

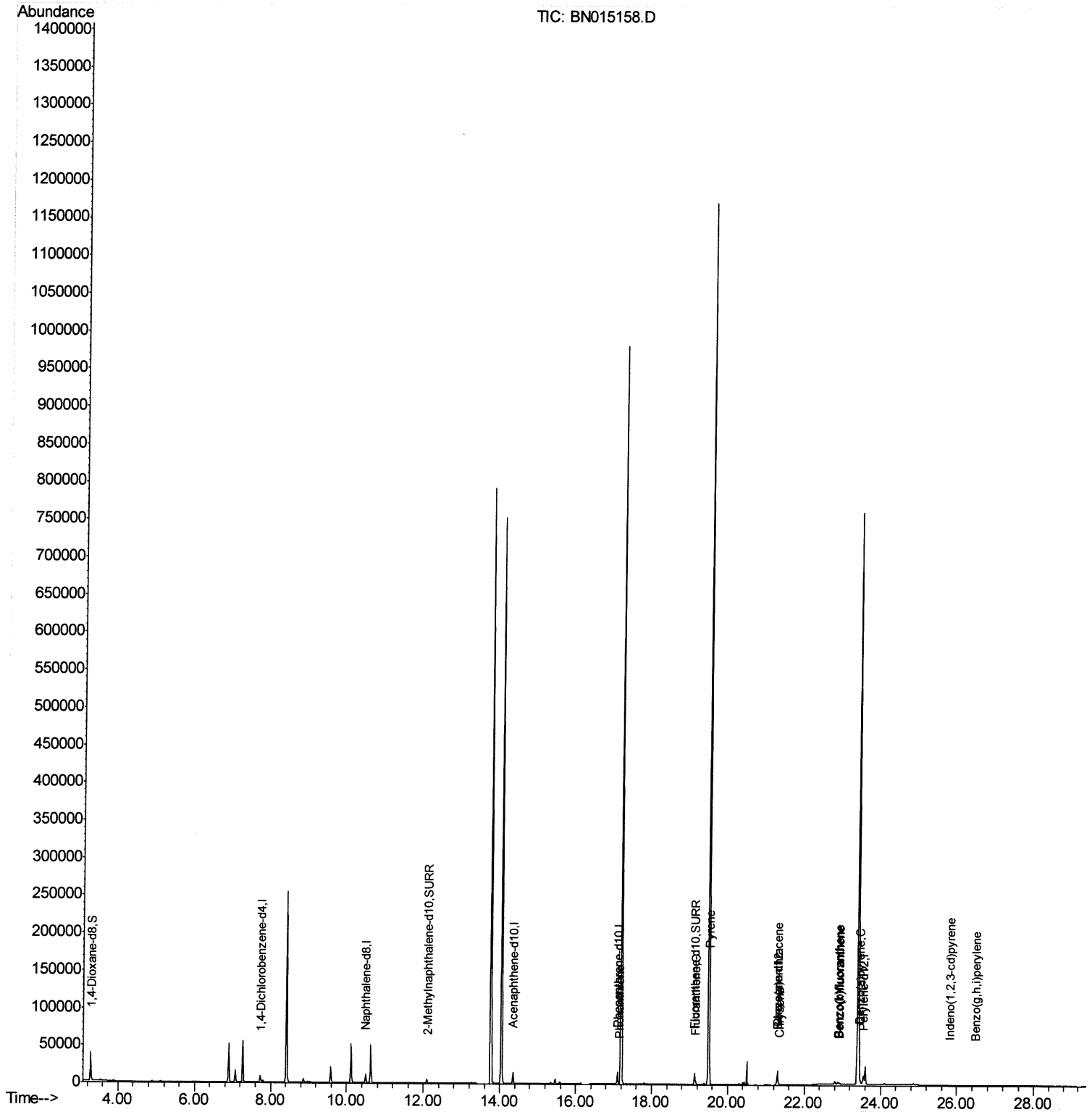
Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN062721\  
Data File : BN015158.D  
Acq On : 27 Jun 2021 18:03  
Operator : CG/JU  
Sample : M2704-10  
Misc :  
ALS Vial : 42 Sample Multiplier: 1

Instrument :  
BNA\_N  
ClientSampleId :  
BGDB9

Quant Time: Jun 28 05:41:32 2021  
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SFAM-EPA-SIM-BN062121.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
QLast Update : Mon Jun 28 05:16:57 2021  
Response via : Initial Calibration

Manual Integrations  
APPROVED

mohammad  
6/28/2021 1:28:57 PM



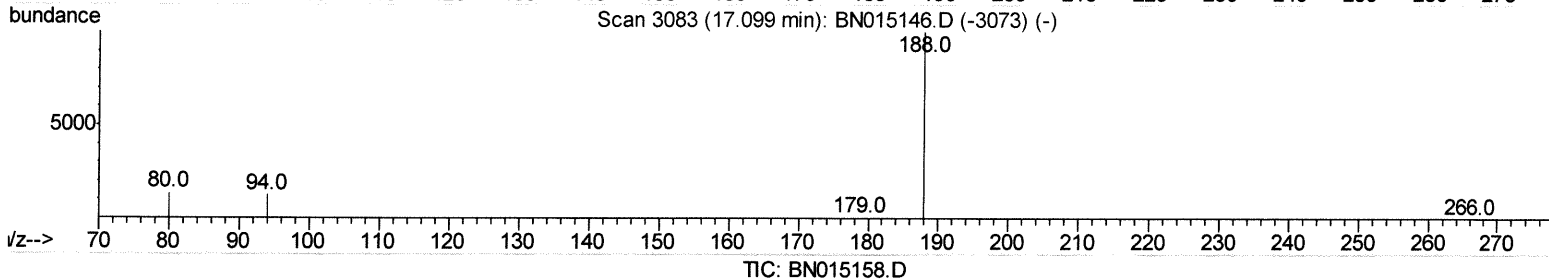
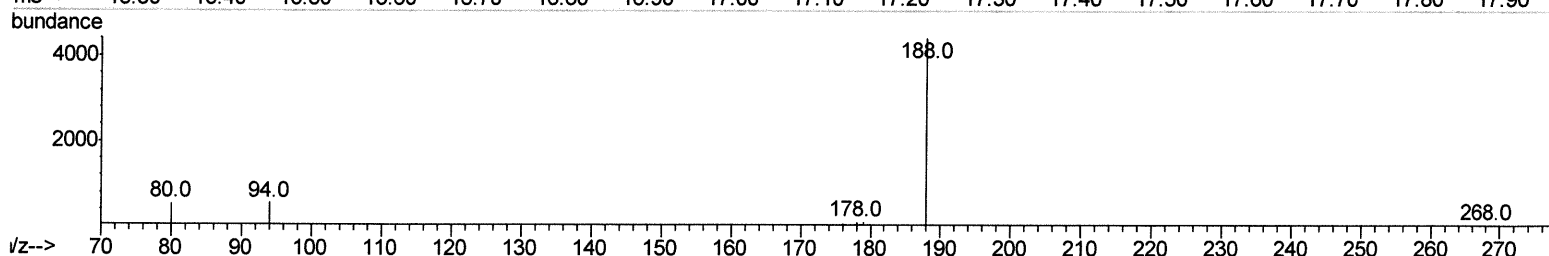
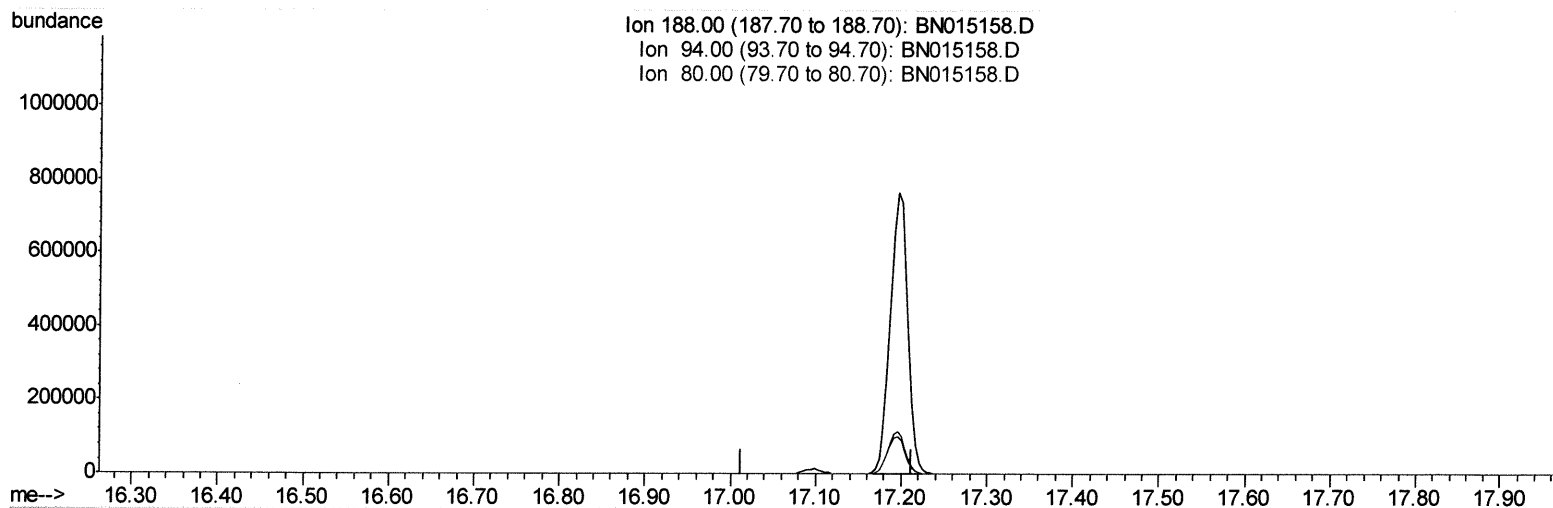
Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN062721\  
 Data File : BN015158.D  
 Acq On : 27 Jun 2021 18:03  
 Operator : CG/JU  
 Sample : M2704-10  
 Misc :  
 ALS Vial : 42 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 BGDB9

Quant Time: Jun 28 05:34:52 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SFAM-EPA-SIM-BN062121.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Jun 28 05:16:57 2021  
 Response via : Initial Calibration

Manual Integrations  
 APPROVED

mohammad  
 6/28/2021 1:28:57 PM



(13) Phenanthrene-d10 (I)

17.113min (-17.113) 0.00ng/ul

response 0

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 188.00 | 100   | 0.00  |
| 94.00  | 14.10 | 0.00# |
| 80.00  | 14.40 | 0.00# |
| 0.00   | 0.00  | 0.00  |

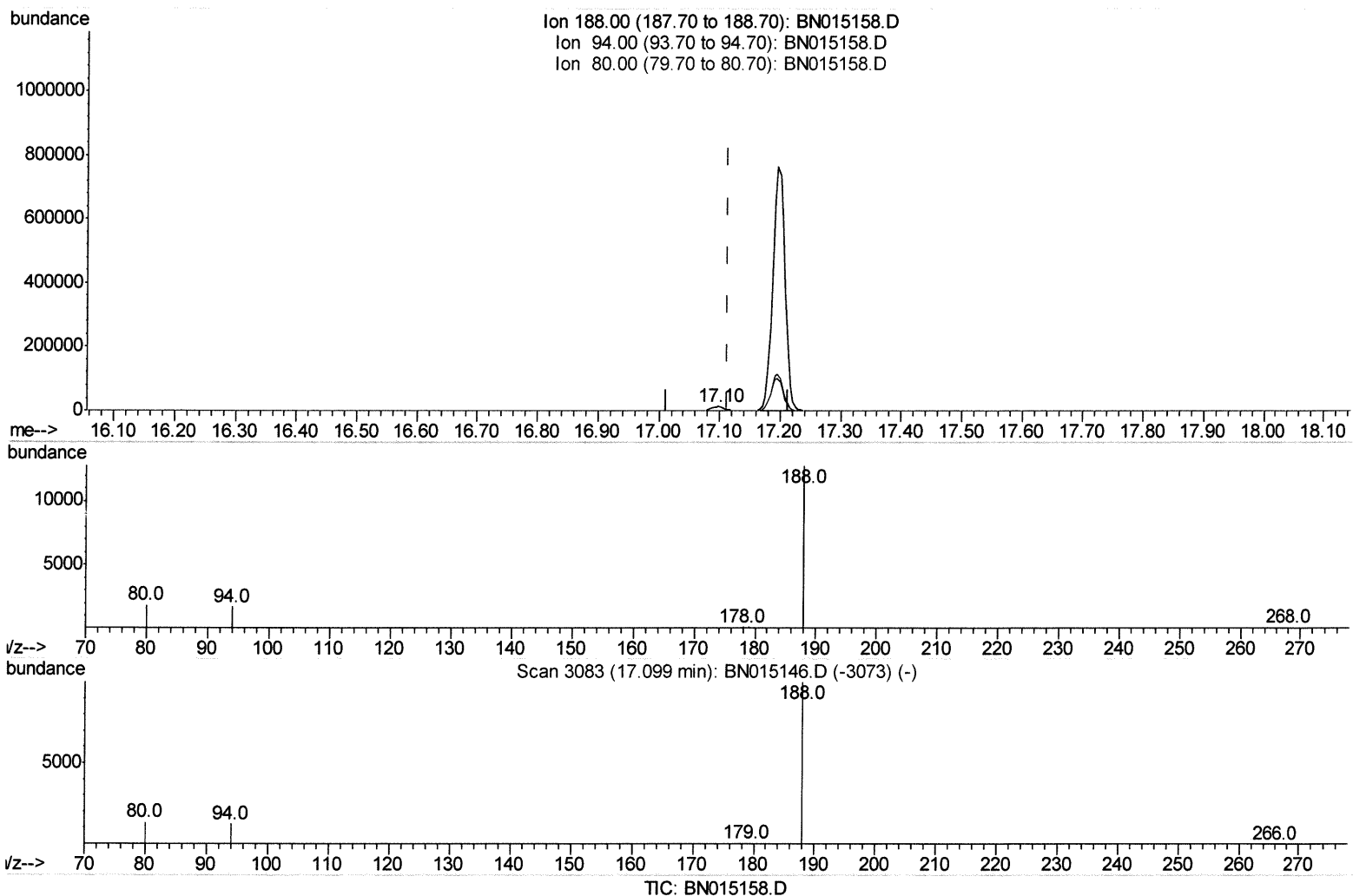
Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN062721\  
 Data File : BN015158.D  
 Acq On : 27 Jun 2021 18:03  
 Operator : CG/JU  
 Sample : M2704-10  
 Misc :  
 ALS Vial : 42 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 BGDB9

Quant Time: Jun 28 05:34:52 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SFAM-EPA-SIM-BN062121.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Jun 28 05:16:57 2021  
 Response via : Initial Calibration

Manual Integrations  
 APPROVED

mohammad  
 6/28/2021 1:28:57 PM



(13) Phenanthrene-d10 (I)

17.098min (-0.015) 0.40ng/ul m

*JU 6/30/21*

response 17728

| Ion    | Exp%  | Act%  |
|--------|-------|-------|
| 188.00 | 100   | 100   |
| 94.00  | 14.10 | 13.17 |
| 80.00  | 14.40 | 14.05 |
| 0.00   | 0.00  | 0.00  |

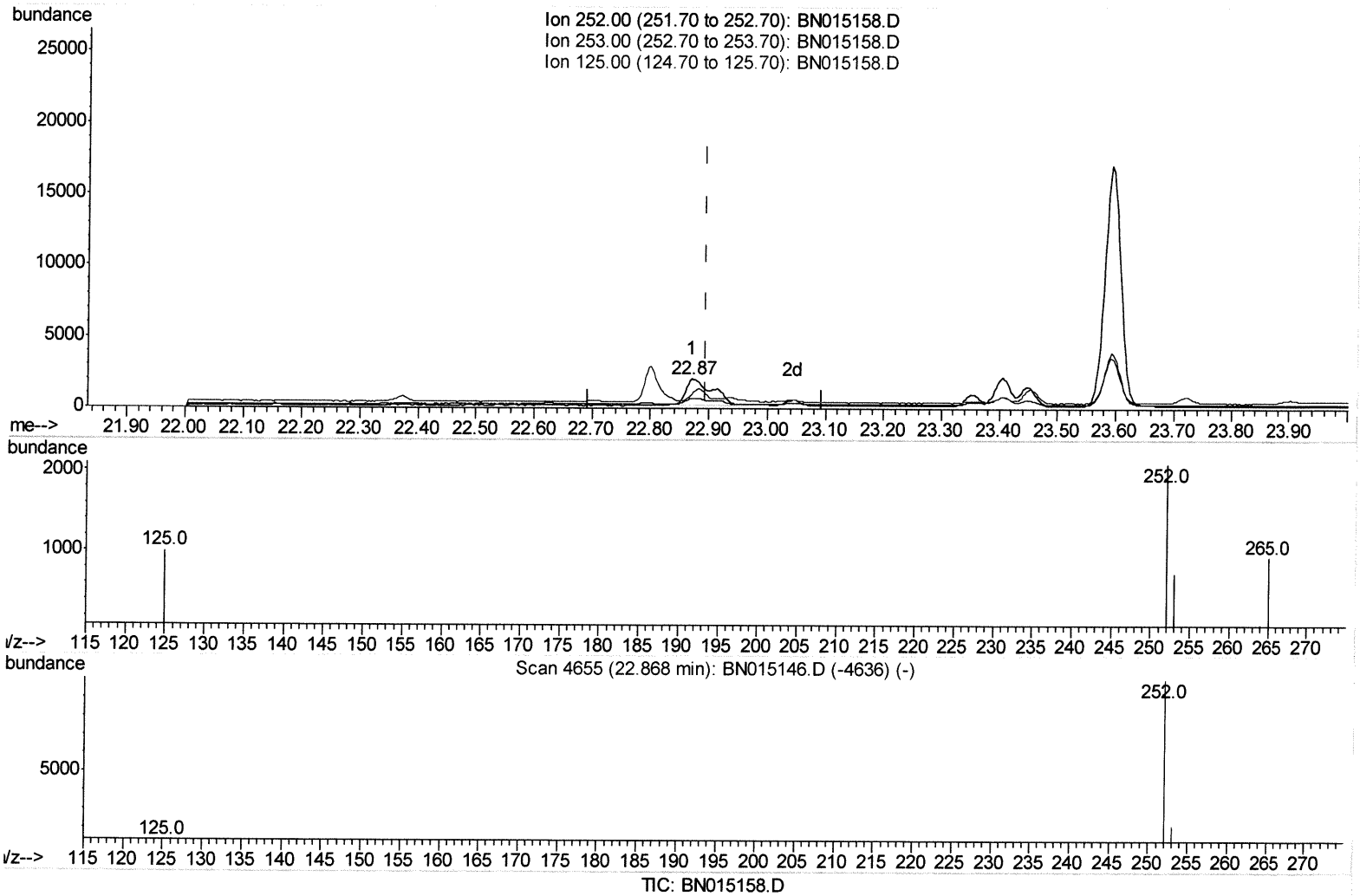
Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN062721\  
 Data File : BN015158.D  
 Acq On : 27 Jun 2021 18:03  
 Operator : CG/JU  
 Sample : M2704-10  
 Misc :  
 ALS Vial : 42 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 BGDB9

Quant Time: Jun 28 05:38:45 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SFAM-EPA-SIM-BN062121.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Jun 28 05:16:57 2021  
 Response via : Initial Calibration

Manual Integrations  
 APPROVED

mohammad  
 6/28/2021 1:28:57 PM



(24) Benzo(b)fluoranthene

22.871min (-0.024) 0.08ng/ul

response 5948

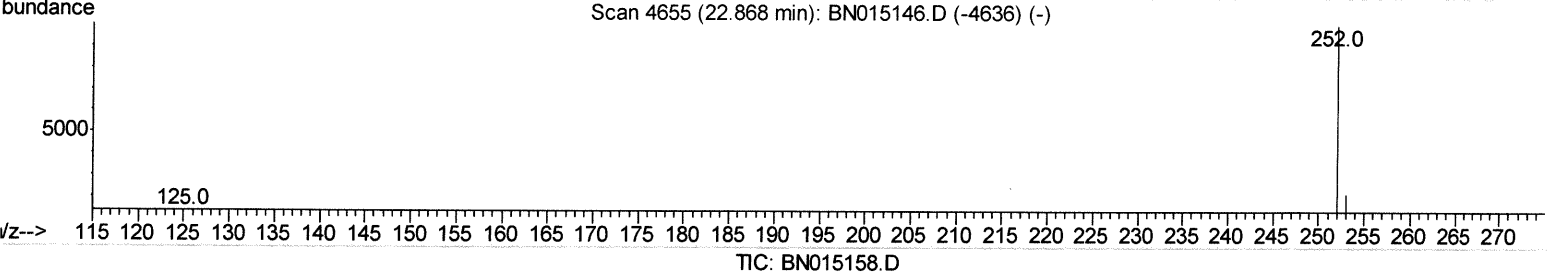
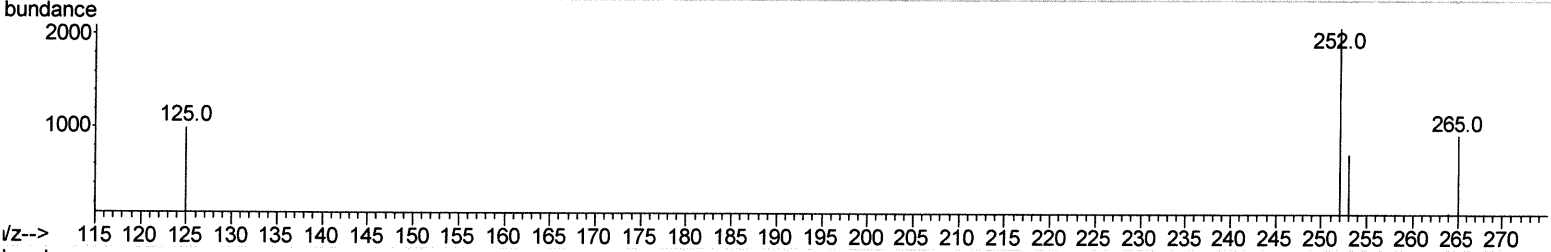
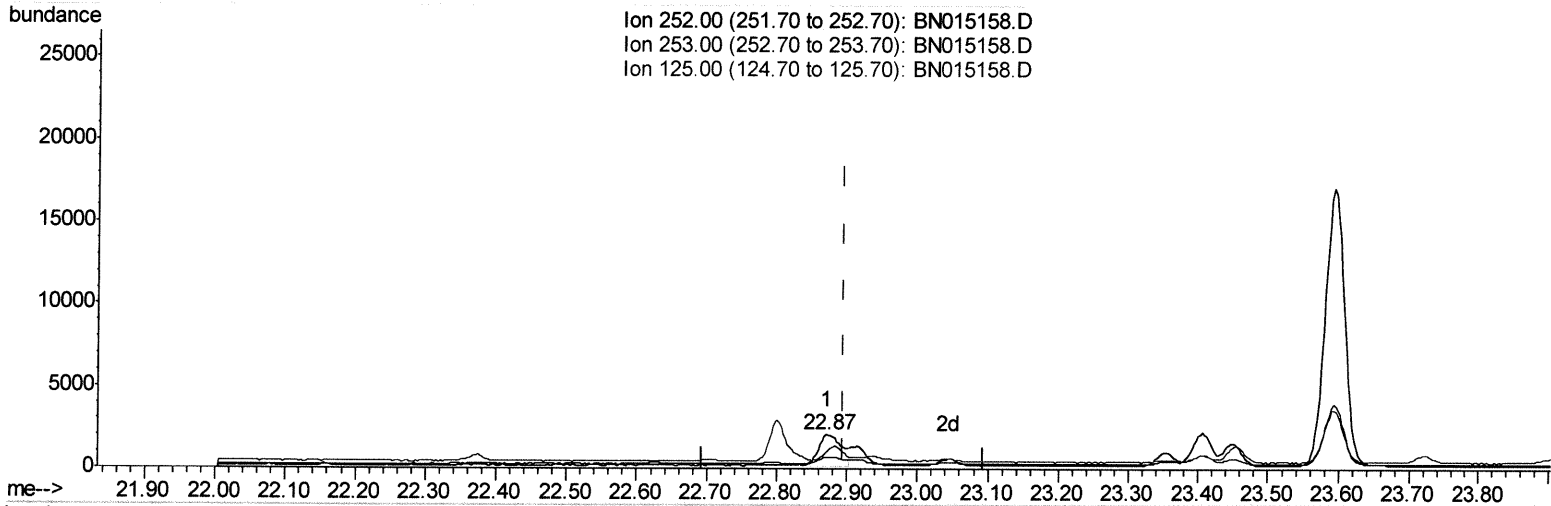
| Ion    | Exp%  | Act%   |
|--------|-------|--------|
| 252.00 | 100   | 100    |
| 253.00 | 22.80 | 35.01  |
| 125.00 | 14.20 | 47.65# |
| 0.00   | 0.00  | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN062721\  
 Data File : BN015158.D  
 Acq On : 27 Jun 2021 18:03  
 Operator : CG/JU  
 Sample : M2704-10  
 Misc :  
 ALS Vial : 42 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampled :  
 BGDB9

Manual Integrations  
 APPROVED  
 mohammad  
 6/28/2021 1:28:57 PM

Quant Time: Jun 28 05:38:45 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SFAM-EPA-SIM-BN062121.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Jun 28 05:16:57 2021  
 Response via : Initial Calibration



(24) Benzo(b)fluoranthene

22.871min (-0.024) 0.06ng/ul m

response 4484

| Ion    | Exp%  | Act%   |
|--------|-------|--------|
| 252.00 | 100   | 100    |
| 253.00 | 22.80 | 35.01  |
| 125.00 | 14.20 | 47.65# |
| 0.00   | 0.00  | 0.00   |

*ju 6/30/21*

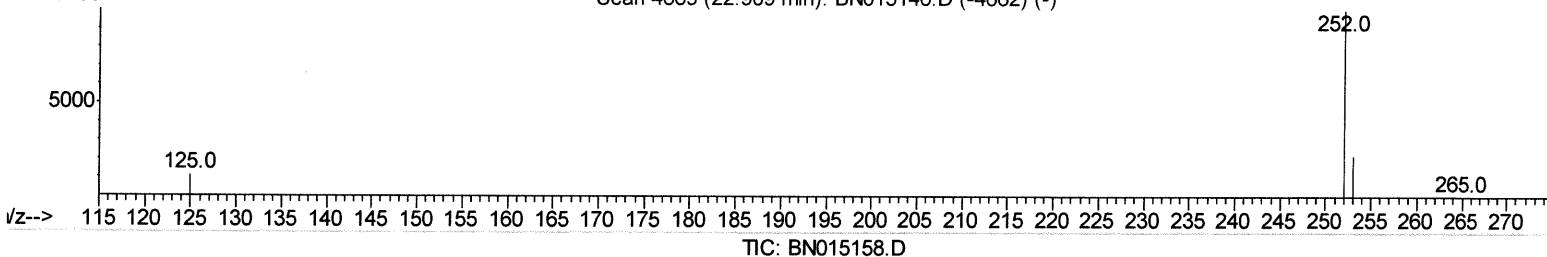
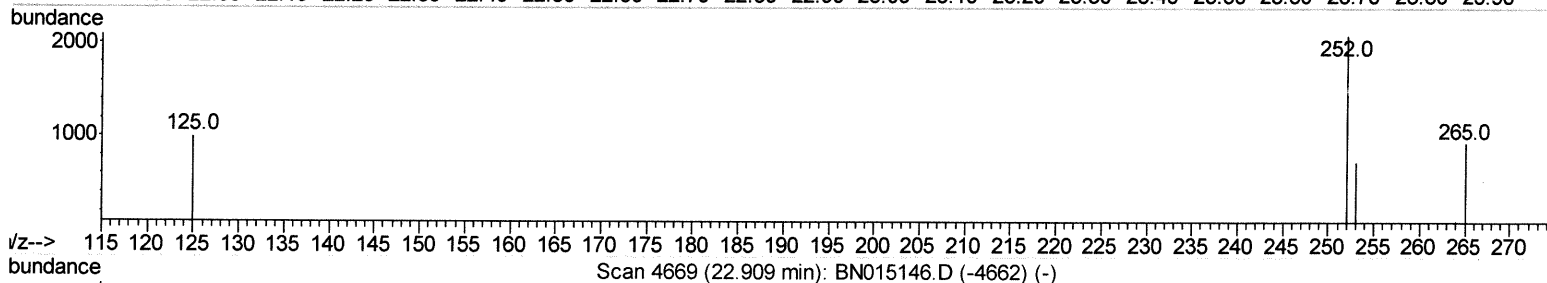
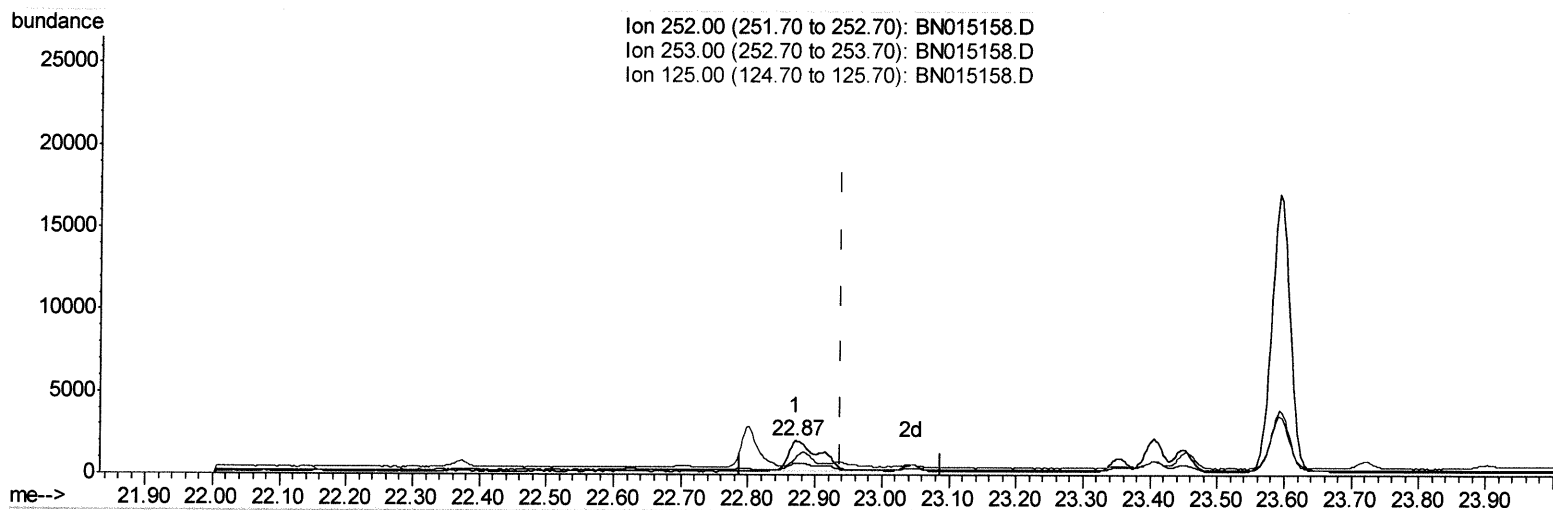
Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN062721\  
 Data File : BN015158.D  
 Acq On : 27 Jun 2021 18:03  
 Operator : CG/JU  
 Sample : M2704-10  
 Misc :  
 ALS Vial : 42 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 BGDB9

Manual Integrations  
 APPROVED

mohammad  
 6/28/2021 1:28:57 PM

Quant Time: Jun 28 05:39:35 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SFAM-EPA-SIM-BN062121.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Jun 28 05:16:57 2021  
 Response via : Initial Calibration



(25) Benzo(k)fluoranthene

22.871min (-0.067) 0.07ng/ul

response 5948

| Ion    | Exp%  | Act%   |
|--------|-------|--------|
| 252.00 | 100   | 100    |
| 253.00 | 22.60 | 35.01# |
| 125.00 | 14.40 | 47.65# |
| 0.00   | 0.00  | 0.00   |

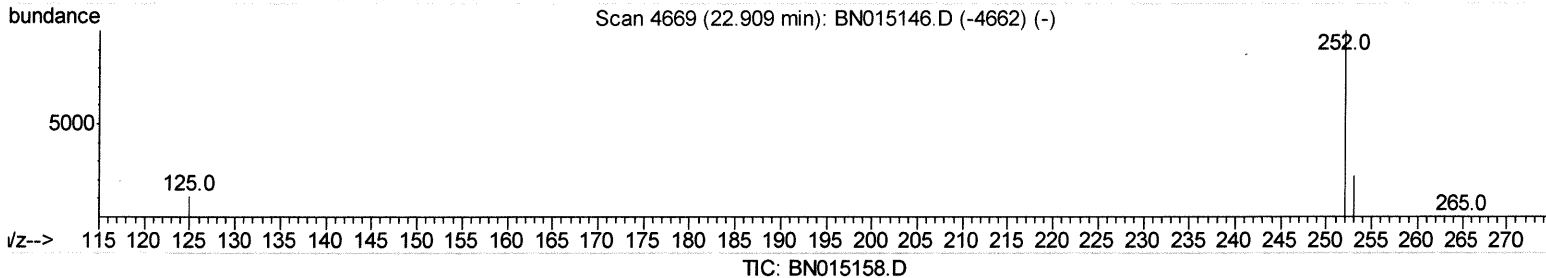
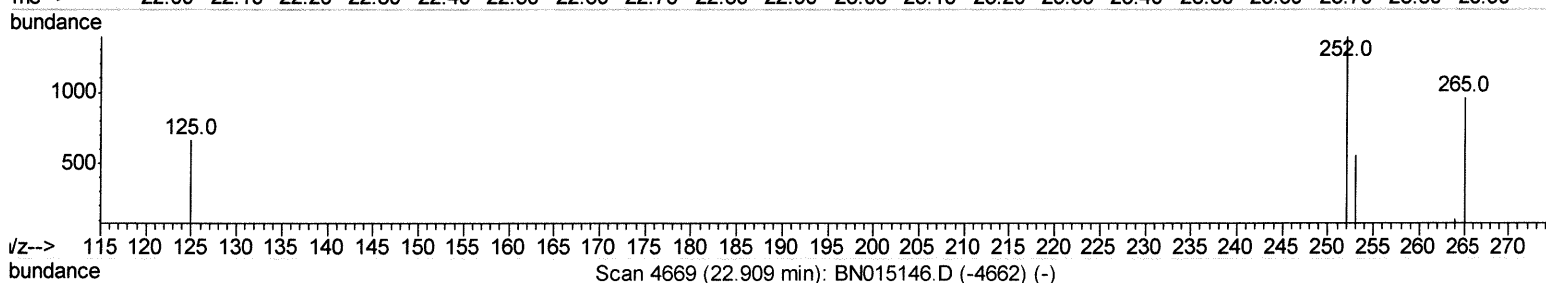
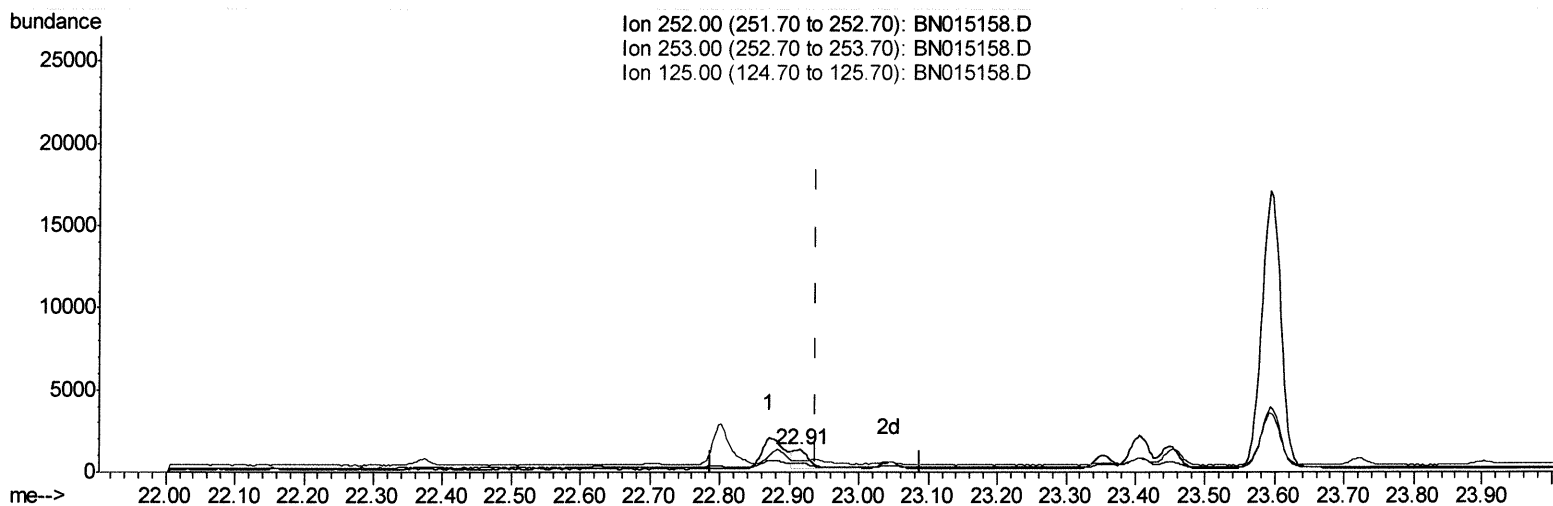
Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN062721\  
 Data File : BN015158.D  
 Acq On : 27 Jun 2021 18:03  
 Operator : CG/JU  
 Sample : M2704-10  
 Misc :  
 ALS Vial : 42 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 BGDB9

Manual Integrations  
 APPROVED

mohammad  
 6/28/2021 1:28:57 PM

Quant Time: Jun 28 05:39:35 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SFAM-EPA-SIM-BN062121.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Jun 28 05:16:57 2021  
 Response via : Initial Calibration



(25) Benzo(k)fluoranthene

22.915min (-0.024) 0.02ng/ul m

response 1823

| Ion    | Exp%  | Act%   |
|--------|-------|--------|
| 252.00 | 100   | 100    |
| 253.00 | 22.60 | 40.19# |
| 125.00 | 14.40 | 47.88# |
| 0.00   | 0.00  | 0.00   |

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN062721\  
 Data File : BN015158.D  
 Acq On : 27 Jun 2021 18:03  
 Operator : CG/JU  
 Sample : M2704-10  
 Misc :  
 ALS Vial : 42 Sample Multiplier: 1

Instrument :  
 BNA\_N  
 ClientSampleId :  
 BGDB9

Quant Time: Jun 28 05:41:32 2021  
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SFAM-EPA-SIM-BN062121.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Jun 28 05:16:57 2021  
 Response via : Initial Calibration

Manual Integrations  
 APPROVED

mohammad  
 6/28/2021 1:28:57 PM

| Internal Standards        | R.T.  | QIon | Response | Conc | Units  | Dev(Min) |
|---------------------------|-------|------|----------|------|--------|----------|
| 1) 1,4-Dichlorobenzene-d4 | 7.73  | 152  | 4163     | 0.40 | ng/ul  | -0.01    |
| 4) Naphthalene-d8         | 10.50 | 136  | 15190    | 0.40 | ng/ul  | -0.01    |
| 9) Acenaphthene-d10       | 14.35 | 164  | 8551     | 0.40 | ng/ul  | -0.01    |
| 13) Phenanthrene-d10      | 17.10 | 188  | 17728m>  | 0.40 | ng/ul> | -0.01    |
| 17) Chrysene-d12          | 21.30 | 240  | 15272    | 0.40 | ng/ul  | -0.01    |
| 23) Perylene-d12          | 23.54 | 264  | 15133    | 0.40 | ng/ul  | -0.03    |

| System Monitoring Compounds | R.T.  | QIon | Response | Conc | Units | Dev(Min) |
|-----------------------------|-------|------|----------|------|-------|----------|
| 3) 1,4-Dioxane-d8           | 3.29  | 96   | 23136    | 4.93 | ng/ul | -0.01    |
| 6) 2-Methylnaphthalene-d10  | 12.09 | 152  | 7158     | 0.30 | ng/ul | -0.01    |
| 18) Fluoranthene-d10        | 19.13 | 212  | 15667    | 0.31 | ng/ul | -0.02    |

| Target Compounds           | R.T.  | QIon | Response | Conc  | Units  | Ovalue |
|----------------------------|-------|------|----------|-------|--------|--------|
| 15) Phenanthrene           | 17.14 | 178  | 2014     | 0.031 | ng/ul# | 90     |
| 19) Fluoranthene           | 19.16 | 202  | 4819     | 0.068 | ng/ul# | 92     |
| 20) Pyrene                 | 19.53 | 202  | 3898     | 0.055 | ng/ul# | 90     |
| 21) Benzo(a)anthracene     | 21.28 | 228  | 3320     | 0.059 | ng/ul# | 88     |
| 22) Chrysene               | 21.33 | 228  | 3276     | 0.048 | ng/ul# | 91     |
| 24) Benzo(b)fluoranthene   | 22.87 | 252  | 4484m{   | 0.063 | ng/ul{ | 91     |
| 25) Benzo(k)fluoranthene   | 22.91 | 252  | 1823m{   | 0.023 | ng/ul{ |        |
| 26) Benzo(a)pyrene         | 23.45 | 252  | 2458     | 0.041 | ng/ul# |        |
| 27) Indeno(1,2,3-cd)pyrene | 25.80 | 276  | 1638     | 0.023 | ng/ul# | 91     |
| 29) Benzo(a,h,i)perylene   | 26.49 | 276  | 1431     | 0.022 | ng/ul# | 68     |

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