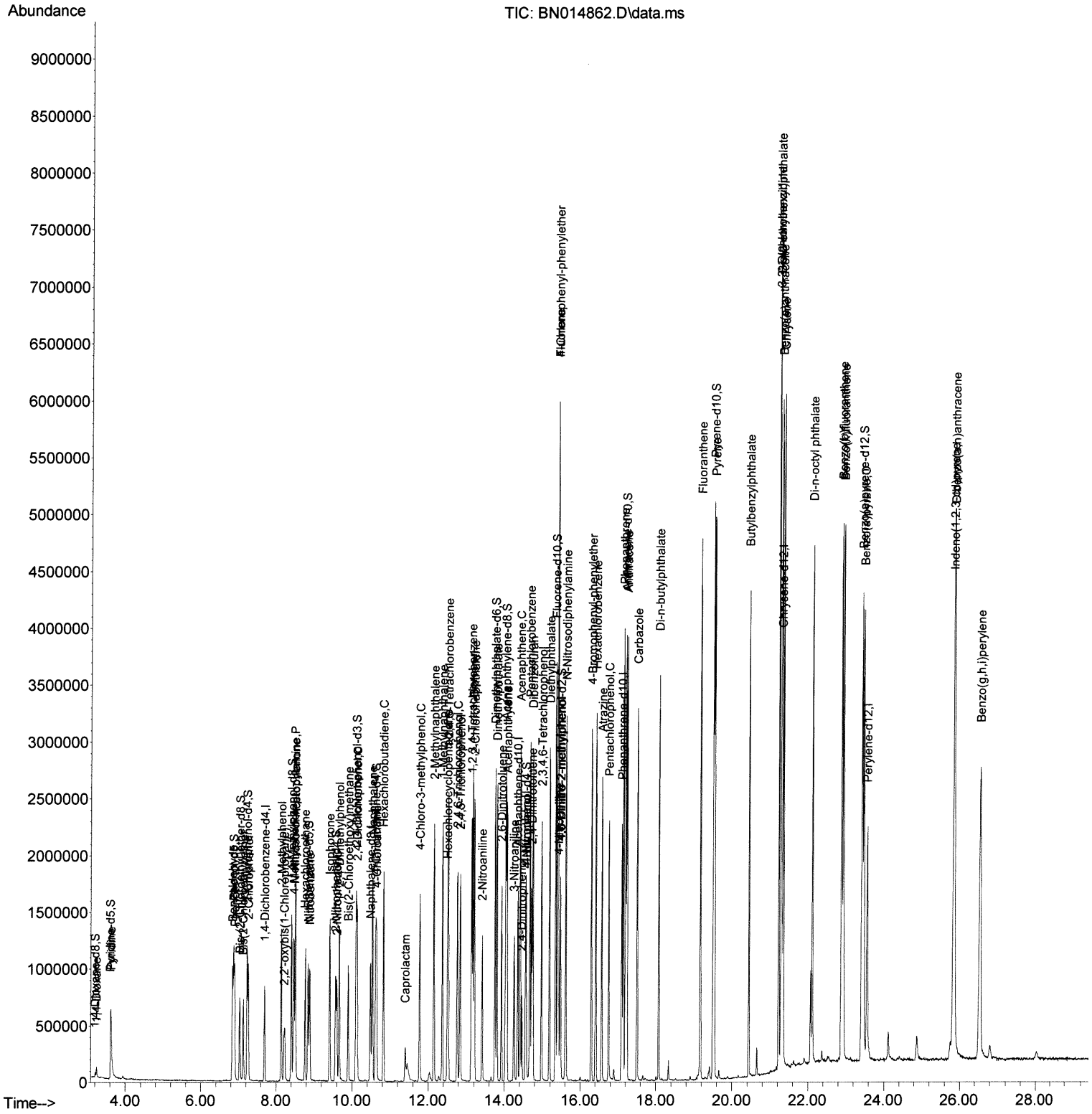
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061021\
Data File : BN014862.D
Acq On : 10 Jun 2021 11:51
Operator : CG/JU
Sample : PB136905BS
Misc :
ALS Vial : 4 Sample Multiplier: 1

Quant Time: Jun 10 12:39:54 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN052521.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jun 10 11:55:23 2021
Response via : Initial Calibration

**Instrument :**
BNA_N
**ClientSampleId :**
SLCS905

**Manual Integrations APPROVED**
mohammad
6/11/2021 2:13:03 PM

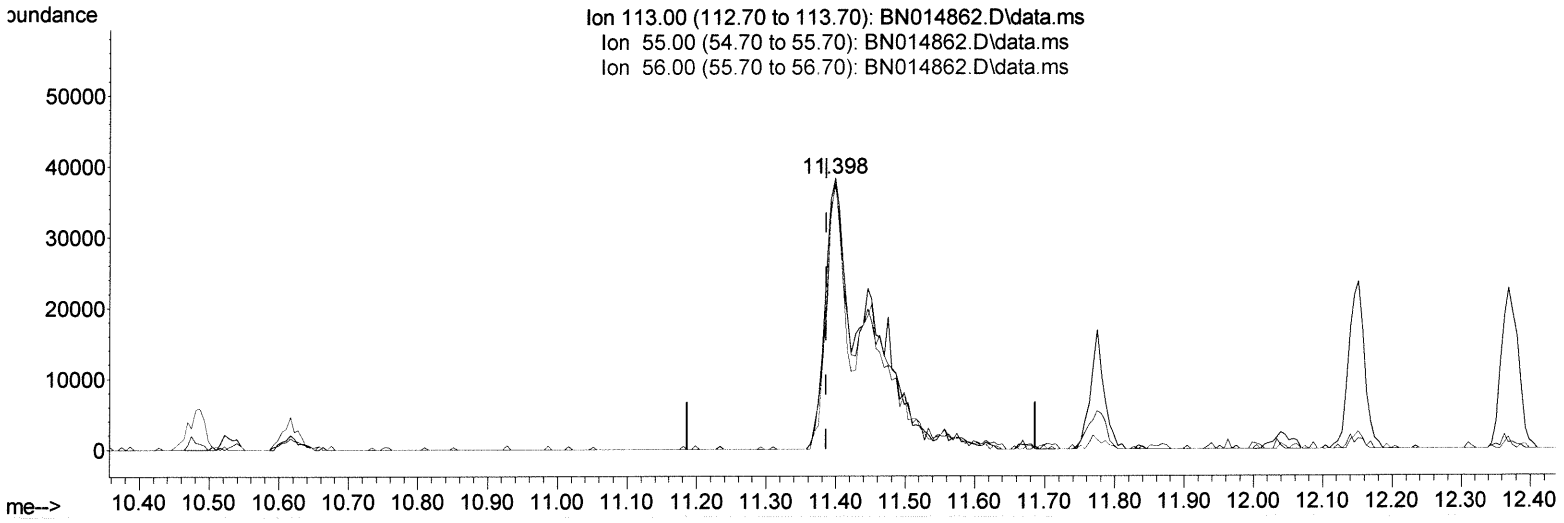


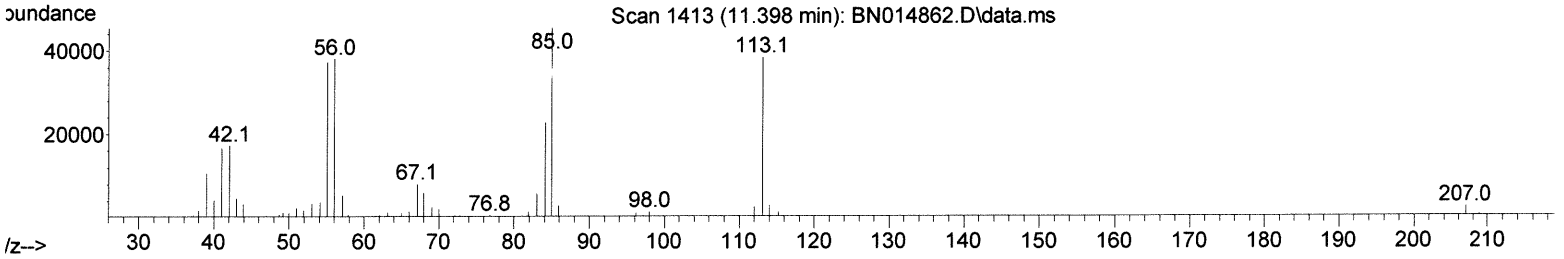
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061021\
Data File : BN014862.D
Acq On : 10 Jun 2021 11:51
Operator : CG/JU
Sample : PB136905BS
Misc :
ALS Vial : 4 Sample Multiplier: 1

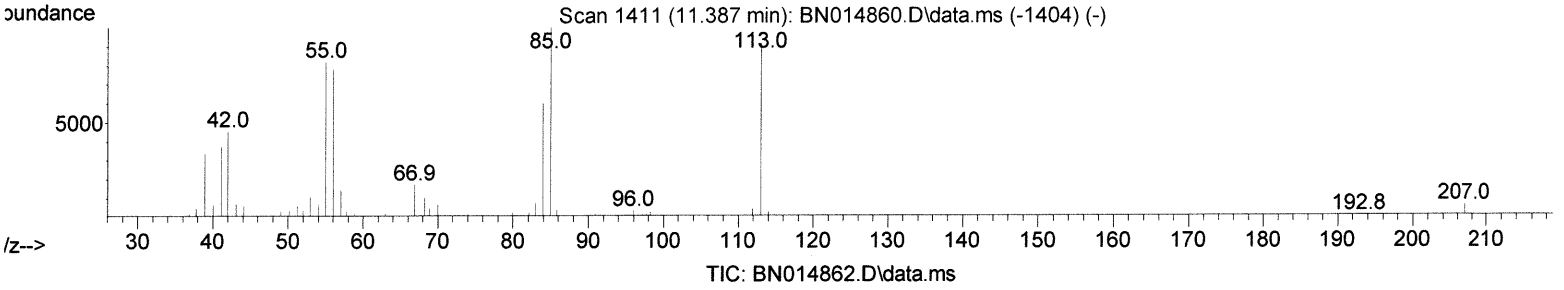
Quant Time: Jun 10 12:39:54 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN052521.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jun 10 11:55:23 2021
Response via : Initial Calibration

**Instrument :**
BNA_N
**ClientSampleId :**
SLCS905

**Manual Integrations APPROVED**
mohammad
6/11/2021 2:13:03 PM




TIC: BN014862.D\data.ms

**(34) Caprolactam**

**11.398min (+ 0.012) 17.08 ng/ul**

**response 76050**

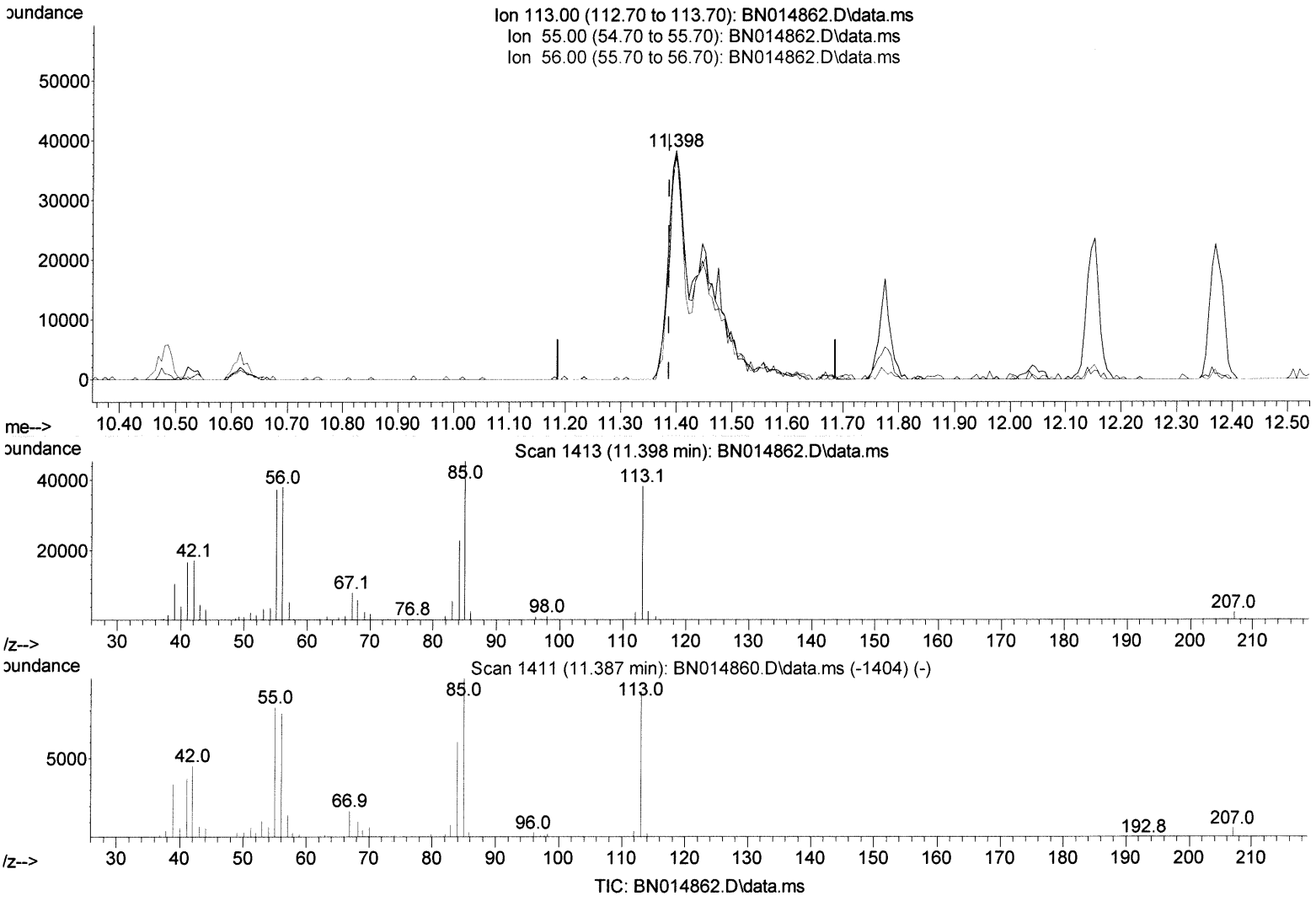
| Ion | Exp% | Act% |
|---|---|---|
| 113.00 | 100.00 | 100.00 |
| 55.00 | 118.50 | 97.43 |
| 56.00 | 73.20 | 99.54# |
| 0.00 | 0.00 | 0.00 |

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061021\
Data File : BN014862.D
Acq On : 10 Jun 2021 11:51
Operator : CG/JU
Sample : PB136905BS
Misc :
ALS Vial : 4 Sample Multiplier: 1

Quant Time: Jun 10 12:39:54 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN052521.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jun 10 11:55:23 2021
Response via : Initial Calibration

**Instrument :**
BNA_N
**ClientSampleId :**
SLCS905

**Manual Integrations APPROVED**
mohammad
6/11/2021 2:13:03 PM




(34) Caprolactam

11.398min (+ 0.012) 34.28 ng/ul m JU 6/16/21

response 152652

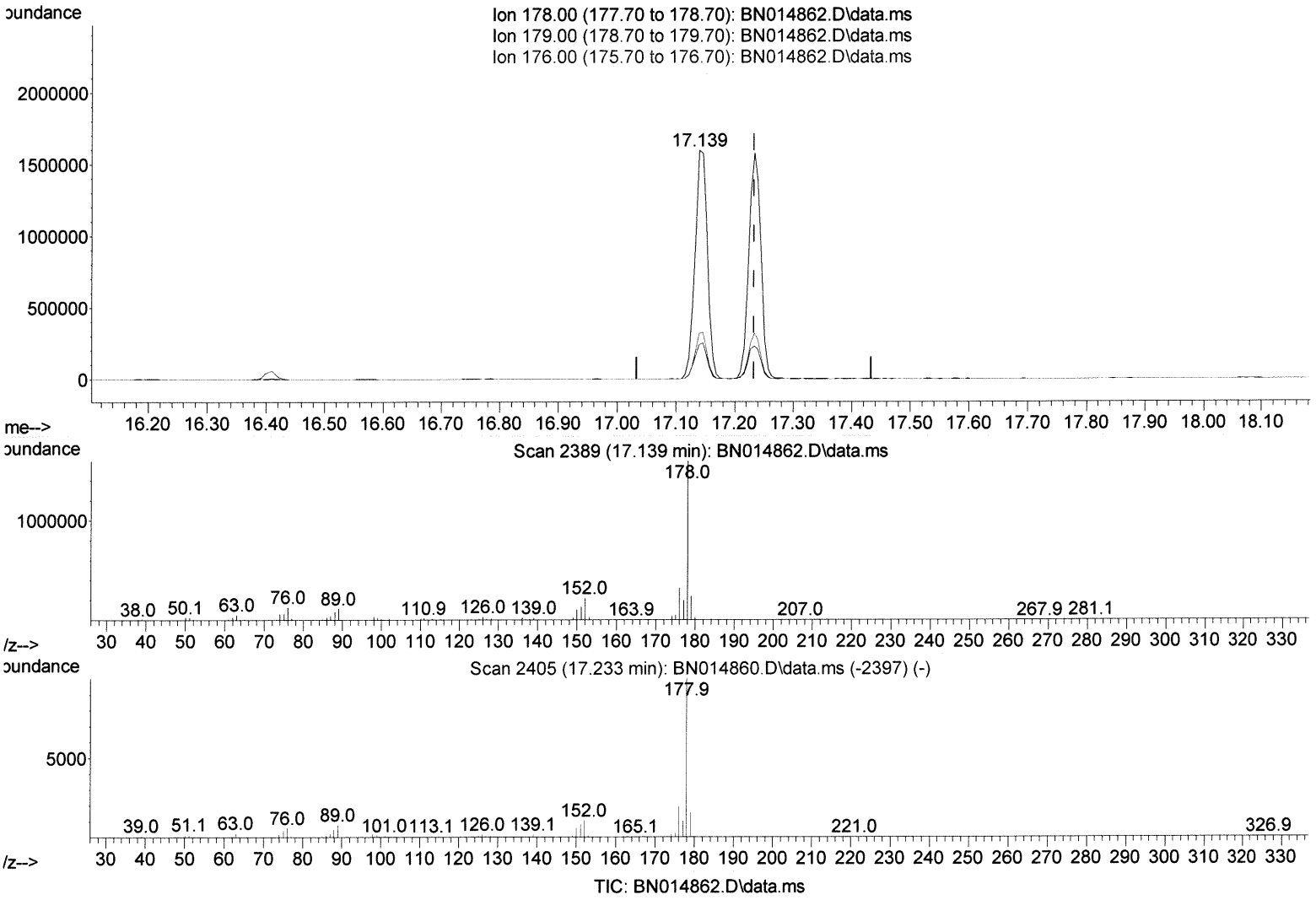
| Ion | Exp% | Act% |
|---|---|---|
| 113.00 | 100.00 | 100.00 |
| 55.00 | 118.50 | 97.43 |
| 56.00 | 73.20 | 99.54# |
| 0.00 | 0.00 | 0.00 |

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061021\
Data File : BN014862.D
Acq On : 10 Jun 2021 11:51
Operator : CG/JU
Sample : PB136905BS
Misc :
ALS Vial : 4 Sample Multiplier: 1

Quant Time: Jun 10 12:39:54 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN052521.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jun 10 11:55:23 2021
Response via : Initial Calibration

**Instrument :**
BNA_N
**ClientSampleId :**
SLCS905

**Manual Integrations APPROVED**
mohammad
6/11/2021 2:13:03 PM




**(74) Anthracene**

**17.139min (-0.094) 35.42 ng/ul**

**response 2329804**

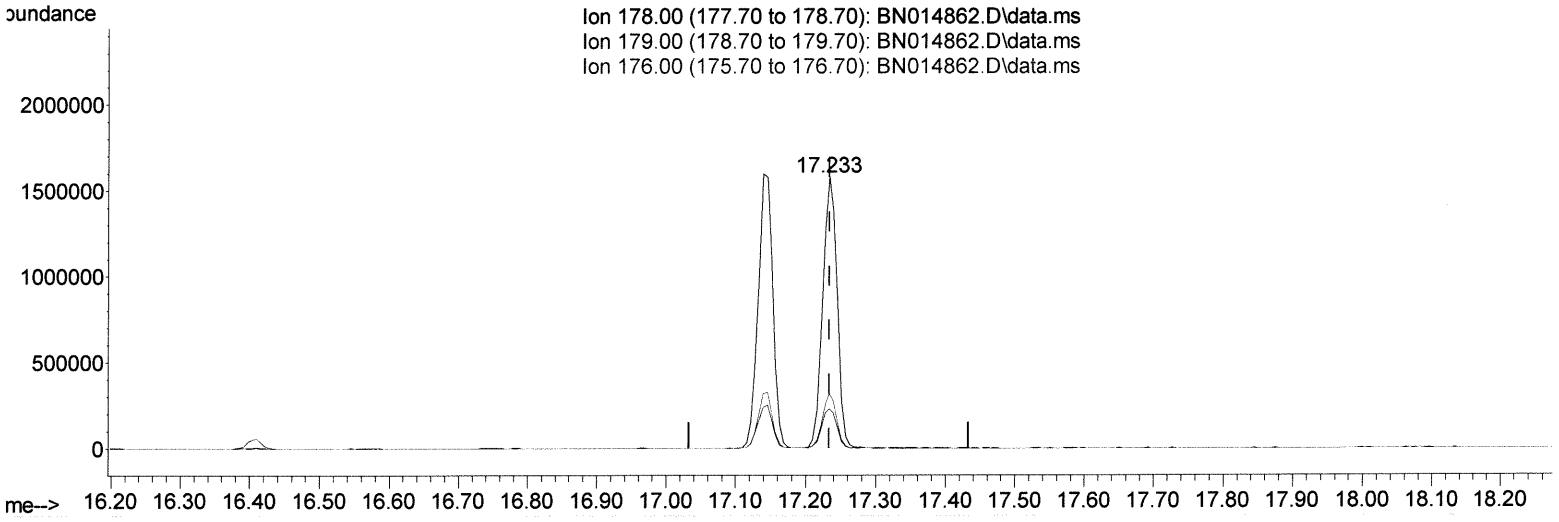
| Ion | Exp% | Act% |
|---|---|---|
| 178.00 | 100.00 | 100.00 |
| 179.00 | 16.10 | 14.92 |
| 176.00 | 19.10 | 19.97 |
| 0.00 | 0.00 | 0.00 |

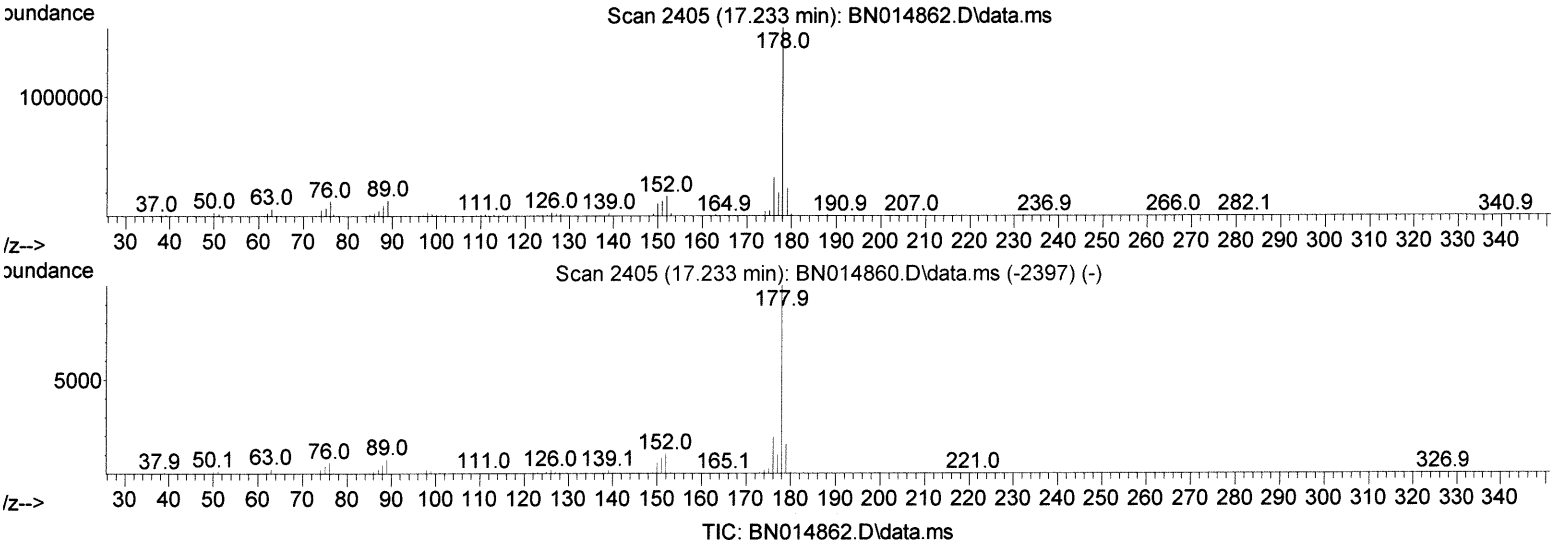
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061021\
Data File : BN014862.D
Acq On : 10 Jun 2021 11:51
Operator : CG/JU
Sample : PB136905BS
Misc :
ALS Vial : 4 Sample Multiplier: 1

Quant Time: Jun 10 12:39:54 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN052521.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jun 10 11:55:23 2021
Response via : Initial Calibration

**Instrument :**
BNA_N
**ClientSampleId :**
SLCS905

**Manual Integrations APPROVED**
mohammad
6/11/2021 2:13:03 PM




(74) Anthracene

17.233min (-0.000) 34.37 ng/ul m

response 2261080

JU 6/14/21

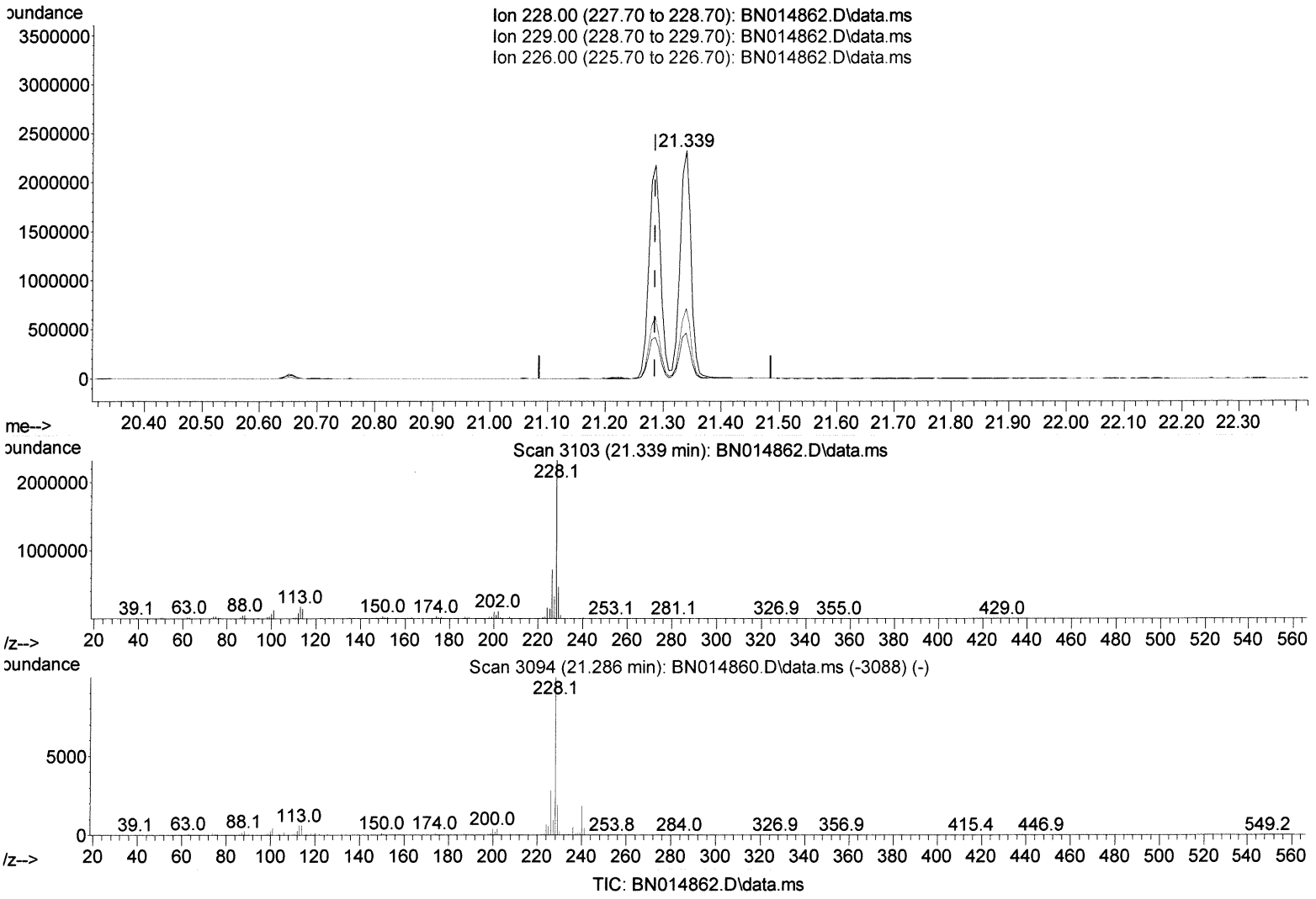
| Ion | Exp% | Act% |
|---|---|---|
| 178.00 | 100.00 | 100.00 |
| 179.00 | 16.10 | 14.51 |
| 176.00 | 19.10 | 20.29 |
| 0.00 | 0.00 | 0.00 |

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061021\
Data File : BN014862.D
Acq On : 10 Jun 2021 11:51
Operator : CG/JU
Sample : PB136905BS
Misc :
ALS Vial : 4 Sample Multiplier: 1

Quant Time: Jun 10 12:39:54 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN052521.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jun 10 11:55:23 2021
Response via : Initial Calibration

**Instrument :**
BNA_N
**ClientSampleId :**
SLCS905

**Manual Integrations APPROVED**
mohammad
6/11/2021 2:13:03 PM




**(85) Benzo(a)anthracene**

**21.339min (+ 0.053) 35.58 ng/ul**

**response 3056503**

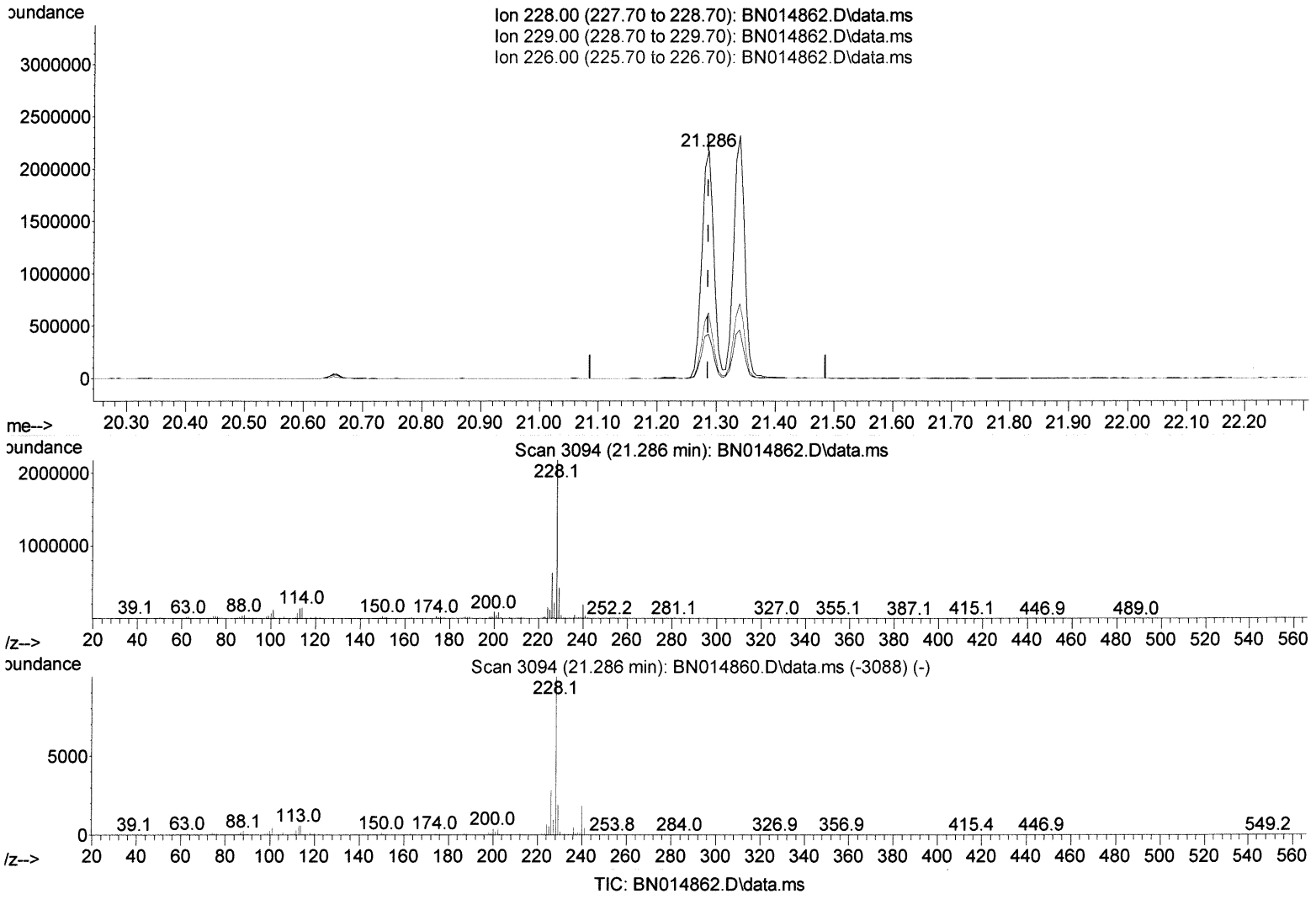
| Ion | Exp% | Act% |
|---|---|---|
| 228.00 | 100.00 | 100.00 |
| 229.00 | 20.30 | 19.95 |
| 226.00 | 27.00 | 30.82 |
| 0.00 | 0.00 | 0.00 |

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061021\
Data File : BN014862.D
Acq On : 10 Jun 2021 11:51
Operator : CG/JU
Sample : PB136905BS
Misc :
ALS Vial : 4 Sample Multiplier: 1

Quant Time: Jun 10 12:39:54 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN052521.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jun 10 11:55:23 2021
Response via : Initial Calibration

**Instrument :**
BNA_N
**ClientSampleId :**
SLCS905

**Manual Integrations APPROVED**
mohammad
6/11/2021 2:13:03 PM




(85) Benzo(a)anthracene

21.286min (-0.000) 35.34 ng/ul m JU 6/14/21

response 3035225

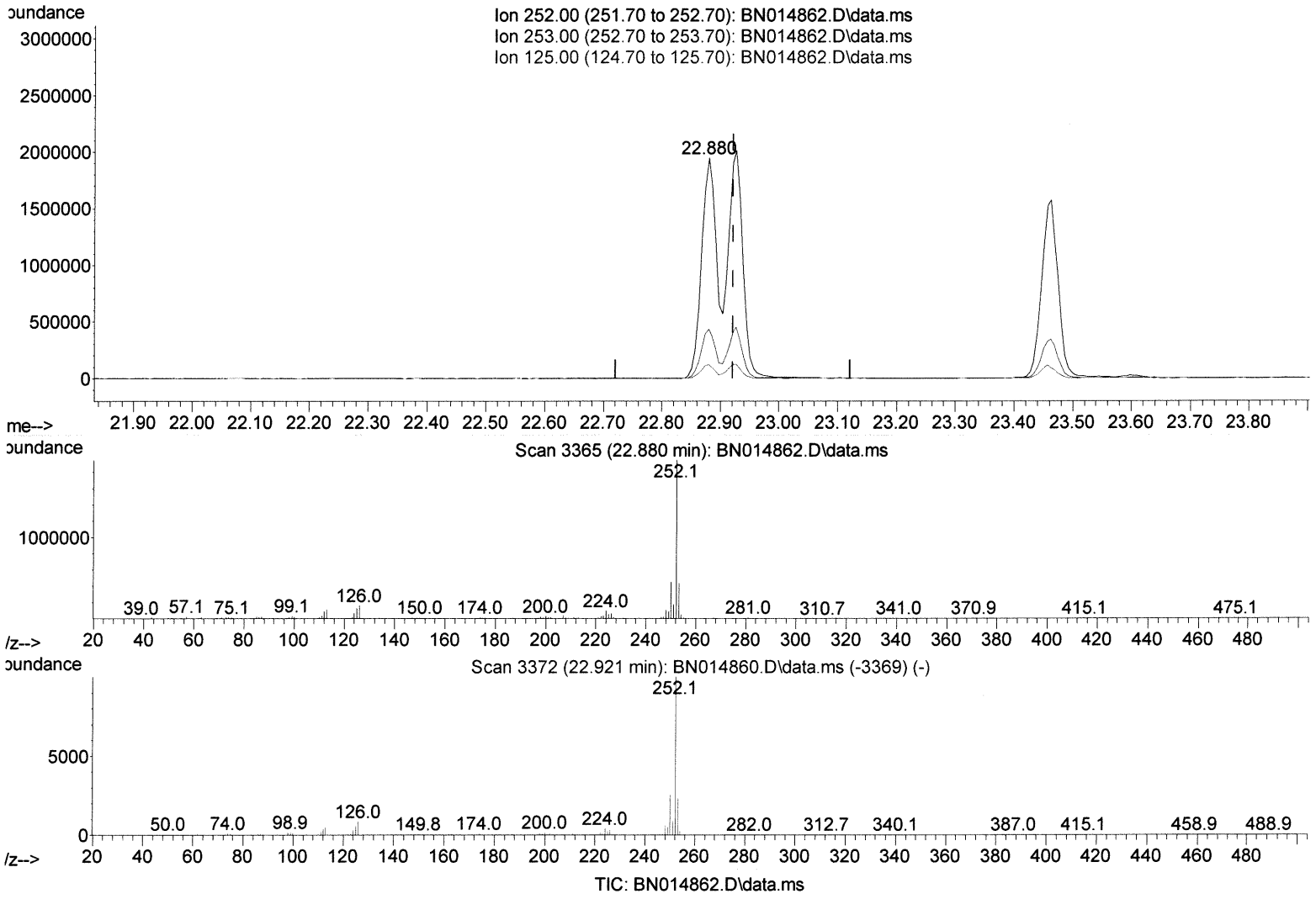
| Ion | Exp% | Act% |
|---|---|---|
| 228.00 | 100.00 | 100.00 |
| 229.00 | 20.30 | 19.39 |
| 226.00 | 27.00 | 28.83 |
| 0.00 | 0.00 | 0.00 |

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061021\
Data File : BN014862.D
Acq On : 10 Jun 2021 11:51
Operator : CG/JU
Sample : PB136905BS
Misc :
ALS Vial : 4 Sample Multiplier: 1

Quant Time: Jun 10 12:39:54 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN052521.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jun 10 11:55:23 2021
Response via : Initial Calibration

**Instrument :**
BNA_N
**ClientSampleId :**
SLCS905

**Manual Integrations APPROVED**
mohammad
6/11/2021 2:13:03 PM




(91) Benzo(k)fluoranthene

22.880min (-0.041) 40.87 ng/ul

response 3506679

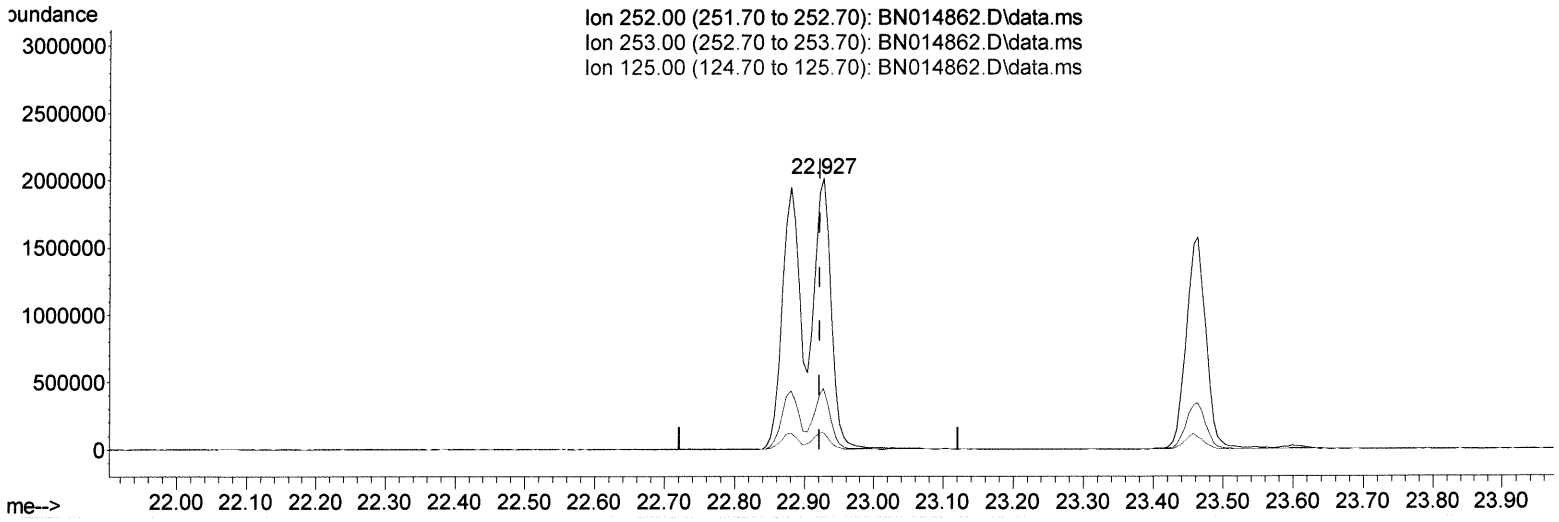
| Ion | Exp% | Act% |
|---|---|---|
| 252.00 | 100.00 | 100.00 |
| 253.00 | 21.20 | 22.37 |
| 125.00 | 6.00 | 6.32 |
| 0.00 | 0.00 | 0.00 |

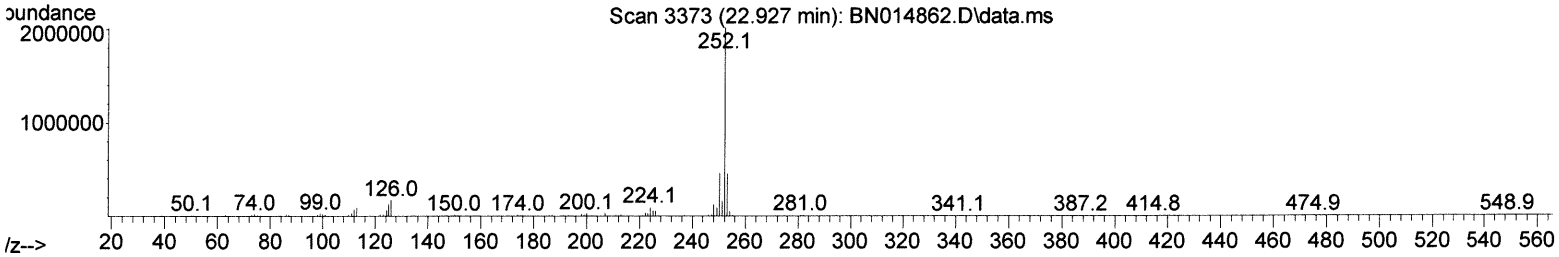
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061021\
Data File : BN014862.D
Acq On : 10 Jun 2021 11:51
Operator : CG/JU
Sample : PB136905BS
Misc :
ALS Vial : 4 Sample Multiplier: 1

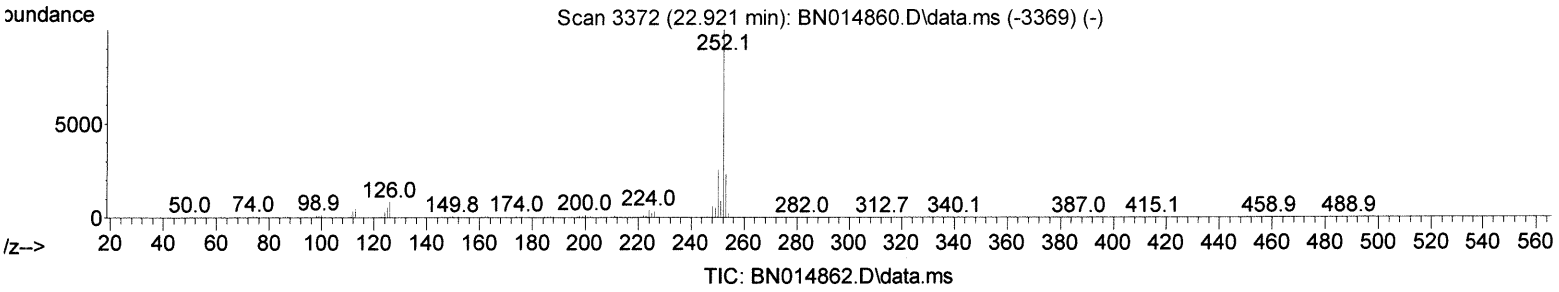
Quant Time: Jun 10 12:39:54 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN052521.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jun 10 11:55:23 2021
Response via : Initial Calibration

**Instrument :**
BNA_N
**ClientSampleId :**
SLCS905

**Manual Integrations APPROVED**
mohammad
6/11/2021 2:13:03 PM




TIC: BN014862.D\data.ms

(91) Benzo(k)fluoranthene

22.927min (+ 0.006) 39.51 ng/ul m JG 6/14/21

response 3389383

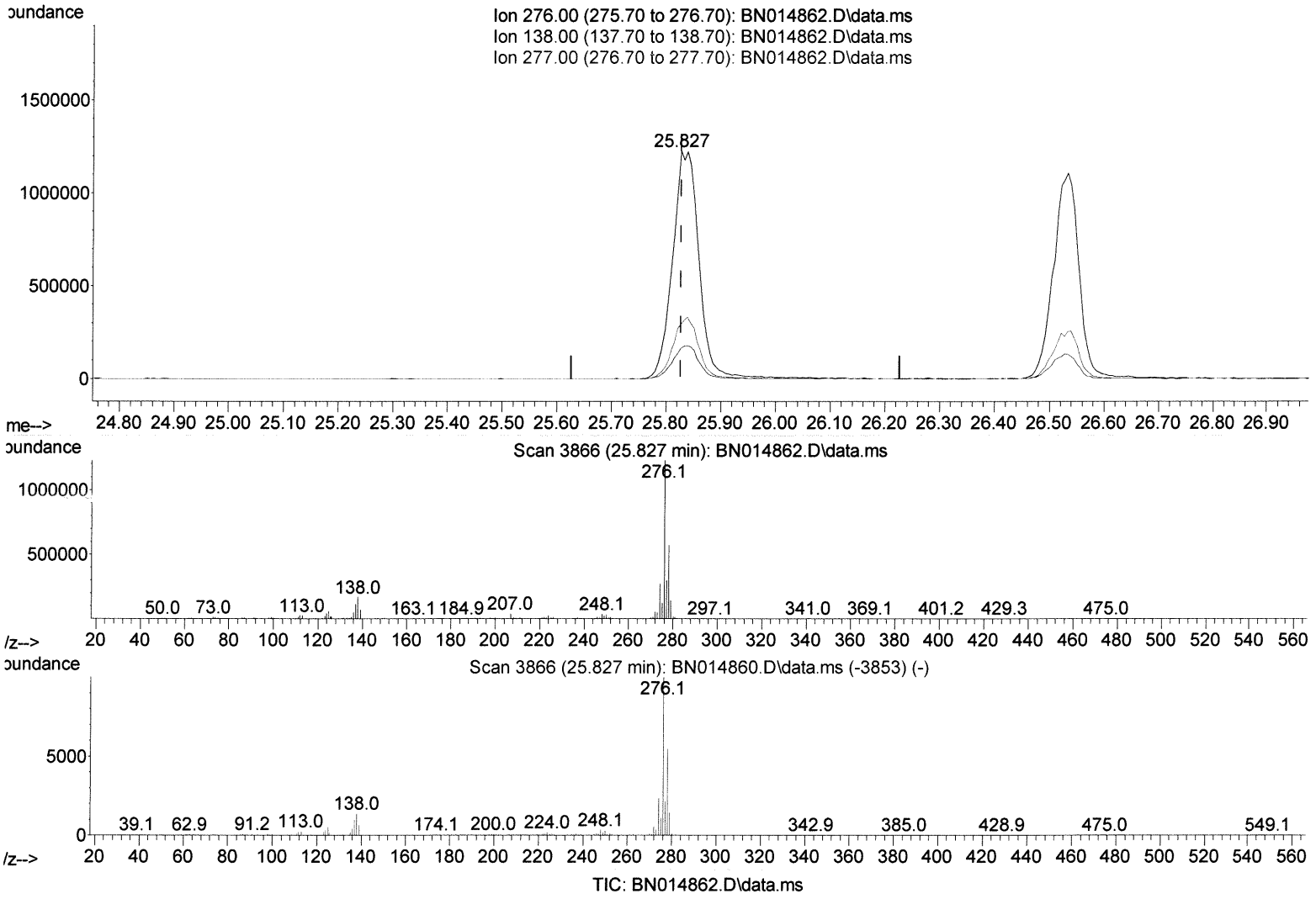
| Ion | Exp% | Act% |
|---|---|---|
| 252.00 | 100.00 | 100.00 |
| 253.00 | 21.20 | 22.52 |
| 125.00 | 6.00 | 6.16 |
| 0.00 | 0.00 | 0.00 |

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061021\
Data File : BN014862.D
Acq On : 10 Jun 2021 11:51
Operator : CG/JU
Sample : PB136905BS
Misc :
ALS Vial : 4 Sample Multiplier: 1

Quant Time: Jun 10 12:39:54 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN052521.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jun 10 11:55:23 2021
Response via : Initial Calibration

**Instrument :**
BNA_N
**ClientSampleId :**
SLCS905

**Manual Integrations APPROVED**
mohammad
6/11/2021 2:13:03 PM




(94) Indeno(1,2,3-cd)pyrene

25.827min (-0.000) 19.27 ng/ul

response 2058104

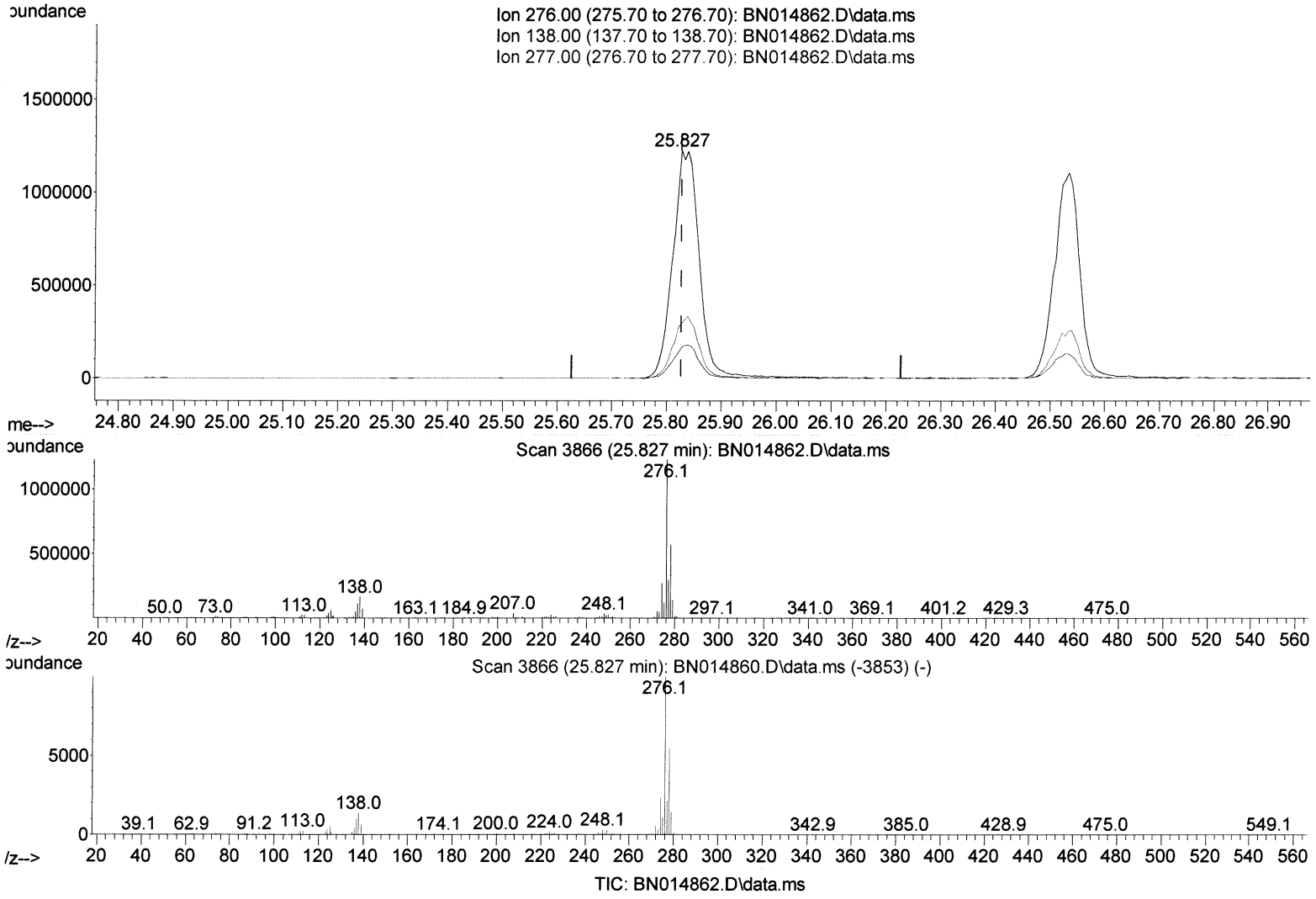
| Ion | Exp% | Act% |
|---|---|---|
| 276.00 | 100.00 | 100.00 |
| 138.00 | 12.50 | 13.44 |
| 277.00 | 24.40 | 24.00 |
| 0.00 | 0.00 | 0.00 |

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061021\
Data File : BN014862.D
Acq On : 10 Jun 2021 11:51
Operator : CG/JU
Sample : PB136905BS
Misc :
ALS Vial : 4 Sample Multiplier: 1

Quant Time: Jun 10 12:39:54 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN052521.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jun 10 11:55:23 2021
Response via : Initial Calibration

**Instrument :**
BNA_N
**ClientSampleId :**
SLCS905

**Manual Integrations APPROVED**
mohammad
6/11/2021 2:13:03 PM




(94) Indeno(1,2,3-cd)pyrene

25.827min (-0.000) 37.73 ng/ul m JG 6/14/21

response 4028730

| Ion | Exp% | Act% |
|---|---|---|
| 276.00 | 100.00 | 100.00 |
| 138.00 | 12.50 | 13.44 |
| 277.00 | 24.40 | 24.00 |
| 0.00 | 0.00 | 0.00 |

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN061021\
Data File : BN014862.D
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Operator : CG/JU
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Misc :
ALS Vial : 4 Sample Multiplier: 1

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QLast Update : Thu Jun 10 11:55:23 2021
Response via : Initial Calibration

**Instrument :** BNA_N
**ClientSampleId :** SLCS905

**Manual Integrations APPROVED**
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| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|---|---|---|---|---|---|---|
| **Internal Standards** | | | | | | |
| 1) 1,4-Dichlorobenzene-d4 | 7.693 | 152 | 197977 | 20.000 | ng/ul | 0.00 |
| 20) Naphthalene-d8 | 10.481 | 136 | 778369 | 20.000 | ng/ul | 0.00 |
| 38) Acenaphthene-d10 | 14.345 | 164 | 555202 | 20.000 | ng/ul | 0.00 |
| 64) Phenanthrene-d10 | 17.098 | 188 | 1249160 | 20.000 | ng/ul | # 0.00 |
| 79) Chrysene-d12 | 21.304 | 240 | 1377956 | 20.000 | ng/ul | 0.00 |
| 88) Perylene-d12 | 23.556 | 264 | 1434873 | 20.000 | ng/ul | 0.00 |
| **System Monitoring Compounds** | | | | | | |
| 3) 1,4-Dioxane-d8 | 3.210 | 96 | 28462 | 5.854 | ng/uL | 0.00 |
| 4) Pyridine-d5 | 3.610 | 84 | 328973 | 26.869 | ng/ul | 0.00 |
| 7) Phenol-d5 | 6.869 | 99 | 615832 | 31.768 | ng/ul | 0.00 |
| 9) Bis-(2-Chloroethyl)eth... | 7.034 | 67 | 293502 | 30.448 | ng/ul | 0.00 |
| 11) 2-Chlorophenol-d4 | 7.228 | 132 | 448733 | 37.376 | ng/ul | 0.00 |
| 15) 4-Methylphenol-d8 | 8.404 | 113 | 498648 | 33.085 | ng/ul | 0.00 |
| 21) Nitrobenzene-d5 | 8.851 | 128 | 231128 | 39.419 | ng/ul | 0.00 |
| 24) 2-Nitrophenol-d4 | 9.563 | 143 | 248006 | 39.351 | ng/ul | 0.00 |
| 28) 2,4-Dichlorophenol-d3 | 10.104 | 165 | 528586 | 36.785 | ng/ul | 0.00 |
| 31) 4-Chloroaniline-d4 | 10.616 | 131 | 568884 | 33.580 | ng/ul | 0.00 |
| 46) Dimethylphthalate-d6 | 13.757 | 166 | 1677843 | 38.038 | ng/ul | 0.00 |
| 49) Acenaphthylene-d8 | 14.033 | 160 | 1909032 | 37.427 | ng/ul | 0.00 |
| 54) 4-Nitrophenol-d4 | 14.557 | 143 | 256931 | 36.666 | ng/ul | 0.00 |
| 60) Fluorene-d10 | 15.345 | 176 | 1401909 | 36.886 | ng/ul | 0.00 |
| 65) 4,6-Dinitro-2-methylph... | 15.469 | 200 | 272601 | 39.899 | ng/ul | 0.00 |
| 73) Anthracene-d10 | 17.198 | 188 | 2188212 | 37.499 | ng/ul | 0.00 |
| 81) Pyrene-d10 | 19.504 | 212 | 2667238 | 41.052 | ng/ul | 0.00 |
| 92) Benzo(a)pyrene-d12 | 23.415 | 264 | 3238875 | 42.377 | ng/ul | 0.00 |
| **Target Compounds** | | | | | | Qvalue |
| 2) 1,4-Dioxane | 3.246 | 88 | 62562 | 10.961 | ng/uL# | 85 |
| 5) Pyridine | 3.628 | 79 | 317518 | 24.748 | ng/ul | 95 |
| 6) Benzaldehyde | 6.840 | 77 | 343780 | 32.932 | ng/ul | 92 |
| 8) Phenol | 6.899 | 94 | 551069 | 27.627 | ng/ul | 97 |
| 10) Bis(2-Chloroethyl)ether | 7.122 | 93 | 419754 | 27.265 | ng/ul | 98 |
| 12) 2-Chlorophenol | 7.263 | 128 | 413194 | 32.882 | ng/ul# | 89 |
| 13) 2-Methylphenol | 8.140 | 108 | 416474 | 28.431 | ng/ul | 98 |
| 14) 2,2'-oxybis(1-Chloropr... | 8.216 | 45 | 262899 | 27.069 | ng/ul | 93 |
| 16) Acetophenone | 8.510 | 105 | 737068 | 29.542 | ng/ul | 89 |
| 17) N-Nitroso-di-n-propyla... | 8.498 | 70 | 330341 | 27.488 | ng/ul# | 93 |
| 18) 4-Methylphenol | 8.463 | 108 | 462828 | 28.969 | ng/ul | 96 |
| 19) Hexachloroethane | 8.763 | 117 | 200211 | 31.823 | ng/ul# | 88 |
| 22) Nitrobenzene | 8.887 | 77 | 530533 | 28.993 | ng/ul | 95 |
| 23) Isophorone | 9.410 | 82 | 986620 | 28.394 | ng/ul | 96 |
| 25) 2-Nitrophenol | 9.598 | 139 | 228815 | 34.164 | ng/ul# | 89 |
| 26) 2,4-Dimethylphenol | 9.663 | 107 | 485752 | 23.561 | ng/ul | 94 |
| 27) Bis(2-Chloroethoxy)met... | 9.898 | 93 | 594566 | 28.161 | ng/ul# | 97 |
| 29) 2,4-Dichlorophenol | 10.128 | 162 | 452973 | 32.868 | ng/ul | 94 |
| 30) Naphthalene | 10.528 | 128 | 1331291 | 31.468 | ng/ul# | 96 |
| 32) 4-Chloroaniline | 10.640 | 127 | 517462 | 28.701 | ng/ul | 95 |
| 33) Hexachlorobutadiene | 10.822 | 225 | 406473 | 32.985 | ng/ul | 89 |
| 34) Caprolactam | 11.398 | 113 | 152652m | 34.284 | ng/ul | 100 |
| 35) 4-Chloro-3-methylphenol | 11.775 | 107 | 630874 | 31.589 | ng/ul | 100 |

> JU 6/14/21

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Data File : BN014862.D
Acq On : 10 Jun 2021 11:51
Operator : CG/JU
Sample : PB136905BS
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
SLCS905

Manual Integrations APPROVED
mohammad
6/11/2021 2:13:03 PM

Quant Time: Jun 10 12:39:54 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-BN052521.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jun 10 11:55:23 2021
Response via : Initial Calibration

| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|---|---|---|---|---|---|---|
| 36) 2-Methylnaphthalene | 12.151 | 142 | 994240 | 31.528 | ng/ul | 99 |
| 37) 1-Methylnaphthalene | 12.369 | 142 | 968819 | 31.248 | ng/ul# | 94 |
| 39) 1,2,4,5-Tetrachloroben... | 12.522 | 216 | 677666 | 32.282 | ng/ul | 95 |
| 40) Hexachlorocyclopentadiene | 12.504 | 237 | 307166 | 22.088 | ng/ul | 98 |
| 41) 2,4,6-Trichlorophenol | 12.769 | 196 | 421905 | 35.791 | ng/ul | 95 |
| 42) 2,4,5-Trichlorophenol | 12.839 | 196 | 454956 | 35.364 | ng/ul | 92 |
| 43) 1,1'-Biphenyl | 13.175 | 154 | 1347180 | 33.919 | ng/ul# | 95 |
| 44) 2-Chloronaphthalene | 13.210 | 162 | 1100221 | 34.572 | ng/ul | 96 |
| 45) 2-Nitroaniline | 13.416 | 65 | 312965 | 32.287 | ng/ul | 91 |
| 47) Dimethylphthalate | 13.804 | 163 | 1519775 | 34.835 | ng/ul | 99 |
| 48) 2,6-Dinitrotoluene | 13.922 | 165 | 312721 | 38.572 | ng/ul | 97 |
| 50) Acenaphthylene | 14.069 | 152 | 1591571 | 33.938 | ng/ul | 96 |
| 51) 3-Nitroaniline | 14.251 | 138 | 273462 | 34.762 | ng/ul | 84 |
| 52) Acenaphthene | 14.410 | 153 | 1141878 | 33.604 | ng/ul | 98 |
| 53) 2,4-Dinitrophenol | 14.457 | 184 | 151956 | 25.815 | ng/ul | 97 |
| 55) 4-Nitrophenol | 14.575 | 109 | 333319 | 24.813 | ng/ul# | 88 |
| 56) Dibenzofuran | 14.745 | 168 | 1679260 | 33.221 | ng/ul# | 87 |
| 57) 2,4-Dinitrotoluene | 14.716 | 165 | 478753 | 37.054 | ng/ul | 94 |
| 58) 2,3,4,6-Tetrachlorophenol | 14.980 | 232 | 436291 | 32.968 | ng/ul# | 93 |
| 59) Diethylphthalate | 15.180 | 149 | 1530131 | 36.482 | ng/ul | 96 |
| 61) Fluorene | 15.398 | 166 | 1278476 | 33.618 | ng/ul | 96 |
| 62) 4-Chlorophenyl-phenyle... | 15.398 | 204 | 719486 | 31.978 | ng/ul# | 85 |
| 63) 4-Nitroaniline | 15.422 | 138 | 279453 | 34.972 | ng/ul | 95 |
| 66) 4,6-Dinitro-2-methylph... | 15.480 | 198 | 255907 | 38.286 | ng/ul# | 93 |
| 67) N-Nitrosodiphenylamine | 15.616 | 169 | 1210191 | 35.969 | ng/ul | 99 |
| 68) 4-Bromophenyl-phenylether | 16.292 | 248 | 574894 | 35.343 | ng/ul# | 82 |
| 69) Hexachlorobenzene | 16.410 | 284 | 701255 | 34.461 | ng/ul | 98 |
| 70) Atrazine | 16.569 | 200 | 498505 | 35.831 | ng/ul | 98 |
| 71) Pentachlorophenol | 16.751 | 266 | 384378 | 32.808 | ng/ul | 96 |
| 72) Phenanthrene | 17.139 | 178 | 2329804 | 35.695 | ng/ul | 98 |
| 74) Anthracene | 17.233 | 178 | 2261080m | 34.372 | ng/ul | JU 6/16/21 |
| 75) 1,2,3,4-Tetrachloroben... | 13.139 | 216 | 709598 | 34.377 | ng/uL | 95 |
| 76) Pentachlorobenzene | 14.669 | 250 | 747552 | 35.016 | ng/uL | 96 |
| 77) Carbazole | 17.504 | 167 | 2021425 | 36.461 | ng/ul | 96 |
| 78) Di-n-butylphthalate | 18.074 | 149 | 2540920 | 38.895 | ng/ul | 99 |
| 80) Fluoranthene | 19.168 | 202 | 2962717 | 38.670 | ng/ul | 96 |
| 82) Pyrene | 19.533 | 202 | 2964502 | 37.626 | ng/ul | 97 |
| 83) Butylbenzylphthalate | 20.439 | 149 | 1155354 | 43.642 | ng/ul | 95 |
| 84) 3,3'-Dichlorobenzidine | 21.221 | 252 | 949348 | 36.542 | ng/ul | 98 |
| 85) Benzo(a)anthracene | 21.286 | 228 | 3035225m | 35.337 | ng/ul | JU 6/16/21 |
| 86) Bis(2-ethylhexyl)phtha... | 21.227 | 149 | 1566629 | 41.055 | ng/ul | 99 |
| 87) Chrysene | 21.339 | 228 | 3056503 | 35.336 | ng/ul | 97 |
| 89) Di-n-octyl phthalate | 22.109 | 149 | 3096154 | 48.772 | ng/ul | 100 |
| 90) Benzo(b)fluoranthene | 22.880 | 252 | 3506679 | 40.722 | ng/ul | 95 |
| 91) Benzo(k)fluoranthene | 22.927 | 252 | 3389383m | 39.506 | ng/ul | JU 6/16/21 |
| 93) Benzo(a)pyrene | 23.462 | 252 | 3001971 | 38.521 | ng/ul | 97 |
| 94) Indeno(1,2,3-cd)pyrene | 25.827 | 276 | 4028730m | 37.731 | ng/ul | JU 6/16/21 |
| 95) Dibenzo(a,h)anthracene | 25.850 | 278 | 3331077 | 37.696 | ng/ul | 100 |
| 96) Benzo(g,h,i)perylene | 26.533 | 276 | 3515979 | 37.206 | ng/ul | 95 |

(#) = qualifier out of range (m) = manual integration (+) = signals summed

**Instrument :**
BNA_N
**ClientSampleId :**
SLCS905

**Manual Integrations APPROVED**

mohammad
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