LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071323\
Data File : BP016208.D
Acq On : 13 Jul 2023 14:57
Operator : MA/JU
Sample : 03385-01ME 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EX7X7ME

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-BP071223.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BP016208.D\data.ms

| peak # | R.T. min | first scan | max scan | last scan | PK TY | peak height | corr. area | corr. % max. | % of total |
|---|---|---|---|---|---|---|---|---|---|
| 1 | 3.722 | 86 | 91 | 106 | rBV3 | 126828 | 405226 | 1.92% | 0.218% |
| 2 | 3.834 | 106 | 110 | 115 | rVV | 252646 | 367934 | 1.74% | 0.198% |
| 3 | 4.058 | 140 | 148 | 166 | rVV3 | 414305 | 1066889 | 5.04% | 0.573% |
| 4 | 4.240 | 173 | 179 | 183 | rBV3 | 286180 | 466858 | 2.21% | 0.251% |
| 5 | 4.293 | 183 | 188 | 206 | rVB | 963800 | 1482553 | 7.01% | 0.797% |
| 6 | 4.446 | 208 | 214 | 229 | rVB | 137620 | 271853 | 1.28% | 0.146% |
| 7 | 4.587 | 230 | 238 | 241 | rBV2 | 107177 | 212337 | 1.00% | 0.114% |
| 8 | 4.628 | 241 | 245 | 252 | rVV2 | 1089013 | 1807344 | 8.54% | 0.971% |
| 9 | 4.693 | 252 | 256 | 257 | rVV | 396248 | 490807 | 2.32% | 0.264% |
| 10 | 4.717 | 257 | 260 | 269 | rVV | 753213 | 1128205 | 5.33% | 0.606% |
| 11 | 5.128 | 322 | 330 | 349 | rVB | 134841 | 326487 | 1.54% | 0.175% |
| 12 | 5.834 | 445 | 450 | 457 | rVV | 444521 | 802788 | 3.79% | 0.431% |
| 13 | 6.246 | 517 | 520 | 531 | rVV2 | 148046 | 298630 | 1.41% | 0.161% |
| 14 | 6.340 | 531 | 536 | 548 | rVB | 166206 | 351674 | 1.66% | 0.189% |
| 15 | 6.587 | 570 | 578 | 595 | rVV | 573698 | 1291114 | 6.10% | 0.694% |
| 16 | 7.705 | 762 | 768 | 779 | rBV4 | 62449 | 217595 | 1.03% | 0.117% |
| 17 | 7.975 | 808 | 814 | 826 | rBV | 285259 | 654863 | 3.10% | 0.352% |
| 18 | 8.187 | 841 | 850 | 855 | rBV2 | 185306 | 375600 | 1.78% | 0.202% |
| 19 | 8.275 | 859 | 865 | 880 | rVB | 1136896 | 2212297 | 10.46% | 1.189% |
| 20 | 8.852 | 958 | 963 | 975 | rVV | 142589 | 455941 | 2.16% | 0.245% |
| 21 | 9.299 | 1029 | 1039 | 1047 | rVV5 | 63539 | 212635 | 1.01% | 0.114% |
| 22 | 10.781 | 1285 | 1291 | 1294 | rBV | 613792 | 1086191 | 5.13% | 0.584% |
| 23 | 10.828 | 1294 | 1299 | 1317 | rVV | 3800009 | 7807730 | 36.91% | 4.196% |
| 24 | 10.969 | 1317 | 1323 | 1335 | rVB2 | 159489 | 375912 | 1.78% | 0.202% |
| 25 | 11.528 | 1411 | 1418 | 1424 | rVV2 | 120316 | 246641 | 1.17% | 0.133% |
| 26 | 12.446 | 1567 | 1574 | 1593 | rVV | 1936716 | 4090341 | 19.33% | 2.198% |
| 27 | 12.663 | 1604 | 1611 | 1630 | rVV | 1003298 | 2091312 | 9.89% | 1.124% |
| 28 | 13.463 | 1741 | 1747 | 1767 | rVV | 640072 | 1212008 | 5.73% | 0.651% |
| 29 | 13.657 | 1775 | 1780 | 1792 | rVB | 234111 | 452674 | 2.14% | 0.243% |
| 30 | 13.945 | 1821 | 1829 | 1834 | rBV | 357702 | 715193 | 3.38% | 0.384% |
| 31 | 13.998 | 1834 | 1838 | 1843 | rVV | 256949 | 442484 | 2.09% | 0.238% |
| 32 | 14.045 | 1843 | 1846 | 1855 | rVB | 114163 | 204537 | 0.97% | 0.110% |
| 33 | 14.187 | 1864 | 1870 | 1880 | rBV | 145900 | 356992 | 1.69% | 0.192% |
| 34 | 14.322 | 1885 | 1893 | 1895 | rVV2 | 125346 | 225595 | 1.07% | 0.121% |
| 35 | 14.345 | 1895 | 1897 | 1911 | rVV2 | 155687 | 252136 | 1.19% | 0.136% |
| 36 | 14.616 | 1930 | 1943 | 1948 | rBV2 | 1225158 | 2078537 | 9.82% | 1.117% |

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071323\
Data File : BP016208.D
Acq On : 13 Jul 2023 14:57
Operator : MA/JU
Sample : 03385-01ME 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EX7X7ME

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-BP071223.MA.M
Title : SVOA CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 37 | 14.681 | 1948 | 1954 | 1971 | rVV | 3809713 | 6068269 | 28.68% | 3.262% |
| 38 | 14.822 | 1971 | 1978 | 1990 | rVV4 | 116901 | 371199 | 1.75% | 0.200% |
| 39 | 15.022 | 2006 | 2012 | 2022 | rVV | 2512000 | 4260654 | 20.14% | 2.290% |
| 40 | 15.104 | 2022 | 2026 | 2031 | rVV | 109100 | 213709 | 1.01% | 0.115% |
| 41 | 15.616 | 2107 | 2113 | 2117 | rVV2 | 203298 | 314244 | 1.49% | 0.169% |
| 42 | 15.675 | 2117 | 2123 | 2136 | rVV | 3448396 | 5646999 | 26.69% | 3.035% |
| 43 | 15.798 | 2136 | 2144 | 2147 | rVV | 185443 | 472547 | 2.23% | 0.254% |
| 44 | 15.834 | 2147 | 2150 | 2156 | rVV | 402753 | 646729 | 3.06% | 0.348% |
| 45 | 15.934 | 2162 | 2167 | 2171 | rVV2 | 217254 | 371411 | 1.76% | 0.200% |
| 46 | 15.981 | 2171 | 2175 | 2184 | rVB | 443933 | 691011 | 3.27% | 0.371% |
| 47 | 16.128 | 2195 | 2200 | 2212 | rBV | 441686 | 1090607 | 5.16% | 0.586% |
| 48 | 16.475 | 2255 | 2259 | 2267 | rVB2 | 161521 | 230830 | 1.09% | 0.124% |
| 49 | 16.687 | 2288 | 2295 | 2300 | rVV2 | 293798 | 523952 | 2.48% | 0.282% |
| 50 | 16.910 | 2324 | 2333 | 2336 | rBV2 | 132207 | 267247 | 1.26% | 0.144% |
| 51 | 16.951 | 2336 | 2340 | 2343 | rVV2 | 134022 | 197440 | 0.93% | 0.106% |
| 52 | 17.104 | 2361 | 2366 | 2377 | rBV4 | 145215 | 372274 | 1.76% | 0.200% |
| 53 | 17.204 | 2377 | 2383 | 2394 | rVB | 1025415 | 1505243 | 7.11% | 0.809% |
| 54 | 17.381 | 2407 | 2413 | 2416 | rBV | 1659970 | 2492444 | 11.78% | 1.340% |
| 55 | 17.428 | 2416 | 2421 | 2428 | rVV | 13644978 | 21156146 | 100.00% | 11.371% |
| 56 | 17.486 | 2428 | 2431 | 2433 | rVV | 195889 | 288393 | 1.36% | 0.155% |
| 57 | 17.528 | 2433 | 2438 | 2449 | rVV | 3020716 | 5096064 | 24.09% | 2.739% |
| 58 | 17.622 | 2450 | 2454 | 2461 | rVV | 177598 | 324877 | 1.54% | 0.175% |
| 59 | 17.810 | 2478 | 2486 | 2497 | rBV | 1162334 | 2083488 | 9.85% | 1.120% |
| 60 | 17.892 | 2497 | 2500 | 2504 | rVB | 165187 | 204667 | 0.97% | 0.110% |
| 61 | 18.245 | 2555 | 2560 | 2565 | rBV | 899586 | 1310227 | 6.19% | 0.704% |
| 62 | 18.304 | 2565 | 2570 | 2579 | rVV | 1015543 | 1718101 | 8.12% | 0.923% |
| 63 | 18.392 | 2579 | 2585 | 2588 | rVV | 327937 | 486537 | 2.30% | 0.261% |
| 64 | 18.439 | 2588 | 2593 | 2597 | rVV | 2052860 | 3050381 | 14.42% | 1.639% |
| 65 | 18.481 | 2597 | 2600 | 2608 | rVV | 496186 | 752319 | 3.56% | 0.404% |
| 66 | 18.733 | 2638 | 2643 | 2650 | rVV | 771810 | 1076283 | 5.09% | 0.578% |
| 67 | 18.816 | 2653 | 2657 | 2660 | rVV2 | 143836 | 219079 | 1.04% | 0.118% |
| 68 | 18.969 | 2679 | 2683 | 2687 | rBV | 163650 | 212922 | 1.01% | 0.114% |
| 69 | 19.128 | 2705 | 2710 | 2715 | rBV | 247253 | 406503 | 1.92% | 0.218% |
| 70 | 19.198 | 2715 | 2722 | 2724 | rBV2 | 135982 | 232354 | 1.10% | 0.125% |
| 71 | 19.263 | 2729 | 2733 | 2735 | rBV2 | 121585 | 200690 | 0.95% | 0.108% |
| 72 | 19.392 | 2749 | 2755 | 2762 | rBV | 11988205 | 16954923 | 80.14% | 9.113% |
| 73 | 19.628 | 2789 | 2795 | 2805 | rBV2 | 333596 | 770938 | 3.64% | 0.414% |
| 74 | 19.745 | 2805 | 2815 | 2823 | rVV2 | 8252372 | 15068281 | 71.22% | 8.099% |
| 75 | 19.822 | 2823 | 2828 | 2833 | rVB | 366832 | 509272 | 2.41% | 0.274% |
| 76 | 19.928 | 2842 | 2846 | 2851 | rBV | 575541 | 711719 | 3.36% | 0.383% |

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071323\
Data File : BP016208.D
Acq On : 13 Jul 2023 14:57
Operator : MA/JU
Sample : 03385-01ME 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EX7X7ME

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-BP071223.MA.M
Title : SVOA CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 77 | 20.063 | 2860 | 2869 | 2878 | rBV3 | 486116 | 1247947 | 5.90% | 0.671% |
| 78 | 20.198 | 2885 | 2892 | 2893 | rVV3 | 295114 | 486967 | 2.30% | 0.262% |
| 79 | 20.228 | 2893 | 2897 | 2905 | rVB | 1609732 | 2299384 | 10.87% | 1.236% |
| 80 | 20.328 | 2907 | 2914 | 2918 | rVV | 1835030 | 2526378 | 11.94% | 1.358% |
| 81 | 20.386 | 2918 | 2924 | 2933 | rVV3 | 591832 | 1582126 | 7.48% | 0.850% |
| 82 | 20.463 | 2934 | 2937 | 2941 | rVV3 | 167958 | 243394 | 1.15% | 0.131% |
| 83 | 20.522 | 2944 | 2947 | 2951 | rVV | 268493 | 396227 | 1.87% | 0.213% |
| 84 | 20.569 | 2951 | 2955 | 2967 | rVB4 | 294434 | 775447 | 3.67% | 0.417% |
| 85 | 21.127 | 3046 | 3050 | 3052 | rBV | 586080 | 746759 | 3.53% | 0.401% |
| 86 | 21.157 | 3052 | 3055 | 3059 | rVV | 682008 | 905686 | 4.28% | 0.487% |
| 87 | 21.204 | 3059 | 3063 | 3074 | rVB3 | 635687 | 1438870 | 6.80% | 0.773% |
| 88 | 21.339 | 3083 | 3086 | 3088 | rBV | 317158 | 392622 | 1.86% | 0.211% |
| 89 | 21.457 | 3099 | 3106 | 3112 | rBV2 | 5188753 | 10933054 | 51.68% | 5.876% |
| 90 | 21.516 | 3112 | 3116 | 3125 | rVB | 3444254 | 5473885 | 25.87% | 2.942% |
| 91 | 21.639 | 3133 | 3137 | 3143 | rVB | 893929 | 1228326 | 5.81% | 0.660% |
| 92 | 22.074 | 3208 | 3211 | 3216 | rVB2 | 454216 | 600214 | 2.84% | 0.323% |
| 93 | 22.251 | 3238 | 3241 | 3244 | rVB | 231185 | 295619 | 1.40% | 0.159% |
| 94 | 22.292 | 3244 | 3248 | 3251 | rBV2 | 295335 | 477959 | 2.26% | 0.257% |
| 95 | 23.204 | 3396 | 3403 | 3409 | rBV | 2977483 | 7293455 | 34.47% | 3.920% |
| 96 | 23.404 | 3433 | 3437 | 3447 | rVB | 417522 | 704869 | 3.33% | 0.379% |
| 97 | 23.745 | 3490 | 3495 | 3502 | rBV | 813260 | 1544827 | 7.30% | 0.830% |
| 98 | 23.863 | 3509 | 3515 | 3526 | rVB | 1373002 | 3241733 | 15.32% | 1.742% |
| 99 | 23.968 | 3527 | 3533 | 3540 | rBV | 1324716 | 2795532 | 13.21% | 1.503% |
| 100 | 26.615 | 3975 | 3983 | 3998 | rVB | 1008900 | 3316925 | 15.68% | 1.783% |

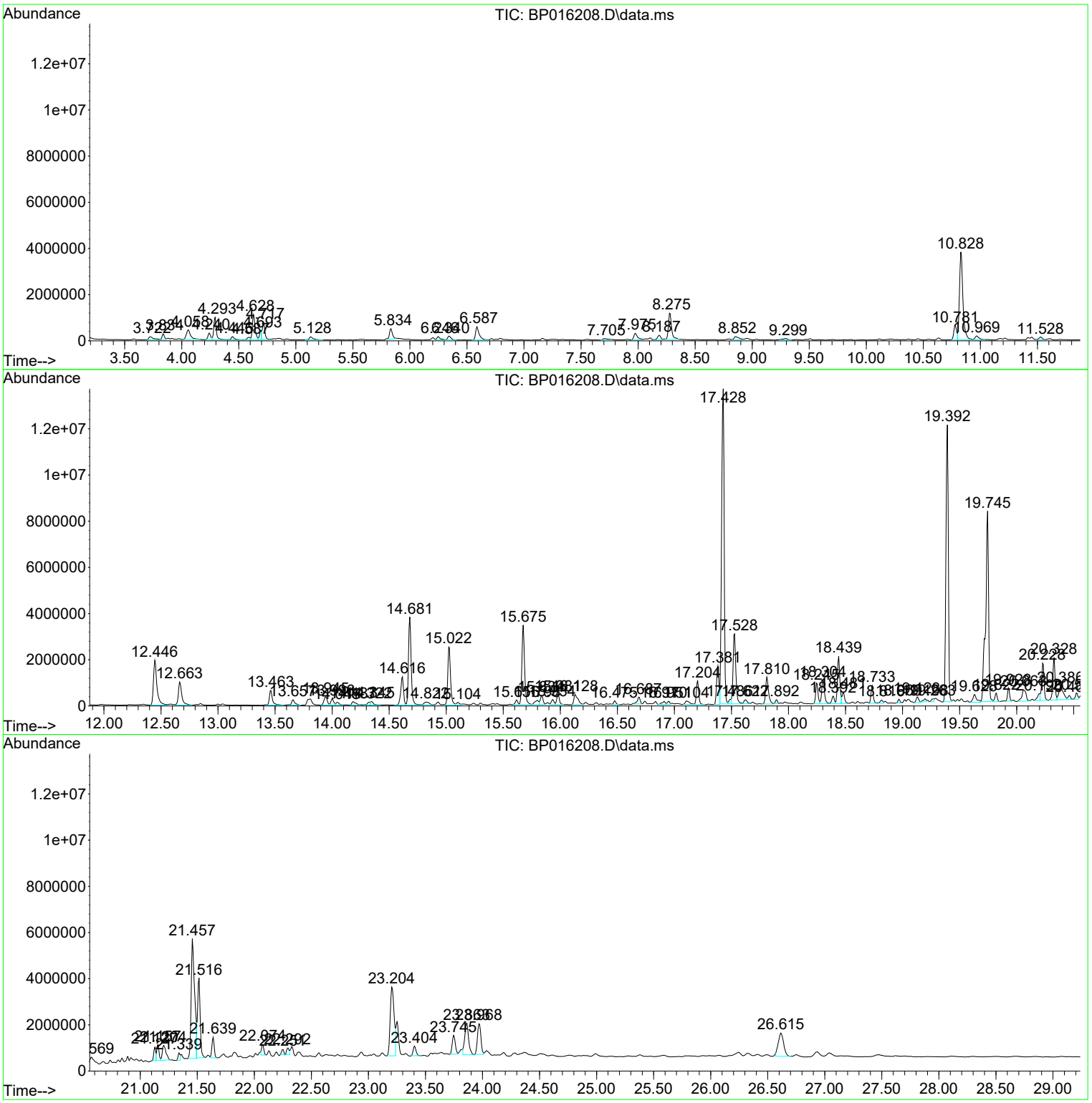
Sum of corrected areas: 186057091

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071323\
Data File : BP016208.D
Acq On : 13 Jul 2023 14:57
Operator : MA/JU
Sample : 03385-01ME 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EX7X7ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.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\BP071323\
Data File : BP016208.D
Acq On : 13 Jul 2023 14:57
Operator : MA/JU
Sample : 03385-01ME 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EX7X7ME

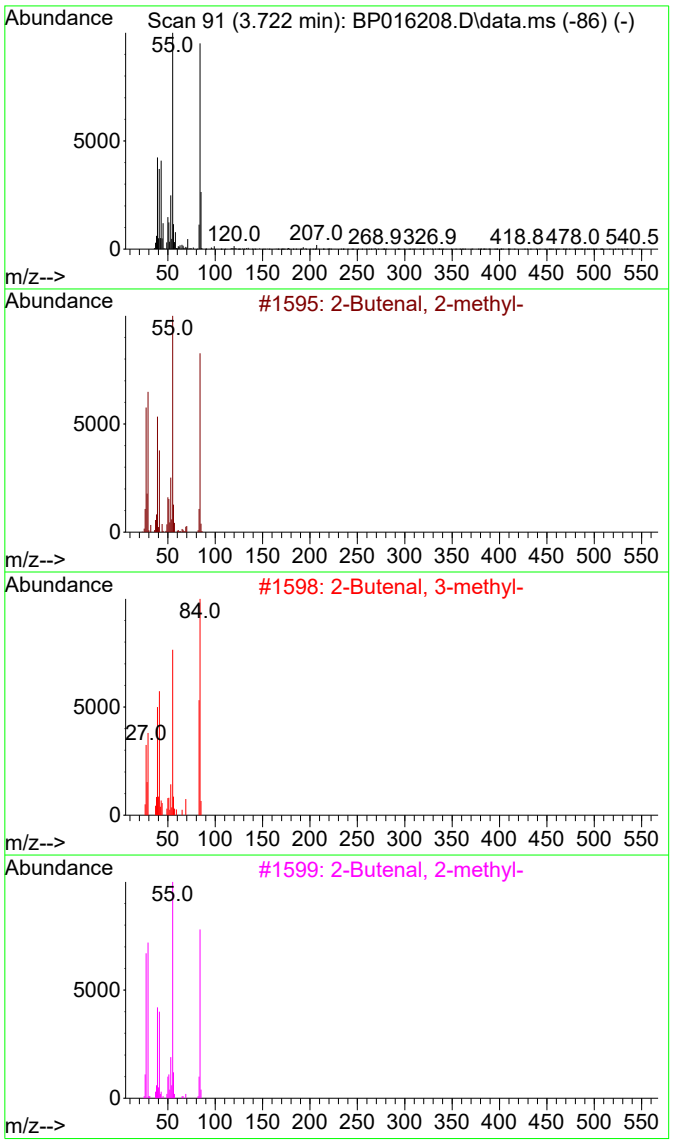
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

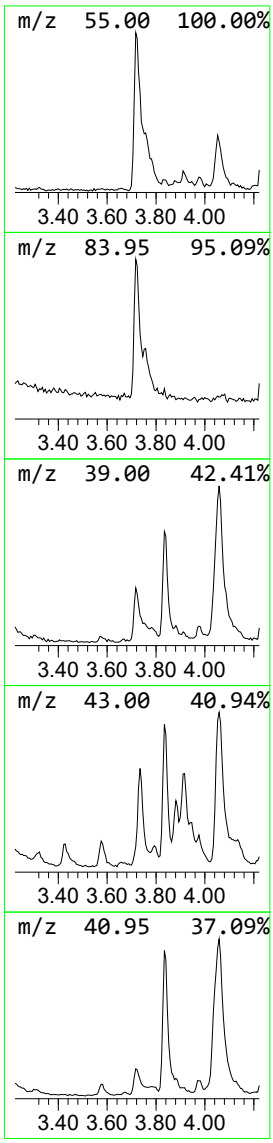
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 1 2-Butenal, 2-methyl- Concentration Rank 12

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 3.722 | 12.38 ng/ul | 405226 | 1,4-Dichlorobenzene-d4 | 7.975 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Butenal, 2-methyl- | 84 | C5H8O | 001115-11-3 | 90 |
| 2 | 2-Butenal, 3-methyl- | 84 | C5H8O | 000107-86-8 | 87 |
| 3 | 2-Butenal, 2-methyl- | 84 | C5H8O | 001115-11-3 | 87 |
| 4 | 2-Pentenal, (E)- | 84 | C5H8O | 001576-87-0 | 72 |
| 5 | 2-Pentenal, (E)- | 84 | C5H8O | 001576-87-0 | 64 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071323\
Data File : BP016208.D
Acq On : 13 Jul 2023 14:57
Operator : MA/JU
Sample : 03385-01ME 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EX7X7ME

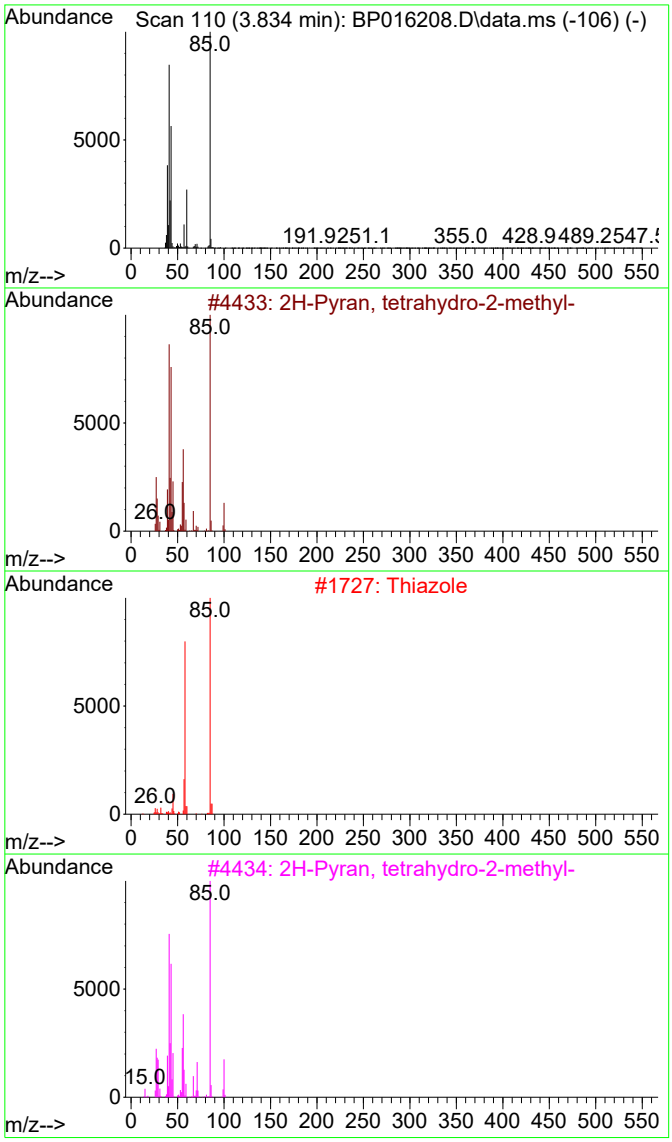
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

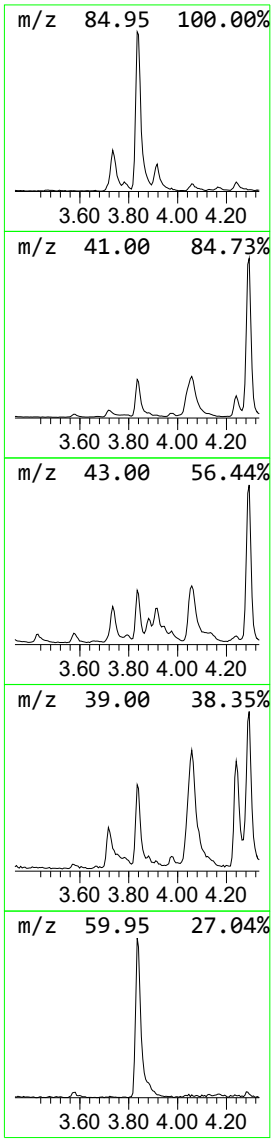
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 2 2H-Pyran, tetrahydro-2-methyl- Concentration Rank 15

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 3.834 | 11.24 ng/ul | 367934 | 1,4-Dichlorobenzene-d4 | 7.975 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2H-Pyran, tetrahydro-2-methyl- | 100 | C6H12O | 010141-72-7 | 50 |
| 2 | Thiazole | 85 | C3H3NS | 000288-47-1 | 35 |
| 3 | 2H-Pyran, tetrahydro-2-methyl- | 100 | C6H12O | 010141-72-7 | 33 |
| 4 | Piperazine, 1-nitroso- | 115 | C4H9N3O | 005632-47-3 | 25 |
| 5 | 2-Pyrrolidinone | 85 | C4H7NO | 000616-45-5 | 23 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071323\
Data File : BP016208.D
Acq On : 13 Jul 2023 14:57
Operator : MA/JU
Sample : 03385-01ME 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EX7X7ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

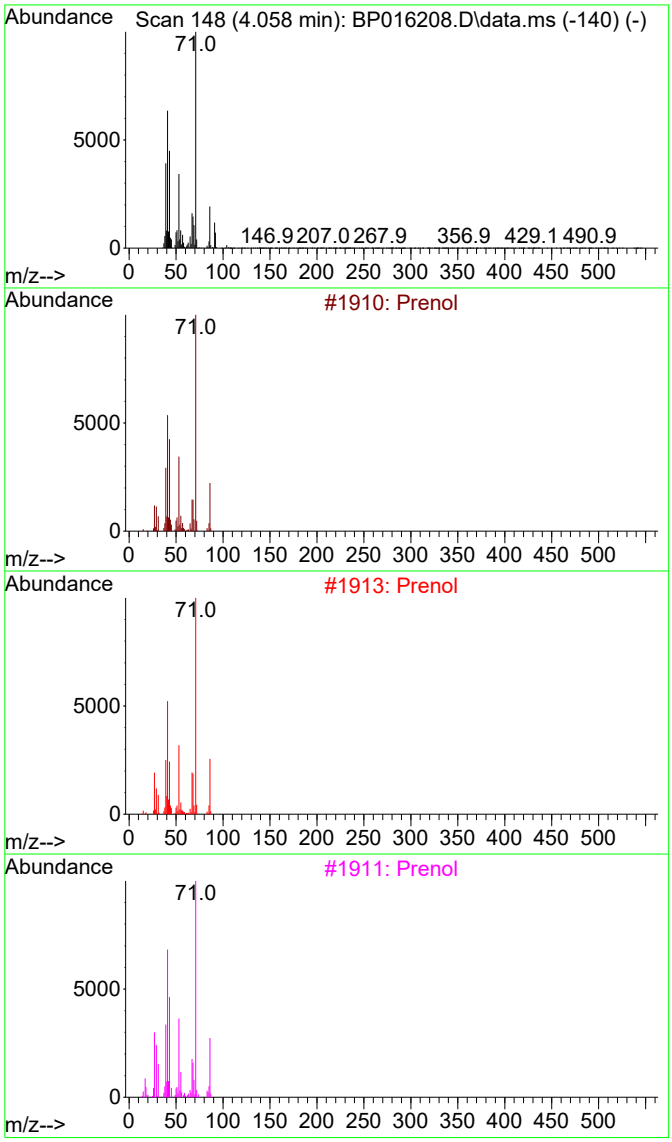
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

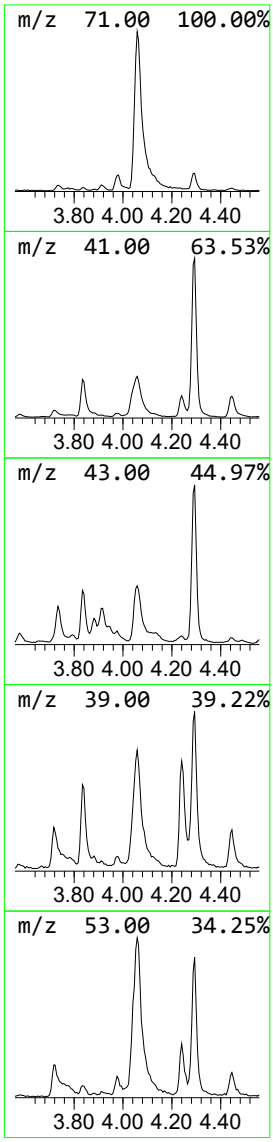
**********

Peak Number 3 Prenol Concentration Rank 6

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 4.058 | 32.58 ng/ul | 1066890 | 1,4-Dichlorobenzene-d4 | 7.975 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Prenol | 86 | C5H10O | 000556-82-1 | 92 |
| 2 | Prenol | 86 | C5H10O | 000556-82-1 | 70 |
| 3 | Prenol | 86 | C5H10O | 000556-82-1 | 58 |
| 4 | 2-Buten-1-ol, 2-methyl- | 86 | C5H10O | 004675-87-0 | 58 |
| 5 | 2-Buten-1-ol, 2-methyl- | 86 | C5H10O | 004675-87-0 | 58 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071323\
Data File : BP016208.D
Acq On : 13 Jul 2023 14:57
Operator : MA/JU
Sample : 03385-01ME 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EX7X7ME

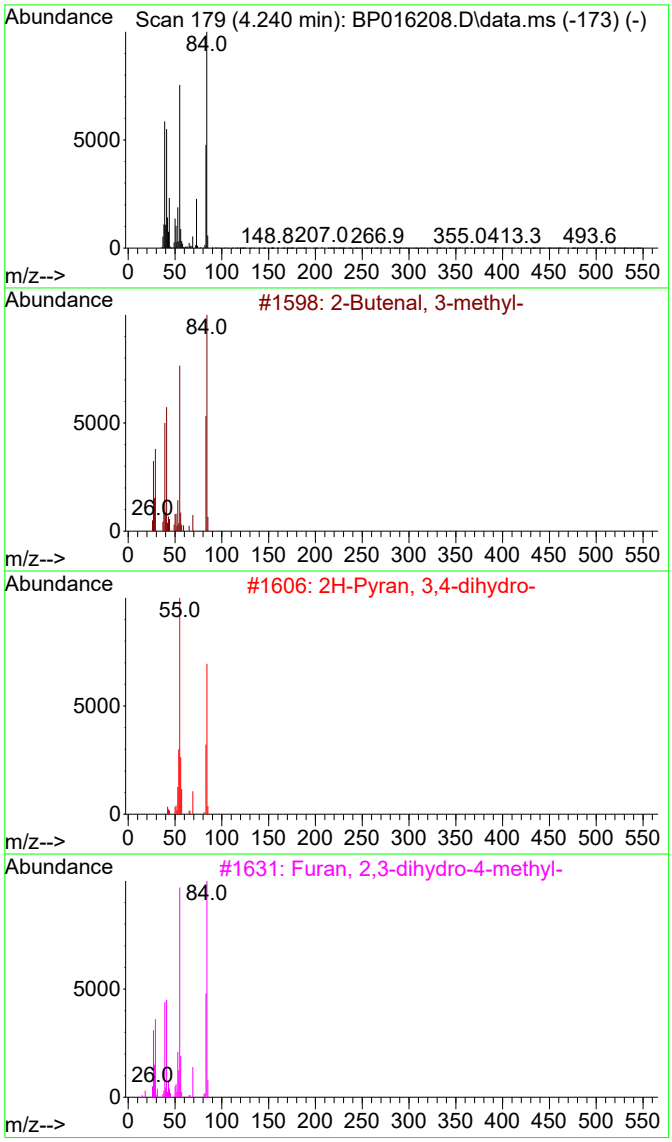
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

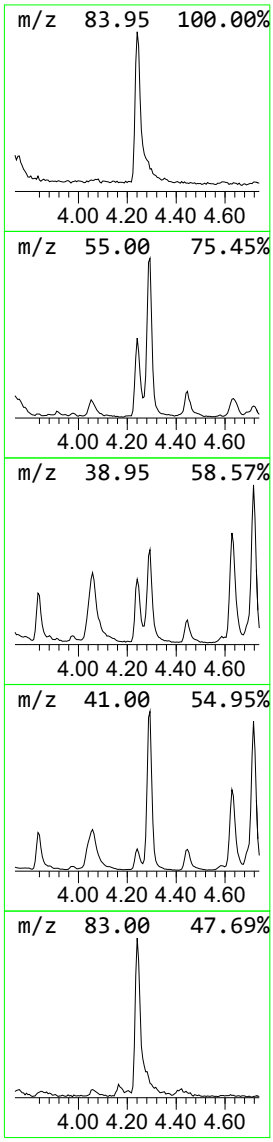
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 4 2-Butenal, 3-methyl- Concentration Rank 10

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 4.240 | 14.26 ng/ul | 466858 | 1,4-Dichlorobenzene-d4 | 7.975 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Butenal, 3-methyl- | 84 | C5H8O | 000107-86-8 | 87 |
| 2 | 2H-Pyran, 3,4-dihydro- | 84 | C5H8O | 000110-87-2 | 64 |
| 3 | Furan, 2,3-dihydro-4-methyl- | 84 | C5H8O | 034314-83-5 | 59 |
| 4 | 2-Butenal, 2-methyl- | 84 | C5H8O | 001115-11-3 | 50 |
| 5 | 2-Ethylacrolein | 84 | C5H8O | 000922-63-4 | 50 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071323\
Data File : BP016208.D
Acq On : 13 Jul 2023 14:57
Operator : MA/JU
Sample : 03385-01ME 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EX7X7ME

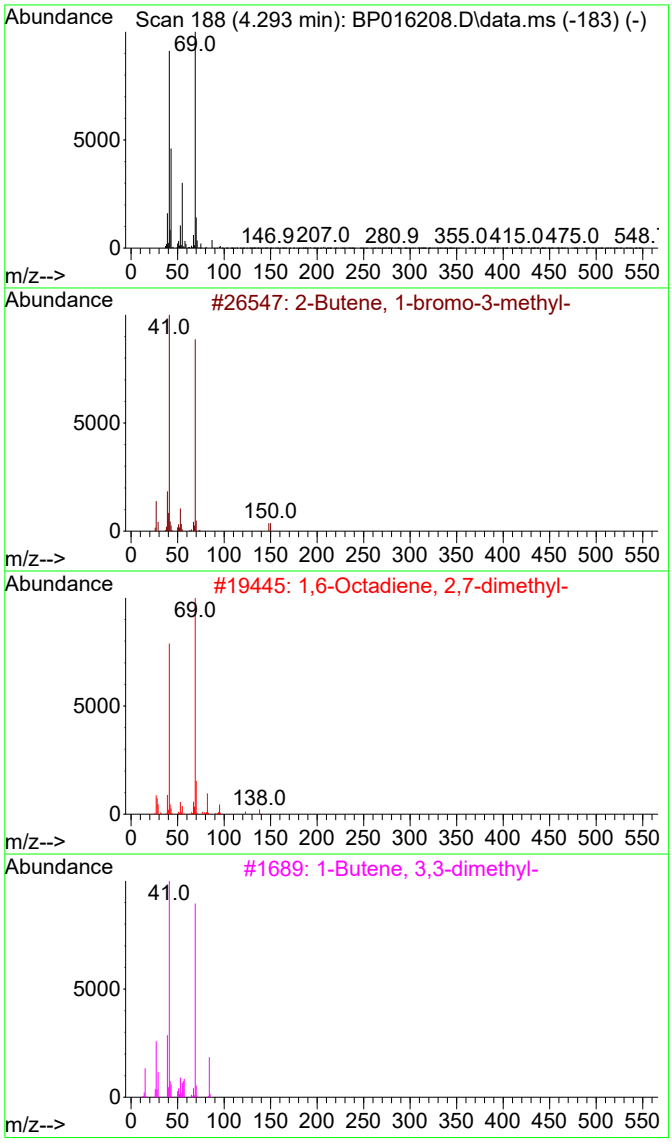
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

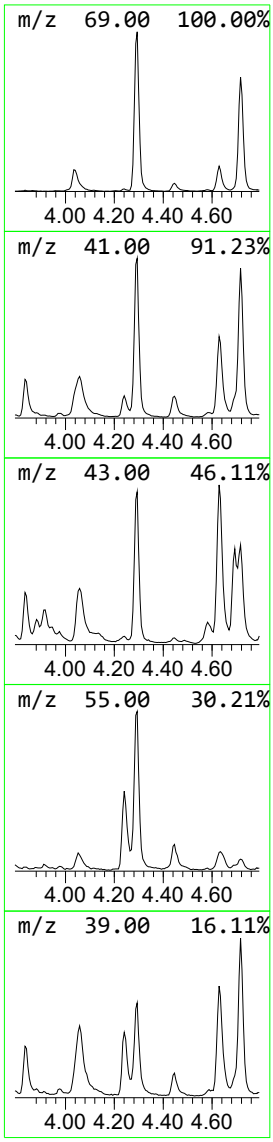
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 5 2-Butene, 1-bromo-3-methyl- Concentration Rank 3

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 4.293 | 45.28 ng/ul | 1482550 | 1,4-Dichlorobenzene-d4 | 7.975 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Butene, 1-bromo-3-methyl- | 148 | C5H9Br | 000870-63-3 | 59 |
| 2 | 1,6-Octadiene, 2,7-dimethyl- | 138 | C10H18 | 040195-09-3 | 56 |
| 3 | 1-Butene, 3,3-dimethyl- | 84 | C6H12 | 000558-37-2 | 39 |
| 4 | 1-Pentene, 3,3-dimethyl- | 98 | C7H14 | 003404-73-7 | 39 |
| 5 | 1-Pentene, 3,3-dimethyl- | 98 | C7H14 | 003404-73-7 | 38 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071323\
Data File : BP016208.D
Acq On : 13 Jul 2023 14:57
Operator : MA/JU
Sample : 03385-01ME 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EX7X7ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

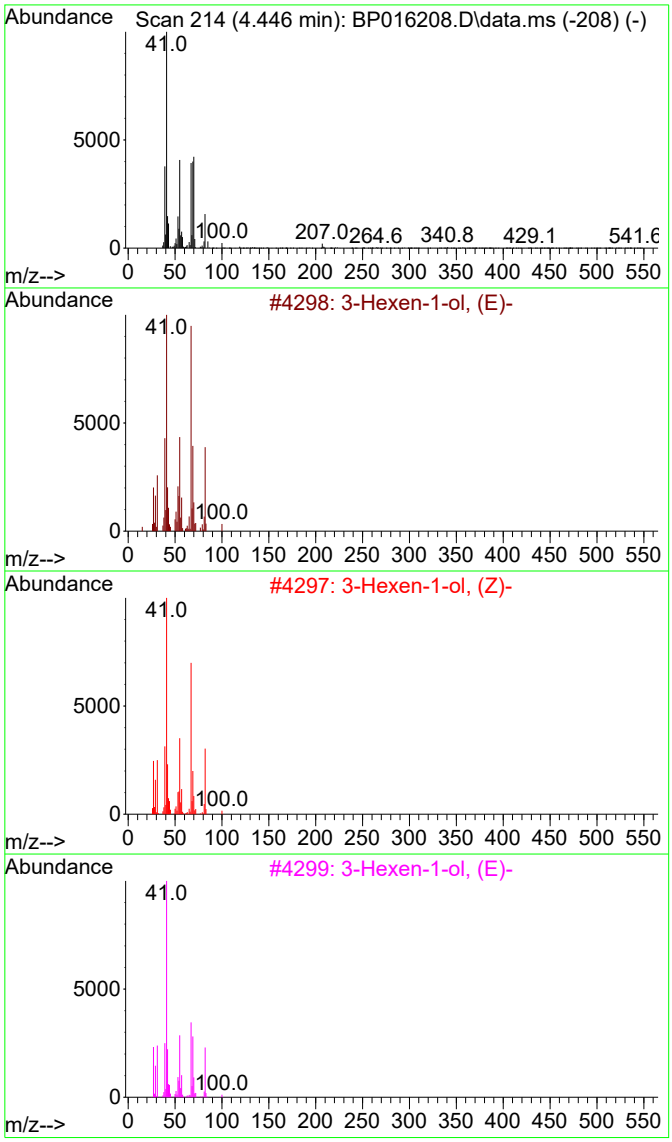
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

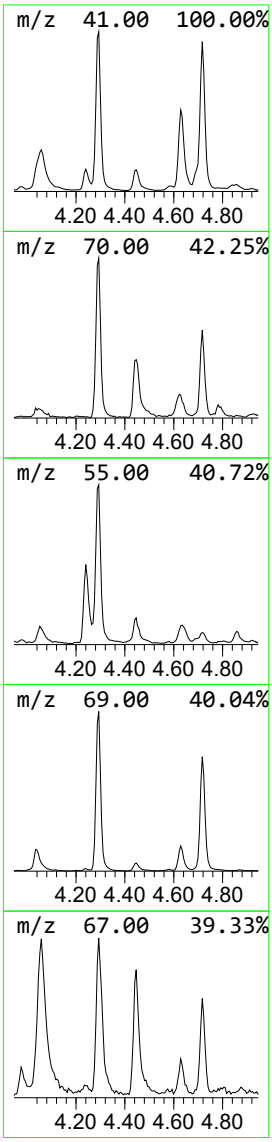
**********

Peak Number 6 unknown-01 Concentration Rank 23

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 4.446 | 8.30 ng/ul | 271853 | 1,4-Dichlorobenzene-d4 | 7.975 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 3-Hexen-1-ol, (E)- | 100 | C6H12O | 000928-97-2 | 47 |
| 2 | 3-Hexen-1-ol, (Z)- | 100 | C6H12O | 000928-96-1 | 43 |
| 3 | 3-Hexen-1-ol, (E)- | 100 | C6H12O | 000928-97-2 | 42 |
| 4 | 2-Hexene, 3,5-dimethyl- | 112 | C8H16 | 003404-79-3 | 38 |
| 5 | 3-Hexen-1-ol | 100 | C6H12O | 000544-12-7 | 38 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071323\
Data File : BP016208.D
Acq On : 13 Jul 2023 14:57
Operator : MA/JU
Sample : 03385-01ME 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EX7X7ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

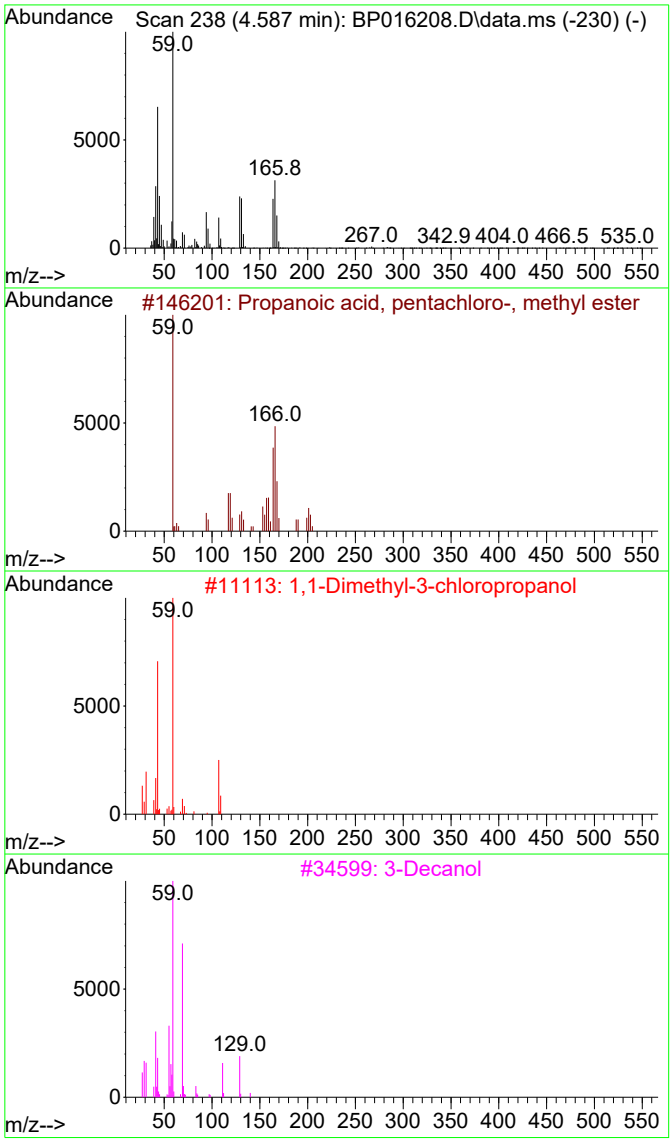
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

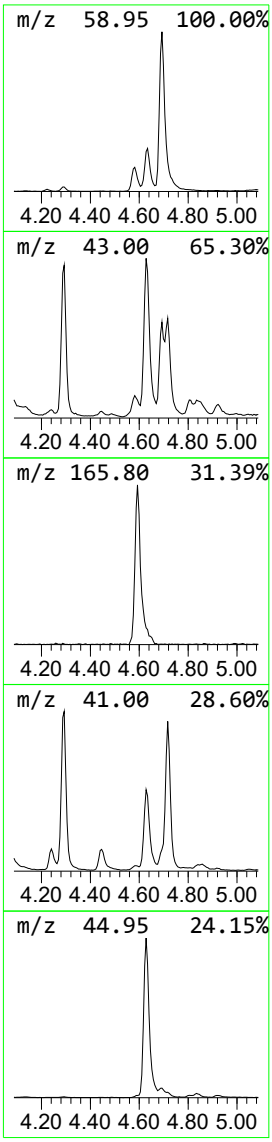
**********

Peak Number 7 unknown-02 Concentration Rank 29

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 4.587 | 6.48 ng/ul | 212337 | 1,4-Dichlorobenzene-d4 | 7.975 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Propanoic acid, pentachloro-, me... | 258 | C4H3Cl5O2 | 000813-46-7 | 36 |
| 2 | 1,1-Dimethyl-3-chloropropanol | 122 | C5H11ClO | 001985-88-2 | 27 |
| 3 | 3-Decanol | 158 | C10H22O | 001565-81-7 | 25 |
| 4 | 1-Propene, 3-[2-(2-methoxyethoxy... | 160 | C8H16O3 | 013752-97-1 | 22 |
| 5 | Guanidine, monothiocyanate | 116 | C2H4N4S | 056960-89-5 | 22 |



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Data File : BP016208.D
Acq On : 13 Jul 2023 14:57
Operator : MA/JU
Sample : 03385-01ME 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EX7X7ME

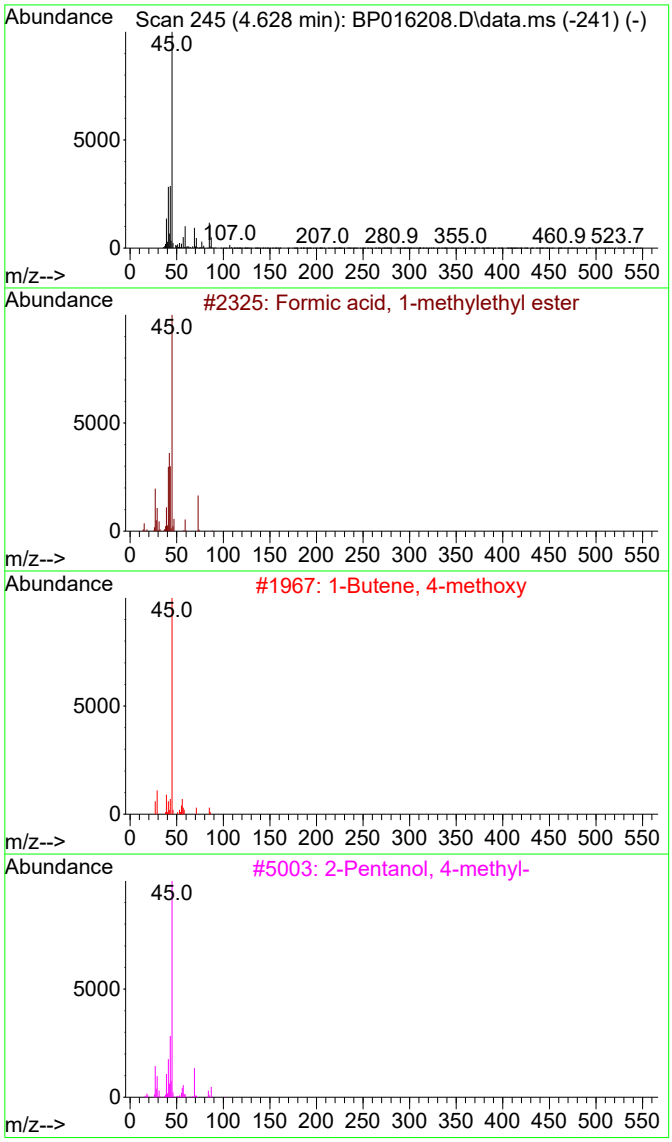
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

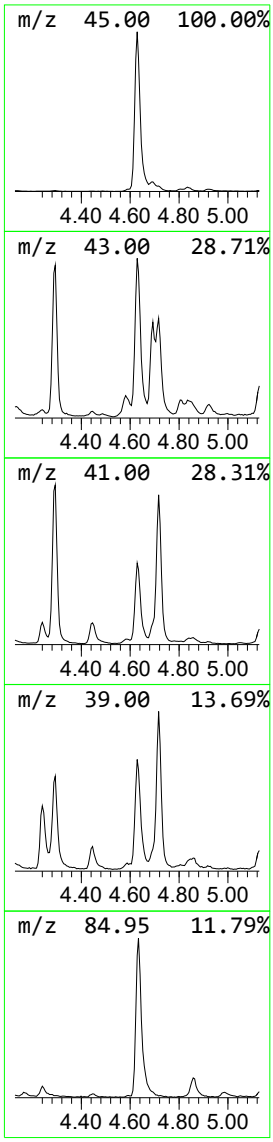
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 8 Formic acid, 1-methylethyl ... Concentration Rank 2

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 4.628 | 55.20 ng/ul | 1807340 | 1,4-Dichlorobenzene-d4 | 7.975 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Formic acid, 1-methylethyl ester | 88 | C4H8O2 | 000625-55-8 | 53 |
| 2 | 1-Butene, 4-methoxy | 86 | C5H10O | 004696-30-4 | 53 |
| 3 | 2-Pentanol, 4-methyl- | 102 | C6H14O | 000108-11-2 | 45 |
| 4 | 3-Ethyl-2-pentanol | 116 | C7H16O | 000609-27-8 | 39 |
| 5 | E-1-Methoxy-3-hexene | 114 | C7H14O | 121441-40-5 | 39 |



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Data File : BP016208.D
Acq On : 13 Jul 2023 14:57
Operator : MA/JU
Sample : 03385-01ME 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EX7X7ME

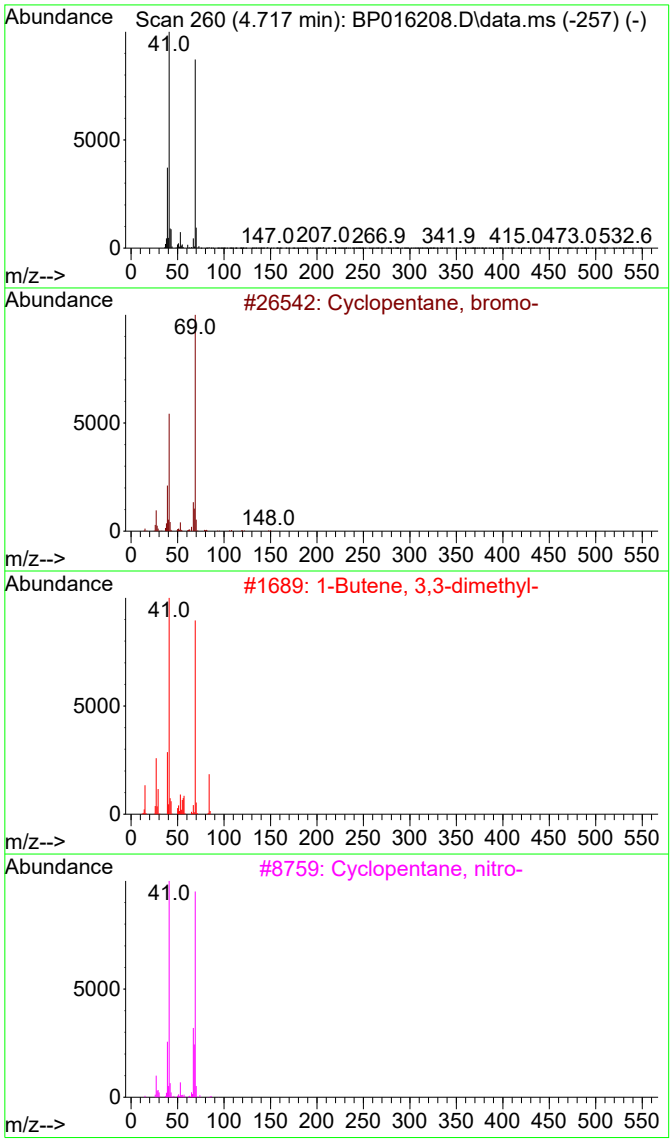
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

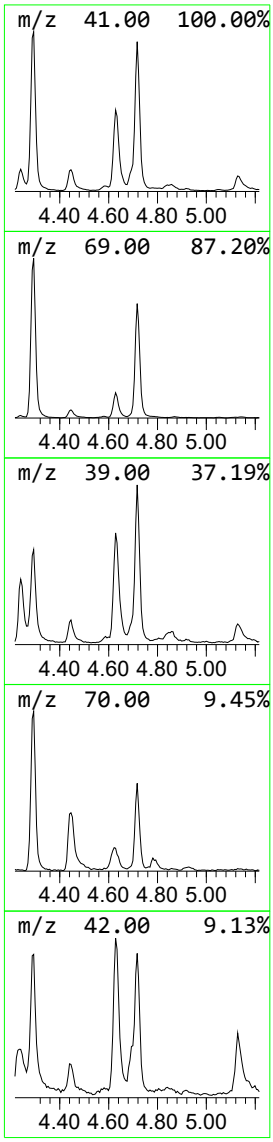
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 10 Cyclopentane, bromo- Concentration Rank 5

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 4.717 | 34.46 ng/ul | 1128210 | 1,4-Dichlorobenzene-d4 | 7.975 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Cyclopentane, bromo- | 148 | C5H9Br | 000137-43-9 | 56 |
| 2 | 1-Butene, 3,3-dimethyl- | 84 | C6H12 | 000558-37-2 | 43 |
| 3 | Cyclopentane, nitro- | 115 | C5H9NO2 | 002562-38-1 | 40 |
| 4 | 3-Cyclopropyl-3-oxopropanenitrile | 109 | C6H7NO | 118431-88-2 | 40 |
| 5 | Zinc, bis(3-methyl-2-butenyl)- | 202 | C10H18Zn | 066094-28-8 | 40 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071323\
Data File : BP016208.D
Acq On : 13 Jul 2023 14:57
Operator : MA/JU
Sample : 03385-01ME 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EX7X7ME

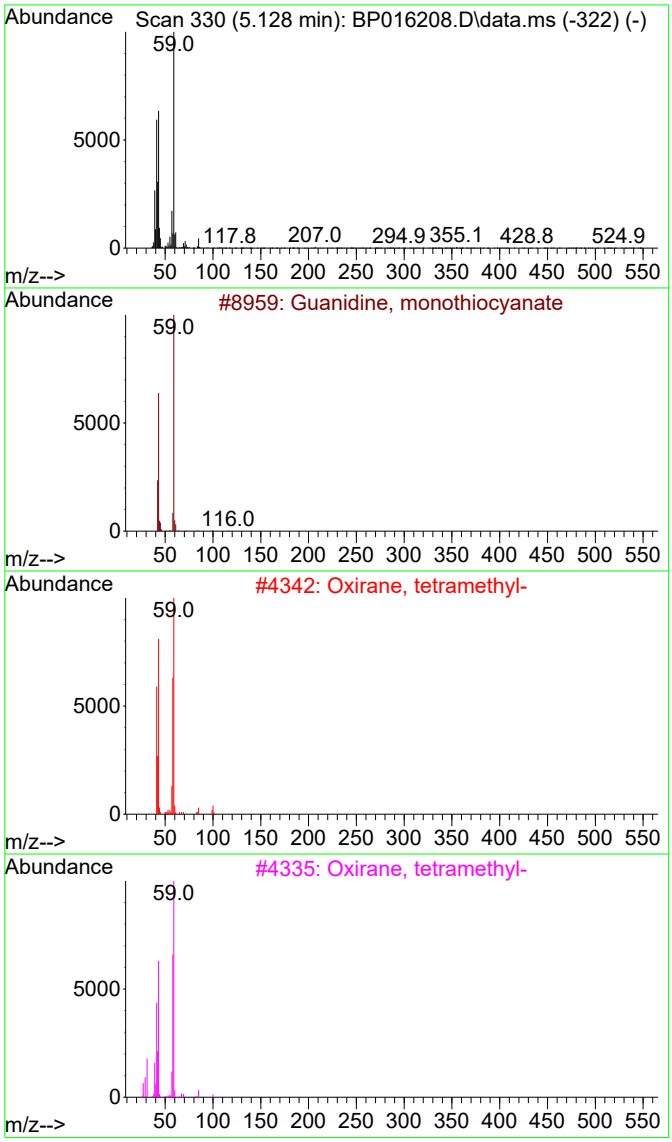
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

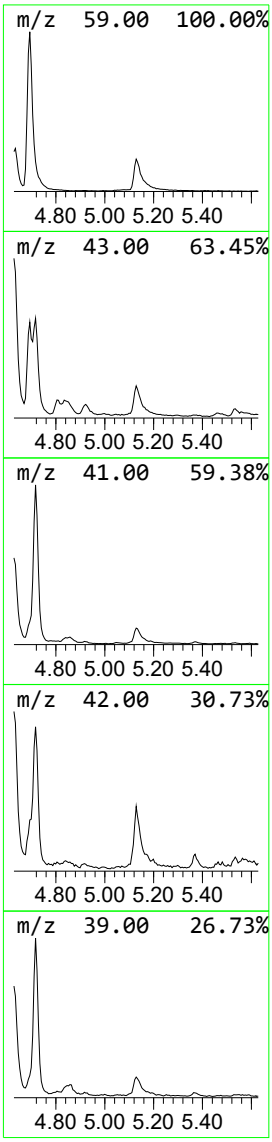
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 11 Guanidine, monothiocyanate Concentration Rank 19

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 5.128 | 9.97 ng/ul | 326487 | 1,4-Dichlorobenzene-d4 | 7.975 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Guanidine, monothiocyanate | 116 | C2H4N4S | 056960-89-5 | 64 |
| 2 | Oxirane, tetramethyl- | 100 | C6H12O | 005076-20-0 | 56 |
| 3 | Oxirane, tetramethyl- | 100 | C6H12O | 005076-20-0 | 56 |
| 4 | Acetamide | 59 | C2H5NO | 000060-35-5 | 53 |
| 5 | Oxirane, tetramethyl- | 100 | C6H12O | 005076-20-0 | 50 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071323\
Data File : BP016208.D
Acq On : 13 Jul 2023 14:57
Operator : MA/JU
Sample : 03385-01ME 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EX7X7ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

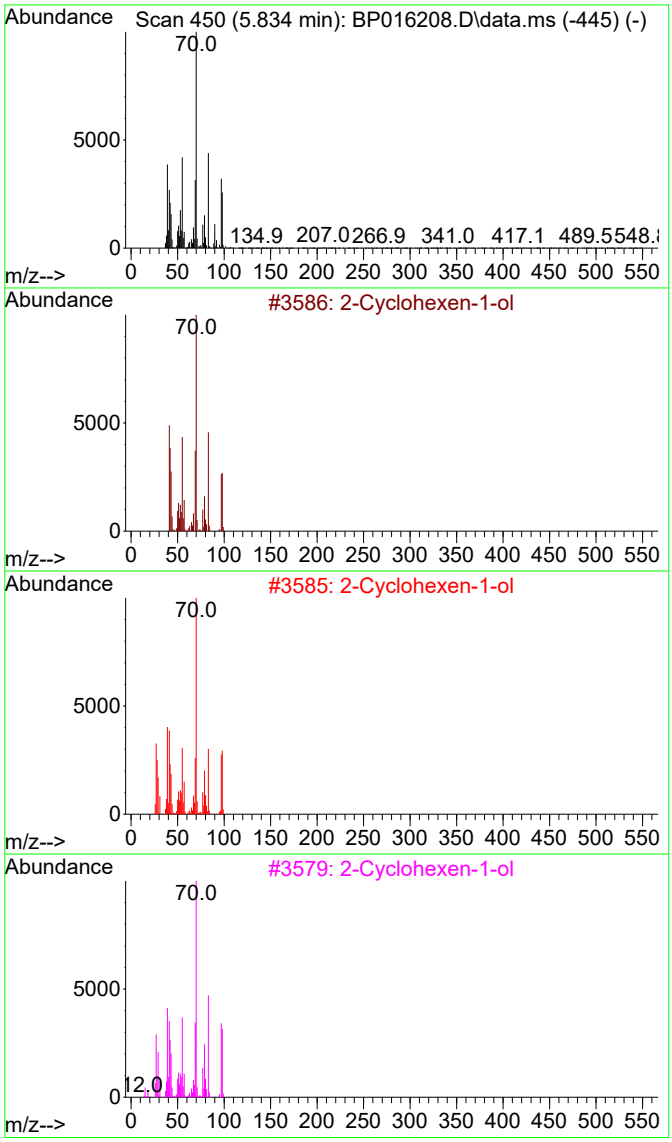
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

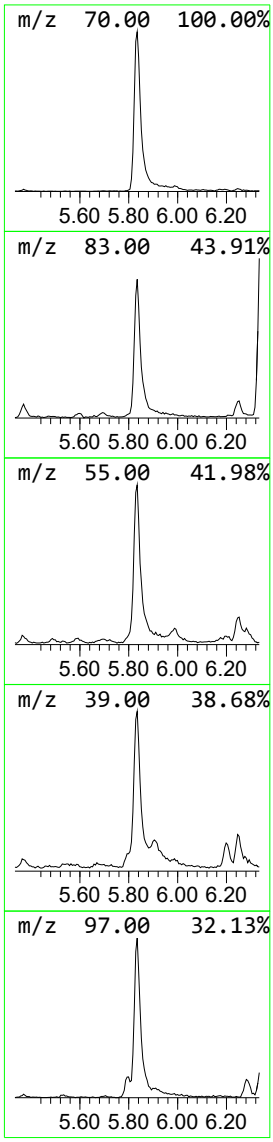
**********

Peak Number 12 2-Cyclohexen-1-ol Concentration Rank 7

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 5.834 | 24.52 ng/ul | 802788 | 1,4-Dichlorobenzene-d4 | 7.975 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Cyclohexen-1-ol | 98 | C6H10O | 000822-67-3 | 74 |
| 2 | 2-Cyclohexen-1-ol | 98 | C6H10O | 000822-67-3 | 72 |
| 3 | 2-Cyclohexen-1-ol | 98 | C6H10O | 000822-67-3 | 46 |
| 4 | 2-Cyclohexen-1-ol | 98 | C6H10O | 000822-67-3 | 45 |
| 5 | Cyclopentane, 1,3-dimethyl- | 98 | C7H14 | 002453-00-1 | 43 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071323\
Data File : BP016208.D
Acq On : 13 Jul 2023 14:57
Operator : MA/JU
Sample : 03385-01ME 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EX7X7ME

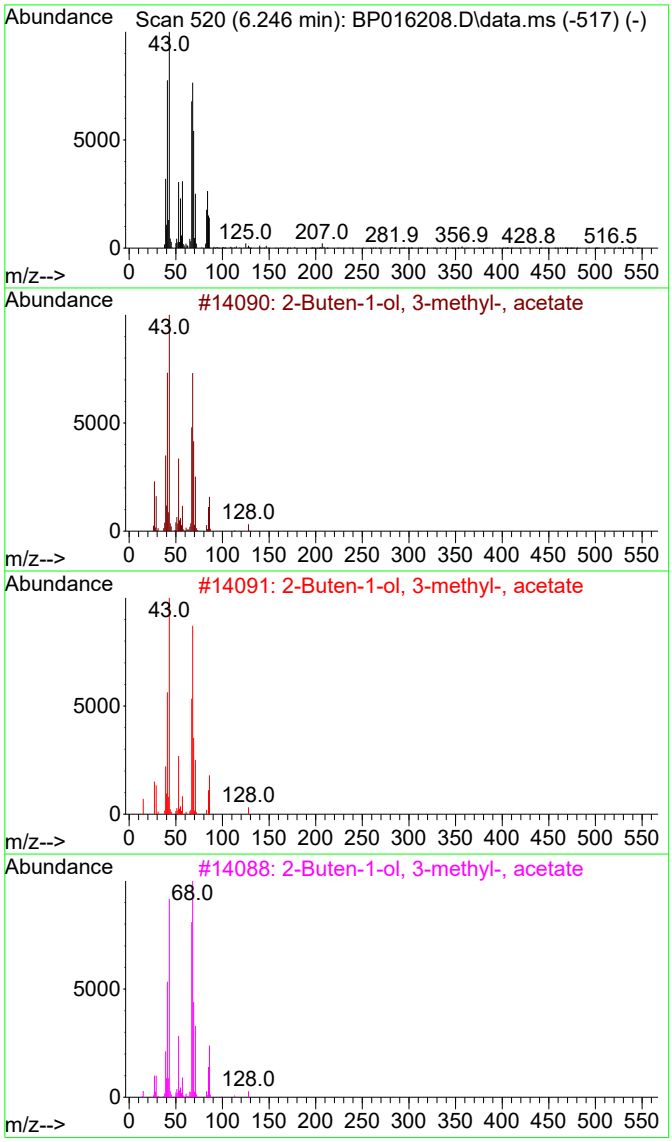
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

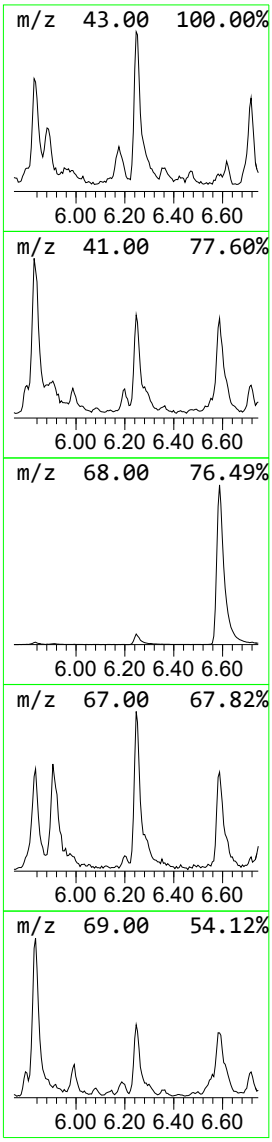
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 13 2-Buten-1-ol, 3-methyl-, ac... Concentration Rank 20

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 6.246 | 9.12 ng/ul | 298630 | 1,4-Dichlorobenzene-d4 | 7.975 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Buten-1-ol, 3-methyl-, acetate | 128 | C7H12O2 | 001191-16-8 | 72 |
| 2 | 2-Buten-1-ol, 3-methyl-, acetate | 128 | C7H12O2 | 001191-16-8 | 64 |
| 3 | 2-Buten-1-ol, 3-methyl-, acetate | 128 | C7H12O2 | 001191-16-8 | 64 |
| 4 | 2-Buten-1-ol, 3-methyl-, acetate | 128 | C7H12O2 | 001191-16-8 | 50 |
| 5 | Bicyclo[3.1.0]hexan-3-one | 96 | C6H8O | 001755-04-0 | 50 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071323\
Data File : BP016208.D
Acq On : 13 Jul 2023 14:57
Operator : MA/JU
Sample : 03385-01ME 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EX7X7ME

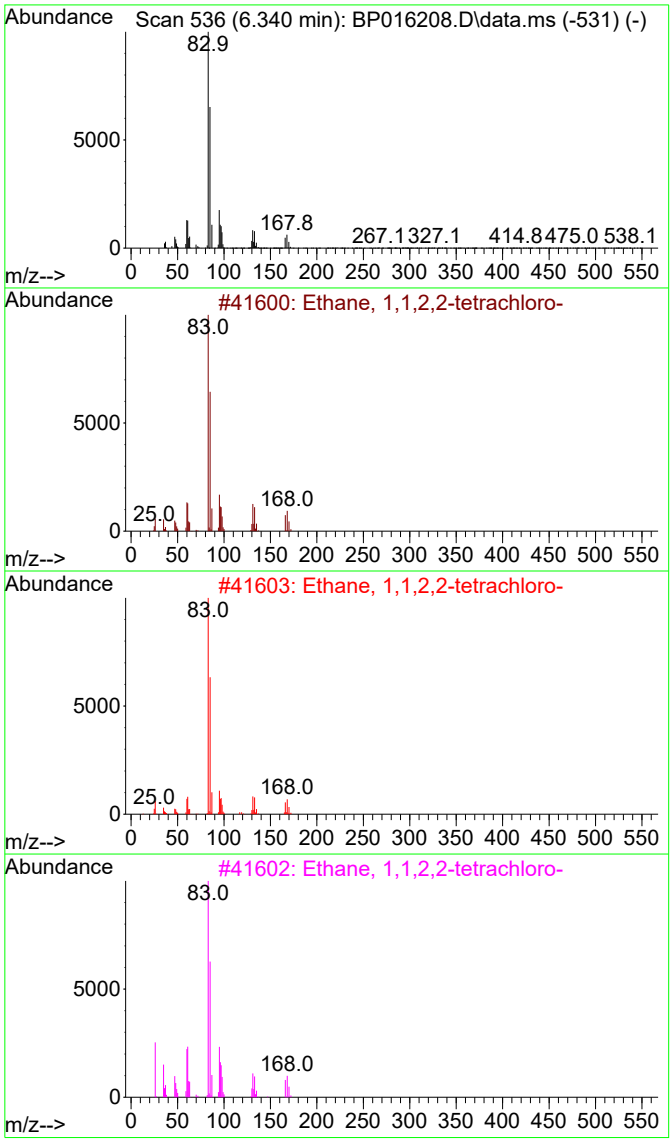
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

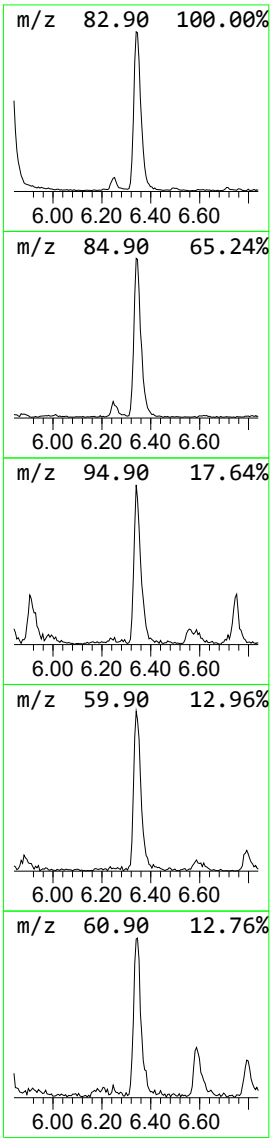
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 14 Ethane, 1,1,2,2-tetrachloro- Concentration Rank 17

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 6.340 | 10.74 ng/ul | 351674 | 1,4-Dichlorobenzene-d4 | 7.975 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Ethane, 1,1,2,2-tetrachloro- | 166 | C2H2Cl4 | 000079-34-5 | 94 |
| 2 | Ethane, 1,1,2,2-tetrachloro- | 166 | C2H2Cl4 | 000079-34-5 | 94 |
| 3 | Ethane, 1,1,2,2-tetrachloro- | 166 | C2H2Cl4 | 000079-34-5 | 91 |
| 4 | Ethane, 1,1,2-trichloro-2-fluoro- | 150 | C2H2Cl3F | 000359-28-4 | 56 |
| 5 | Ethane, 1,2,2-trichloro-1,1-difl... | 168 | C2HCl3F2 | 000354-21-2 | 53 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071323\
Data File : BP016208.D
Acq On : 13 Jul 2023 14:57
Operator : MA/JU
Sample : 03385-01ME 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EX7X7ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

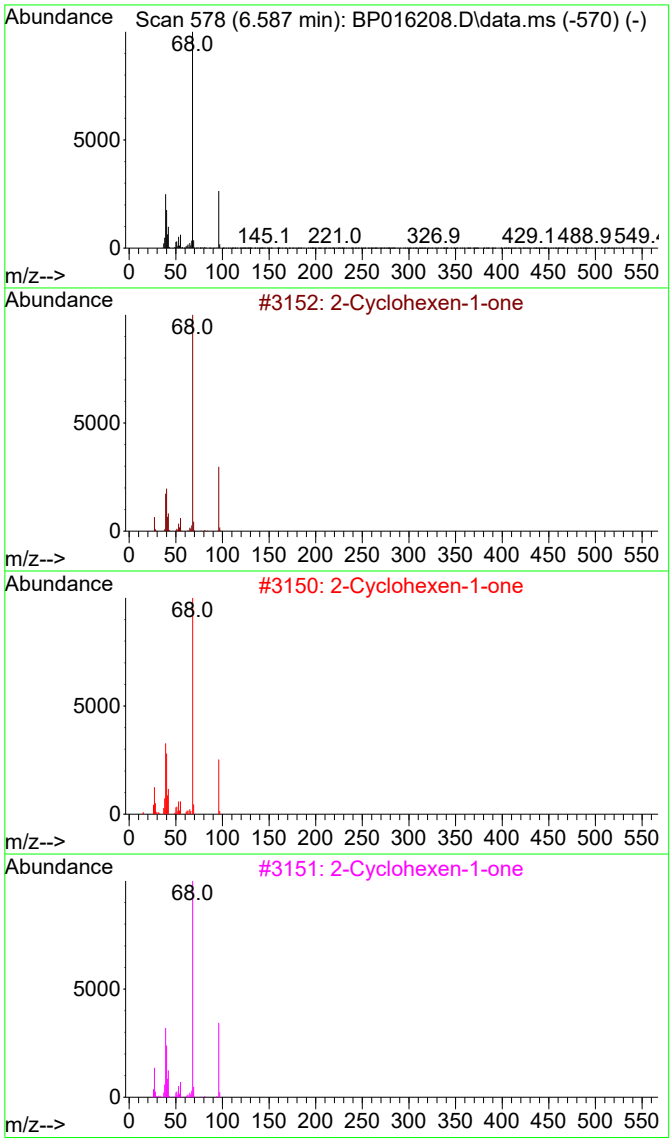
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

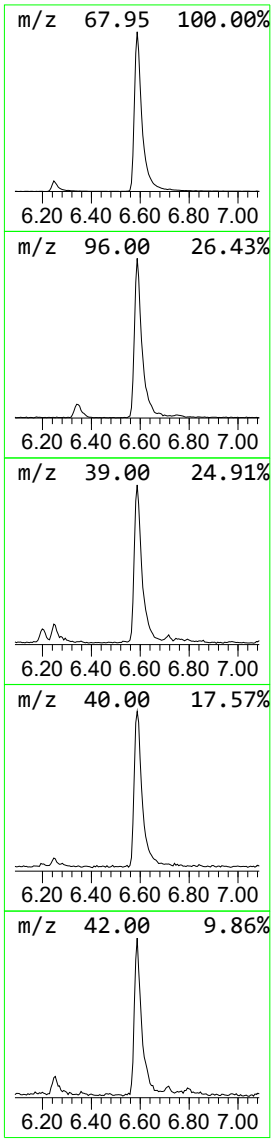
**********

Peak Number 15 2-Cyclohexen-1-one Concentration Rank 4

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 6.587 | 39.43 ng/ul | 1291110 | 1,4-Dichlorobenzene-d4 | 7.975 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Cyclohexen-1-one | 96 | C6H8O | 000930-68-7 | 91 |
| 2 | 2-Cyclohexen-1-one | 96 | C6H8O | 000930-68-7 | 91 |
| 3 | 2-Cyclohexen-1-one | 96 | C6H8O | 000930-68-7 | 91 |
| 4 | 2-Cyclohexen-1-one | 96 | C6H8O | 000930-68-7 | 91 |
| 5 | Bicyclo[3.1.0]hexan-2-one | 96 | C6H8O | 004160-49-0 | 12 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071323\
Data File : BP016208.D
Acq On : 13 Jul 2023 14:57
Operator : MA/JU
Sample : 03385-01ME 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EX7X7ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

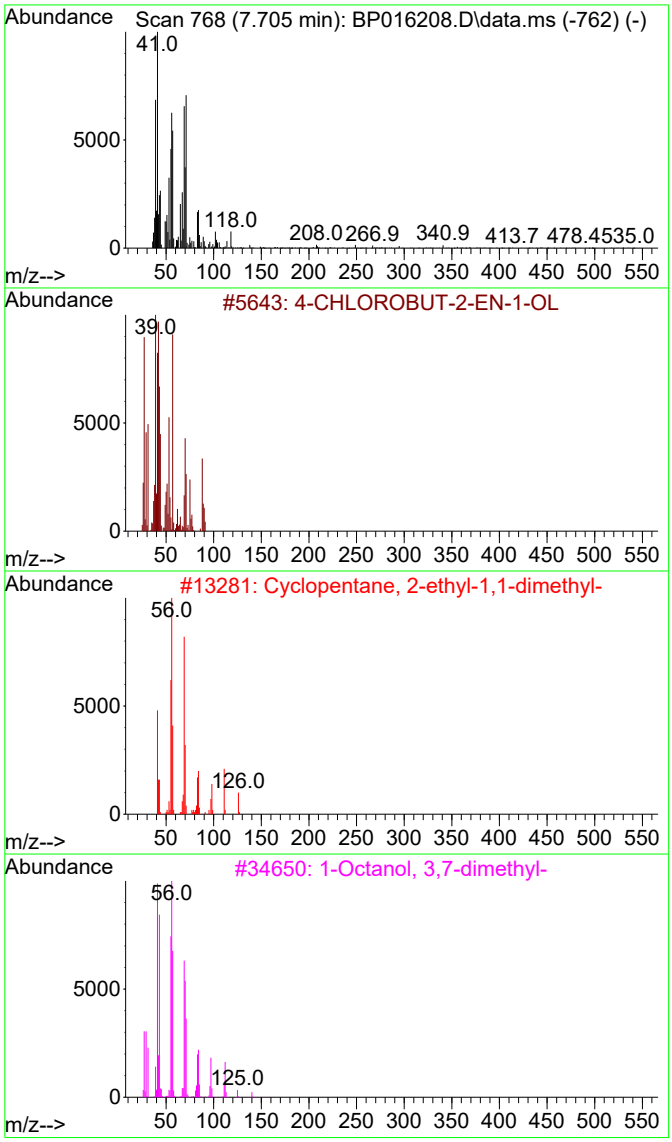
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

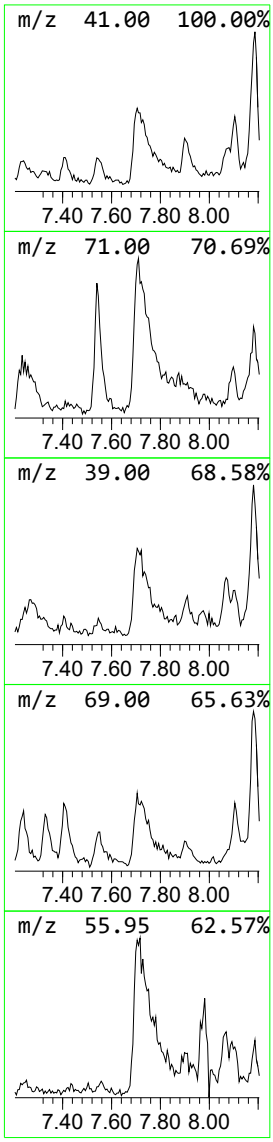
**********

Peak Number 16 unknown-03 Concentration Rank 27

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 7.705 | 6.65 ng/ul | 217595 | 1,4-Dichlorobenzene-d4 | 7.975 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 4-CHLOROBUT-2-EN-1-OL | 106 | C4H7ClO | 007523-44-6 | 25 |
| 2 | Cyclopentane, 2-ethyl-1,1-dimethyl- | 126 | C9H18 | 054549-80-3 | 22 |
| 3 | 1-Octanol, 3,7-dimethyl- | 158 | C10H22O | 000106-21-8 | 22 |
| 4 | Cyclopropane, 1,1-dimethyl-2-nonyl- | 196 | C14H28 | 041977-38-2 | 22 |
| 5 | 1,2-Dimethyl cyclopropene | 68 | C5H8 | 014309-32-1 | 18 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071323\
Data File : BP016208.D
Acq On : 13 Jul 2023 14:57
Operator : MA/JU
Sample : 03385-01ME 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EX7X7ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

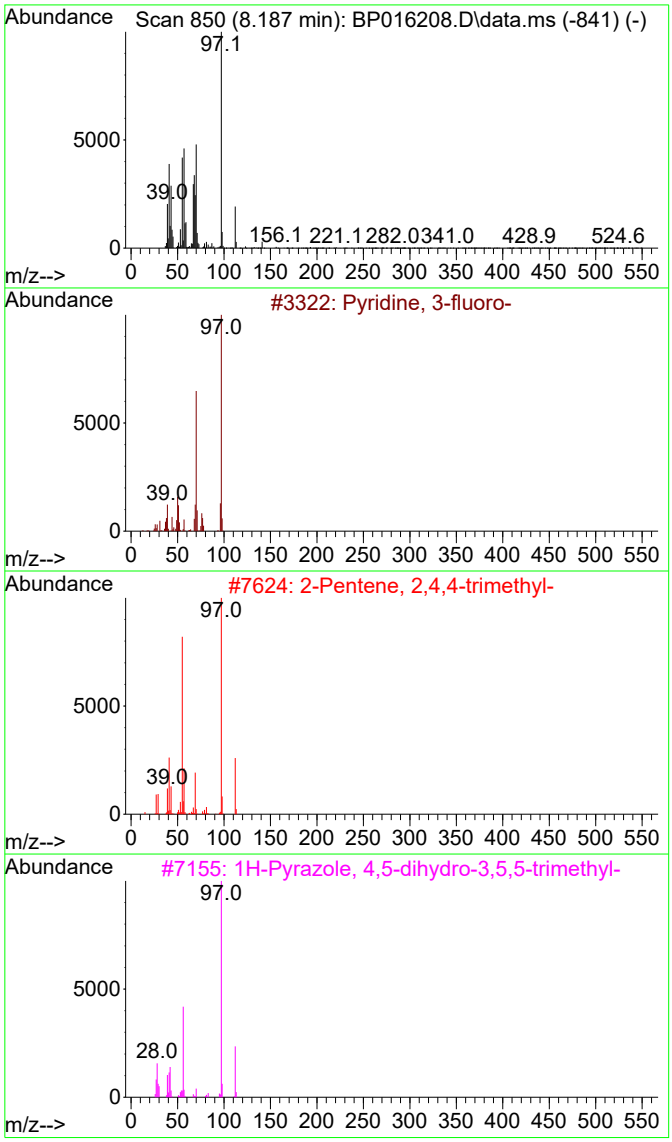
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

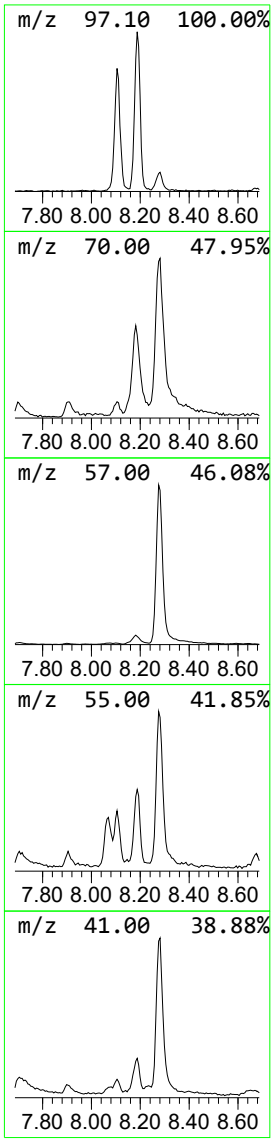
**********

Peak Number 17 unknown-04 Concentration Rank 14

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 8.187 | 11.47 ng/ul | 375600 | 1,4-Dichlorobenzene-d4 | 7.975 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Pyridine, 3-fluoro- | 97 | C5H4FN | 000372-47-4 | 46 |
| 2 | 2-Pentene, 2,4,4-trimethyl- | 112 | C8H16 | 000107-40-4 | 46 |
| 3 | 1H-Pyrazole, 4,5-dihydro-3,5,5-t... | 112 | C6H12N2 | 003975-85-7 | 43 |
| 4 | 1H-Pyrazole, 4,5-dihydro-3,5,5-t... | 112 | C6H12N2 | 003975-85-7 | 43 |
| 5 | Cyclohexane, 1,3-dimethyl- | 112 | C8H16 | 000591-21-9 | 43 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071323\
Data File : BP016208.D
Acq On : 13 Jul 2023 14:57
Operator : MA/JU
Sample : 03385-01ME 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EX7X7ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

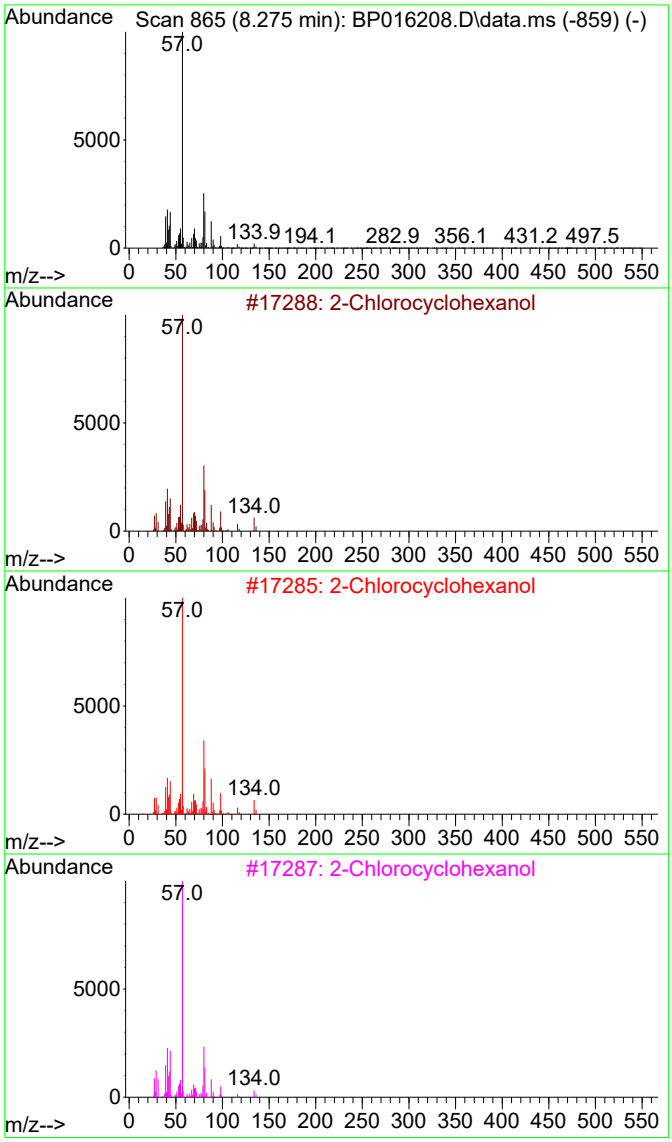
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

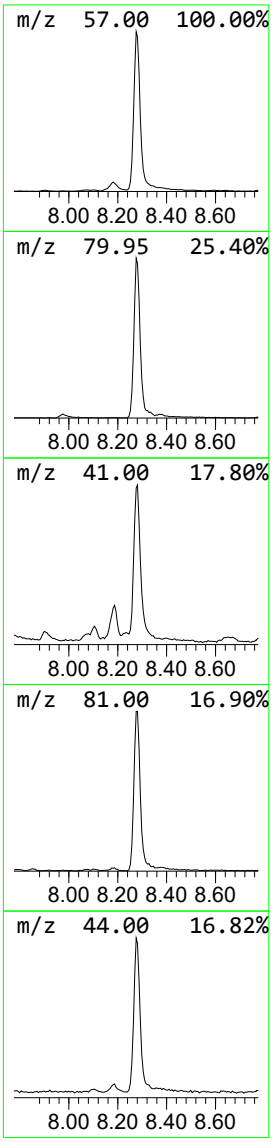
**********

Peak Number 18 2-Chlorocyclohexanol Concentration Rank 1

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 8.275 | 67.57 ng/ul | 2212300 | 1,4-Dichlorobenzene-d4 | 7.975 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2-Chlorocyclohexanol | 134 | C6H11ClO | 001561-86-0 | 93 |
| 2 | 2-Chlorocyclohexanol | 134 | C6H11ClO | 001561-86-0 | 93 |
| 3 | 2-Chlorocyclohexanol | 134 | C6H11ClO | 001561-86-0 | 87 |
| 4 | 2-Chlorocyclohexanol | 134 | C6H11ClO | 001561-86-0 | 64 |
| 5 | 2-Chlorocyclohexanol | 134 | C6H11ClO | 001561-86-0 | 52 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071323\
Data File : BP016208.D
Acq On : 13 Jul 2023 14:57
Operator : MA/JU
Sample : 03385-01ME 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EX7X7ME

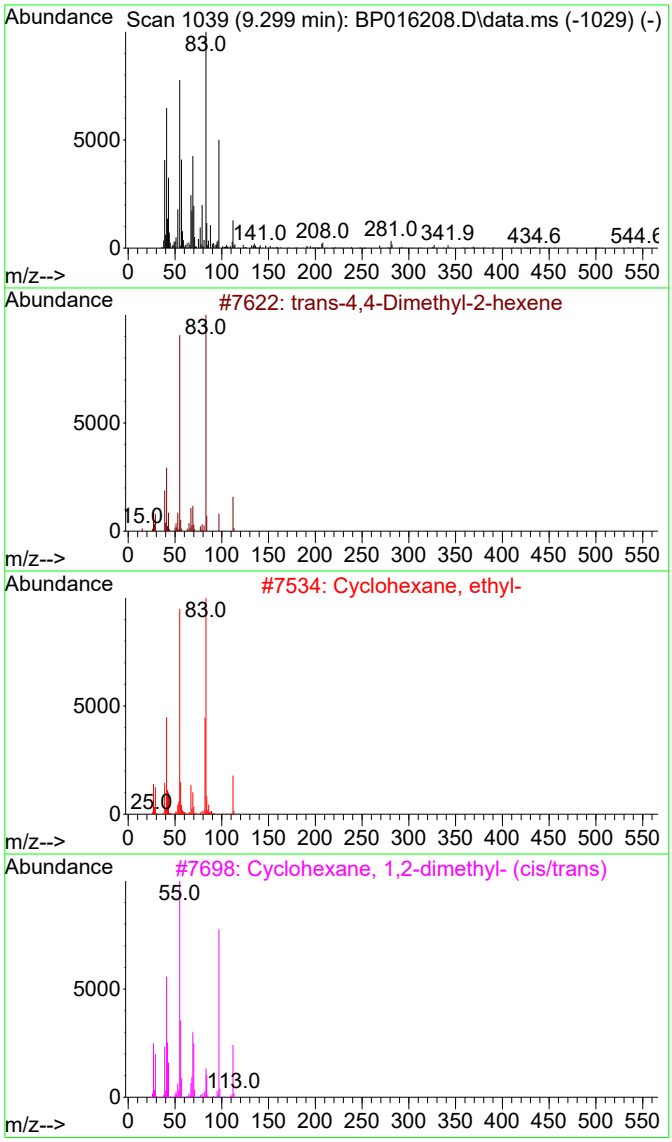
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

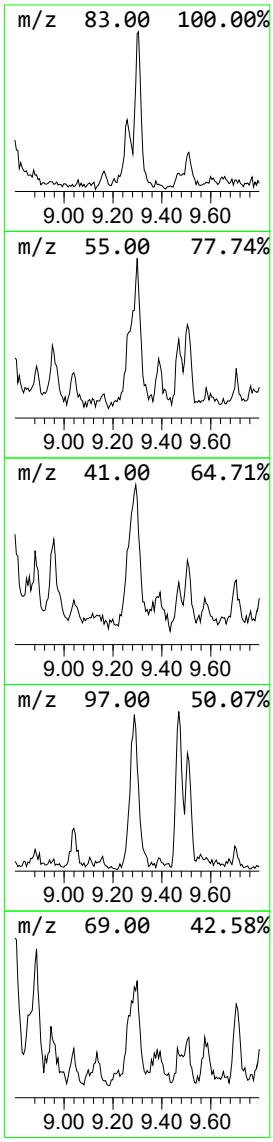
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 28

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 9.299 | 6.49 ng/ul | 212635 | 1,4-Dichlorobenzene-d4 | 7.975 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | trans-4,4-Dimethyl-2-hexene | 112 | C8H16 | 019550-83-5 | 43 |
| 2 | Cyclohexane, ethyl- | 112 | C8H16 | 001678-91-7 | 43 |
| 3 | Cyclohexane, 1,2-dimethyl- (cis/... | 112 | C8H16 | 000583-57-3 | 38 |
| 4 | Cyclohexanone, 3,3,5-trimethyl- | 140 | C9H16O | 000873-94-9 | 38 |
| 5 | 2H-Pyran-2-carboxaldehyde, 3,4-d... | 112 | C6H8O2 | 000100-73-2 | 38 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071323\
Data File : BP016208.D
Acq On : 13 Jul 2023 14:57
Operator : MA/JU
Sample : 03385-01ME 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EX7X7ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

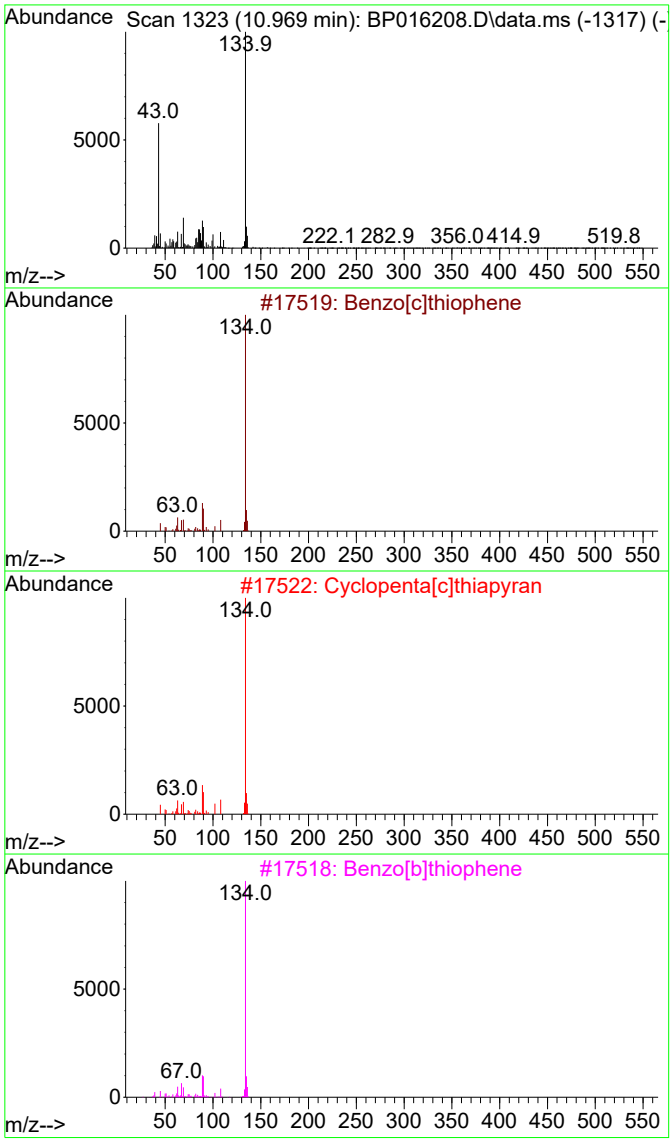
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

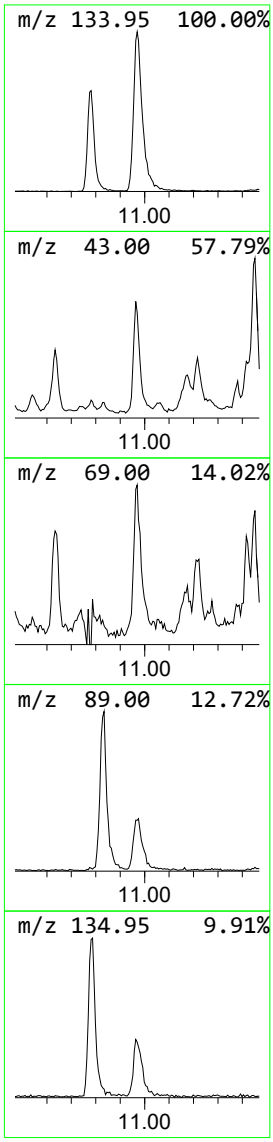
**********

Peak Number 20 Benzo[c]thiophene Concentration Rank 24

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 10.969 | 6.92 ng/ul | 375912 | Naphthalene-d8 | 10.781 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzo[c]thiophene | 134 | C8H6S | 000270-82-6 | 87 |
| 2 | Cyclopenta[c]thiapyran | 134 | C8H6S | 000270-63-3 | 87 |
| 3 | Benzo[b]thiophene | 134 | C8H6S | 000095-15-8 | 87 |
| 4 | Benzo[b]thiophene | 134 | C8H6S | 000095-15-8 | 87 |
| 5 | Benzo[b]thiophene | 134 | C8H6S | 000095-15-8 | 87 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071323\
Data File : BP016208.D
Acq On : 13 Jul 2023 14:57
Operator : MA/JU
Sample : 03385-01ME 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EX7X7ME

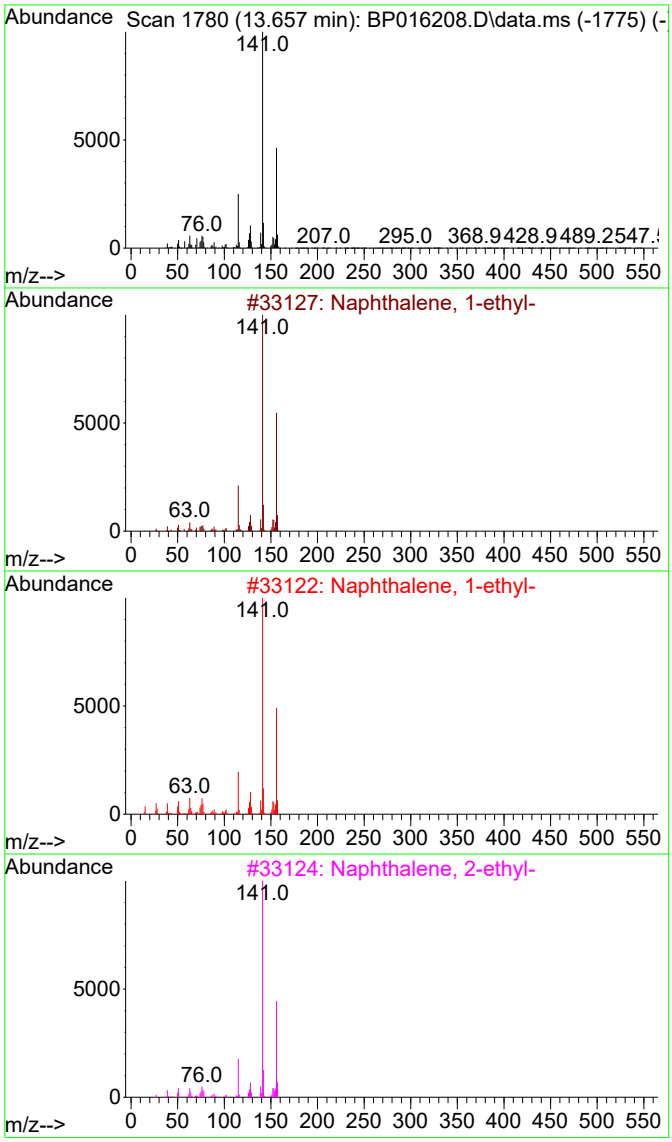
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

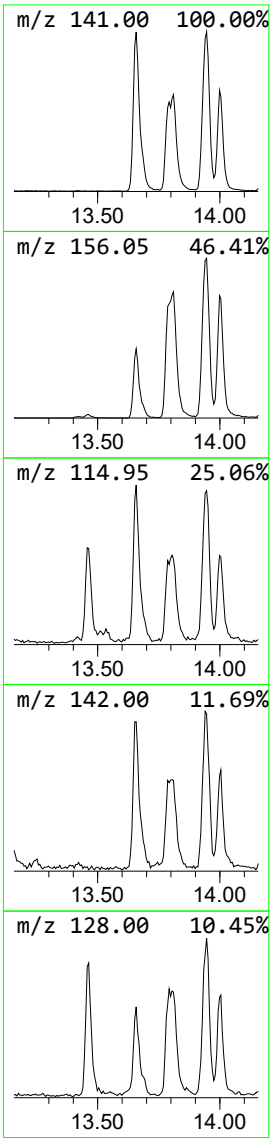
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 22 Naphthalene, 1-ethyl- Concentration Rank 36

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.657 | 4.36 ng/ul | 452674 | Acenaphthene-d10 | 14.616 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Naphthalene, 1-ethyl- | 156 | C12H12 | 001127-76-0 | 96 |
| 2 | Naphthalene, 1-ethyl- | 156 | C12H12 | 001127-76-0 | 95 |
| 3 | Naphthalene, 2-ethyl- | 156 | C12H12 | 000939-27-5 | 94 |
| 4 | Naphthalene, 2-ethyl- | 156 | C12H12 | 000939-27-5 | 94 |
| 5 | Naphthalene, 2-ethyl- | 156 | C12H12 | 000939-27-5 | 94 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071323\
Data File : BP016208.D
Acq On : 13 Jul 2023 14:57
Operator : MA/JU
Sample : 03385-01ME 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EX7X7ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

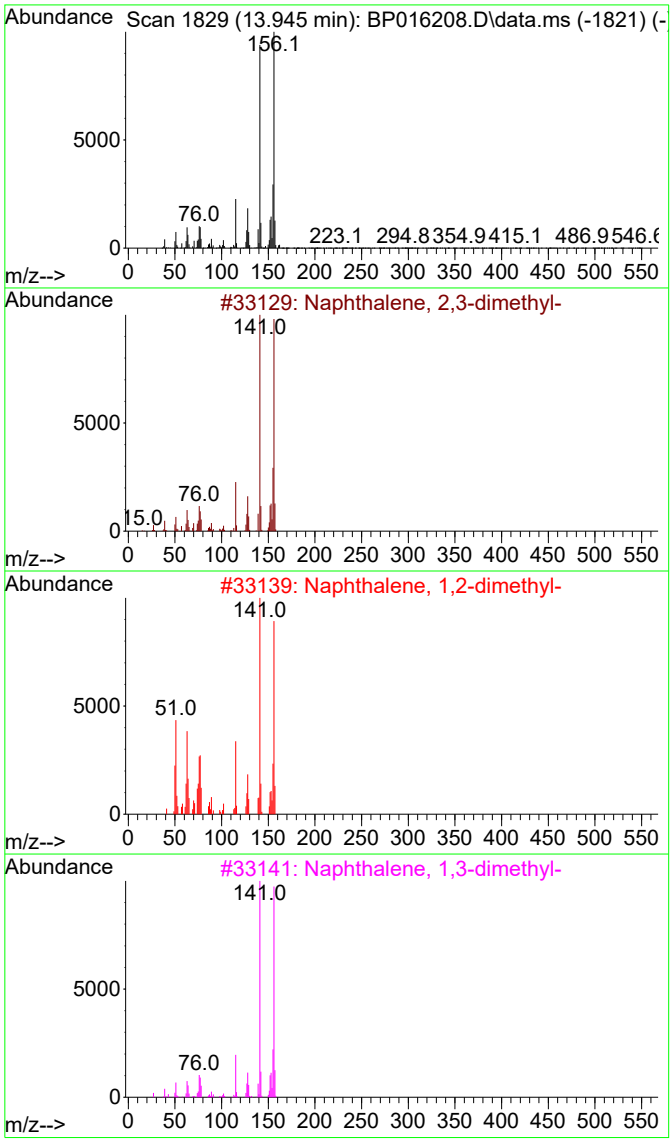
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

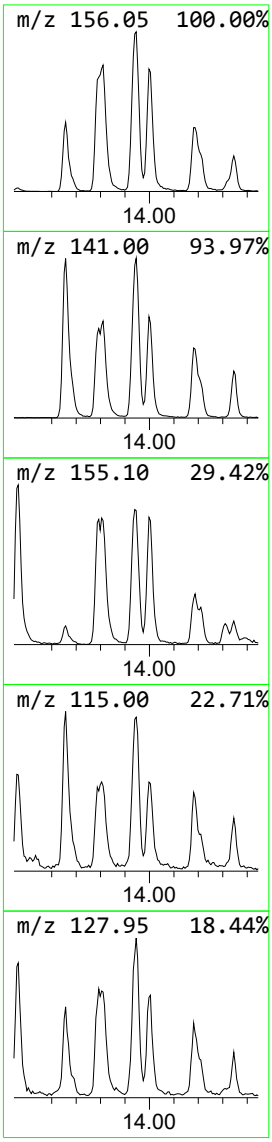
**********

Peak Number 23 Naphthalene, 2,3-dimethyl- Concentration Rank 25

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 13.945 | 6.88 ng/ul | 715193 | Acenaphthene-d10 | 14.616 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Naphthalene, 2,3-dimethyl- | 156 | C12H12 | 000581-40-8 | 98 |
| 2 | Naphthalene, 1,2-dimethyl- | 156 | C12H12 | 000573-98-8 | 97 |
| 3 | Naphthalene, 1,3-dimethyl- | 156 | C12H12 | 000575-41-7 | 97 |
| 4 | Naphthalene, 1,7-dimethyl- | 156 | C12H12 | 000575-37-1 | 96 |
| 5 | Naphthalene, 1,7-dimethyl- | 156 | C12H12 | 000575-37-1 | 96 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071323\
Data File : BP016208.D
Acq On : 13 Jul 2023 14:57
Operator : MA/JU
Sample : 03385-01ME 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EX7X7ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

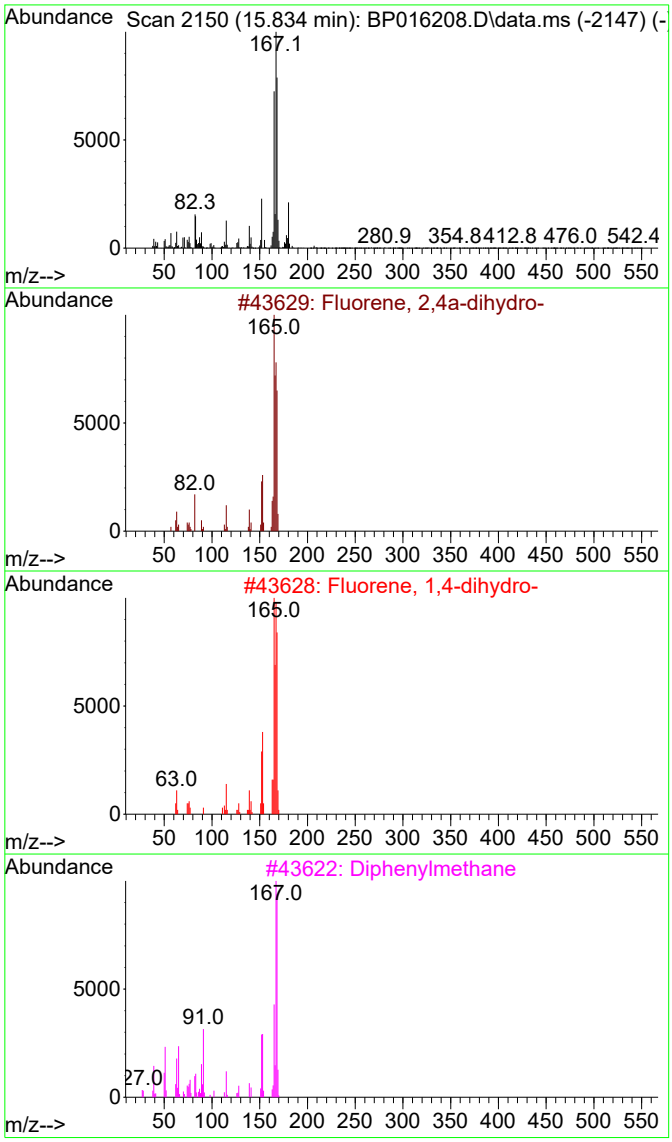
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

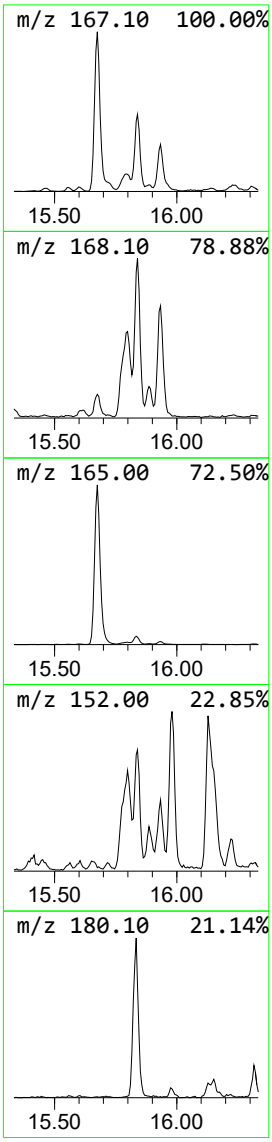
**********

Peak Number 28 Fluorene, 2,4a-dihydro- Concentration Rank 30

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.834 | 6.22 ng/ul | 646729 | Acenaphthene-d10 | 14.616 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Fluorene, 2,4a-dihydro- | 168 | C13H12 | 059247-36-8 | 96 |
| 2 | Fluorene, 1,4-dihydro- | 168 | C13H12 | 041593-21-9 | 93 |
| 3 | Diphenylmethane | 168 | C13H12 | 000101-81-5 | 89 |
| 4 | Benzeneacetamide, .alpha.-phenyl- | 211 | C14H13NO | 004695-13-0 | 80 |
| 5 | Nordiphenamid | 225 | C15H15NO | 000954-21-2 | 80 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071323\
Data File : BP016208.D
Acq On : 13 Jul 2023 14:57
Operator : MA/JU
Sample : 03385-01ME 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EX7X7ME

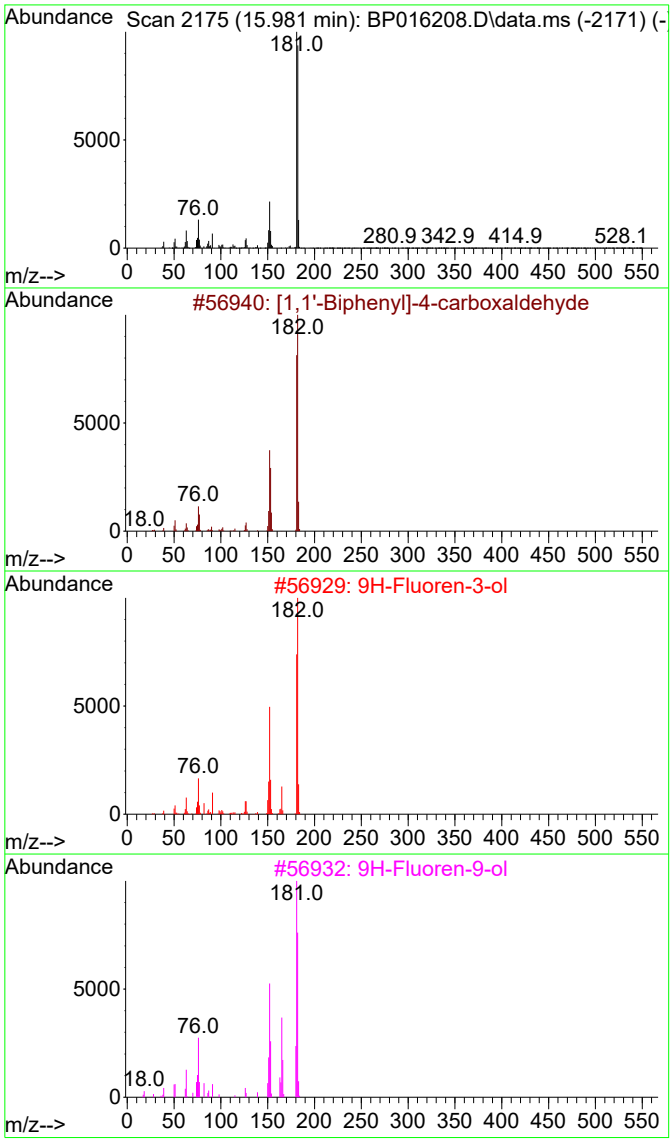
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

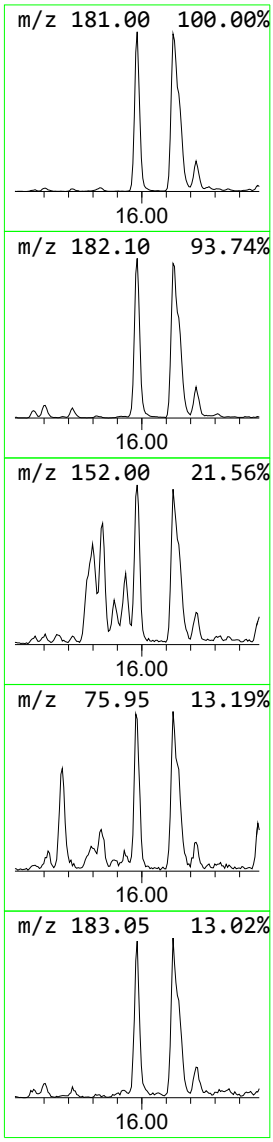
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 30 [1,1'-Biphenyl]-4-carboxald... Concentration Rank 26

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.981 | 6.65 ng/ul | 691011 | Acenaphthene-d10 | 14.616 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | [1,1'-Biphenyl]-4-carboxaldehyde | 182 | C13H10O | 003218-36-8 | 87 |
| 2 | 9H-Fluoren-3-ol | 182 | C13H10O | 006344-67-8 | 81 |
| 3 | 9H-Fluoren-9-ol | 182 | C13H10O | 001689-64-1 | 78 |
| 4 | [1,1'-Biphenyl]-4-carboxaldehyde | 182 | C13H10O | 003218-36-8 | 78 |
| 5 | 9H-Fluoren-9-ol | 182 | C13H10O | 001689-64-1 | 78 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071323\
Data File : BP016208.D
Acq On : 13 Jul 2023 14:57
Operator : MA/JU
Sample : 03385-01ME 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EX7X7ME

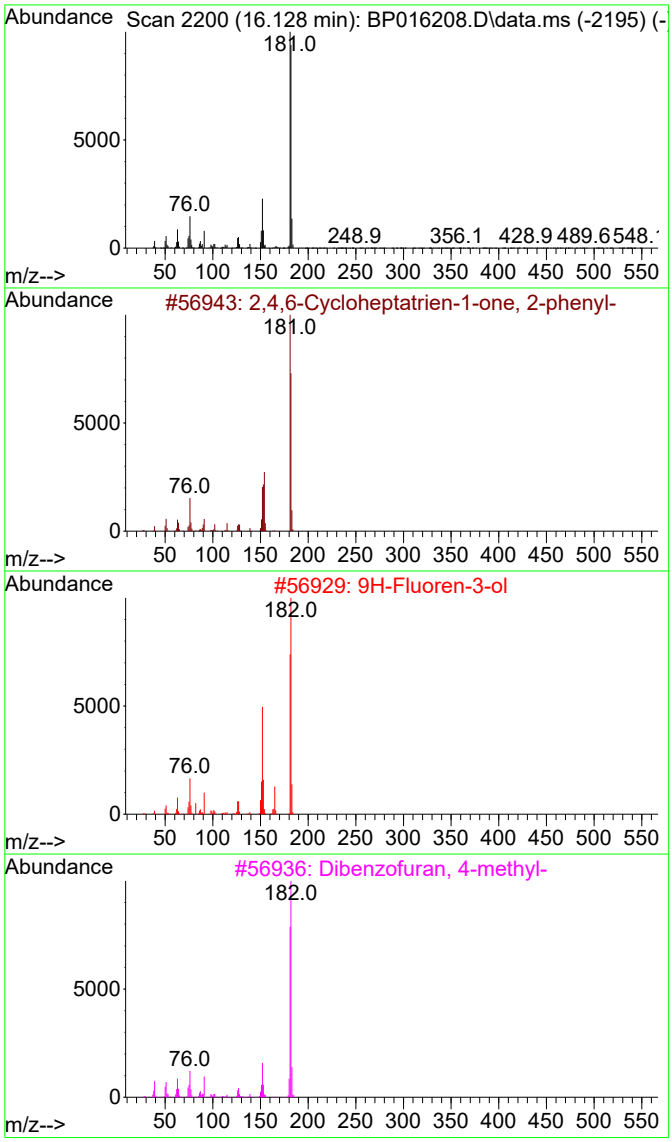
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

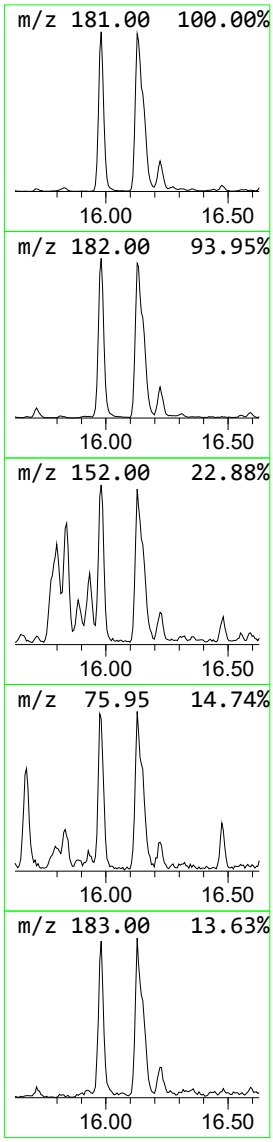
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 31 2,4,6-Cycloheptatrien-1-one... Concentration Rank 21

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 16.128 | 8.75 ng/ul | 1090610 | Phenanthrene-d10 | 17.381 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2,4,6-Cycloheptatrien-1-one, 2-p... | 182 | C13H10O | 014562-09-5 | 91 |
| 2 | 9H-Fluoren-3-ol | 182 | C13H10O | 006344-67-8 | 91 |
| 3 | Dibenzofuran, 4-methyl- | 182 | C13H10O | 007320-53-8 | 89 |
| 4 | [1,1'-Biphenyl]-4-carboxaldehyde | 182 | C13H10O | 003218-36-8 | 87 |
| 5 | 9H-Fluoren-9-ol | 182 | C13H10O | 001689-64-1 | 78 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071323\
Data File : BP016208.D
Acq On : 13 Jul 2023 14:57
Operator : MA/JU
Sample : 03385-01ME 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EX7X7ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

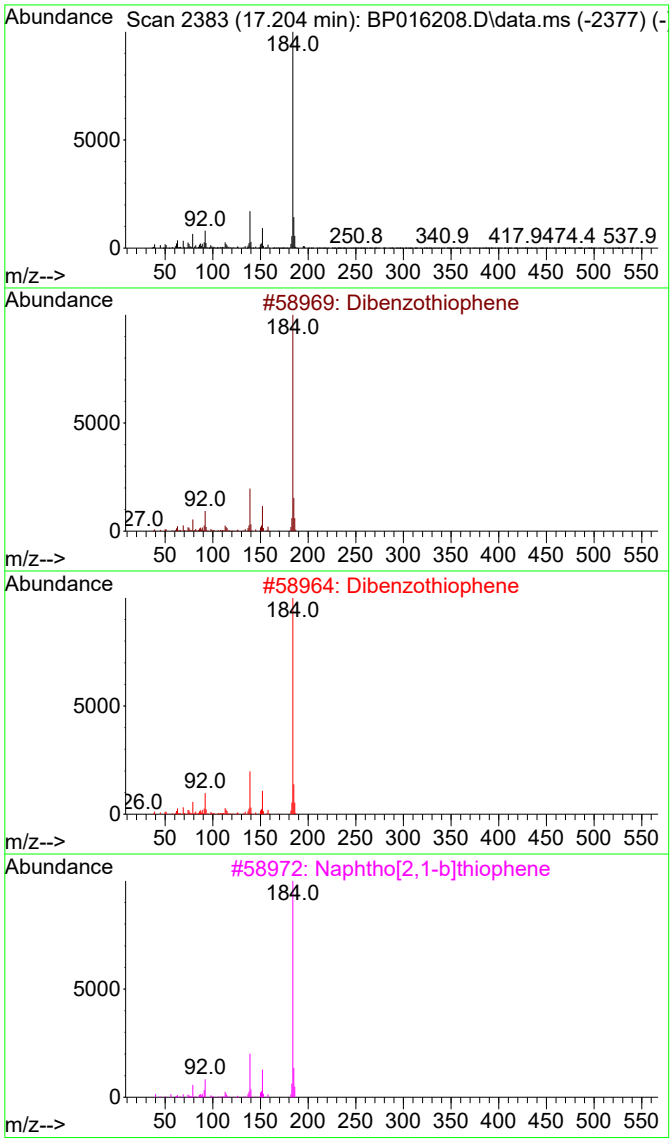
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

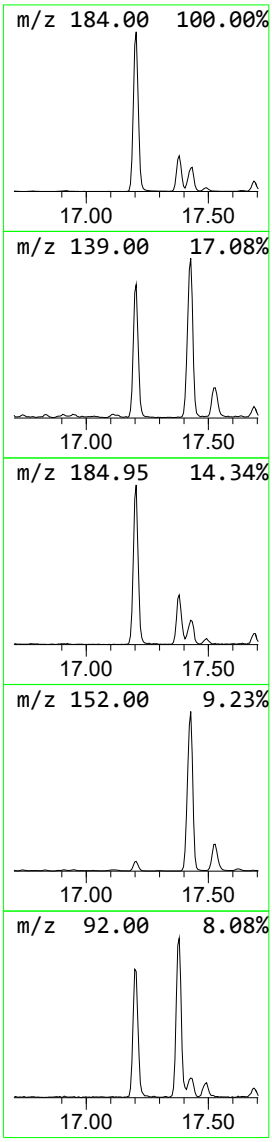
**********

Peak Number 35 Dibenzothiophene Concentration Rank 13

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 17.204 | 12.08 ng/ul | 1505240 | Phenanthrene-d10 | 17.381 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Dibenzothiophene | 184 | C12H8S | 000132-65-0 | 97 |
| 2 | Dibenzothiophene | 184 | C12H8S | 000132-65-0 | 97 |
| 3 | Naphtho[2,1-b]thiophene | 184 | C12H8S | 000233-02-3 | 97 |
| 4 | Dibenzothiophene | 184 | C12H8S | 000132-65-0 | 96 |
| 5 | Dibenzothiophene | 184 | C12H8S | 000132-65-0 | 95 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071323\
Data File : BP016208.D
Acq On : 13 Jul 2023 14:57
Operator : MA/JU
Sample : 03385-01ME 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EX7X7ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

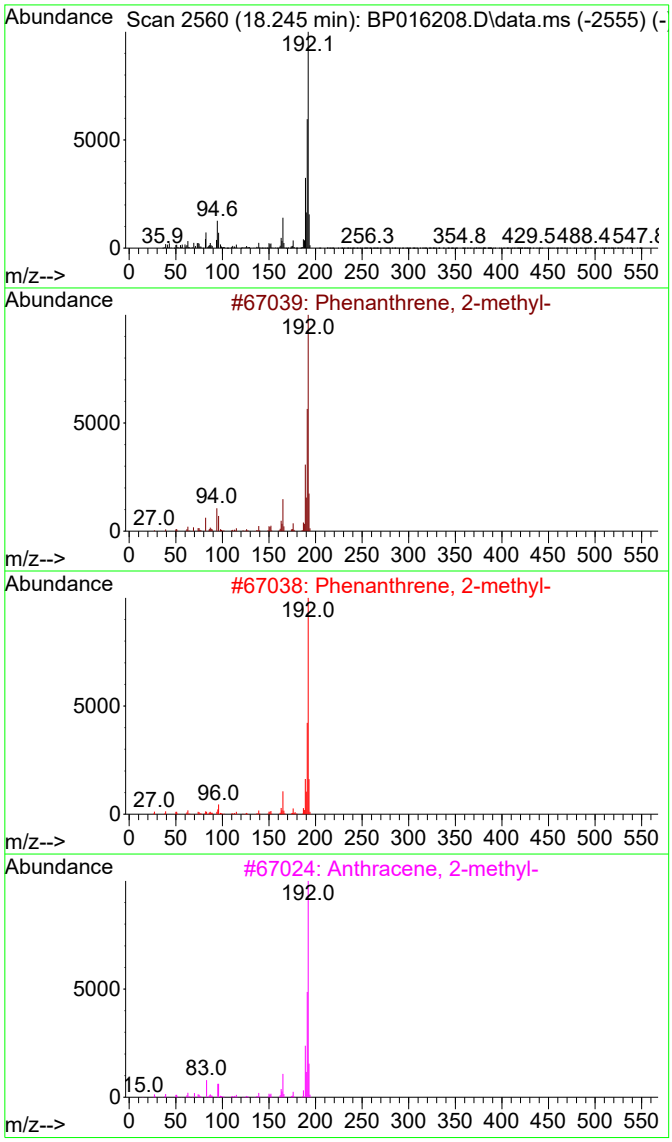
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

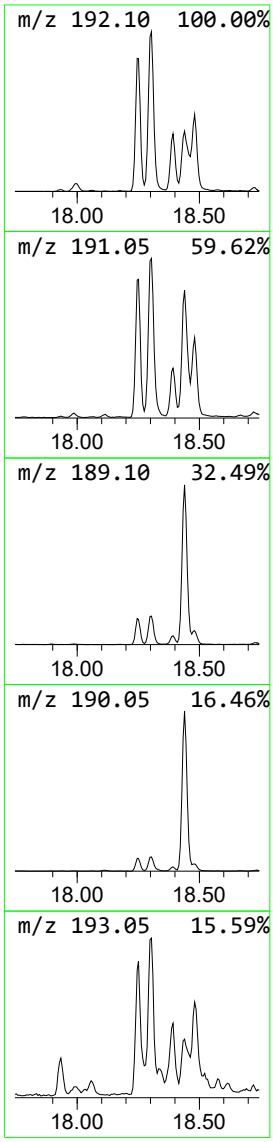
**********

Peak Number 37 Phenanthrene, 2-methyl- Concentration Rank 18

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.245 | 10.51 ng/ul | 1310230 | Phenanthrene-d10 | 17.381 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenanthrene, 2-methyl- | 192 | C15H12 | 002531-84-2 | 97 |
| 2 | Phenanthrene, 2-methyl- | 192 | C15H12 | 002531-84-2 | 96 |
| 3 | Anthracene, 2-methyl- | 192 | C15H12 | 000613-12-7 | 96 |
| 4 | Phenanthrene, 1-methyl- | 192 | C15H12 | 000832-69-9 | 96 |
| 5 | Anthracene, 1-methyl- | 192 | C15H12 | 000610-48-0 | 96 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071323\
Data File : BP016208.D
Acq On : 13 Jul 2023 14:57
Operator : MA/JU
Sample : 03385-01ME 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EX7X7ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

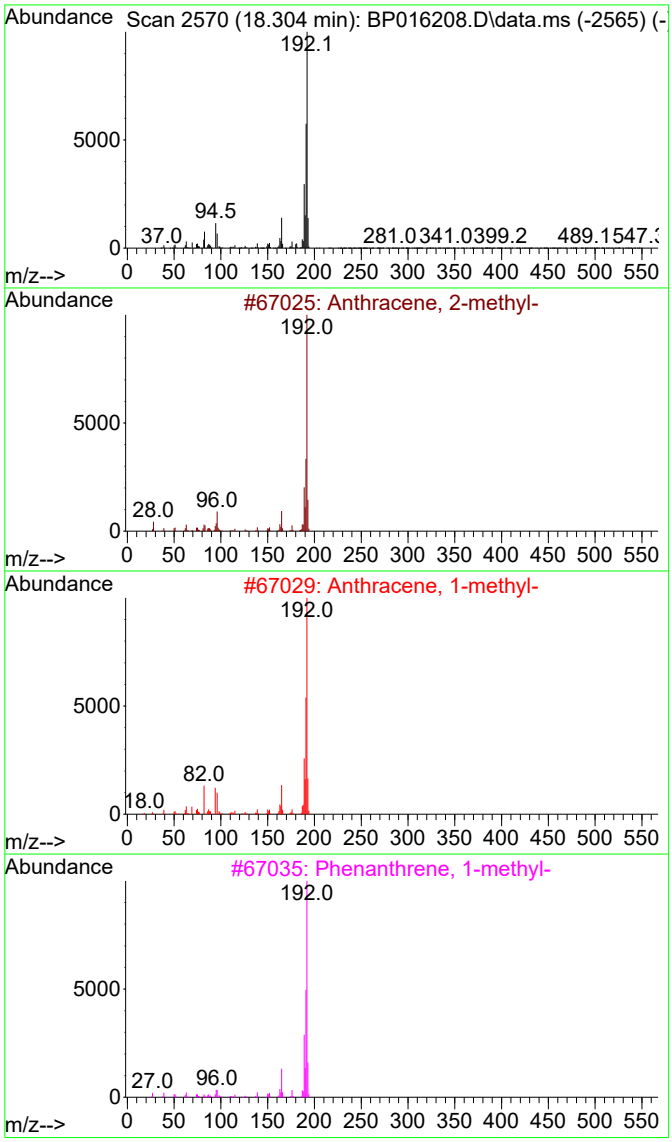
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

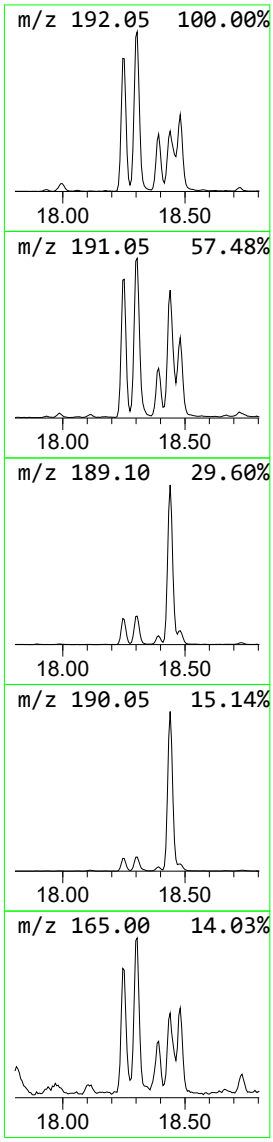
**********

Peak Number 38 Anthracene, 2-methyl- Concentration Rank 11

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.304 | 13.79 ng/ul | 1718100 | Phenanthrene-d10 | 17.381 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Anthracene, 2-methyl- | 192 | C15H12 | 000613-12-7 | 95 |
| 2 | Anthracene, 1-methyl- | 192 | C15H12 | 000610-48-0 | 93 |
| 3 | Phenanthrene, 1-methyl- | 192 | C15H12 | 000832-69-9 | 93 |
| 4 | Phenanthrene, 2-methyl- | 192 | C15H12 | 002531-84-2 | 93 |
| 5 | Anthracene, 9-methyl- | 192 | C15H12 | 000779-02-2 | 91 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071323\
Data File : BP016208.D
Acq On : 13 Jul 2023 14:57
Operator : MA/JU
Sample : 03385-01ME 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EX7X7ME

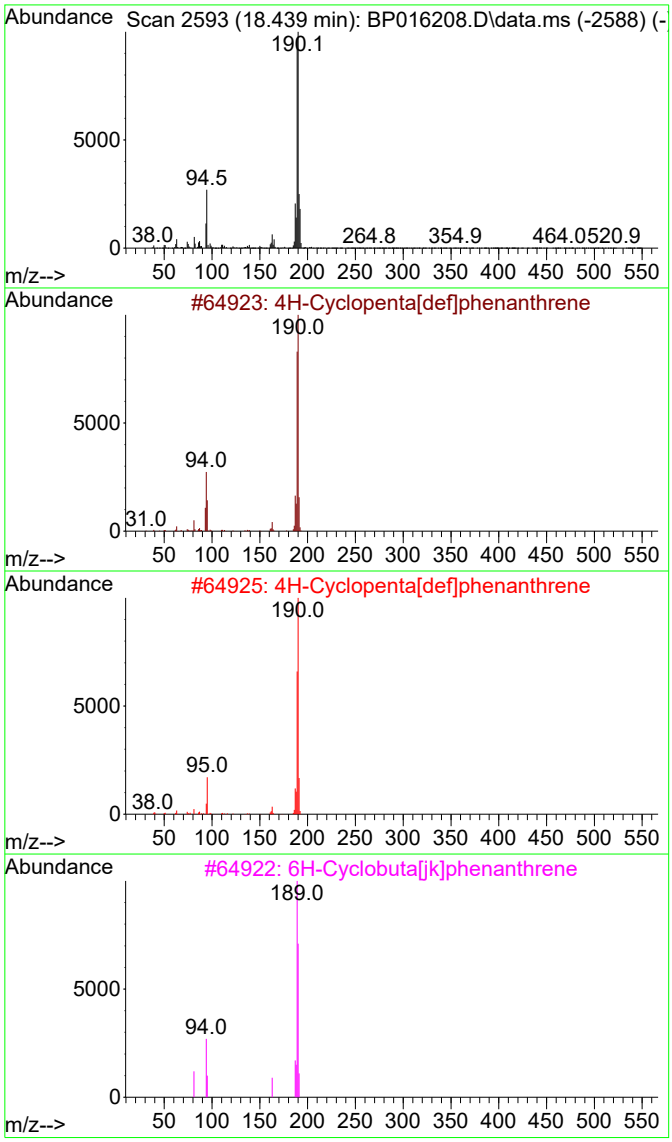
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

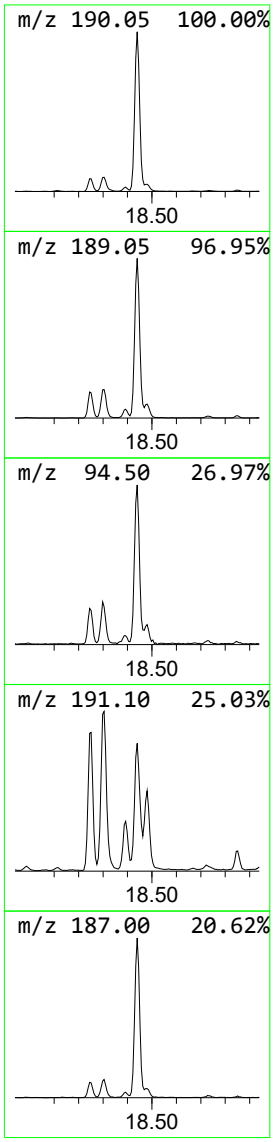
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 40 4H-Cyclopenta[def]phenanthrene Concentration Rank 8

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.439 | 24.48 ng/ul | 3050380 | Phenanthrene-d10 | 17.381 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 4H-Cyclopenta[def]phenanthrene | 190 | C15H10 | 000203-64-5 | 93 |
| 2 | 4H-Cyclopenta[def]phenanthrene | 190 | C15H10 | 000203-64-5 | 68 |
| 3 | 6H-Cyclobuta[jk]phenanthrene | 190 | C15H10 | 083469-43-6 | 64 |
| 4 | 4H-Cyclopenta[def]phenanthrene | 190 | C15H10 | 000203-64-5 | 55 |
| 5 | Methyl diselenide | 190 | C2H6Se2 | 007101-31-7 | 41 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071323\
Data File : BP016208.D
Acq On : 13 Jul 2023 14:57
Operator : MA/JU
Sample : 03385-01ME 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EX7X7ME

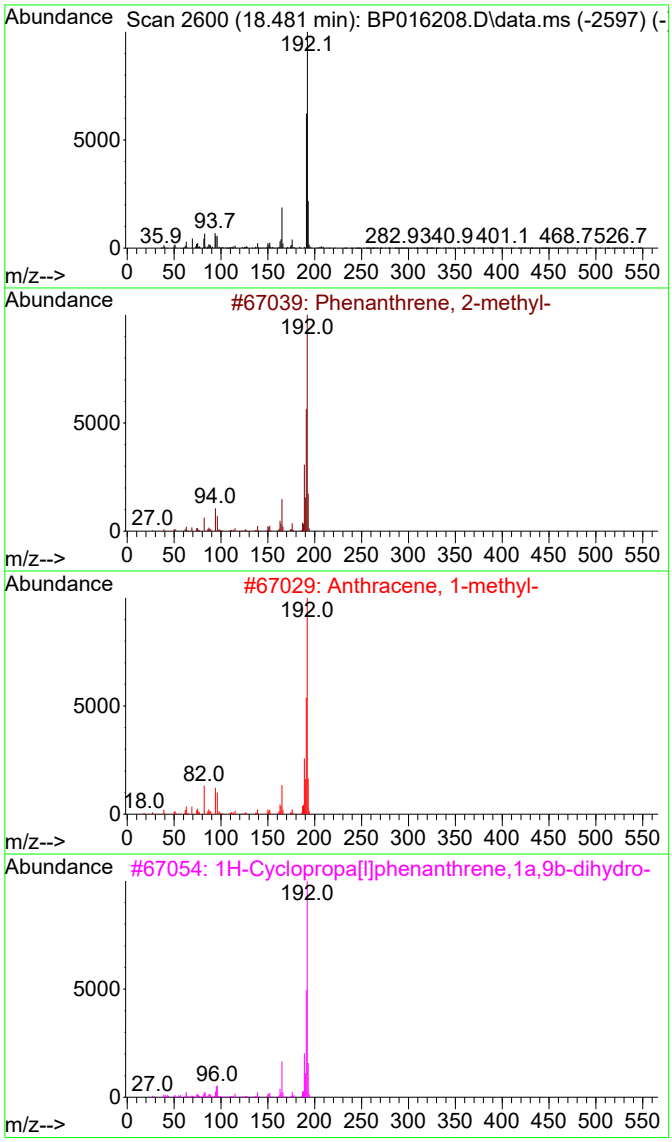
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

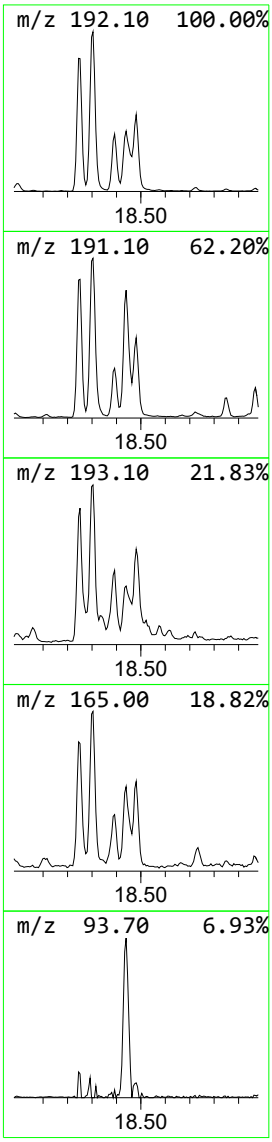
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 41 1H-Cyclopropa[l]phenanthren... Concentration Rank 31

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.480 | 6.04 ng/ul | 752319 | Phenanthrene-d10 | 17.381 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Phenanthrene, 2-methyl- | 192 | C15H12 | 002531-84-2 | 90 |
| 2 | Anthracene, 1-methyl- | 192 | C15H12 | 000610-48-0 | 90 |
| 3 | 1H-Cyclopropa[l]phenanthrene,1a,... | 192 | C15H12 | 000949-41-7 | 90 |
| 4 | Phenanthrene, 1-methyl- | 192 | C15H12 | 000832-69-9 | 90 |
| 5 | 1H-Indene, 2-phenyl- | 192 | C15H12 | 004505-48-0 | 87 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071323\
Data File : BP016208.D
Acq On : 13 Jul 2023 14:57
Operator : MA/JU
Sample : 03385-01ME 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EX7X7ME

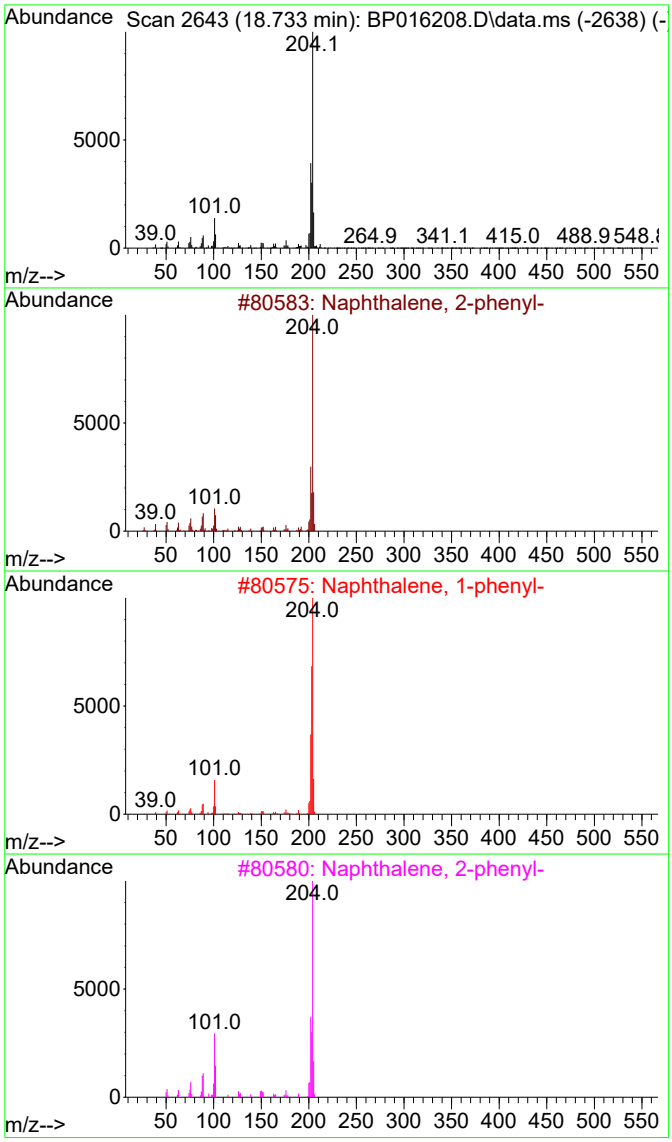
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

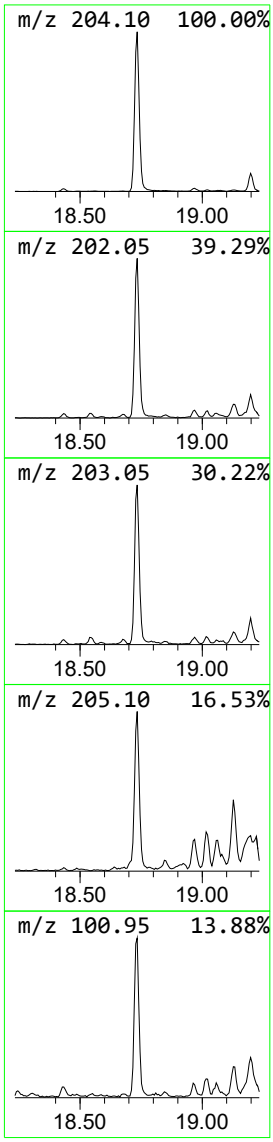
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 42 Naphthalene, 2-phenyl- Concentration Rank 22

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.733 | 8.64 ng/ul | 1076280 | Phenanthrene-d10 | 17.381 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Naphthalene, 2-phenyl- | 204 | C16H12 | 000612-94-2 | 83 |
| 2 | Naphthalene, 1-phenyl- | 204 | C16H12 | 000605-02-7 | 78 |
| 3 | Naphthalene, 2-phenyl- | 204 | C16H12 | 000612-94-2 | 64 |
| 4 | Tricyclo[8.2.2.2(4,7)]hexadeca-2... | 204 | C16H12 | 006572-60-7 | 62 |
| 5 | Naphthalene, 2-phenyl- | 204 | C16H12 | 000612-94-2 | 60 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071323\
Data File : BP016208.D
Acq On : 13 Jul 2023 14:57
Operator : MA/JU
Sample : 03385-01ME 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EX7X7ME

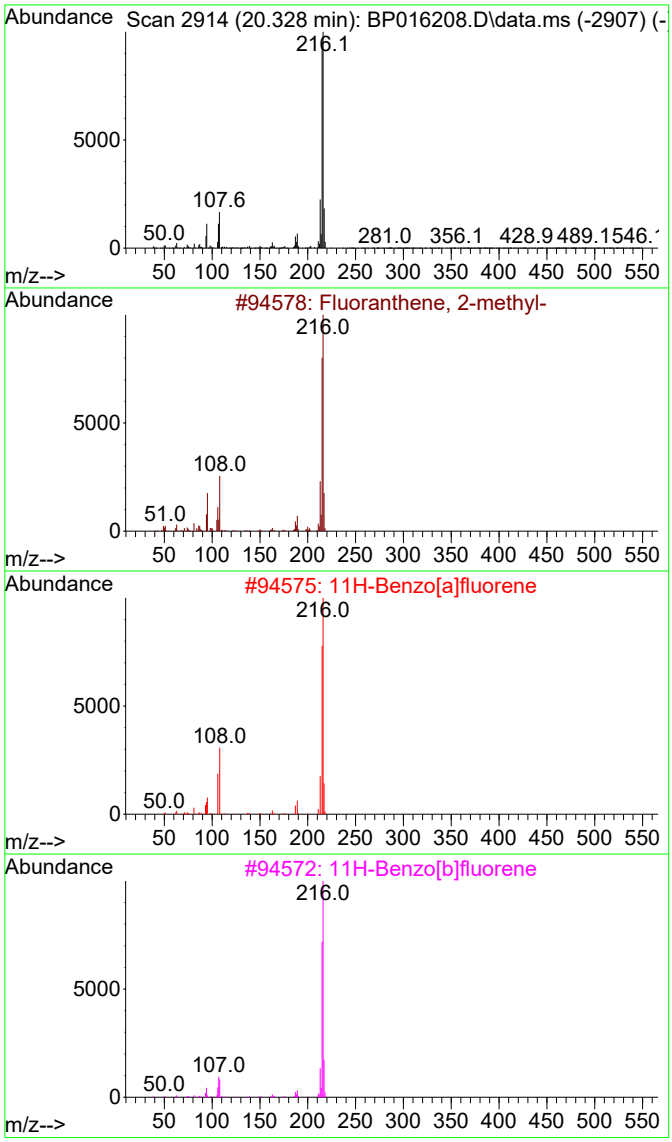
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION

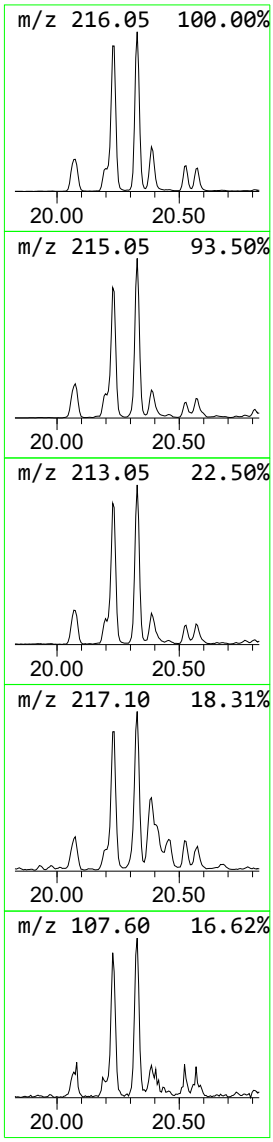
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 46 Fluoranthene, 2-methyl- Concentration Rank 33

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 20.328 | 4.62 ng/ul | 2526380 | Chrysene-d12 | 21.474 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Fluoranthene, 2-methyl- | 216 | C17H12 | 033543-31-6 | 94 |
| 2 | 11H-Benzo[a]fluorene | 216 | C17H12 | 000238-84-6 | 91 |
| 3 | 11H-Benzo[b]fluorene | 216 | C17H12 | 000243-17-4 | 90 |
| 4 | 11H-Benzo[a]fluorene | 216 | C17H12 | 000238-84-6 | 81 |
| 5 | Fluoranthene, 2-methyl- | 216 | C17H12 | 033543-31-6 | 76 |



Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071323\
Data File : BP016208.D
Acq On : 13 Jul 2023 14:57
Operator : MA/JU
Sample : 03385-01ME 5X
Misc :
ALS Vial : 10 Sample Multiplier: 1

**Instrument :**
BNA_P
**ClientSampleId :**
EX7X7ME

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.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 |
| 2-Butenal, 2-me... | 3.722 | 12.4 | ng/ul | 405226 | 1 | 7.975 | 654863 | 20.0 |
| 2H-Pyran, tetra... | 3.834 | 11.2 | ng/ul | 367934 | 1 | 7.975 | 654863 | 20.0 |
| Prenol | 4.058 | 32.6 | ng/ul | 1066890 | 1 | 7.975 | 654863 | 20.0 |
| 2-Butenal, 3-me... | 4.240 | 14.3 | ng/ul | 466858 | 1 | 7.975 | 654863 | 20.0 |
| 2-Butene, 1-bro... | 4.293 | 45.3 | ng/ul | 1482550 | 1 | 7.975 | 654863 | 20.0 |
| unknown-01 | 4.446 | 8.3 | ng/ul | 271853 | 1 | 7.975 | 654863 | 20.0 |
| unknown-02 | 4.587 | 6.5 | ng/ul | 212337 | 1 | 7.975 | 654863 | 20.0 |
| Formic acid, 1-... | 4.628 | 55.2 | ng/ul | 1807340 | 1 | 7.975 | 654863 | 20.0 |
| Cyclopentane, b... | 4.717 | 34.5 | ng/ul | 1128210 | 1 | 7.975 | 654863 | 20.0 |
| Guanidine, mono... | 5.128 | 10.0 | ng/ul | 326487 | 1 | 7.975 | 654863 | 20.0 |
| 2-Cyclohexen-1-ol | 5.834 | 24.5 | ng/ul | 802788 | 1 | 7.975 | 654863 | 20.0 |
| 2-Buten-1-ol, 3... | 6.246 | 9.1 | ng/ul | 298630 | 1 | 7.975 | 654863 | 20.0 |
| Ethane, 1,1,2,2... | 6.340 | 10.7 | ng/ul | 351674 | 1 | 7.975 | 654863 | 20.0 |
| 2-Cyclohexen-1-one | 6.587 | 39.4 | ng/ul | 1291110 | 1 | 7.975 | 654863 | 20.0 |
| unknown-03 | 7.705 | 6.7 | ng/ul | 217595 | 1 | 7.975 | 654863 | 20.0 |
| unknown-04 | 8.187 | 11.5 | ng/ul | 375600 | 1 | 7.975 | 654863 | 20.0 |
| 2-Chlorocyclohe... | 8.275 | 67.6 | ng/ul | 2212300 | 1 | 7.975 | 654863 | 20.0 |
| (DEL) Alkane: S... | 9.299 | 6.5 | ng/ul | 212635 | 1 | 7.975 | 654863 | 20.0 |
| Benzo[c]thiophene | 10.969 | 6.9 | ng/ul | 375912 | 2 | 10.781 | 1086190 | 20.0 |
| Naphthalene, 1-... | 13.657 | 4.4 | ng/ul | 452674 | 3 | 14.616 | 2078540 | 20.0 |
| Naphthalene, 2,... | 13.945 | 6.9 | ng/ul | 715193 | 3 | 14.616 | 2078540 | 20.0 |
| Fluorene, 2,4a-... | 15.834 | 6.2 | ng/ul | 646729 | 3 | 14.616 | 2078540 | 20.0 |
| [1,1'-Biphenyl]... | 15.981 | 6.7 | ng/ul | 691011 | 3 | 14.616 | 2078540 | 20.0 |
| 2,4,6-Cycloheptt... | 16.128 | 8.8 | ng/ul | 1090610 | 4 | 17.381 | 2492440 | 20.0 |
| Dibenzothiophene | 17.204 | 12.1 | ng/ul | 1505240 | 4 | 17.381 | 2492440 | 20.0 |
| Phenanthrene, 2... | 18.245 | 10.5 | ng/ul | 1310230 | 4 | 17.381 | 2492440 | 20.0 |
| Anthracene, 2-m... | 18.304 | 13.8 | ng/ul | 1718100 | 4 | 17.381 | 2492440 | 20.0 |
| 4H-Cyclopenta[d... | 18.439 | 24.5 | ng/ul | 3050380 | 4 | 17.381 | 2492440 | 20.0 |
| 1H-Cyclopropa[l... | 18.480 | 6.0 | ng/ul | 752319 | 4 | 17.381 | 2492440 | 20.0 |
| Naphthalene, 2-... | 18.733 | 8.6 | ng/ul | 1076280 | 4 | 17.381 | 2492440 | 20.0 |
| Fluoranthene, 2... | 20.328 | 4.6 | ng/ul | 2526380 | 5 | 21.474 | 10933100 | 20.0 |