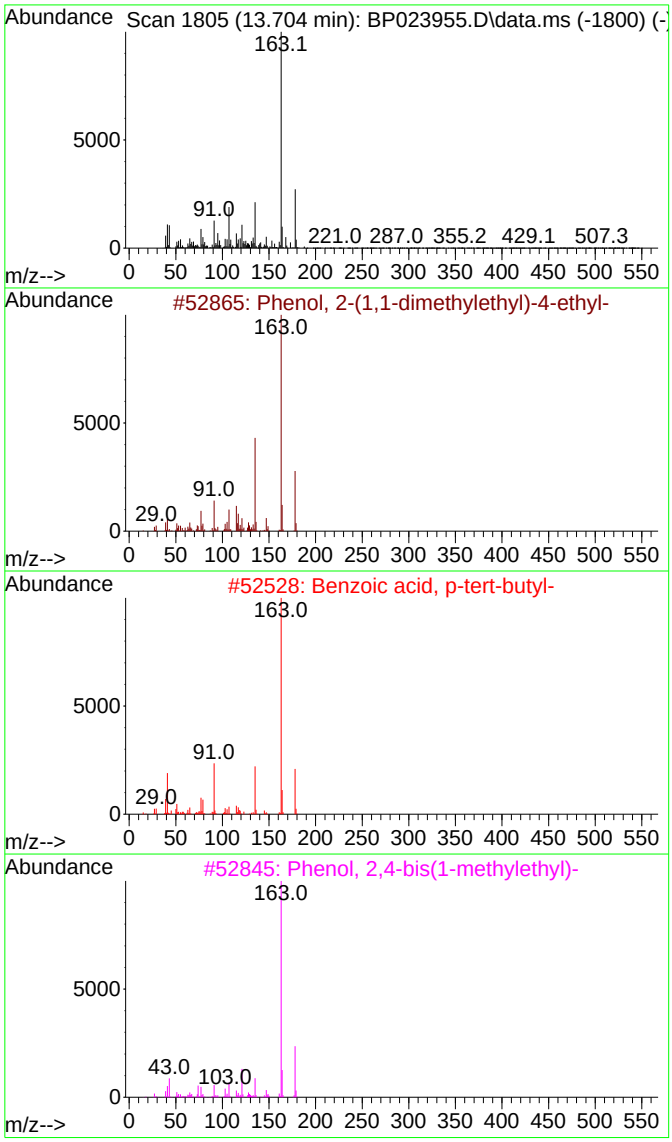
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

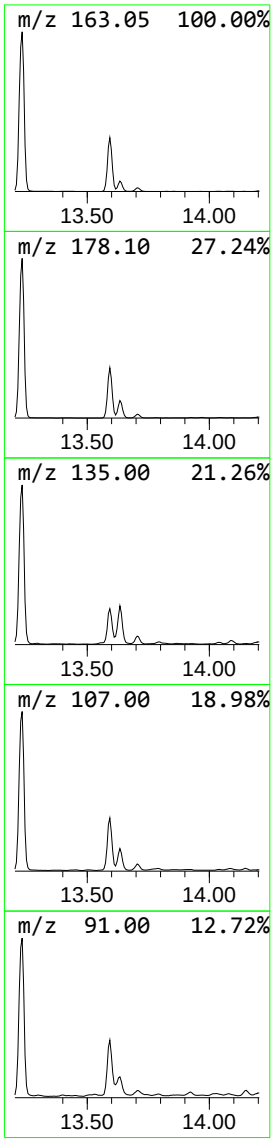
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 27 Phenol, 2-(1,1-dimethylethy... Concentration Rank 24

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.704 | 4.01 ng/ul | 1351180 | Acenaphthene-d10 | 14.387 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, 2-(1,1-dimethylethyl)-4-... | 178 | C12H18O | 000096-70-8 | 87 |
| 2 | Benzoic acid, p-tert-butyl- | 178 | C11H14O2 | 000098-73-7 | 74 |
| 3 | Phenol, 2,4-bis(1-methylethyl)- | 178 | C12H18O | 002934-05-6 | 74 |
| 4 | Phenol, 2,4-bis(1-methylethyl)- | 178 | C12H18O | 002934-05-6 | 74 |
| 5 | Propofol | 178 | C12H18O | 002078-54-8 | 74 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023955.D
Acq On : 06 Feb 2025 04:32
Operator : RC/JU
Sample : Q1202-18
Misc :
ALS Vial : 28 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
E29B1

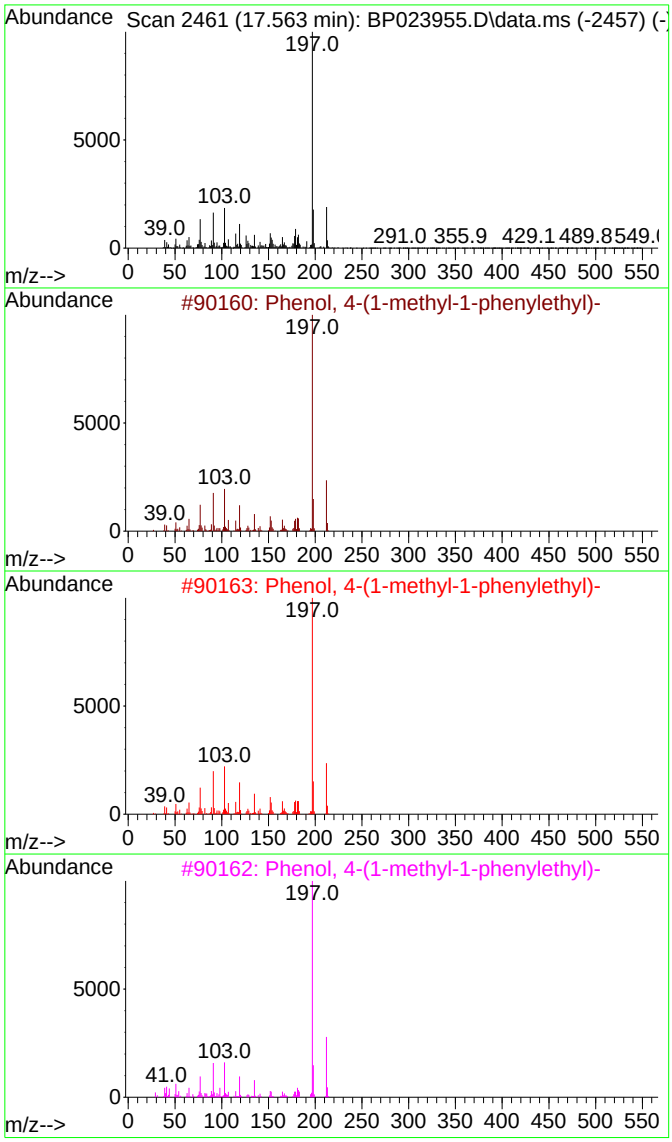
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

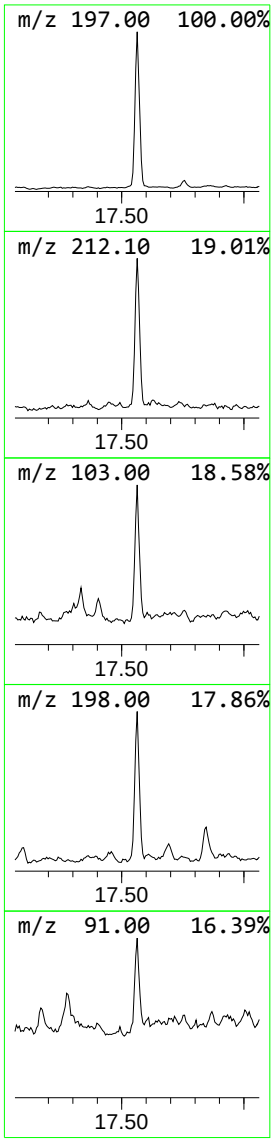
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 35 Phenol, 4-(1-methyl-1-pheny... Concentration Rank 20

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 17.563 | 4.65 ng/ul | 1590390 | Phenanthrene-d10 | 17.175 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenol, 4-(1-methyl-1-phenylethyl)- | 212 | C15H16O | 000599-64-4 | 98 |
| 2 | Phenol, 4-(1-methyl-1-phenylethyl)- | 212 | C15H16O | 000599-64-4 | 96 |
| 3 | Phenol, 4-(1-methyl-1-phenylethyl)- | 212 | C15H16O | 000599-64-4 | 91 |
| 4 | Dimethyl diphenylsilane | 212 | C14H16Si | 000778-24-5 | 81 |
| 5 | Dimethyl diphenylsilane | 212 | C14H16Si | 000778-24-5 | 81 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023955.D
Acq On : 06 Feb 2025 04:32
Operator : RC/JU
Sample : Q1202-18
Misc :
ALS Vial : 28 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
E29B1

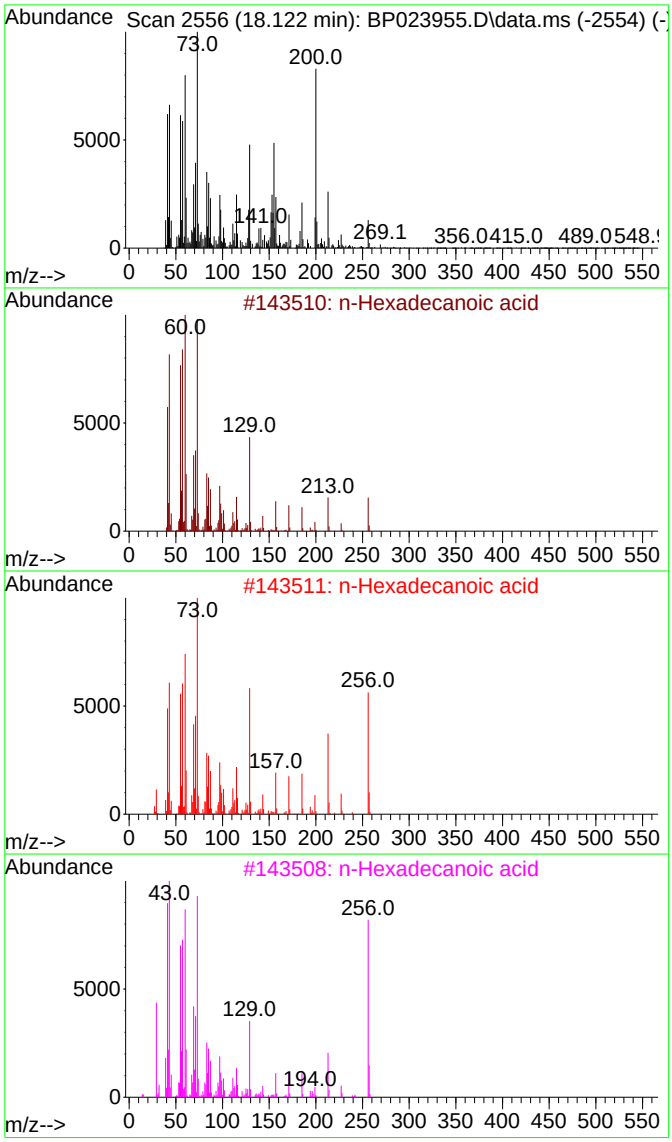
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA CALIBRATION

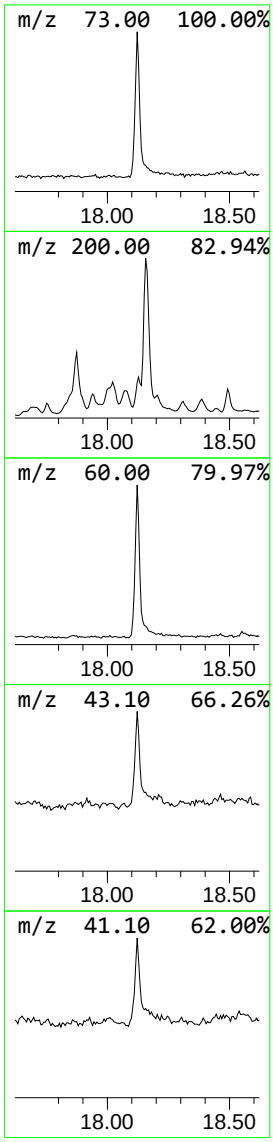
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 36 n-Hexadecanoic acid Concentration Rank 37

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.122 | 2.09 ng/ul | 714798 | Phenanthrene-d10 | 17.175 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 96 |
| 2 | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 90 |
| 3 | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 87 |
| 4 | Dodecanoic acid | 200 | C12H24O2 | 000143-07-7 | 53 |
| 5 | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 52 |



Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020525\
Data File : BP023955.D
Acq On : 06 Feb 2025 04:32
Operator : RC/JU
Sample : Q1202-18
Misc :
ALS Vial : 28 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
E29B1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP012925.MA.M
Quant Title : SVOA 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 |
| 4-Penten-1-ol, ... | 6.852 | 2.2 | ng/ul | 304562 | 1 | 7.769 | 2750790 | 20.0 |
| Hexanoic acid, ... | 9.099 | 15.1 | ng/ul | 2082230 | 1 | 7.769 | 2750790 | 20.0 |
| Phenol, 2,6-dim... | 9.181 | 7.1 | ng/ul | 1402620 | 2 | 10.546 | 3960830 | 20.0 |
| Phenol, 2-ethyl- | 9.522 | 2.3 | ng/ul | 458790 | 2 | 10.546 | 3960830 | 20.0 |
| Phenol, 3-ethyl- | 9.993 | 17.3 | ng/ul | 3432170 | 2 | 10.546 | 3960830 | 20.0 |
| Phenol, 3,5-dim... | 10.034 | 18.7 | ng/ul | 3709150 | 2 | 10.546 | 3960830 | 20.0 |
| Phenol, 3,4-dim... | 10.422 | 6.8 | ng/ul | 1343510 | 2 | 10.546 | 3960830 | 20.0 |
| Phenol, 2-(1-me... | 10.475 | 217.5 | ng/ul | 43074100 | 2 | 10.546 | 3960830 | 20.0 |
| p-Cumenol | 10.899 | 59.3 | ng/ul | 11746200 | 2 | 10.546 | 3960830 | 20.0 |
| Phenol, 4-ethyl... | 11.087 | 4.1 | ng/ul | 806508 | 2 | 10.546 | 3960830 | 20.0 |
| Phenol, 2-ethyl... | 11.163 | 3.8 | ng/ul | 753716 | 2 | 10.546 | 3960830 | 20.0 |
| Phenol, 3,4,5-t... | 11.622 | 9.1 | ng/ul | 1807620 | 2 | 10.546 | 3960830 | 20.0 |
| Phenol, m-tert-... | 11.904 | 340.2 | ng/ul | 67382400 | 2 | 10.546 | 3960830 | 20.0 |
| Phenol, 2,4-bis... | 13.234 | 102.1 | ng/ul | 34418200 | 3 | 14.387 | 6742750 | 20.0 |
| Propofol | 13.592 | 34.3 | ng/ul | 11564800 | 3 | 14.387 | 6742750 | 20.0 |
| Phenol, 3,5-bis... | 13.634 | 10.6 | ng/ul | 3586330 | 3 | 14.387 | 6742750 | 20.0 |
| Phenol, 2-(1,1-... | 13.704 | 4.0 | ng/ul | 1351180 | 3 | 14.387 | 6742750 | 20.0 |
| Phenol, 4-(1-me... | 17.563 | 4.7 | ng/ul | 1590390 | 4 | 17.175 | 6847490 | 20.0 |
| n-Hexadecanoic ... | 18.122 | 2.1 | ng/ul | 714798 | 4 | 17.175 | 6847490 | 20.0 |