LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092325\
Data File : BG064437.D
Acq On : 23 Sep 2025 13:18
Operator : RC/JU
Sample : Q3155-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
COMP-1

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082825.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BG064437.D\data.ms

| peak # | R.T. min | first scan | max scan | last scan | PK TY | peak height | corr. area | corr. % max. | % of total |
|---|---|---|---|---|---|---|---|---|---|
| 1 | 5.432 | 392 | 399 | 406 | rBV | 241866 | 388046 | 32.10% | 4.121% |
| 2 | 7.013 | 658 | 668 | 678 | rBV | 272381 | 463141 | 38.31% | 4.918% |
| 3 | 7.836 | 802 | 808 | 817 | rBV2 | 80328 | 134673 | 11.14% | 1.430% |
| 4 | 8.987 | 998 | 1004 | 1014 | rBV | 168376 | 293392 | 24.27% | 3.115% |
| 5 | 10.620 | 1276 | 1282 | 1290 | rBV | 102940 | 185866 | 15.37% | 1.974% |
| 6 | 13.082 | 1694 | 1701 | 1709 | rBV | 434606 | 651841 | 53.91% | 6.922% |
| 7 | 14.463 | 1929 | 1936 | 1942 | rBV | 182160 | 267612 | 22.13% | 2.842% |
| 8 | 15.820 | 2161 | 2167 | 2173 | rVB | 62744 | 96029 | 7.94% | 1.020% |
| 9 | 15.950 | 2182 | 2189 | 2199 | rBV | 440269 | 663300 | 54.86% | 7.043% |
| 10 | 17.107 | 2382 | 2386 | 2393 | rBV3 | 22956 | 33900 | 2.80% | 0.360% |
| 11 | 17.201 | 2393 | 2402 | 2406 | rVV | 246381 | 379895 | 31.42% | 4.034% |
| 12 | 17.242 | 2406 | 2409 | 2416 | rVV | 99014 | 138412 | 11.45% | 1.470% |
| 13 | 17.336 | 2420 | 2425 | 2430 | rVB | 15391 | 22899 | 1.89% | 0.243% |
| 14 | 17.606 | 2463 | 2471 | 2474 | rBV4 | 9654 | 17149 | 1.42% | 0.182% |
| 15 | 17.977 | 2528 | 2534 | 2541 | rBV5 | 20288 | 44663 | 3.69% | 0.474% |
| 16 | 18.041 | 2541 | 2545 | 2548 | rVV4 | 12800 | 18314 | 1.51% | 0.194% |
| 17 | 18.106 | 2548 | 2556 | 2565 | rVB3 | 153666 | 255781 | 21.16% | 2.716% |
| 18 | 18.270 | 2579 | 2584 | 2588 | rBV3 | 16815 | 30389 | 2.51% | 0.323% |
| 19 | 18.400 | 2602 | 2606 | 2611 | rBV5 | 21242 | 34969 | 2.89% | 0.371% |
| 20 | 18.611 | 2639 | 2642 | 2647 | rVB3 | 13397 | 17558 | 1.45% | 0.186% |
| 21 | 19.128 | 2727 | 2730 | 2736 | rVB3 | 15818 | 25038 | 2.07% | 0.266% |
| 22 | 19.257 | 2744 | 2752 | 2761 | rBV | 227406 | 377103 | 31.19% | 4.004% |
| 23 | 19.375 | 2766 | 2772 | 2780 | rVV7 | 52081 | 92473 | 7.65% | 0.982% |
| 24 | 19.616 | 2803 | 2813 | 2821 | rBV | 210056 | 341550 | 28.25% | 3.627% |
| 25 | 19.680 | 2822 | 2824 | 2831 | rVB5 | 9637 | 16866 | 1.39% | 0.179% |
| 26 | 19.821 | 2840 | 2848 | 2853 | rBV | 949059 | 1209041 | 100.00% | 12.838% |
| 27 | 19.933 | 2863 | 2867 | 2870 | rBV5 | 13637 | 20723 | 1.71% | 0.220% |
| 28 | 20.109 | 2894 | 2897 | 2901 | rVB2 | 16529 | 22293 | 1.84% | 0.237% |
| 29 | 20.262 | 2918 | 2923 | 2927 | rVV5 | 17626 | 31784 | 2.63% | 0.338% |
| 30 | 20.297 | 2927 | 2929 | 2935 | rVB7 | 11715 | 17353 | 1.44% | 0.184% |
| 31 | 20.456 | 2952 | 2956 | 2960 | rVV5 | 14731 | 21927 | 1.81% | 0.233% |
| 32 | 20.856 | 3020 | 3024 | 3031 | rVB | 30292 | 39938 | 3.30% | 0.424% |
| 33 | 21.032 | 3048 | 3054 | 3057 | rVV3 | 16936 | 27120 | 2.24% | 0.288% |
| 34 | 21.073 | 3059 | 3061 | 3063 | rVV3 | 12372 | 15318 | 1.27% | 0.163% |
| 35 | 21.108 | 3063 | 3067 | 3074 | rVV4 | 75402 | 136194 | 11.26% | 1.446% |
| 36 | 21.173 | 3074 | 3078 | 3086 | rVB3 | 31181 | 48426 | 4.01% | 0.514% |

LSC Area Percent Report

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Acq On : 23 Sep 2025 13:18
Operator : RC/JU
Sample : Q3155-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
COMP-1

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082825.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

| | | | | | | | | | |
|---|---|---|---|---|---|---|---|---|---|
| 37 | 21.337 | 3102 | 3106 | 3110 | rBV | 63498 | 74338 | 6.15% | 0.789% |
| 38 | 21.425 | 3114 | 3121 | 3126 | rVV2 | 319438 | 610863 | 50.52% | 6.487% |
| 39 | 21.472 | 3126 | 3129 | 3138 | rVB | 85547 | 142313 | 11.77% | 1.511% |
| 40 | 21.625 | 3151 | 3155 | 3160 | rVB6 | 17485 | 28113 | 2.33% | 0.299% |
| 41 | 21.813 | 3181 | 3187 | 3193 | rBV9 | 15471 | 34350 | 2.84% | 0.365% |
| 42 | 22.119 | 3235 | 3239 | 3243 | rVB4 | 16564 | 27839 | 2.30% | 0.296% |
| 43 | 22.195 | 3249 | 3252 | 3259 | rVB7 | 19638 | 38331 | 3.17% | 0.407% |
| 44 | 22.377 | 3278 | 3283 | 3286 | rBV7 | 13411 | 25103 | 2.08% | 0.267% |
| 45 | 22.542 | 3309 | 3311 | 3315 | rVB5 | 14222 | 16477 | 1.36% | 0.175% |
| 46 | 22.642 | 3320 | 3328 | 3336 | rBV3 | 58787 | 133054 | 11.00% | 1.413% |
| 47 | 22.812 | 3353 | 3357 | 3367 | rVB3 | 28767 | 69963 | 5.79% | 0.743% |
| 48 | 23.018 | 3387 | 3392 | 3398 | rVB4 | 40066 | 70917 | 5.87% | 0.753% |
| 49 | 23.159 | 3410 | 3416 | 3423 | rVB5 | 52945 | 107315 | 8.88% | 1.140% |
| 50 | 23.488 | 3463 | 3472 | 3477 | rBV2 | 75087 | 234930 | 19.43% | 2.495% |
| 51 | 23.635 | 3492 | 3497 | 3507 | rBV4 | 59307 | 130496 | 10.79% | 1.386% |
| 52 | 24.146 | 3578 | 3584 | 3595 | rVB2 | 41631 | 108827 | 9.00% | 1.156% |
| 53 | 24.281 | 3599 | 3607 | 3620 | rVB2 | 59323 | 178069 | 14.73% | 1.891% |
| 54 | 24.428 | 3623 | 3632 | 3641 | rBV | 170500 | 441961 | 36.55% | 4.693% |
| 55 | 24.651 | 3668 | 3670 | 3679 | rVB10 | 7956 | 16367 | 1.35% | 0.174% |
| 56 | 24.927 | 3713 | 3717 | 3727 | rVB7 | 35884 | 87767 | 7.26% | 0.932% |
| 57 | 25.468 | 3807 | 3809 | 3810 | rBV2 | 15228 | 12712 | 1.05% | 0.135% |
| 58 | 25.603 | 3824 | 3832 | 3833 | rBV8 | 17734 | 38136 | 3.15% | 0.405% |
| 59 | 27.783 | 4195 | 4203 | 4204 | rBV7 | 24858 | 51288 | 4.24% | 0.545% |
| 60 | 28.817 | 4376 | 4379 | 4380 | rBV3 | 12269 | 13575 | 1.12% | 0.144% |
| 61 | 31.255 | 4785 | 4794 | 4795 | rBV9 | 16925 | 32498 | 2.69% | 0.345% |
| 62 | 31.766 | 4869 | 4881 | 4894 | rVB9 | 23305 | 110732 | 9.16% | 1.176% |

Sum of corrected areas: 9417310

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092325\
Data File : BG064437.D
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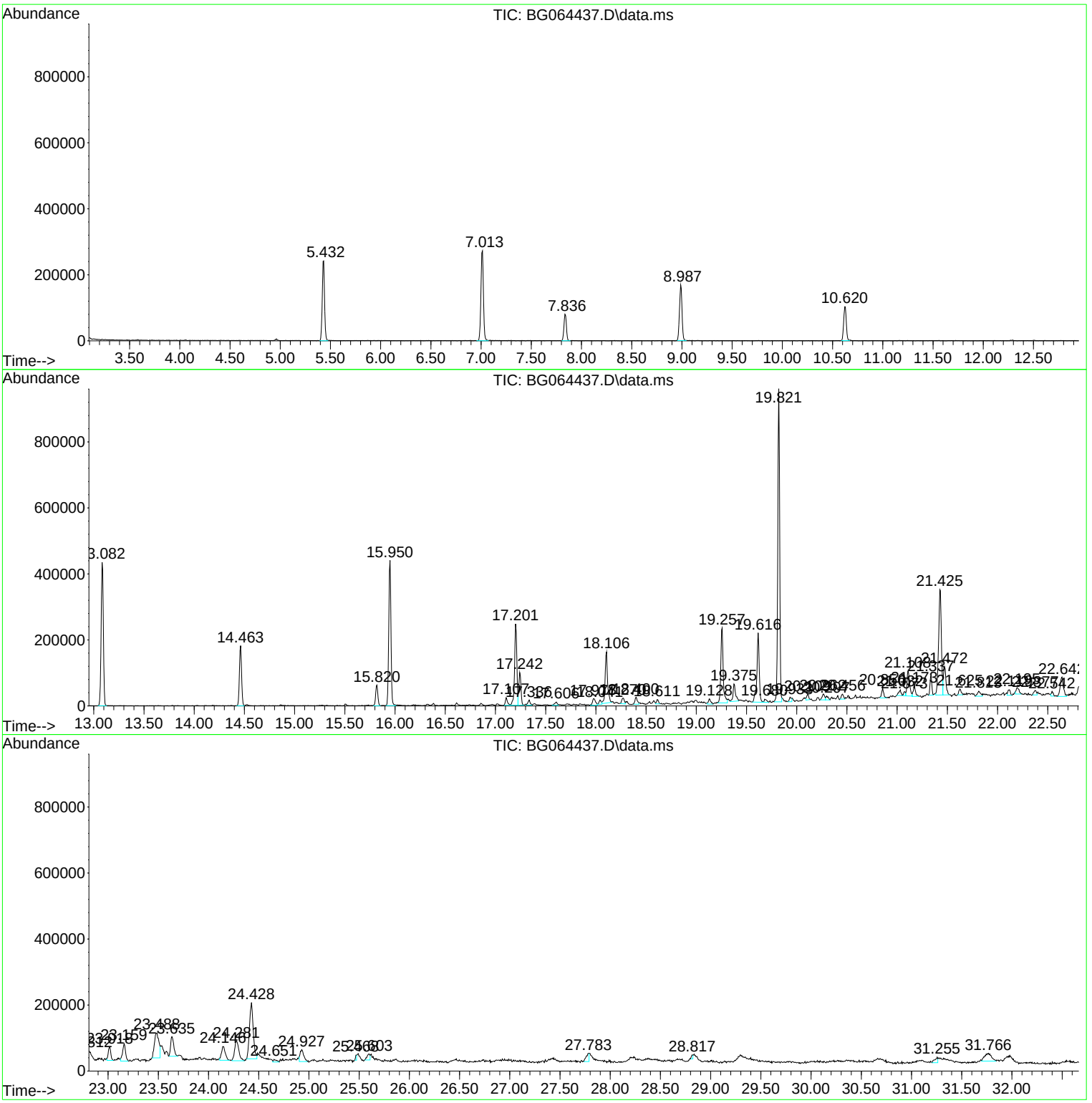
Sample : Q3155-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
COMP-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082825.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092325\
Data File : BG064437.D
Acq On : 23 Sep 2025 13:18
Operator : RC/JU
Sample : Q3155-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
COMP-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082825.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

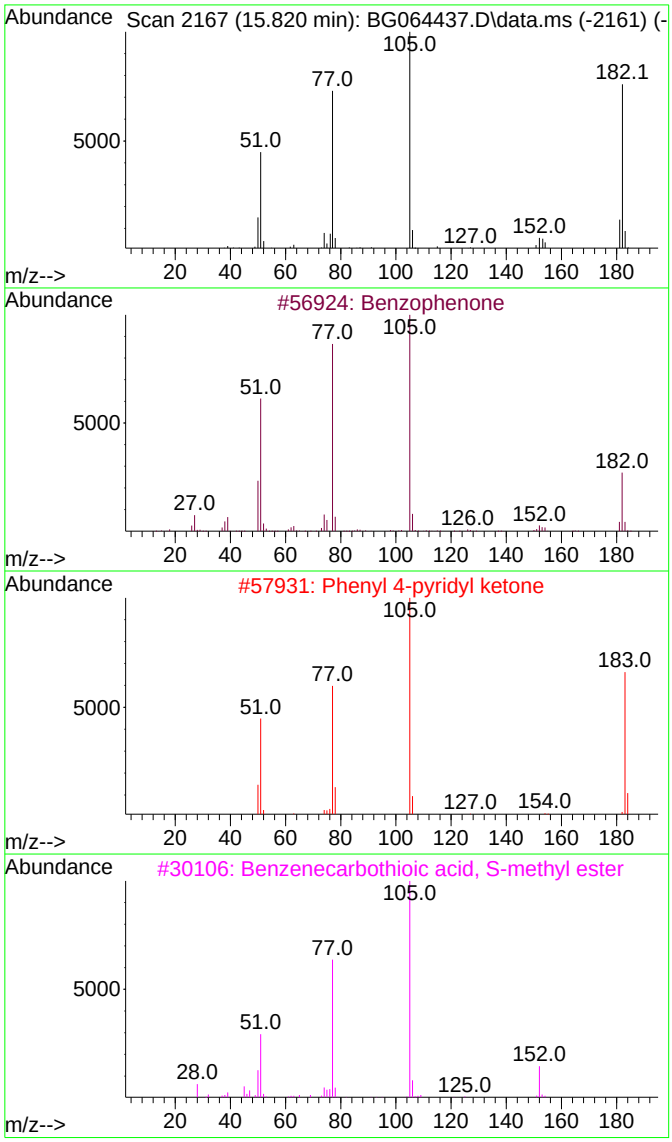
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

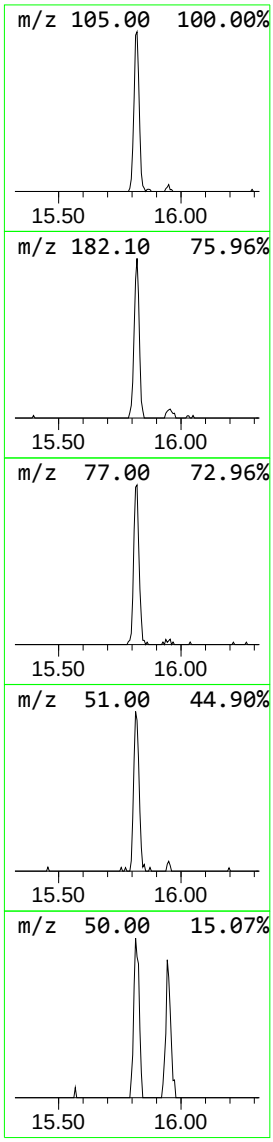
**********

Peak Number 1 Benzophenone Concentration Rank 2

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 15.820 | 7.18 ng | 96029 | Acenaphthene-d10 | 14.463 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzophenone | 182 | C13H10O | 000119-61-9 | 97 |
| 2 | Phenyl 4-pyridyl ketone | 183 | C12H9NO | 014548-46-0 | 55 |
| 3 | Benzenecarbothioic acid, S-methy... | 152 | C8H8OS | 005925-68-8 | 45 |
| 4 | Benzoylformic acid | 150 | C8H6O3 | 000611-73-4 | 38 |
| 5 | Benzenepropanenitrile, .beta.-oxo- | 145 | C9H7NO | 000614-16-4 | 38 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092325\
Data File : BG064437.D
Acq On : 23 Sep 2025 13:18
Operator : RC/JU
Sample : Q3155-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
COMP-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082825.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

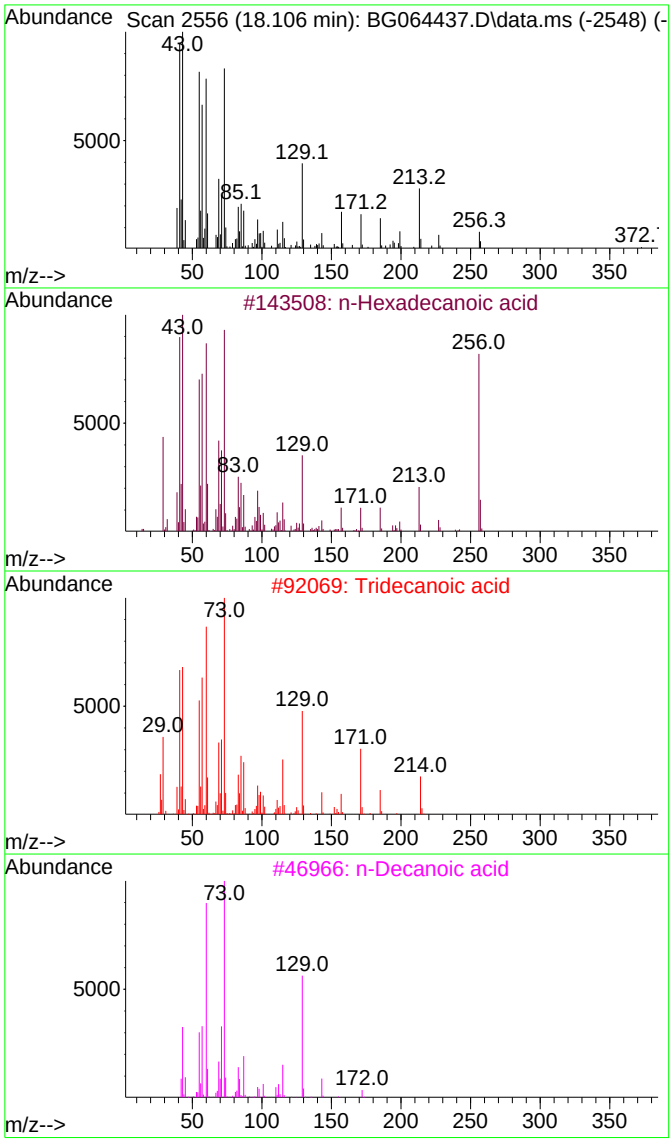
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

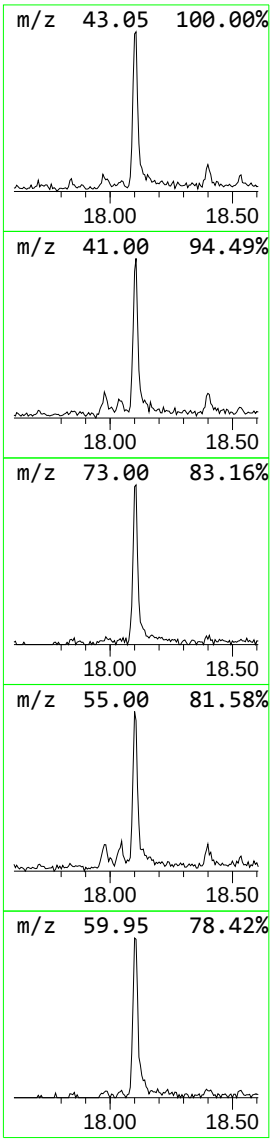
**********

Peak Number 3 n-Hexadecanoic acid Concentration Rank 1

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 18.106 | 13.47 ng | 255781 | Phenanthrene-d10 | 17.201 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | n-Hexadecanoic acid | 256 | C16H32O2 | 000057-10-3 | 94 |
| 2 | Tridecanoic acid | 214 | C13H26O2 | 000638-53-9 | 87 |
| 3 | n-Decanoic acid | 172 | C10H20O2 | 000334-48-5 | 78 |
| 4 | Tetradecanoic acid | 228 | C14H28O2 | 000544-63-8 | 74 |
| 5 | Pentadecanoic acid | 242 | C15H30O2 | 001002-84-2 | 72 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092325\
Data File : BG064437.D
Acq On : 23 Sep 2025 13:18
Operator : RC/JU
Sample : Q3155-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
COMP-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082825.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

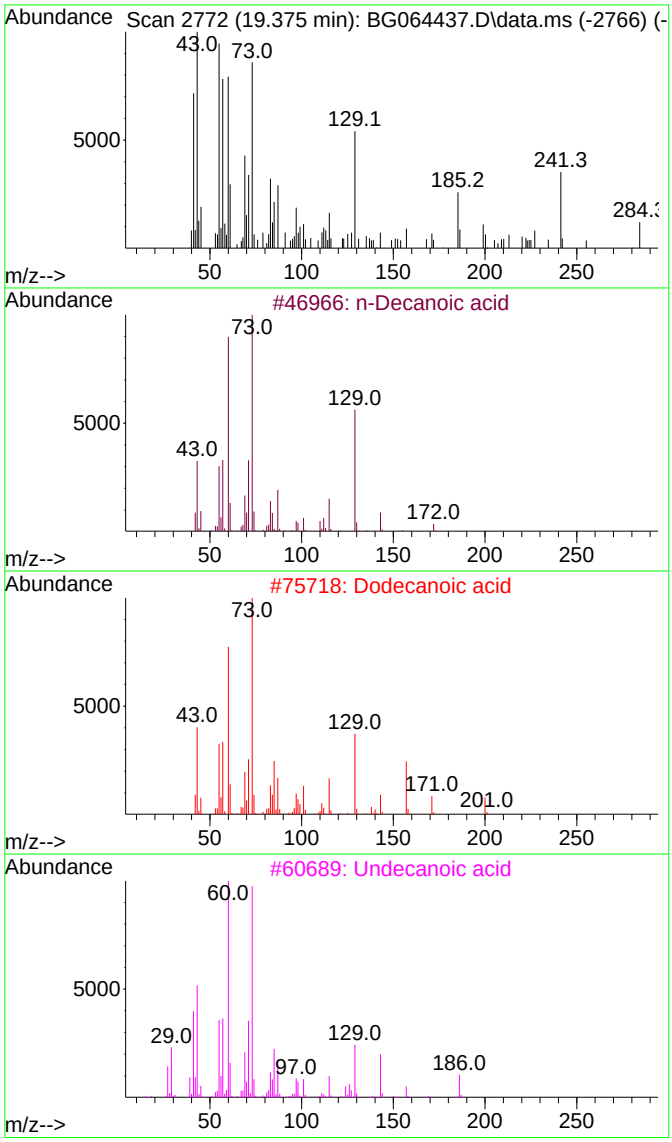
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

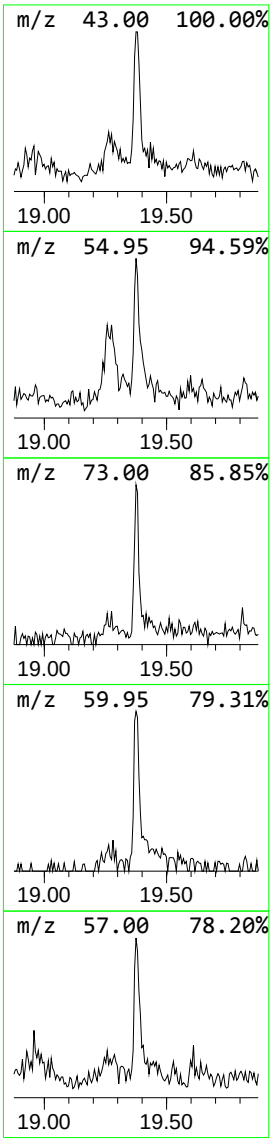
**********

Peak Number 4 n-Decanoic acid Concentration Rank 11

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 19.375 | 3.03 ng | 92473 | Chrysene-d12 | 21.431 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | n-Decanoic acid | 172 | C10H20O2 | 000334-48-5 | 90 |
| 2 | Dodecanoic acid | 200 | C12H24O2 | 000143-07-7 | 78 |
| 3 | Undecanoic acid | 186 | C11H22O2 | 000112-37-8 | 60 |
| 4 | Pentadecanoic acid | 242 | C15H30O2 | 001002-84-2 | 59 |
| 5 | Tetradecanoic acid | 228 | C14H28O2 | 000544-63-8 | 53 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092325\
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Acq On : 23 Sep 2025 13:18
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Misc :
ALS Vial : 4 Sample Multiplier: 1

**Instrument :**
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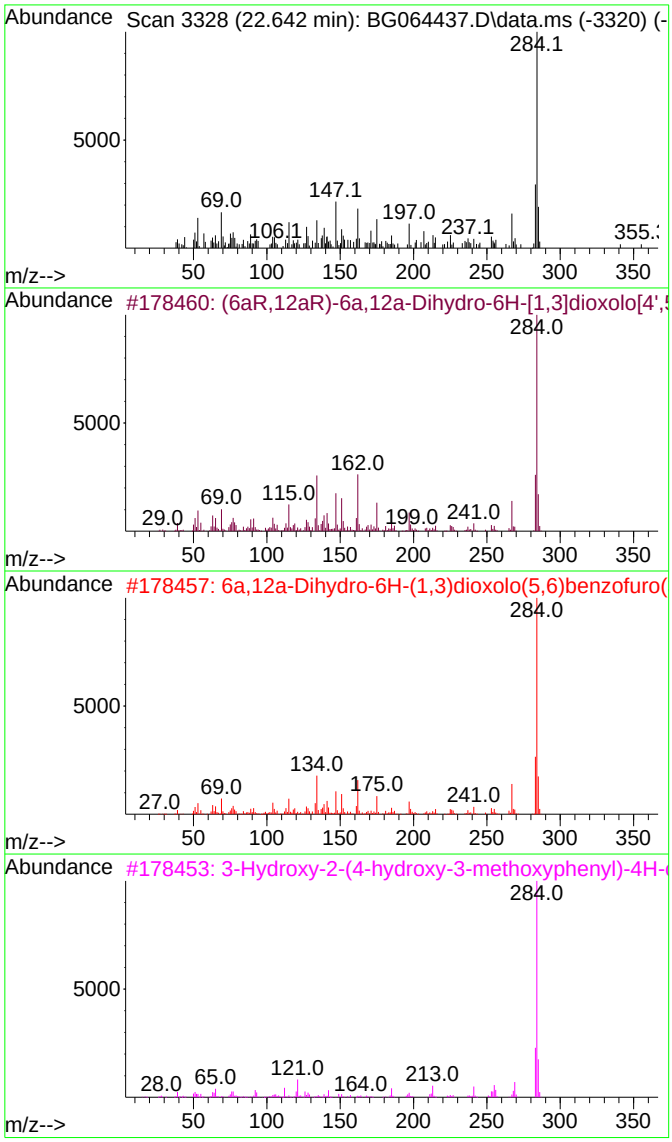
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082825.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

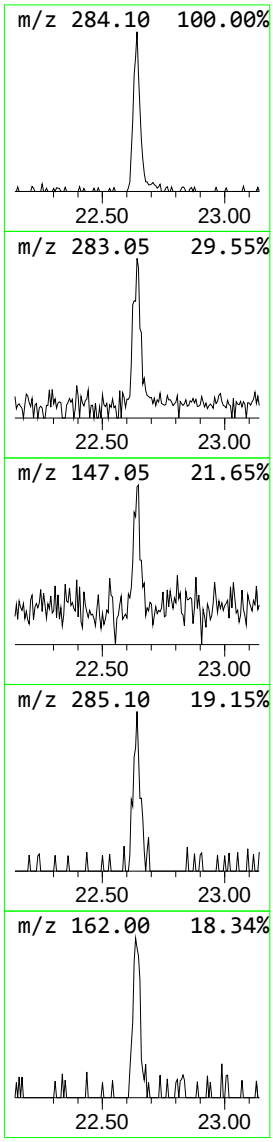
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 6 (6aR,12aR)-6a,12a-Dihydro-6... Concentration Rank 8

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 22.642 | 4.36 ng | 133054 | Chrysene-d12 | 21.431 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | (6aR,12aR)-6a,12a-Dihydro-6H-[1,... | 284 | C16H12O5 | 002035-15-6 | 93 |
| 2 | 6a,12a-Dihydro-6H-(1,3)dioxolo(5... | 284 | C16H12O5 | 076496-57-6 | 93 |
| 3 | 3-Hydroxy-2-(4-hydroxy-3-methoxy... | 284 | C16H12O5 | 159184-95-9 | 83 |
| 4 | Maackiain, acetate | 326 | C18H14O6 | 002035-16-7 | 81 |
| 5 | Coumaran-6-ol-3-one, 2-[4-hydrox... | 284 | C16H12O5 | 032396-80-8 | 70 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092325\
Data File : BG064437.D
Acq On : 23 Sep 2025 13:18
Operator : RC/JU
Sample : Q3155-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

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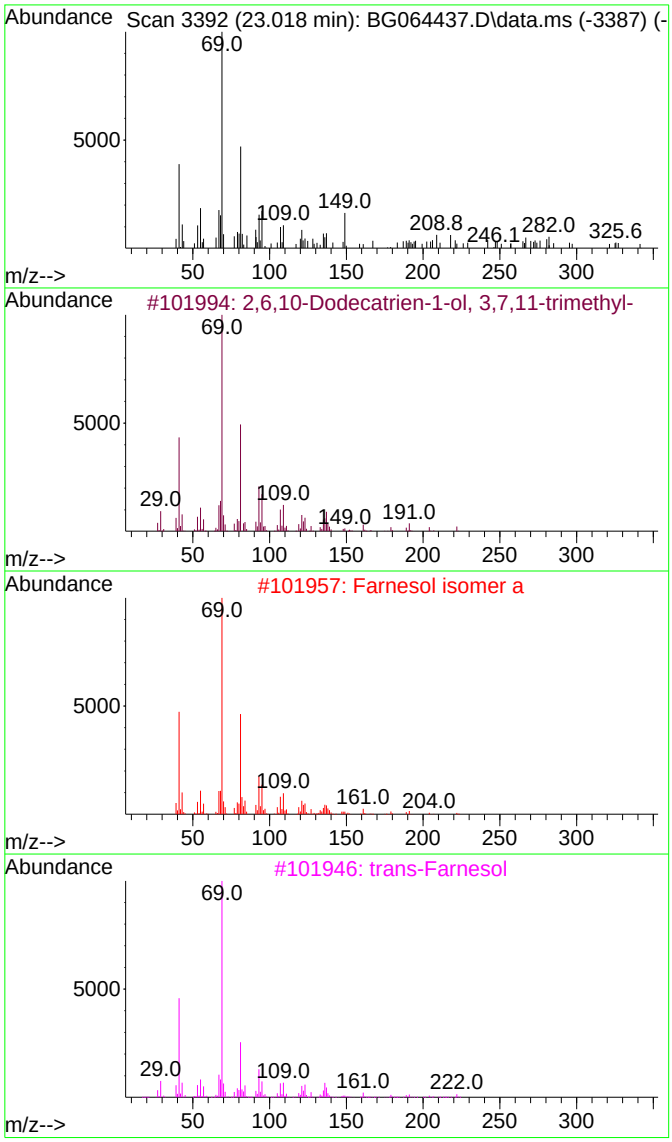
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082825.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

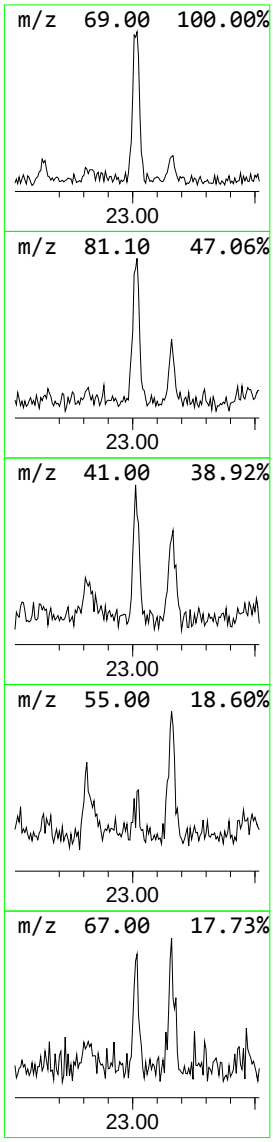
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

**********
Peak Number 8 2,6,10-Dodecatrien-1-ol, 3,... Concentration Rank 10

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 23.018 | 3.21 ng | 70917 | Perylene-d12 | 24.428 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 2,6,10-Dodecatrien-1-ol, 3,7,11-... | 222 | C15H26O | 004602-84-0 | 97 |
| 2 | Farnesol isomer a | 222 | C15H26O | 1000108-92-4 | 93 |
| 3 | trans-Farnesol | 222 | C15H26O | 000106-28-5 | 70 |
| 4 | 2,6,10-Dodecatrien-1-ol, 3,7,11-... | 222 | C15H26O | 003790-71-4 | 70 |
| 5 | 1,5,9-Undecatriene, 2,6,10-trime... | 192 | C14H24 | 062951-96-6 | 58 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092325\
Data File : BG064437.D
Acq On : 23 Sep 2025 13:18
Operator : RC/JU
Sample : Q3155-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

**Instrument :**
BNA_G
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082825.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

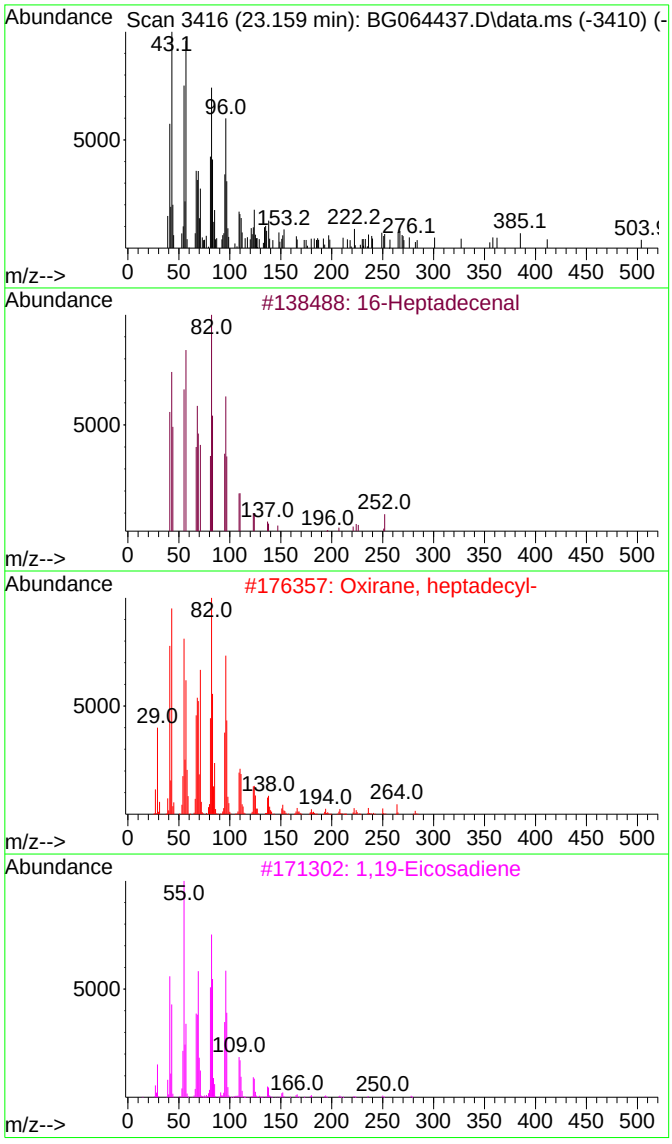
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

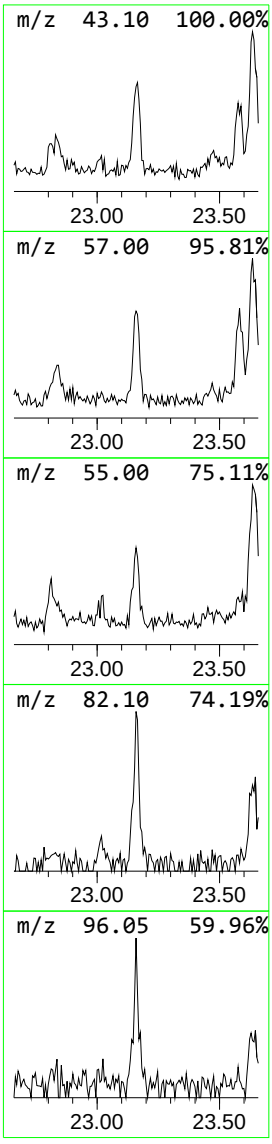
**********

Peak Number 9 16-Heptadecenal Concentration Rank 6

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 23.159 | 4.86 ng | 107315 | Perylene-d12 | 24.428 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | 16-Heptadecenal | 252 | C17H32O | 1010144-57-9 | 91 |
| 2 | Oxirane, heptadecyl- | 282 | C19H38O | 067860-04-2 | 91 |
| 3 | 1,19-Eicosadiene | 278 | C20H38 | 014811-95-1 | 87 |
| 4 | 1,16-Hexadecanediol | 258 | C16H34O2 | 007735-42-4 | 87 |
| 5 | Pentadecanal- | 226 | C15H30O | 002765-11-9 | 83 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092325\
Data File : BG064437.D
Acq On : 23 Sep 2025 13:18
Operator : RC/JU
Sample : Q3155-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082825.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

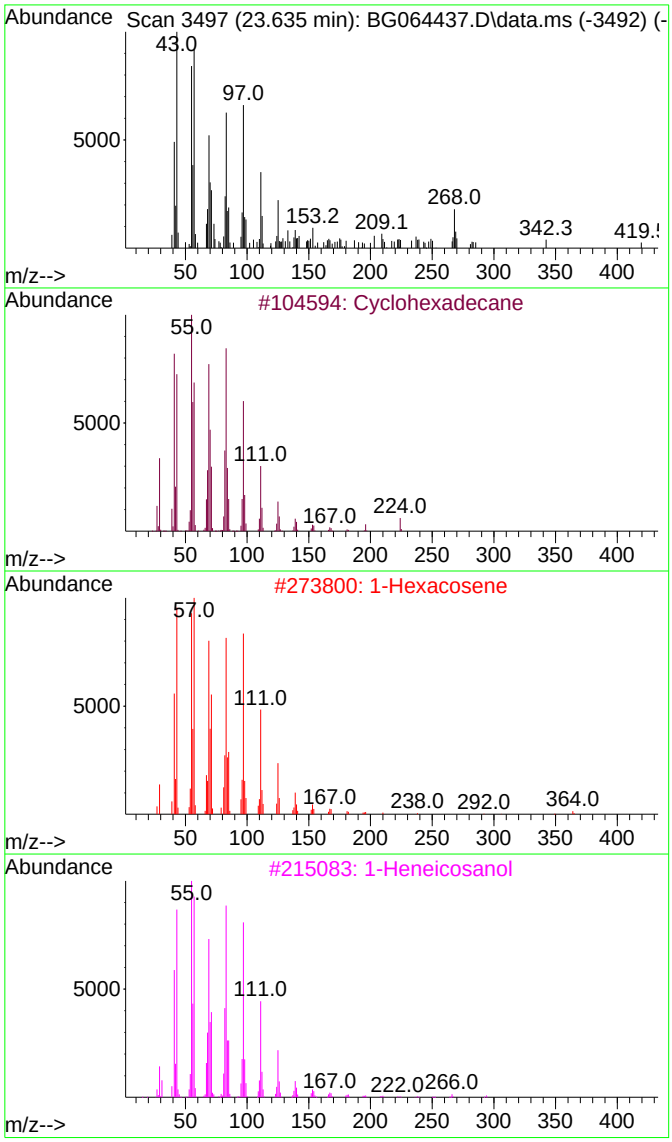
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

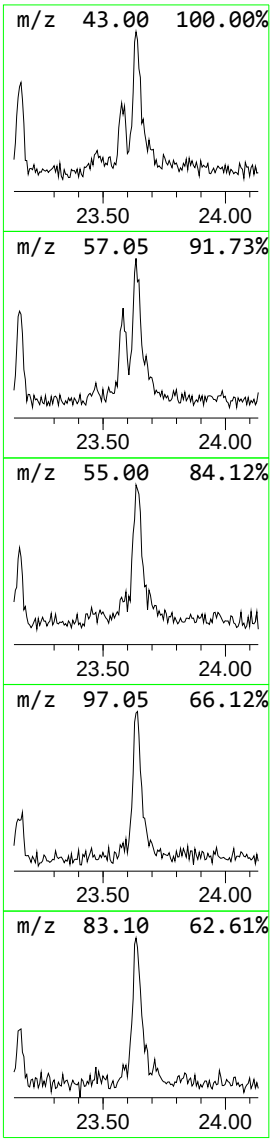
**********

Peak Number 10 Cyclohexadecane Concentration Rank 3

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 23.635 | 5.91 ng | 130496 | Perylene-d12 | 24.428 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Cyclohexadecane | 224 | C16H32 | 000295-65-8 | 95 |
| 2 | 1-Hexacosene | 364 | C26H52 | 018835-33-1 | 91 |
| 3 | 1-Heneicosanol | 312 | C21H44O | 015594-90-8 | 87 |
| 4 | Z-5-Nonadecene | 266 | C19H38 | 1000131-11-8 | 81 |
| 5 | 1-Docosene | 308 | C22H44 | 001599-67-3 | 81 |



Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092325\
Data File : BG064437.D
Acq On : 23 Sep 2025 13:18
Operator : RC/JU
Sample : Q3155-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
COMP-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082825.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

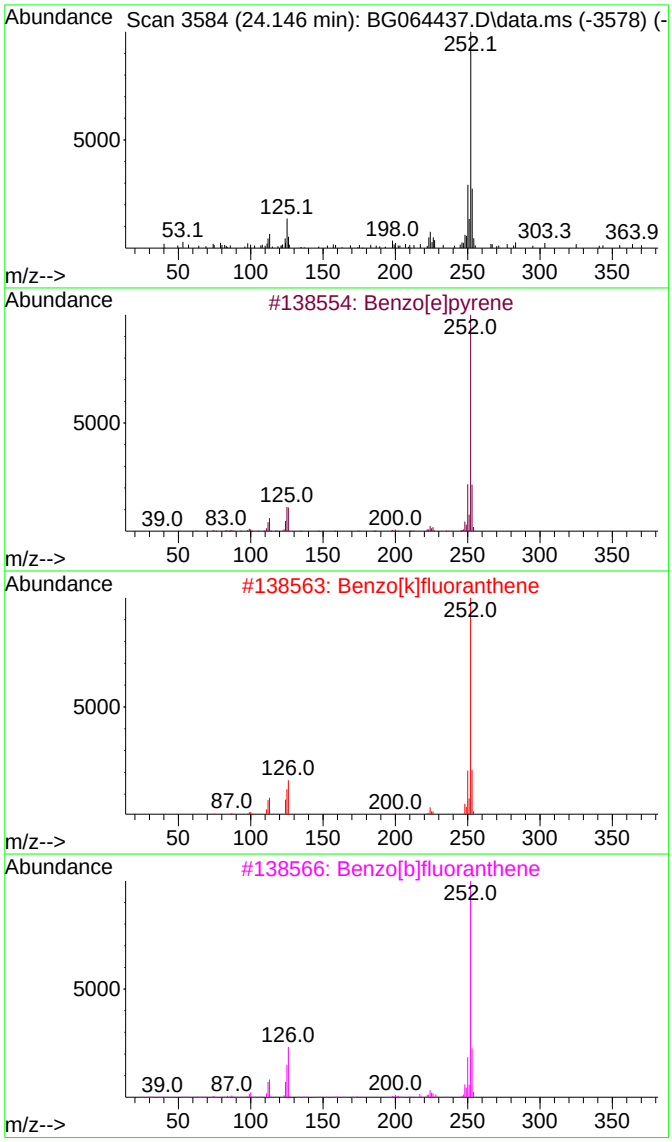
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

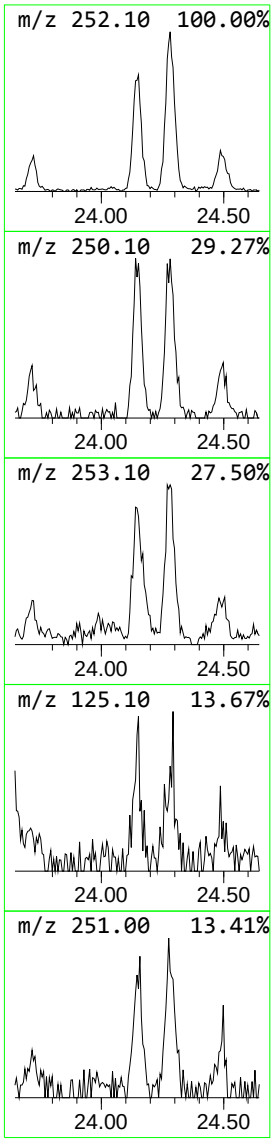
**********

Peak Number 11 Benzo[e]pyrene Concentration Rank 5

| R.T. | EstConc | Area | Relative to ISTD | R.T. |
|---|---|---|---|---|
| 24.146 | 4.92 ng | 108827 | Perylene-d12 | 24.428 |

| Hit# of 5 | Tentative ID | MW | MolForm | CAS# | Qual |
|---|---|---|---|---|---|
| 1 | Benzo[e]pyrene | 252 | C20H12 | 000192-97-2 | 90 |
| 2 | Benzo[k]fluoranthene | 252 | C20H12 | 000207-08-9 | 90 |
| 3 | Benzo[b]fluoranthene | 252 | C20H12 | 000205-99-2 | 89 |
| 4 | Benzo[a]pyrene | 252 | C20H12 | 000050-32-8 | 81 |
| 5 | Perylene | 252 | C20H12 | 000198-55-0 | 81 |



Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092325\
Data File : BG064437.D
Acq On : 23 Sep 2025 13:18
Operator : RC/JU
Sample : Q3155-01
Misc :
ALS Vial : 4 Sample Multiplier: 1

**Instrument :**
BNA_G
**ClientSampleId :**
COMP-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG082825.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

| | | | | | --Internal Standard--- | | | |
|---|---|---|---|---|---|---|---|---|
| TIC Top Hit name | RT | EstConc | Units | Response | # | RT | Resp | Conc |
| Benzophenone | 15.820 | 7.2 | ng | 96029 | 3 | 14.463 | 267612 | 20.0 |
| n-Hexadecanoic ... | 18.106 | 13.5 | ng | 255781 | 4 | 17.201 | 379895 | 20.0 |
| n-Decanoic acid | 19.375 | 3.0 | ng | 92473 | 5 | 21.431 | 610863 | 20.0 |
| (6aR,12aR)-6a,1... | 22.642 | 4.4 | ng | 133054 | 5 | 21.431 | 610863 | 20.0 |
| 2,6,10-Dodecatr... | 23.018 | 3.2 | ng | 70917 | 6 | 24.428 | 441961 | 20.0 |
| 16-Heptadecenal | 23.159 | 4.9 | ng | 107315 | 6 | 24.428 | 441961 | 20.0 |
| Cyclohexadecane | 23.635 | 5.9 | ng | 130496 | 6 | 24.428 | 441961 | 20.0 |
| Benzo[e]pyrene | 24.146 | 4.9 | ng | 108827 | 6 | 24.428 | 441961 | 20.0 |