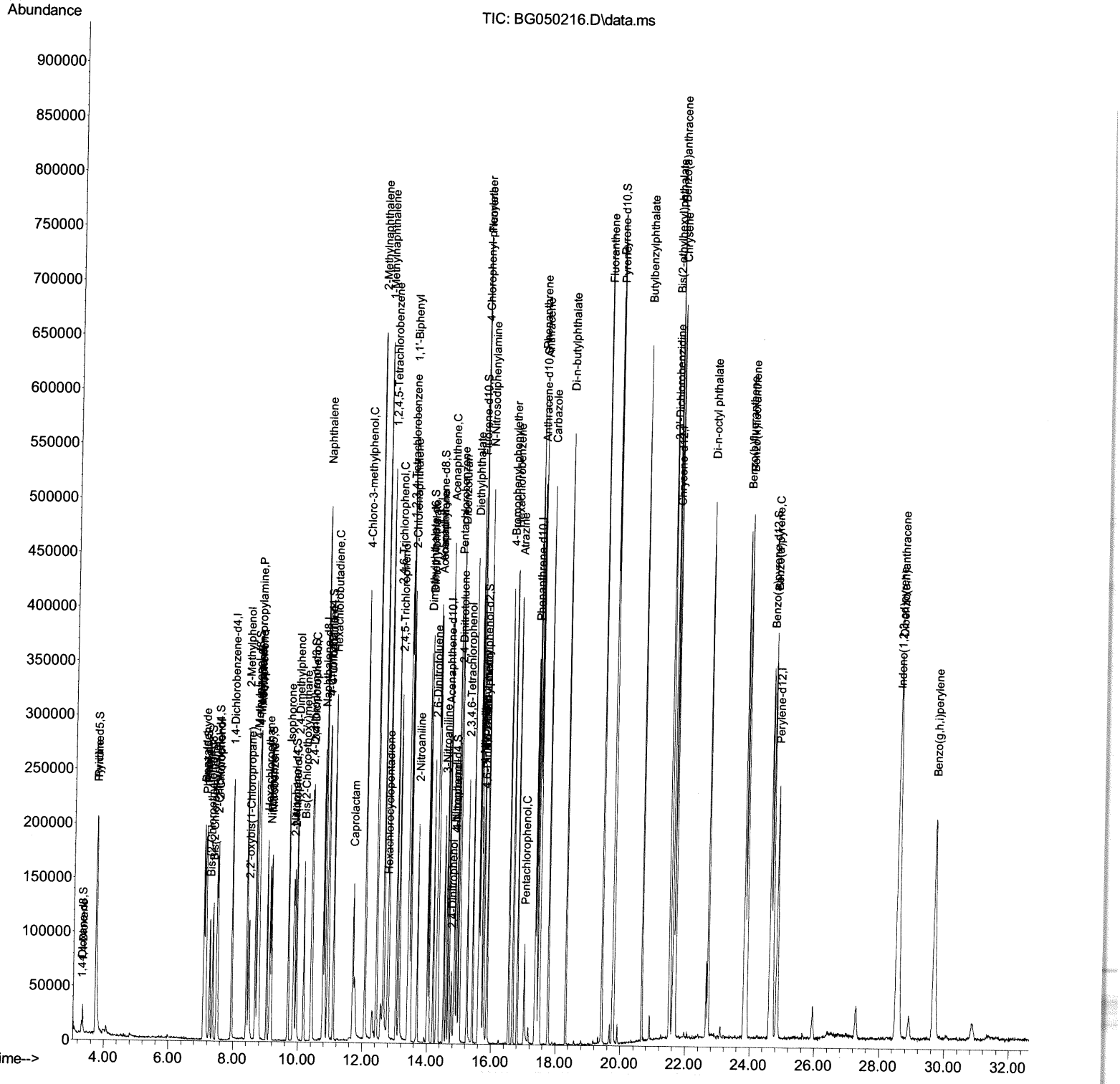
Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050216.D
Acq On : 9 Sep 2021 2:59
Operator : CG/JU
Sample : PB138845BS
Misc :
ALS Vial : 44 Sample Multiplier: 1

Quant Time: Sep 09 05:21:21 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Sep 07 18:32:38 2021
Response via : Initial Calibration

**Instrument :**
BNA_G
**ClientSampleId :**
SLCS845

**Manual Integrations APPROVED**
mohammad
9/9/2021 11:47:10 AM



Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050216.D
Acq On : 9 Sep 2021 2:59
Operator : CG/JU
Sample : PB138845BS
Misc :
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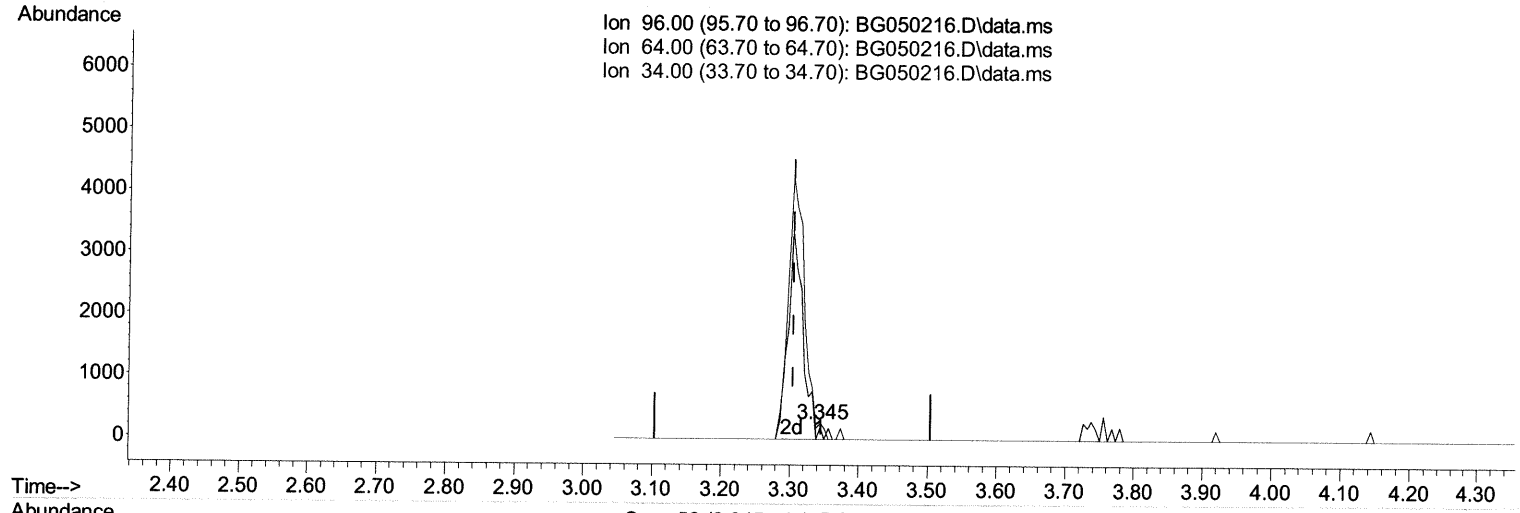
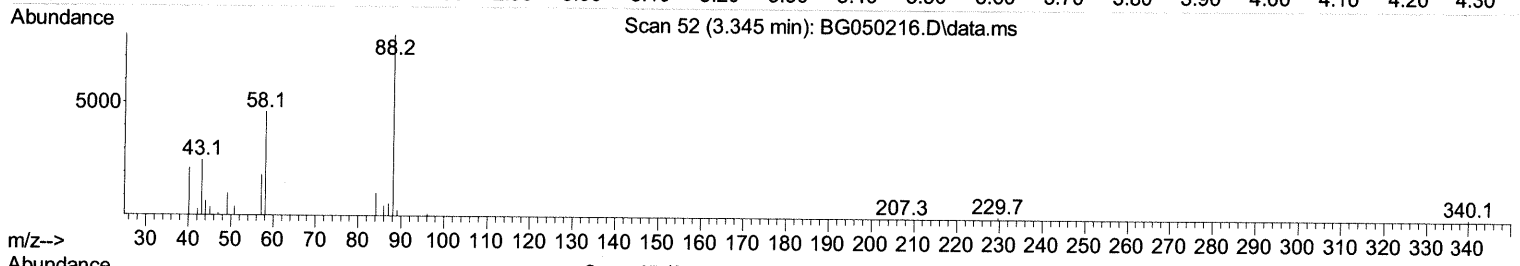
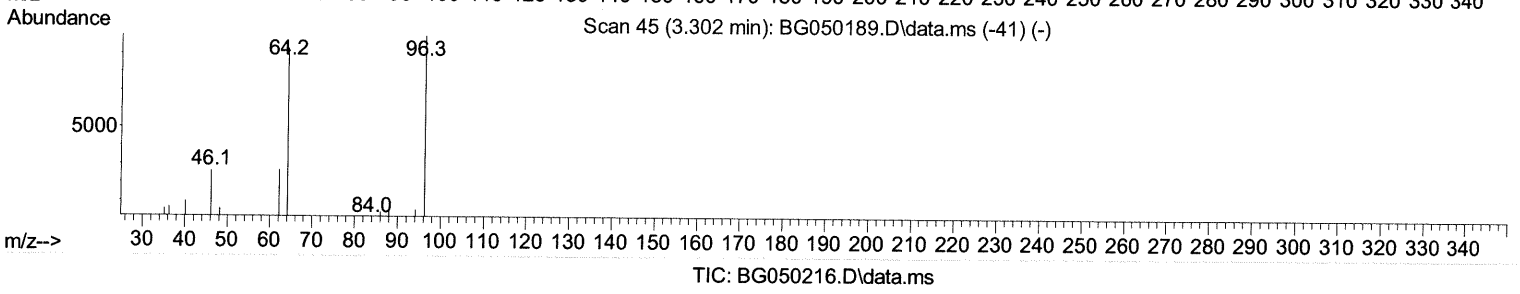
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Sep 07 18:32:38 2021
Response via : Initial Calibration

**Instrument :**
BNA_G
**ClientSampleId :**
SLCS845

**Manual Integrations APPROVED**
mohammad
9/9/2021 11:47:10 AM




TIC: BG050216.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.345min (+ 0.041) 0.11 ng/uL

response 152

| Ion | Exp% | Act% |
|---|---|---|
| 96.00 | 100.00 | 100.00 |
| 64.00 | 63.40 | 75.89 |
| 34.00 | 0.00 | 0.00 |
| 0.00 | 0.00 | 0.00 |

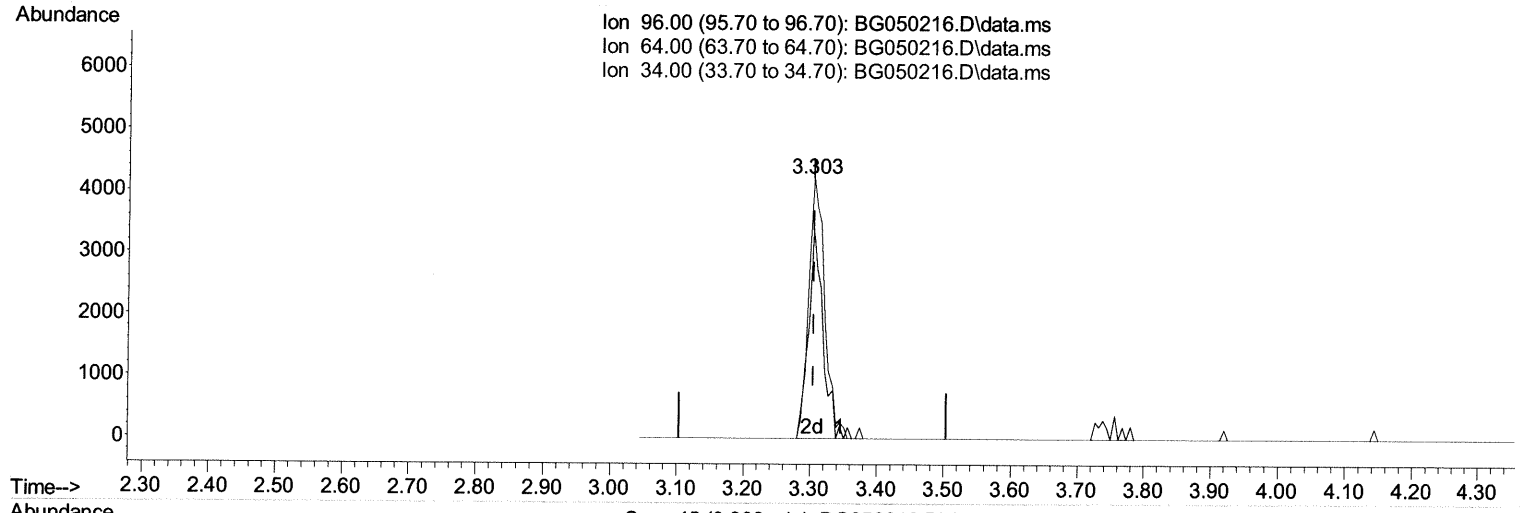
Quantitation Report (Qedit)

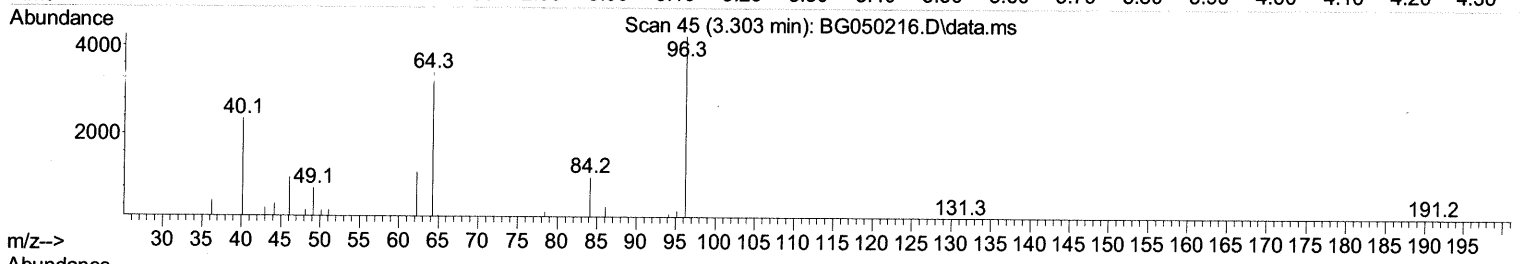
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050216.D
Acq On : 9 Sep 2021 2:59
Operator : CG/JU
Sample : PB138845BS
Misc :
ALS Vial : 44 Sample Multiplier: 1

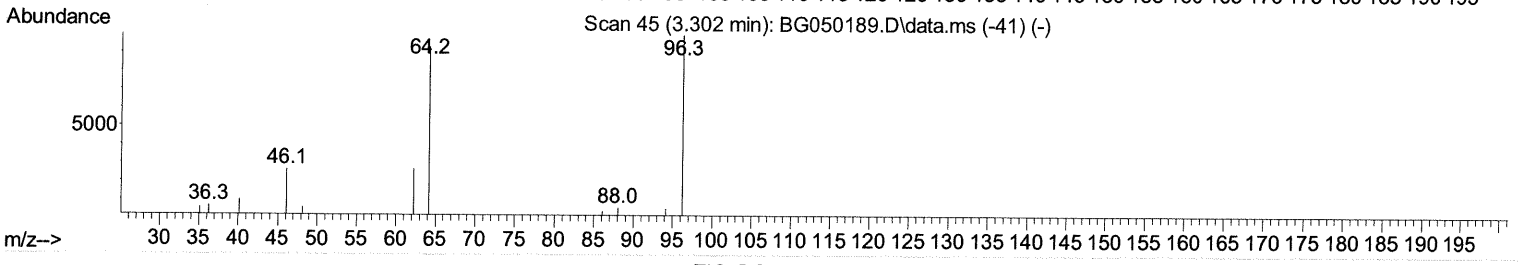
Quant Time: Sep 09 05:21:21 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Sep 07 18:32:38 2021
Response via : Initial Calibration

**Instrument :**
BNA_G
**ClientSampleId :**
SLCS845

**Manual Integrations APPROVED**
mohammad
9/9/2021 11:47:10 AM




TIC: BG050216.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.303min (-0.001) 5.07 ng/uL m 09/10/21 JU

response 7148

| Ion | Exp% | Act% |
|---|---|---|
| 96.00 | 100.00 | 100.00 |
| 64.00 | 63.40 | 80.26# |
| 34.00 | 0.00 | 0.00 |
| 0.00 | 0.00 | 0.00 |

Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050216.D
Acq On : 9 Sep 2021 2:59
Operator : CG/JU
Sample : PB138845BS
Misc :
ALS Vial : 44 Sample Multiplier: 1

Quant Time: Sep 09 05:21:21 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Sep 07 18:32:38 2021
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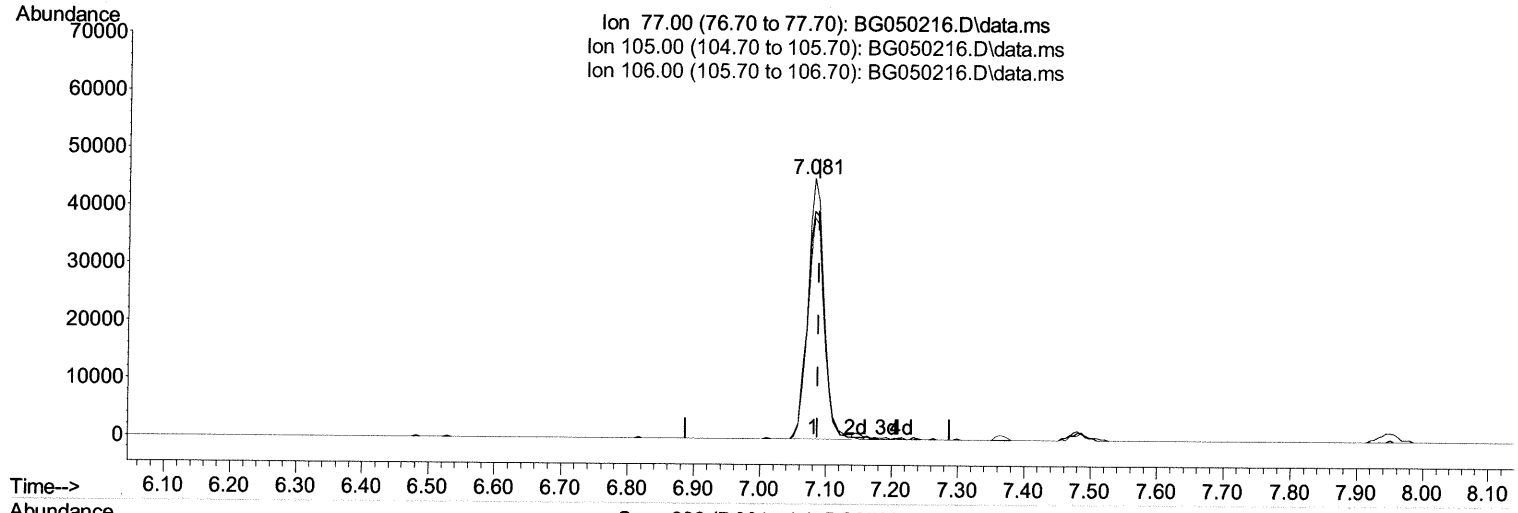
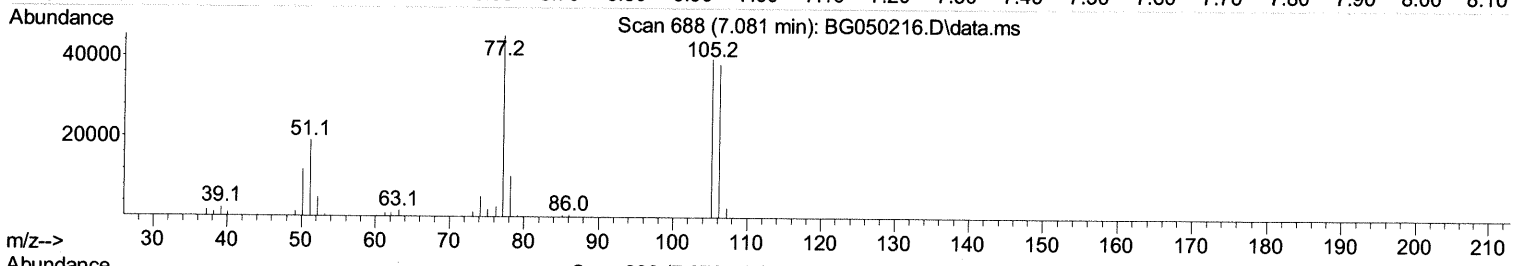
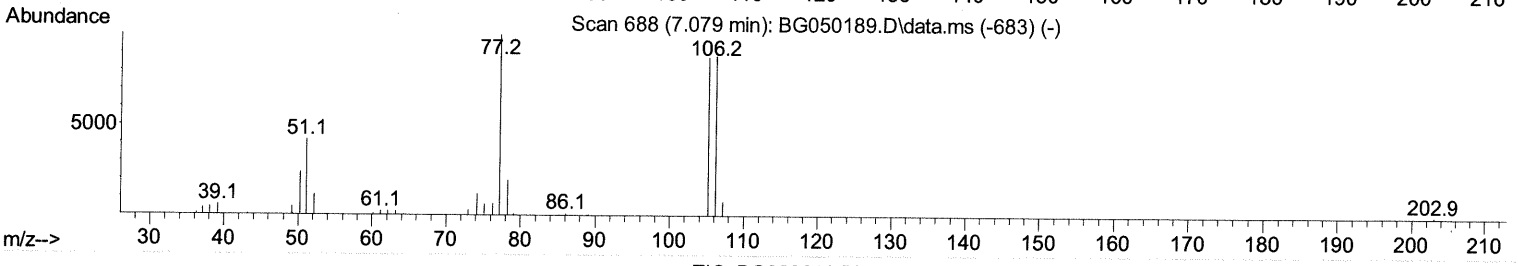
**Instrument :**
BNA_G
**ClientSampleId :**
SLCS845

**Manual Integrations APPROVED**
mohammad
9/9/2021 11:47:10 AM




TIC: BG050216.D\data.ms

**(6) Benzaldehyde**

**7.081min (-0.006) 25.99 ng/ul**

**response 77956**

| Ion | Exp% | Act% |
|---|---|---|
| 77.00 | 100.00 | 100.00 |
| 105.00 | 90.40 | 87.44 |
| 106.00 | 90.40 | 84.98 |
| 0.00 | 0.00 | 0.00 |

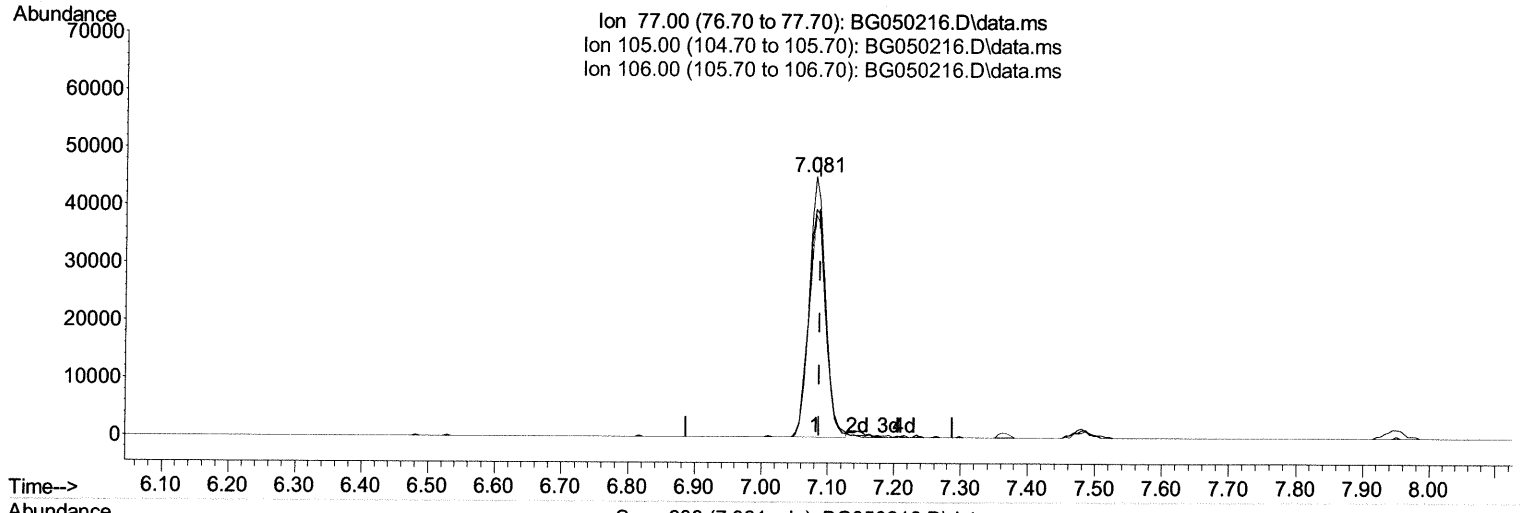
Quantitation Report (Qedit)

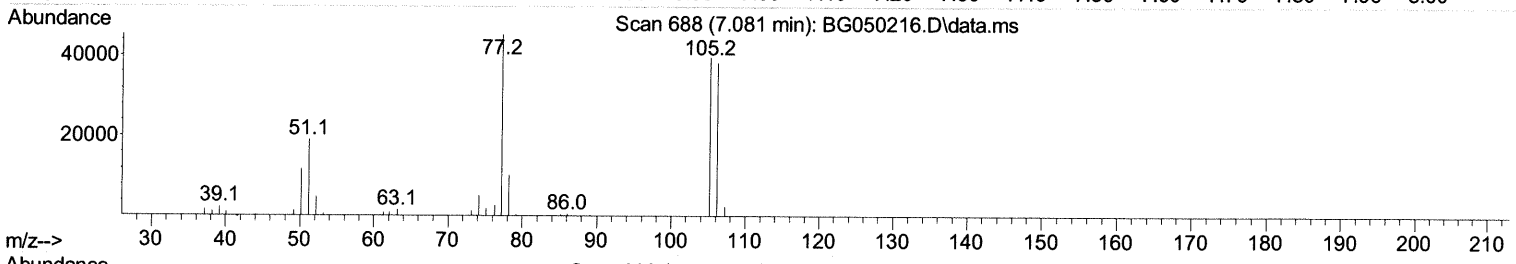
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050216.D
Acq On : 9 Sep 2021 2:59
Operator : CG/JU
Sample : PB138845BS
Misc :
ALS Vial : 44 Sample Multiplier: 1

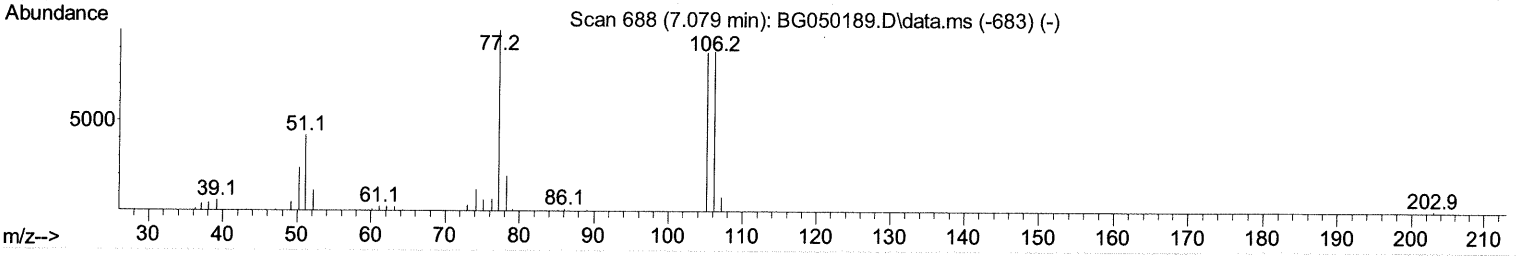
Quant Time: Sep 09 05:21:21 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Sep 07 18:32:38 2021
Response via : Initial Calibration

**Instrument :**
BNA_G
**ClientSampleId :**
SLCS845

**Manual Integrations APPROVED**
mohammad
9/9/2021 11:47:10 AM




TIC: BG050216.D\data.ms

**(6) Benzaldehyde**

**7.081min (-0.006) 25.82 ng/ul m** 09/10/2021

**response 77460**

| Ion | Exp% | Act% |
|---|---|---|
| 77.00 | 100.00 | 100.00 |
| 105.00 | 90.40 | 87.44 |
| 106.00 | 90.40 | 84.98 |
| 0.00 | 0.00 | 0.00 |

Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050216.D
Acq On : 9 Sep 2021 2:59
Operator : CG/JU
Sample : PB138845BS
Misc :
ALS Vial : 44 Sample Multiplier: 1

Quant Time: Sep 09 05:21:21 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
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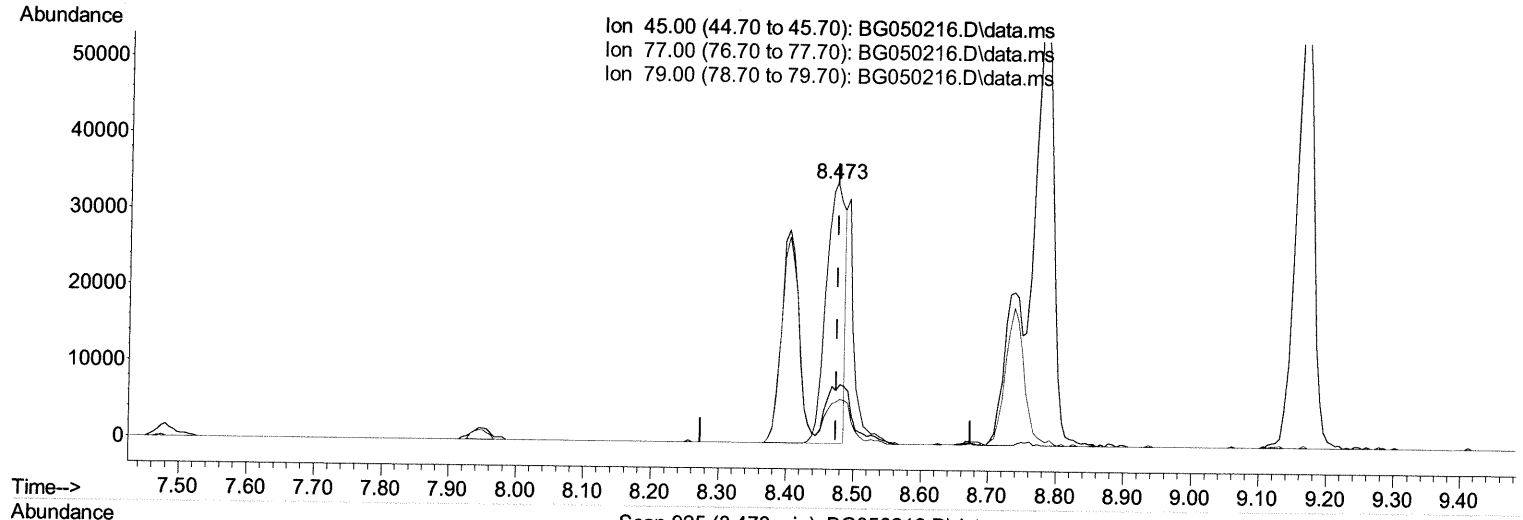
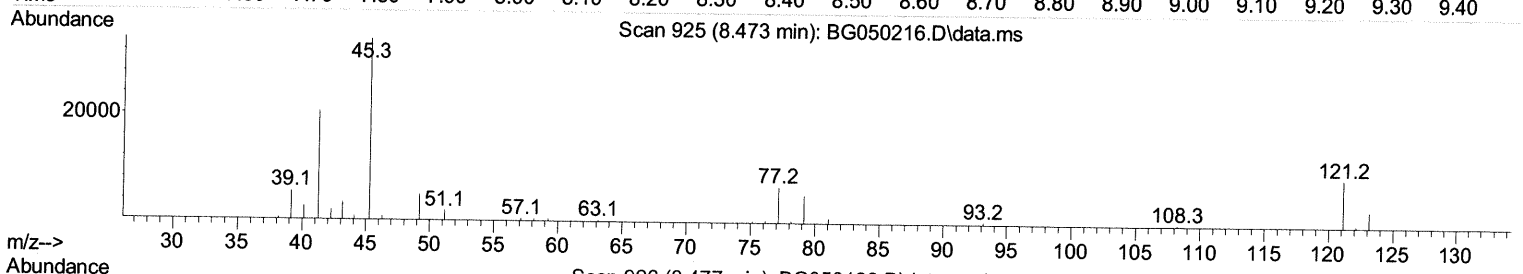
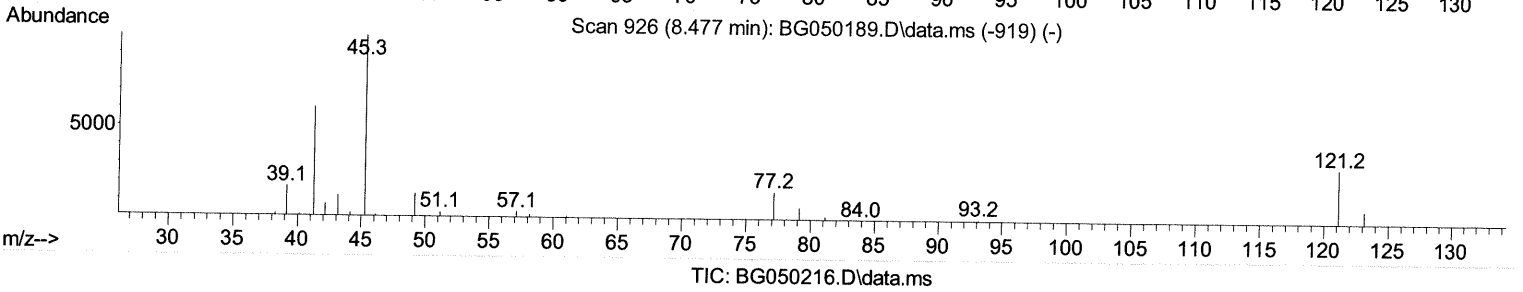
QLast Update : Tue Sep 07 18:32:38 2021
Response via : Initial Calibration

**Instrument :**
BNA_G
**ClientSampleId :**
SLCS845

**Manual Integrations APPROVED**
mohammad
9/9/2021 11:47:10 AM




TIC: BG050216.D\data.ms

**(14) 2,2'-oxybis(1-Chloropropane)**

**8.473min (-0.001) 16.81 ng/ul**

**response 67036**

| Ion | Exp% | Act% |
|---|---|---|
| 45.00 | 100.00 | 100.00 |
| 77.00 | 21.10 | 20.22 |
| 79.00 | 13.70 | 15.67 |
| 0.00 | 0.00 | 0.00 |

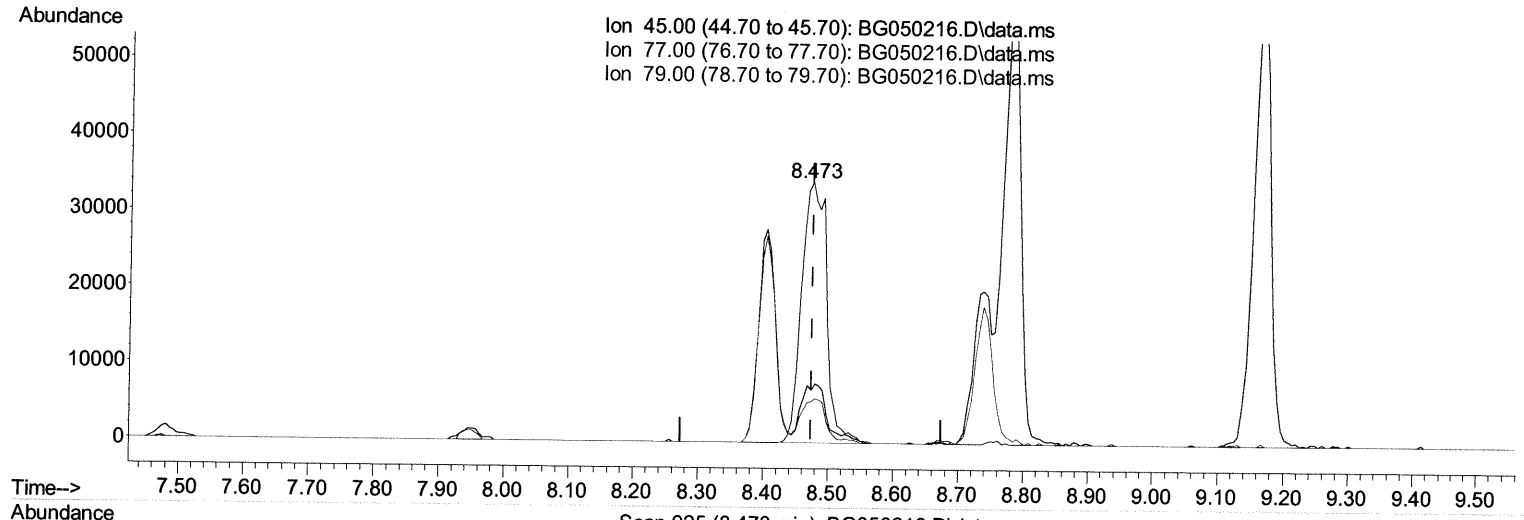
Quantitation Report (Qedit)

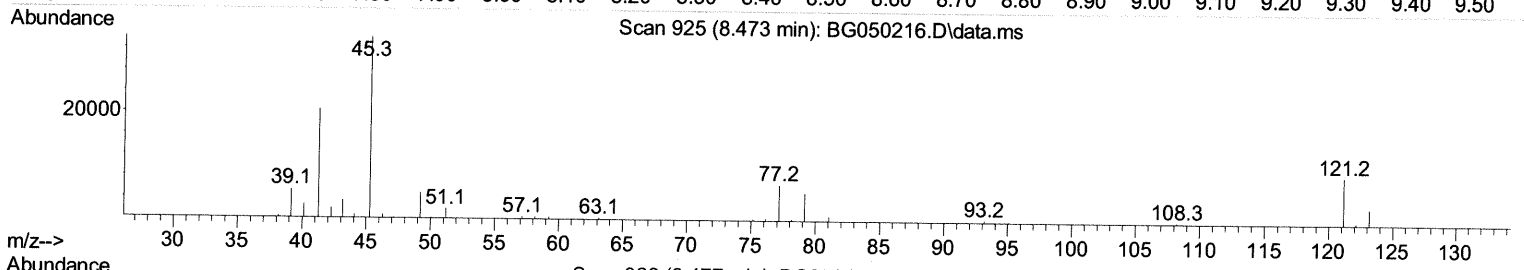
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050216.D
Acq On : 9 Sep 2021 2:59
Operator : CG/JU
Sample : PB138845BS
Misc :
ALS Vial : 44 Sample Multiplier: 1

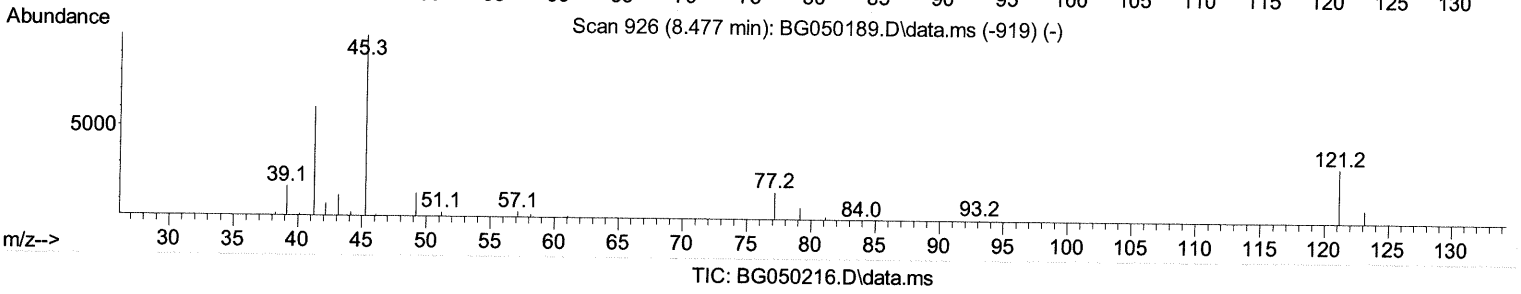
Quant Time: Sep 09 05:21:21 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Sep 07 18:32:38 2021
Response via : Initial Calibration

**Instrument :**
BNA_G
**ClientSampleId :**
SLCS845

**Manual Integrations APPROVED**
mohammad
9/9/2021 11:47:10 AM




TIC: BG050216.D\data.ms

**(14) 2,2'-oxybis(1-Chloropropane)**

**8.473min (-0.001) 22.83 ng/ul m** 6-9/10/2021

**response 91020**

| Ion | Exp% | Act% |
|---|---|---|
| 45.00 | 100.00 | 100.00 |
| 77.00 | 21.10 | 20.22 |
| 79.00 | 13.70 | 15.67 |
| 0.00 | 0.00 | 0.00 |

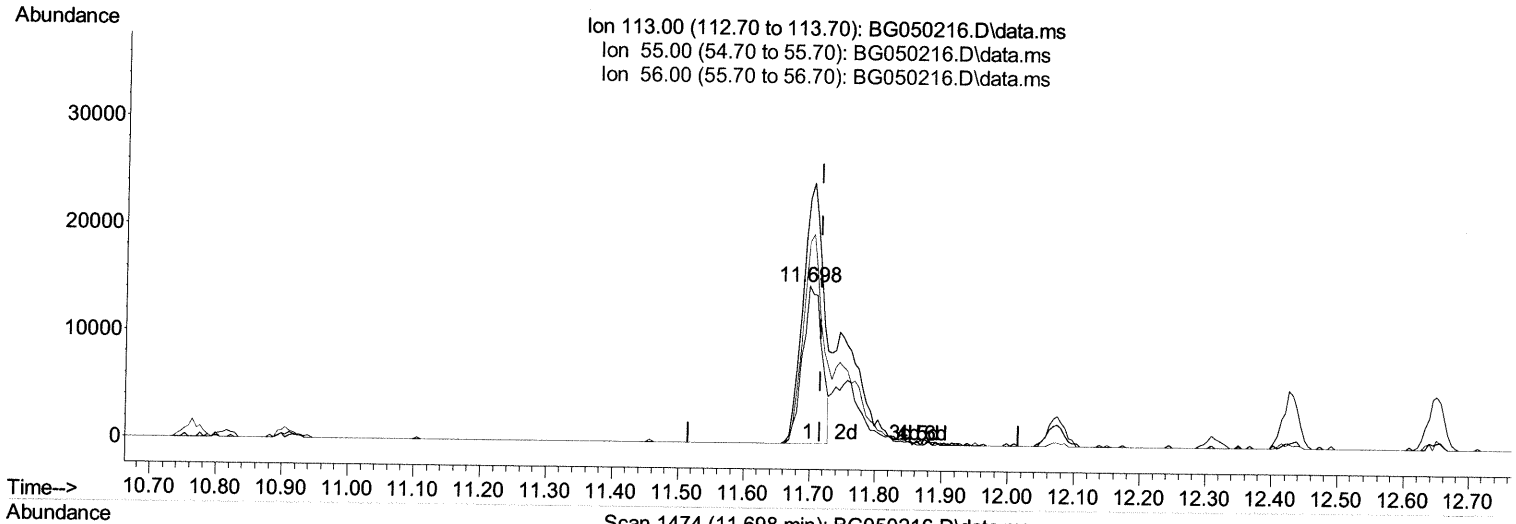
Quantitation Report (Qedit)

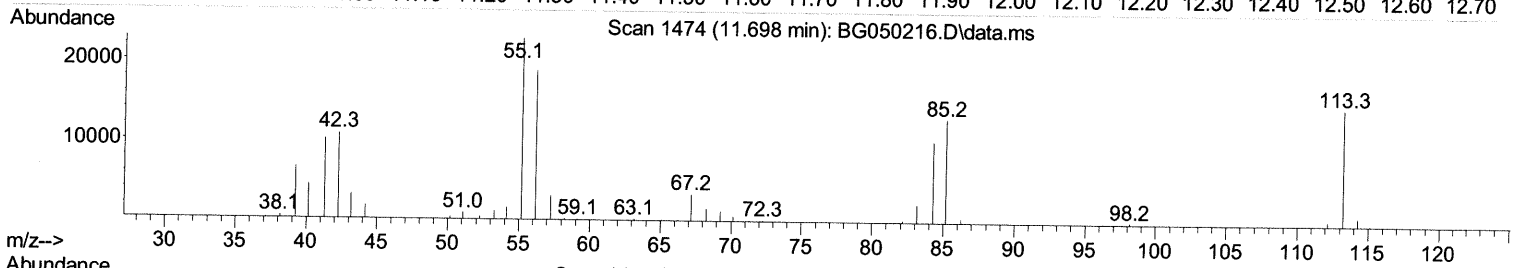
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050216.D
Acq On : 9 Sep 2021 2:59
Operator : CG/JU
Sample : PB138845BS
Misc :
ALS Vial : 44 Sample Multiplier: 1

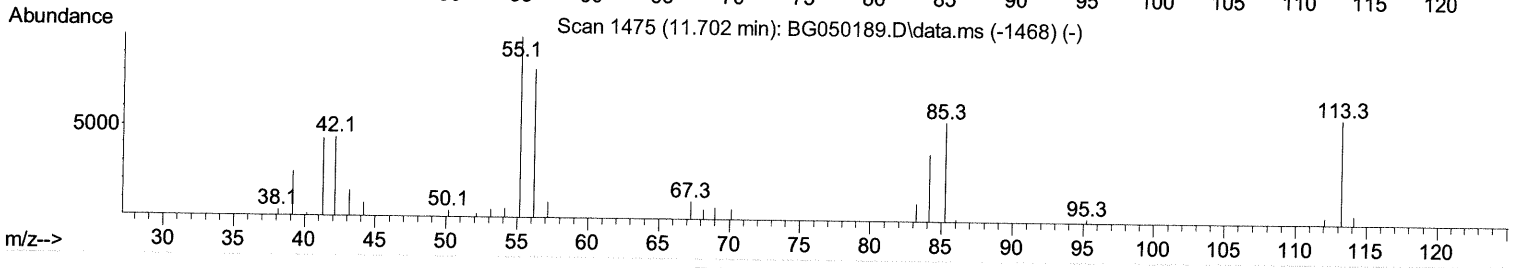
Quant Time: Sep 09 05:21:21 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Sep 07 18:32:38 2021
Response via : Initial Calibration

**Instrument :**
BNA_G
**ClientSampleId :**
SLCS845

**Manual Integrations APPROVED**
mohammad
9/9/2021 11:47:10 AM




TIC: BG050216.D\data.ms

**(34) Caprolactam**

**11.698min (-0.018) 32.35 ng/ul**

**response 30331**

| Ion | Exp% | Act% |
|---|---|---|
| 113.00 | 100.00 | 100.00 |
| 55.00 | 142.50 | 155.36 |
| 56.00 | 134.20 | 128.14 |
| 0.00 | 0.00 | 0.00 |

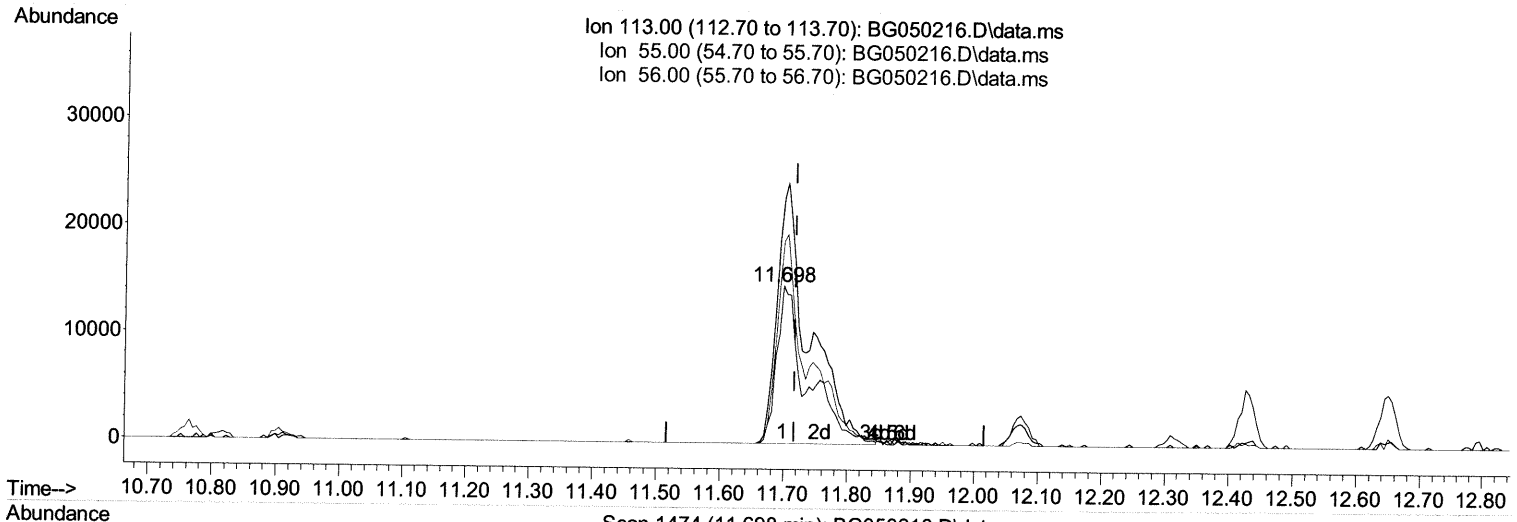
Quantitation Report (Qedit)

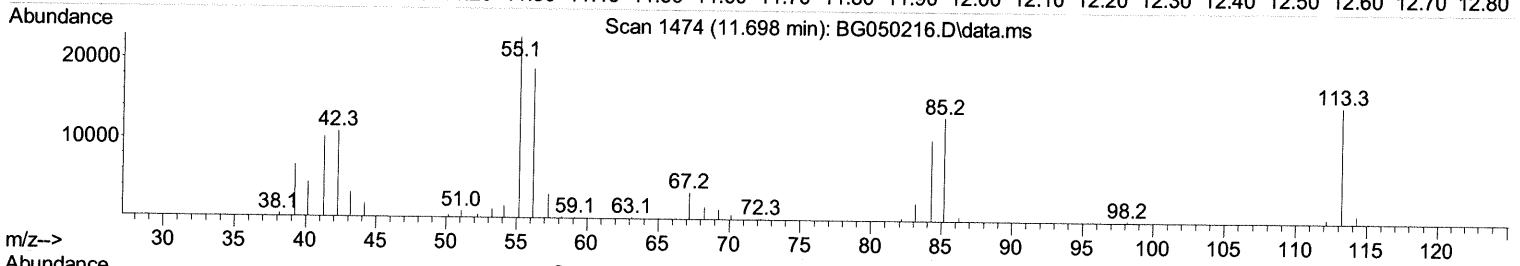
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050216.D
Acq On : 9 Sep 2021 2:59
Operator : CG/JU
Sample : PB138845BS
Misc :
ALS Vial : 44 Sample Multiplier: 1

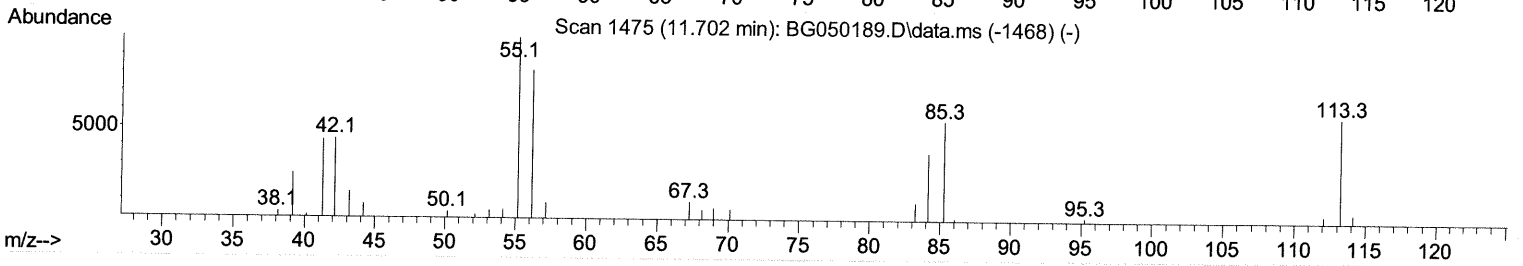
Quant Time: Sep 09 05:21:21 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Sep 07 18:32:38 2021
Response via : Initial Calibration

**Instrument :**
BNA_G
**ClientSampleId :**
SLCS845

**Manual Integrations APPROVED**
mohammad
9/9/2021 11:47:10 AM




TIC: BG050216.D\data.ms

**(34) Caprolactam**

**11.698min (-0.018) 51.64 ng/ul m** 09/10/2021

**response 48423**

| Ion | Exp% | Act% |
|---|---|---|
| 113.00 | 100.00 | 100.00 |
| 55.00 | 142.50 | 155.36 |
| 56.00 | 134.20 | 128.14 |
| 0.00 | 0.00 | 0.00 |

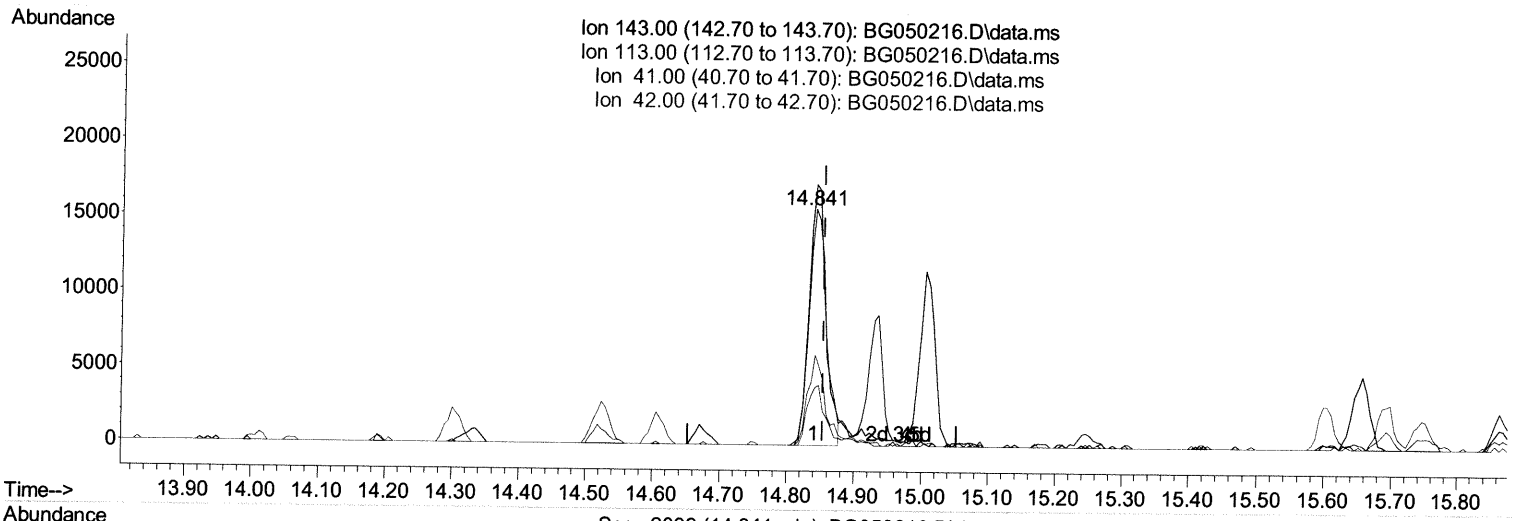
Quantitation Report (Qedit)

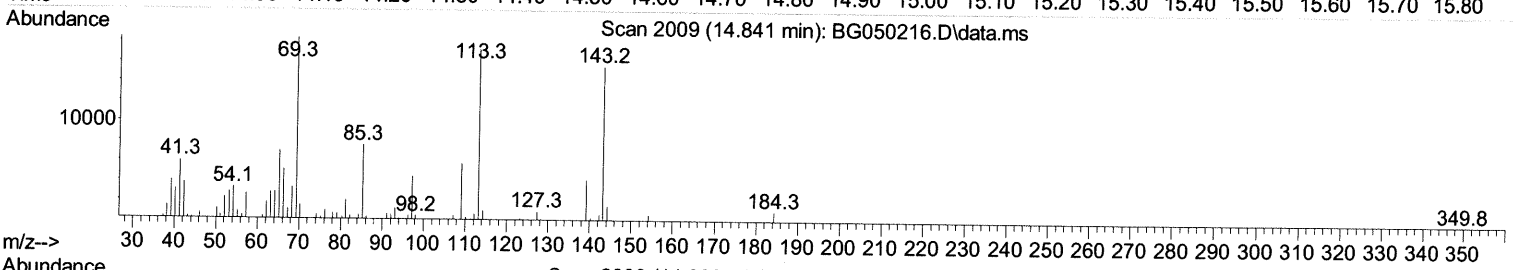
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050216.D
Acq On : 9 Sep 2021 2:59
Operator : CG/JU
Sample : PB138845BS
Misc :
ALS Vial : 44 Sample Multiplier: 1

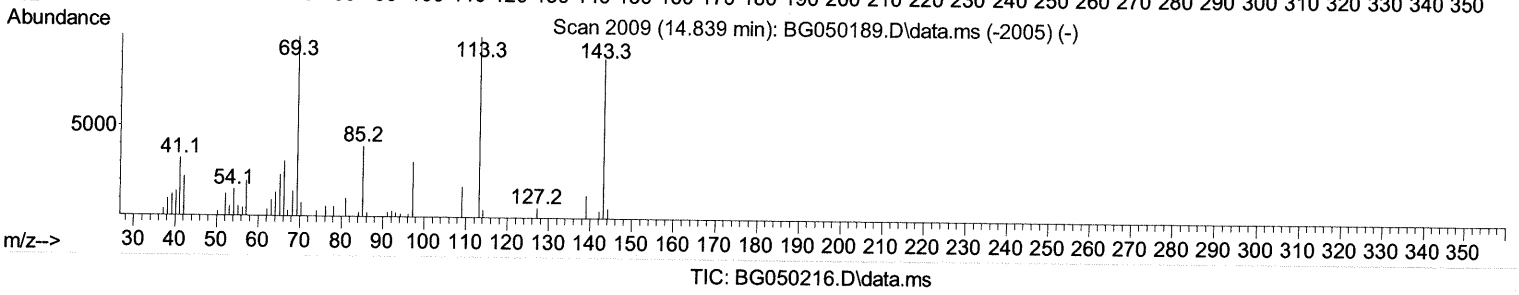
Quant Time: Sep 09 05:21:21 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Sep 07 18:32:38 2021
Response via : Initial Calibration

**Instrument :**
BNA_G
**ClientSampleId :**
SLCS845

**Manual Integrations APPROVED**
mohammad
9/9/2021 11:47:10 AM




TIC: BG050216.D\data.ms

**(54) 4-Nitrophenol-d4 (S)**

**14.841min (-0.012) 30.00 ng/ul**

**response 28175**

| Ion | Exp% | Act% |
|---|---|---|
| 143.00 | 100.00 | 100.00 |
| 113.00 | 121.00 | 110.42 |
| 41.00 | 45.30 | 38.24 |
| 42.00 | 25.60 | 24.47 |

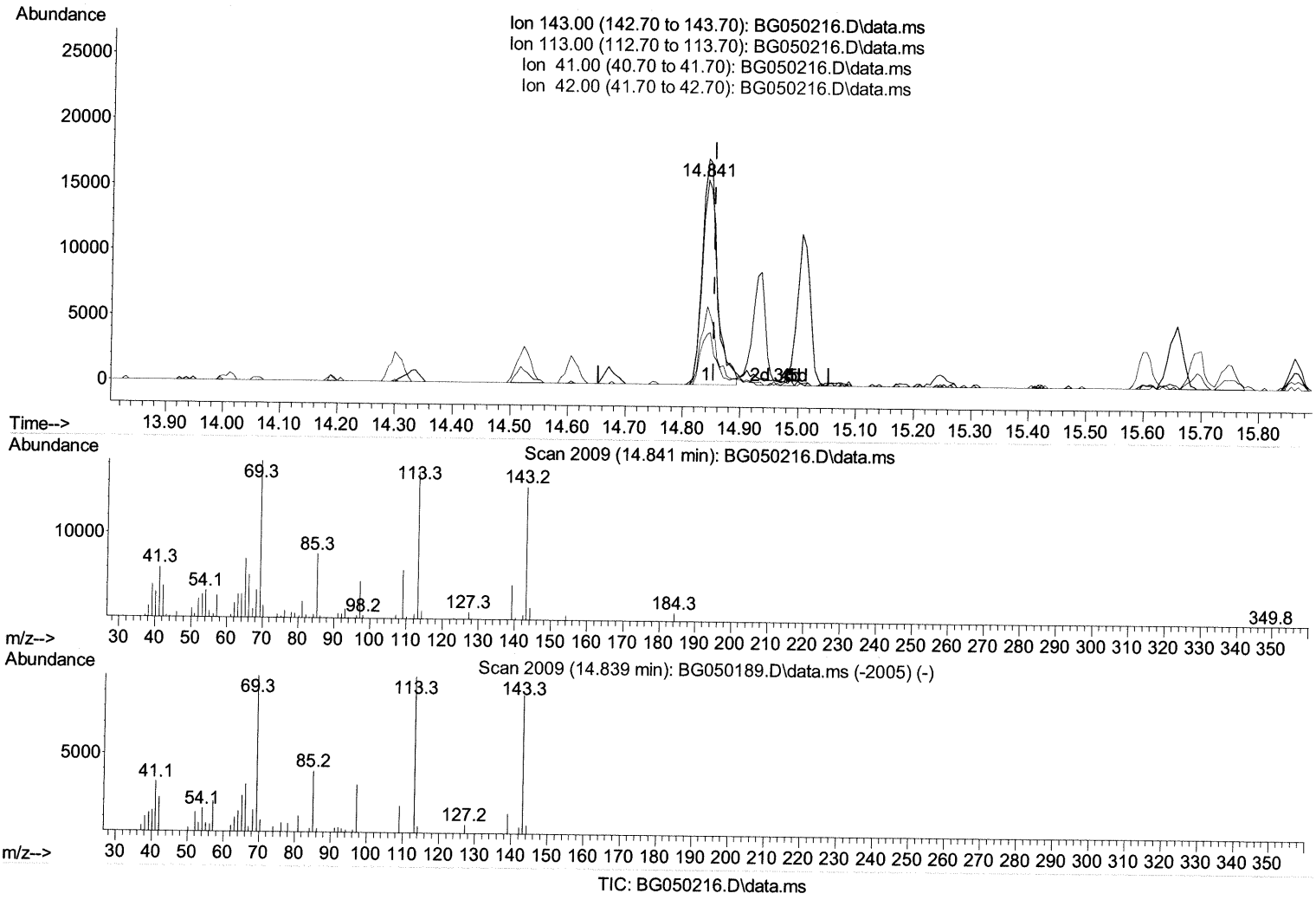
Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050216.D
Acq On : 9 Sep 2021 2:59
Operator : CG/JU
Sample : PB138845BS
Misc :
ALS Vial : 44 Sample Multiplier: 1

Quant Time: Sep 09 05:21:21 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Sep 07 18:32:38 2021
Response via : Initial Calibration

**Instrument :**
BNA_G
**ClientSampleId :**
SLCS845

**Manual Integrations APPROVED**
mohammad
9/9/2021 11:47:10 AM




TIC: BG050216.D\data.ms

**(54) 4-Nitrophenol-d4 (S)**

**14.841min (-0.012) 31.31 ng/ul m** 09/16/21 JU

**response 29413**

| Ion | Exp% | Act% |
|---|---|---|
| 143.00 | 100.00 | 100.00 |
| 113.00 | 121.00 | 110.42 |
| 41.00 | 45.30 | 38.24 |
| 42.00 | 25.60 | 24.47 |

Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050216.D
Acq On : 9 Sep 2021 2:59
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Response via : Initial Calibration

**Instrument :** BNA_G
**ClientSampleId :** SLCS845

**Manual Integrations APPROVED**
mohammad
9/9/2021 11:47:10 AM

| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|---|---|---|---|---|---|---|
| **Internal Standards** | | | | | | |
| 1) 1,4-Dichlorobenzene-d4 | 7.950 | 152 | 69701 | 20.000 | ng/ul | 0.00 |
| 20) Naphthalene-d8 | 10.764 | 136 | 211531 | 20.000 | ng/ul | -0.01 |
| 38) Acenaphthene-d10 | 14.606 | 164 | 103096 | 20.000 | ng/ul | -0.01 |
| 64) Phenanthrene-d10 | 17.367 | 188 | 231417 | 20.000 | ng/ul | 0.00 |
| 79) Chrysene-d12 | 21.637 | 240 | 236381 | 20.000 | ng/ul | 0.00 |
| 88) Perylene-d12 | 24.845 | 264 | 239393 | 20.000 | ng/ul | -0.01 |
| **System Monitoring Compounds** | | | | | | |
| 3) 1,4-Dioxane-d8 | 3.303 | 96 | 7148m | 5.070 | ng/uL | 0.00 |
| 4) Pyridine-d5 | 3.732 | 84 | 106652 | 21.311 | ng/ul | 0.00 |
| 7) Phenol-d5 | 7.122 | 99 | 101017 | 18.666 | ng/ul | 0.00 |
| 9) Bis-(2-Chloroethyl)eth... | 7.269 | 67 | 54733 | 17.317 | ng/ul | -0.01 |
| 11) 2-Chlorophenol-d4 | 7.480 | 132 | 78641 | 17.509 | ng/ul | 0.00 |
| 15) 4-Methylphenol-d8 | 8.673 | 113 | 90318 | 21.226 | ng/ul | 0.00 |
| 21) Nitrobenzene-d5 | 9.119 | 128 | 38260 | 26.351 | ng/ul | 0.00 |
| 24) 2-Nitrophenol-d4 | 9.847 | 143 | 44863 | 25.411 | ng/ul | 0.00 |
| 28) 2,4-Dichlorophenol-d3 | 10.394 | 165 | 82487 | 24.480 | ng/ul | -0.01 |
| 31) 4-Chloroaniline-d4 | 10.911 | 131 | 124731 | 30.045 | ng/ul | 0.00 |
| 46) Dimethylphthalate-d6 | 14.012 | 166 | 241680 | 31.046 | ng/ul | 0.00 |
| 49) Acenaphthylene-d8 | 14.306 | 160 | 295316 | 31.860 | ng/ul | 0.00 |
| 54) 4-Nitrophenol-d4 | 14.841 | 143 | 29413m | 31.314 | ng/ul | -0.01 |
| 60) Fluorene-d10 | 15.604 | 176 | 204552 | 31.646 | ng/ul | 0.00 |
| 65) 4,6-Dinitro-2-methylph... | 15.745 | 200 | 41037 | 29.397 | ng/ul | 0.00 |
| 73) Anthracene-d10 | 17.461 | 188 | 315742 | 30.393 | ng/ul | 0.00 |
| 81) Pyrene-d10 | 19.758 | 212 | 396886 | 33.708 | ng/ul | 0.00 |
| 92) Benzo(a)pyrene-d12 | 24.633 | 264 | 392361 | 30.961 | ng/ul | 0.00 |
| **Target Compounds** | | | | | | Qvalue |
| 2) 1,4-Dioxane | 3.339 | 88 | 16371 | 10.197 | ng/uL | 95 |
| 5) Pyridine | 3.750 | 79 | 113864 | 21.646 | ng/ul# | 83 |
| 6) Benzaldehyde | 7.081 | 77 | 77460m | 25.820 | ng/ul | |
| 8) Phenol | 7.145 | 94 | 106826 | 20.045 | ng/ul | 97 |
| 10) Bis(2-Chloroethyl)ether | 7.363 | 93 | 75738 | 19.168 | ng/ul | 97 |
| 12) 2-Chlorophenol | 7.509 | 128 | 80107 | 18.304 | ng/ul# | 91 |
| 13) 2-Methylphenol | 8.402 | 108 | 100483 | 23.426 | ng/ul | 94 |
| 14) 2,2'-oxybis(1-Chloropr... | 8.473 | 45 | 91020m | 22.825 | ng/ul | |
| 16) Acetophenone | 8.778 | 105 | 132433 | 21.154 | ng/ul | 96 |
| 17) N-Nitroso-di-n-propyla... | 8.761 | 70 | 67915 | 20.343 | ng/ul | 94 |
| 18) 4-Methylphenol | 8.737 | 108 | 92664 | 22.745 | ng/ul | 81 |
| 19) Hexachloroethane | 9.025 | 117 | 35426 | 19.405 | ng/ul | 99 |
| 22) Nitrobenzene | 9.166 | 77 | 105108 | 27.797 | ng/ul | 99 |
| 23) Isophorone | 9.683 | 82 | 178286 | 24.691 | ng/ul | 95 |
| 25) 2-Nitrophenol | 9.877 | 139 | 45375 | 26.052 | ng/ul# | 91 |
| 26) 2,4-Dimethylphenol | 9.941 | 107 | 91453 | 22.065 | ng/ul | 91 |
| 27) Bis(2-Chloroethoxy)met... | 10.165 | 93 | 98749 | 22.546 | ng/ul | 99 |
| 29) 2,4-Dichlorophenol | 10.423 | 162 | 81475 | 26.436 | ng/ul | 98 |
| 30) Naphthalene | 10.817 | 128 | 415618 | 42.781 | ng/ul | 98 |
| 32) 4-Chloroaniline | 10.928 | 127 | 123963 | 31.700 | ng/ul | 93 |
| 33) Hexachlorobutadiene | 11.093 | 225 | 77014 | 31.525 | ng/ul | 97 |
| 34) Caprolactam | 11.698 | 113 | 48423m | 51.643 | ng/ul | |
| 35) 4-Chloro-3-methylphenol | 12.074 | 107 | 163914 | 47.188 | ng/ul | 98 |

Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090721\
Data File : BG050216.D
Acq On : 9 Sep 2021 2:59
Operator : CG/JU
Sample : PB138845BS
Misc :
ALS Vial : 44 Sample Multiplier: 1

Quant Time: Sep 09 05:21:21 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\SFAM-EPA-BG090721.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Sep 07 18:32:38 2021
Response via : Initial Calibration

**Instrument :** BNA_G
**ClientSampleId :** SLCS845

**Manual Integrations APPROVED**
mohammad
9/9/2021 11:47:10 AM

| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|---|---|---|---|---|---|---|
| 36) 2-Methylnaphthalene | 12.432 | 142 | 318366 | 44.315 | ng/ul | 97 |
| 37) 1-Methylnaphthalene | 12.650 | 142 | 299572 | 43.427 | ng/ul | 99 |
| 39) 1,2,4,5-Tetrachloroben... | 12.802 | 216 | 166295 | 47.690 | ng/ul# | 97 |
| 40) Hexachlorocyclopentadiene | 12.773 | 237 | 24846 | 26.278 | ng/ul# | 95 |
| 41) 2,4,6-Trichlorophenol | 13.049 | 196 | 94600 | 48.051 | ng/ul | 88 |
| 42) 2,4,5-Trichlorophenol | 13.131 | 196 | 85158 | 41.362 | ng/ul | 93 |
| 43) 1,1'-Biphenyl | 13.437 | 154 | 303562 | 40.786 | ng/ul | 99 |
| 44) 2-Chloronaphthalene | 13.484 | 162 | 213363 | 36.453 | ng/ul | 98 |
| 45) 2-Nitroaniline | 13.695 | 65 | 59854 | 32.301 | ng/ul | 98 |
| 47) Dimethylphthalate | 14.059 | 163 | 240968 | 32.431 | ng/ul | 99 |
| 48) 2,6-Dinitrotoluene | 14.189 | 165 | 52753 | 33.345 | ng/ul | 91 |
| 50) Acenaphthylene | 14.330 | 152 | 280997 | 32.437 | ng/ul | 99 |
| 51) 3-Nitroaniline | 14.524 | 138 | 50957 | 35.380 | ng/ul | 99 |
| 52) Acenaphthene | 14.670 | 153 | 203522 | 32.971 | ng/ul | 96 |
| 53) 2,4-Dinitrophenol | 14.753 | 184 | 18518 | 28.581 | ng/ul# | 77 |
| 55) 4-Nitrophenol | 14.858 | 109 | 34142 | 32.662 | ng/ul | 92 |
| 56) Dibenzofuran | 15.011 | 168 | 277415 | 32.666 | ng/ul | 100 |
| 57) 2,4-Dinitrotoluene | 14.988 | 165 | 76373 | 33.783 | ng/ul | 91 |
| 58) 2,3,4,6-Tetrachlorophenol | 15.246 | 232 | 57572 | 36.920 | ng/ul# | 93 |
| 59) Diethylphthalate | 15.422 | 149 | 246130 | 32.774 | ng/ul | 95 |
| 61) Fluorene | 15.657 | 166 | 233389 | 32.779 | ng/ul | 96 |
| 62) 4-Chlorophenyl-phenyle... | 15.646 | 204 | 121905 | 32.156 | ng/ul | 98 |
| 63) 4-Nitroaniline | 15.693 | 138 | 53496 | 39.881 | ng/ul | 93 |
| 66) 4,6-Dinitro-2-methylph... | 15.763 | 198 | 37752 | 30.451 | ng/ul | 99 |
| 67) N-Nitrosodiphenylamine | 15.863 | 169 | 197029 | 31.487 | ng/ul | 99 |
| 68) 4-Bromophenyl-phenylether | 16.544 | 248 | 91397 | 33.328 | ng/ul | 93 |
| 69) Hexachlorobenzene | 16.674 | 284 | 100642 | 32.369 | ng/ul | 93 |
| 70) Atrazine | 16.815 | 200 | 87846 | 33.481 | ng/ul | 99 |
| 71) Pentachlorophenol | 17.032 | 266 | 22155 | 26.013 | ng/ul | 90 |
| 72) Phenanthrene | 17.408 | 178 | 383111 | 32.459 | ng/ul | 98 |
| 74) Anthracene | 17.496 | 178 | 377304 | 31.526 | ng/ul | 95 |
| 75) 1,2,3,4-Tetrachloroben... | 13.407 | 216 | 134068 | 37.383 | ng/uL | 90 |
| 76) Pentachlorobenzene | 14.935 | 250 | 113402 | 30.625 | ng/uL | 96 |
| 77) Carbazole | 17.772 | 167 | 338753 | 32.106 | ng/ul | 99 |
| 78) Di-n-butylphthalate | 18.313 | 149 | 419501 | 32.877 | ng/ul | 98 |
| 80) Fluoranthene | 19.423 | 202 | 482589 | 32.977 | ng/ul# | 98 |
| 82) Pyrene | 19.781 | 202 | 474640 | 33.068 | ng/ul | 98 |
| 83) Butylbenzylphthalate | 20.656 | 149 | 190563 | 34.715 | ng/ul | 96 |
| 84) 3,3'-Dichlorobenzidine | 21.526 | 252 | 174689 | 34.261 | ng/ul# | 97 |
| 85) Benzo(a)anthracene | 21.614 | 228 | 465022 | 32.134 | ng/ul | 98 |
| 86) Bis(2-ethylhexyl)phtha... | 21.502 | 149 | 268832 | 32.792 | ng/ul | 100 |
| 87) Chrysene | 21.684 | 228 | 442527 | 32.804 | ng/ul | 96 |
| 89) Di-n-octyl phthalate | 22.701 | 149 | 457497 | 33.021 | ng/ul | 100 |
| 90) Benzo(b)fluoranthene | 23.823 | 252 | 493022 | 32.908 | ng/ul# | 99 |
| 91) Benzo(k)fluoranthene | 23.893 | 252 | 468673 | 32.772 | ng/ul | 100 |
| 93) Benzo(a)pyrene | 24.698 | 252 | 432123 | 32.992 | ng/ul | 98 |
| 94) Indeno(1,2,3-cd)pyrene | 28.540 | 276 | 556288 | 32.414 | ng/ul# | 96 |
| 95) Dibenzo(a,h)anthracene | 28.581 | 278 | 461946 | 32.447 | ng/ul | 98 |
| 96) Benzo(g,h,i)perylene | 29.685 | 276 | 452190 | 32.886 | ng/ul | 97 |

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