Quantitation Report (Qedit)

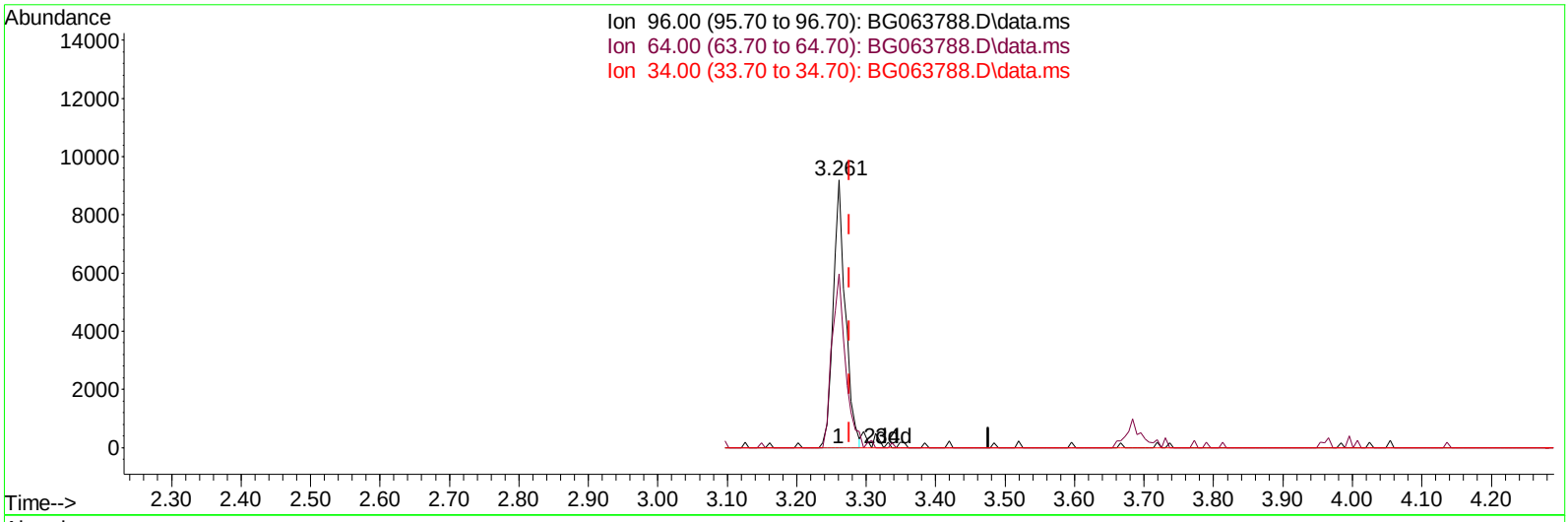
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122324\
Data File : BG063788.D
Acq On : 23 Dec 2024 17:37
Operator : RC/JU
Sample : P5302-04MS
Misc :
ALS Vial : 11 Sample Multiplier: 1

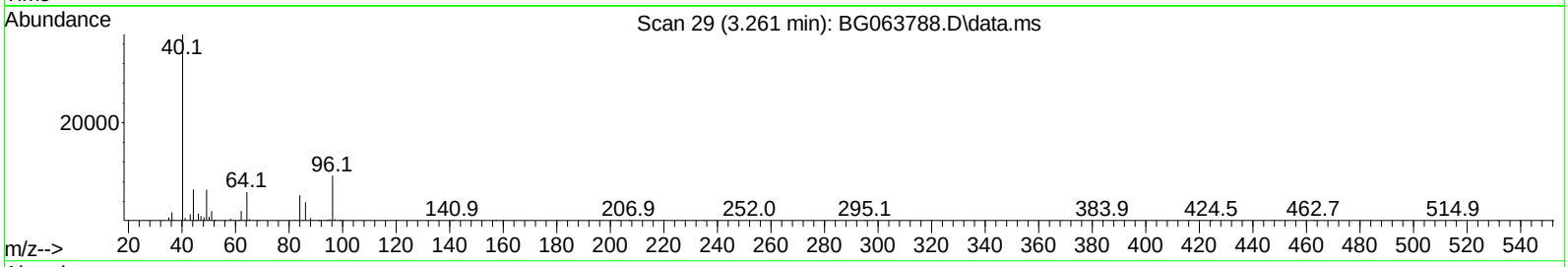
Quant Time: Dec 23 23:53:50 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121124.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Dec 11 23:59:23 2024
Response via : Initial Calibration

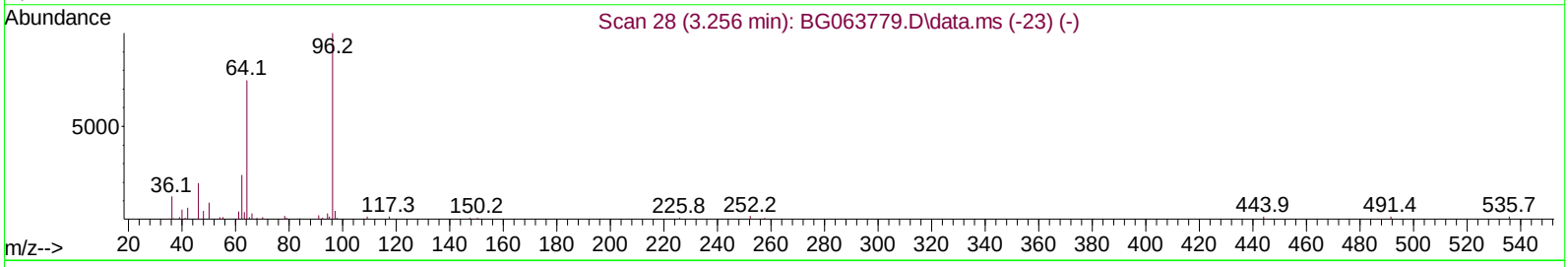
**Instrument :**
BNA_G
**ClientSampleId :**
E2AR6MS

**Manual IntegrationsAPPROVED**

Reviewed By :Yogesh Patel 12/24/2024
Supervised By :mohammad ahmed 12/26/2024




TIC: BG063788.D\data.ms

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

**3.261min (-0.015) 4.07 ng/uL**

**response 11243**

| Ion | Exp% | Act% |
|---|---|---|
| 96.00 | 100.00 | 100.00 |
| 64.00 | 82.80 | 64.90# |
| 34.00 | 0.00 | 0.00 |
| 0.00 | 0.00 | 0.00 |

Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122324\
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Acq On : 23 Dec 2024 17:37
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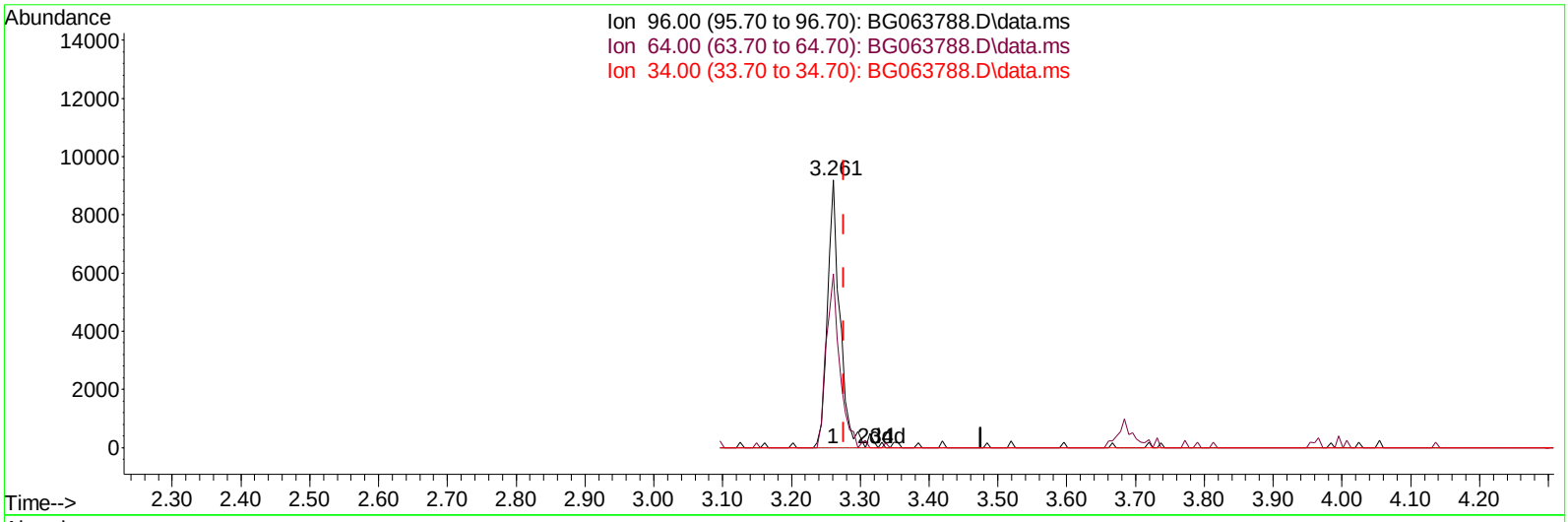
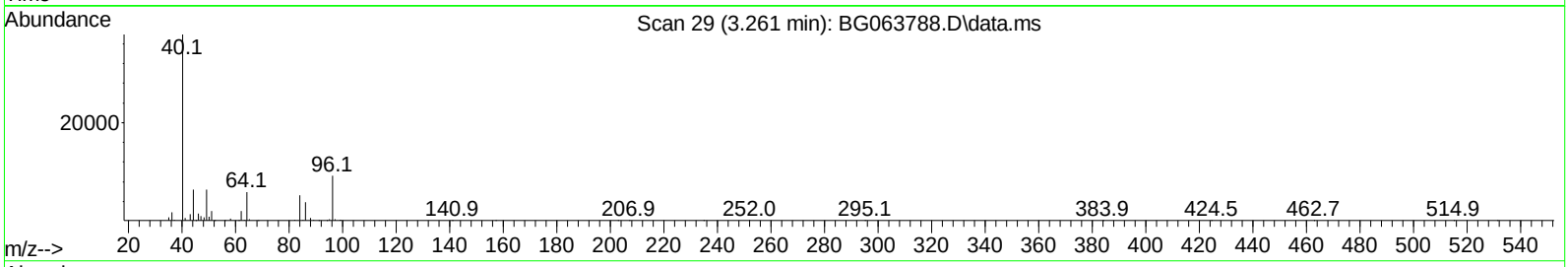
QLast Update : Wed Dec 11 23:59:23 2024
Response via : Initial Calibration

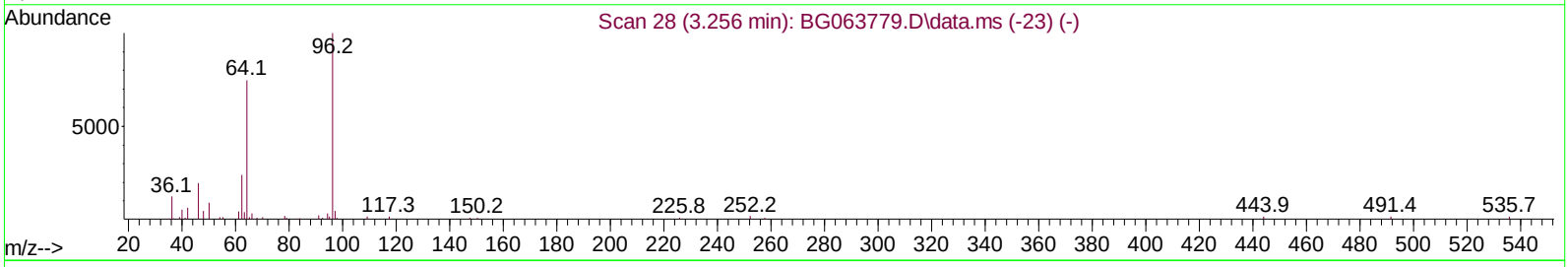
**Instrument :**
BNA_G
**ClientSampleId :**
E2AR6MS

**Manual IntegrationsAPPROVED**

Reviewed By :Yogesh Patel 12/24/2024
Supervised By :mohammad ahmed 12/26/2024




TIC: BG063788.D\data.ms

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

**3.261min (-0.015) 4.17 ng/uL m**

**response 11524**

| Ion | Exp% | Act% |
|---|---|---|
| 96.00 | 100.00 | 100.00 |
| 64.00 | 82.80 | 64.90# |
| 34.00 | 0.00 | 0.00 |
| 0.00 | 0.00 | 0.00 |

Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122324\
Data File : BG063788.D
Acq On : 23 Dec 2024 17:37
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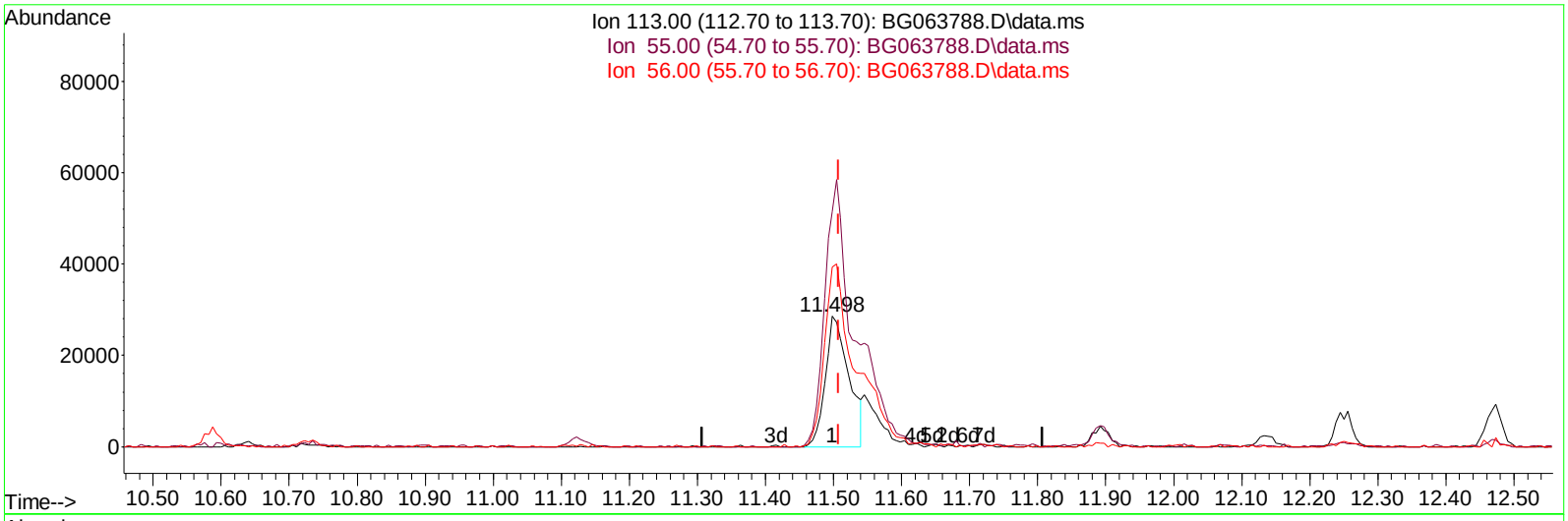
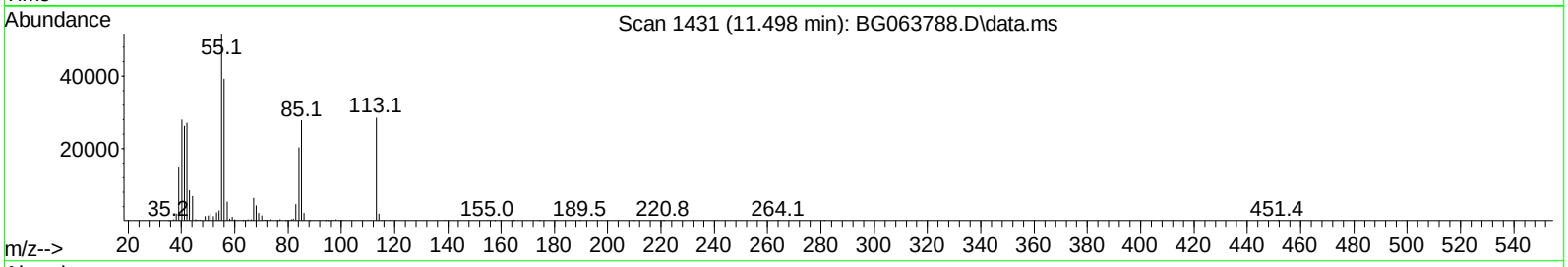
Quant Title : SVOA CALIBRATION
QLast Update : Wed Dec 11 23:59:23 2024
Response via : Initial Calibration

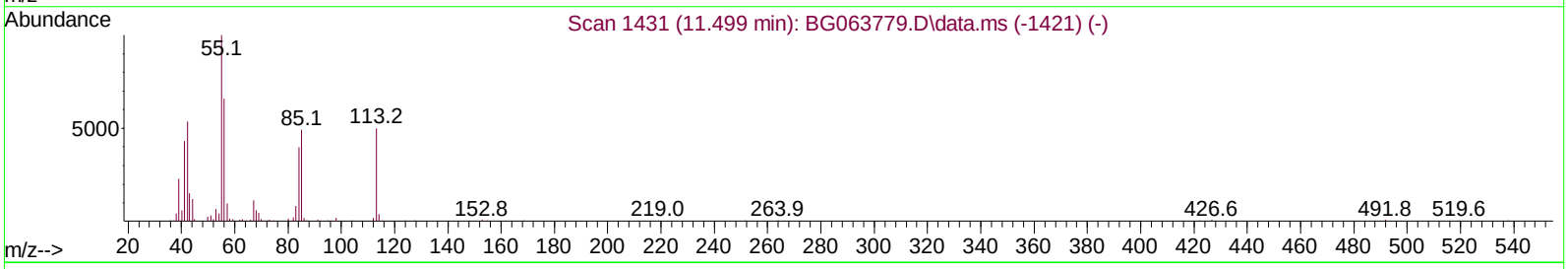
**Instrument :**
BNA_G
**ClientSampleId :**
E2AR6MS

**Manual IntegrationsAPPROVED**

Reviewed By :Yogesh Patel 12/24/2024
Supervised By :mohammad ahmed 12/26/2024




TIC: BG063788.D\data.ms

**(34) Caprolactam**

**11.498min (-0.009) 29.02 ng/ul**

**response 68738**

| Ion | Exp% | Act% |
|---|---|---|
| 113.00 | 100.00 | 100.00 |
| 55.00 | 202.90 | 179.99 |
| 56.00 | 133.90 | 137.55 |
| 0.00 | 0.00 | 0.00 |

Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122324\
Data File : BG063788.D
Acq On : 23 Dec 2024 17:37
Operator : RC/JU
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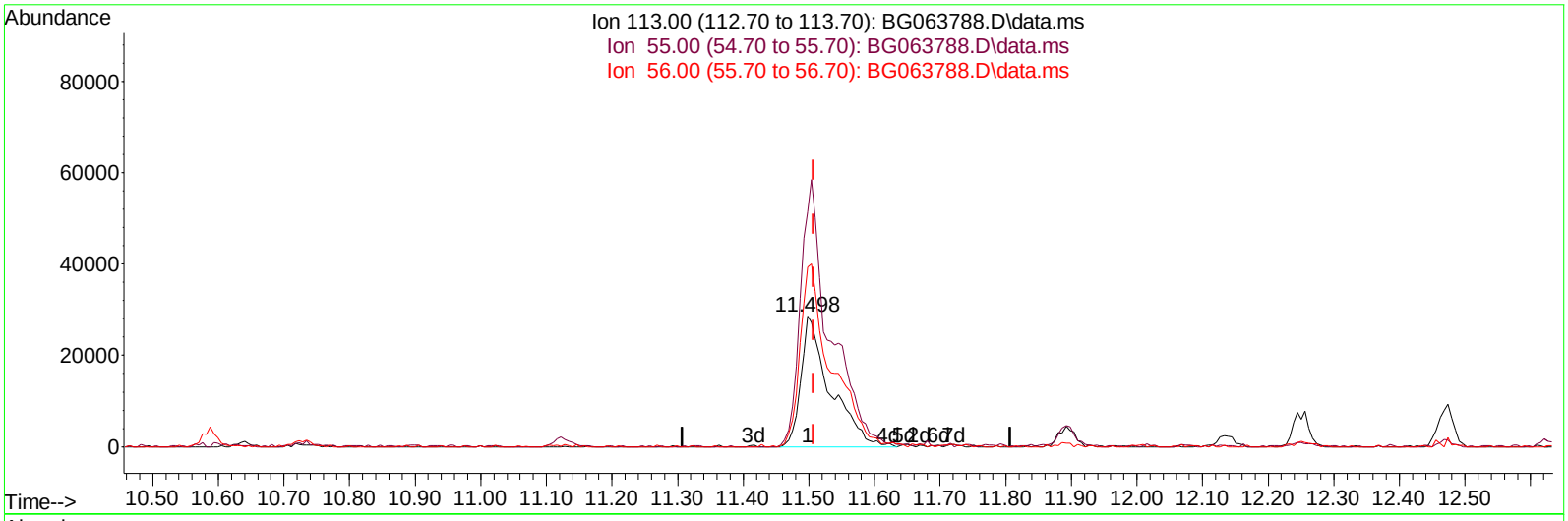
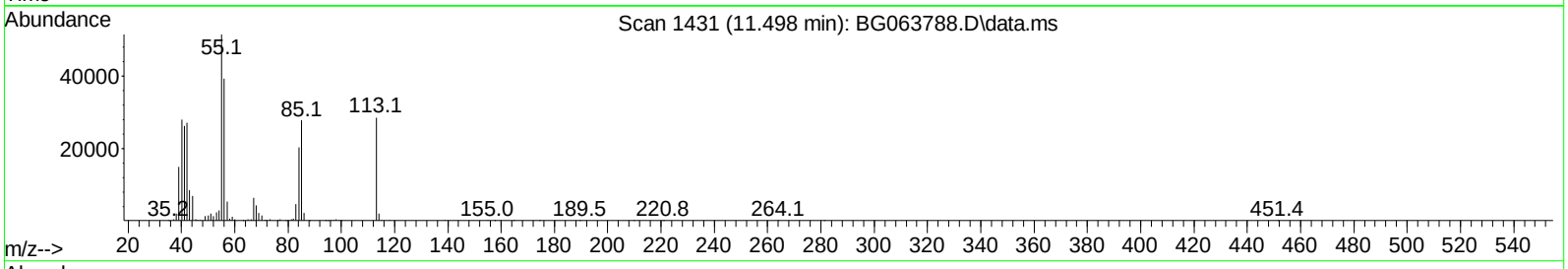
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121124.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Dec 11 23:59:23 2024
Response via : Initial Calibration

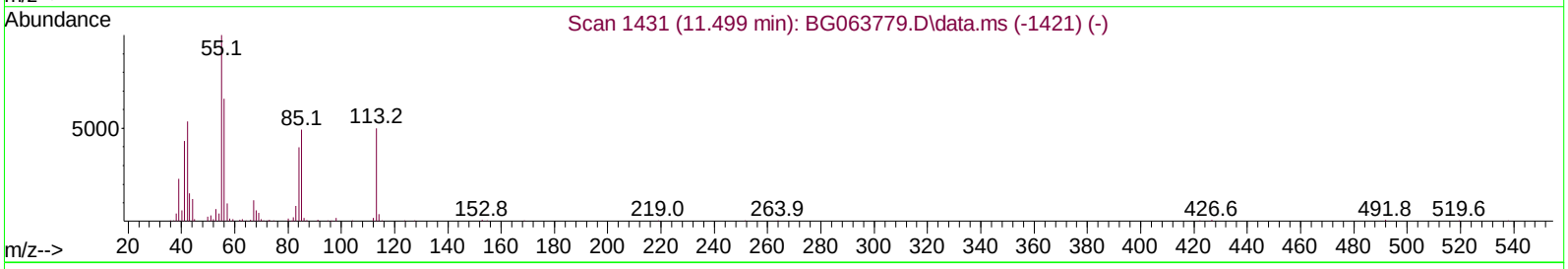
**Instrument :**
BNA_G
**ClientSampleId :**
E2AR6MS

**Manual IntegrationsAPPROVED**

Reviewed By :Yogesh Patel 12/24/2024
Supervised By :mohammad ahmed 12/26/2024




TIC: BG063788.D\data.ms

**(34) Caprolactam**

**11.498min (-0.009) 37.61 ng/ul m**

**response 89092**

| Ion | Exp% | Act% |
|---|---|---|
| 113.00 | 100.00 | 100.00 |
| 55.00 | 202.90 | 179.99 |
| 56.00 | 133.90 | 137.55 |
| 0.00 | 0.00 | 0.00 |

Quantitation Report (Qedit)

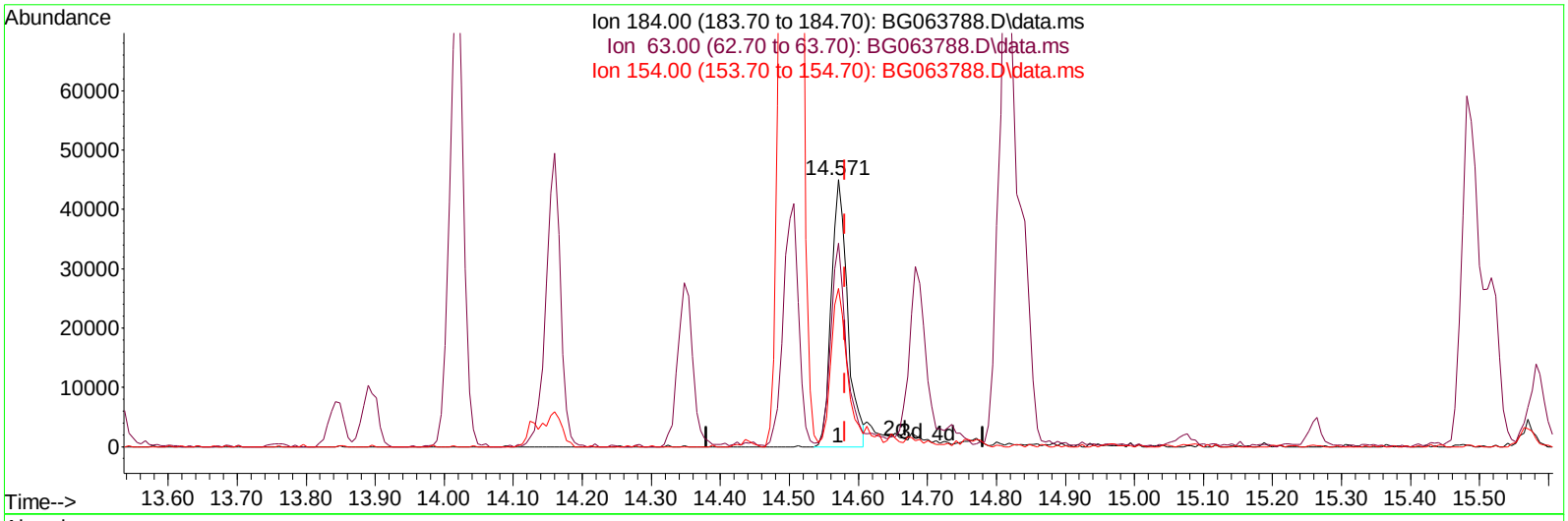
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122324\
Data File : BG063788.D
Acq On : 23 Dec 2024 17:37
Operator : RC/JU
Sample : P5302-04MS
Misc :
ALS Vial : 11 Sample Multiplier: 1

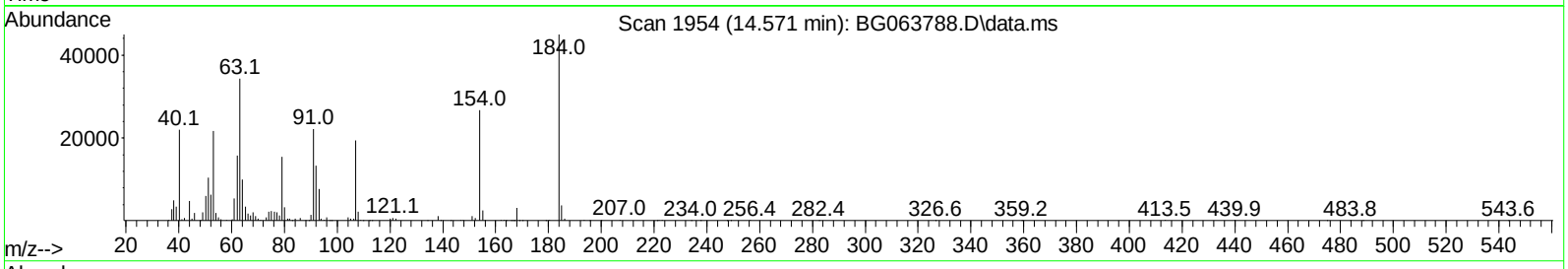
Quant Time: Dec 23 23:53:50 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121124.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Dec 11 23:59:23 2024
Response via : Initial Calibration

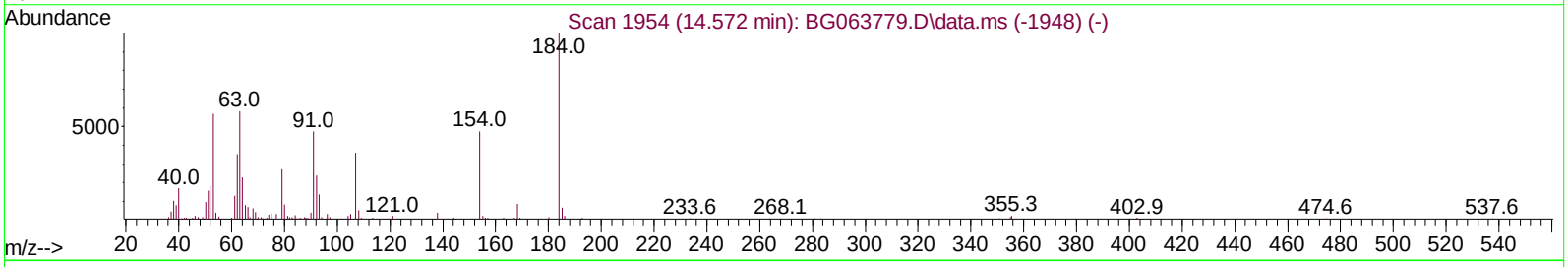
**Instrument :**
BNA_G
**ClientSampleId :**
E2AR6MS

**Manual IntegrationsAPPROVED**

Reviewed By :Yogesh Patel 12/24/2024
Supervised By :mohammad ahmed 12/26/2024




TIC: BG063788.D\data.ms

**(53) 2,4-Dinitrophenol**

**14.571min (-0.009) 31.14 ng/ul**

**response 74152**

| Ion | Exp% | Act% |
|---|---|---|
| 184.00 | 100.00 | 100.00 |
| 63.00 | 65.10 | 76.30 |
| 154.00 | 59.70 | 59.31 |
| 0.00 | 0.00 | 0.00 |

Quantitation Report (Qedit)

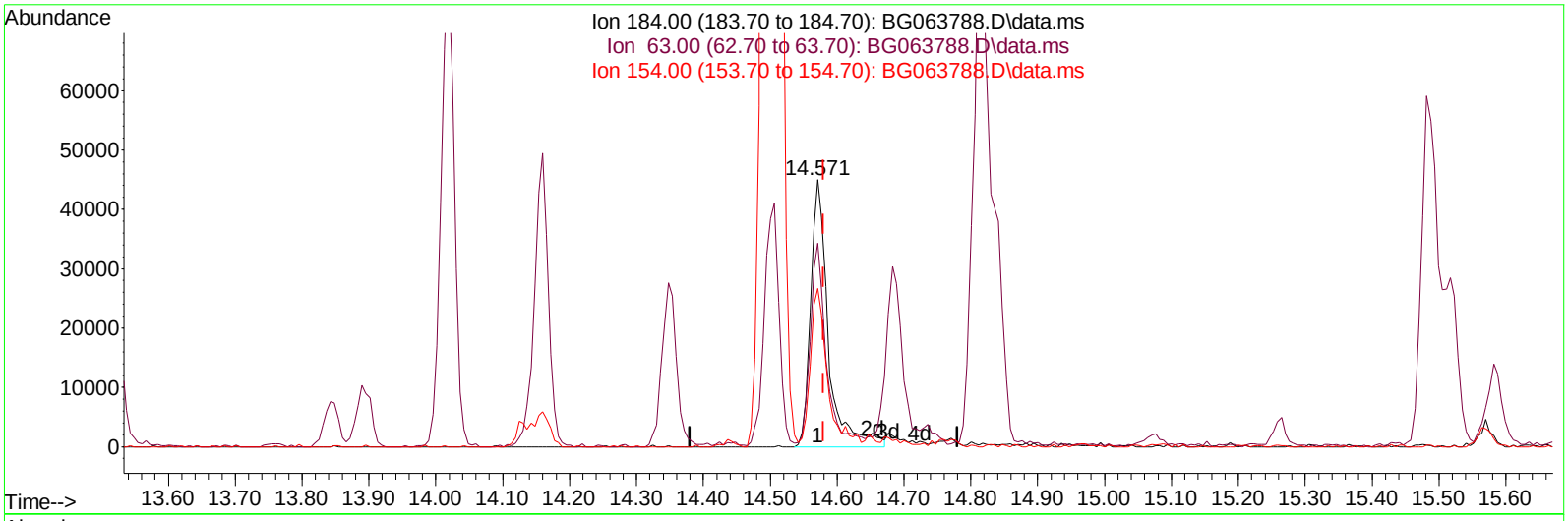
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122324\
Data File : BG063788.D
Acq On : 23 Dec 2024 17:37
Operator : RC/JU
Sample : P5302-04MS
Misc :
ALS Vial : 11 Sample Multiplier: 1

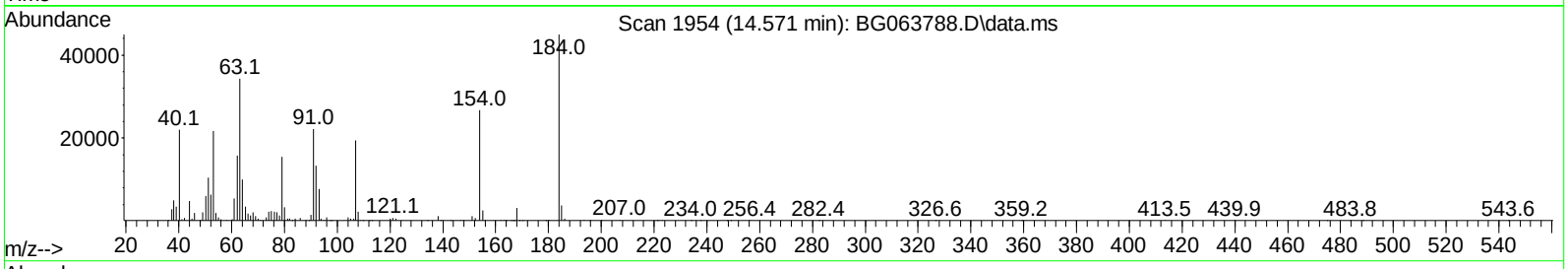
Quant Time: Dec 23 23:53:50 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121124.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Dec 11 23:59:23 2024
Response via : Initial Calibration

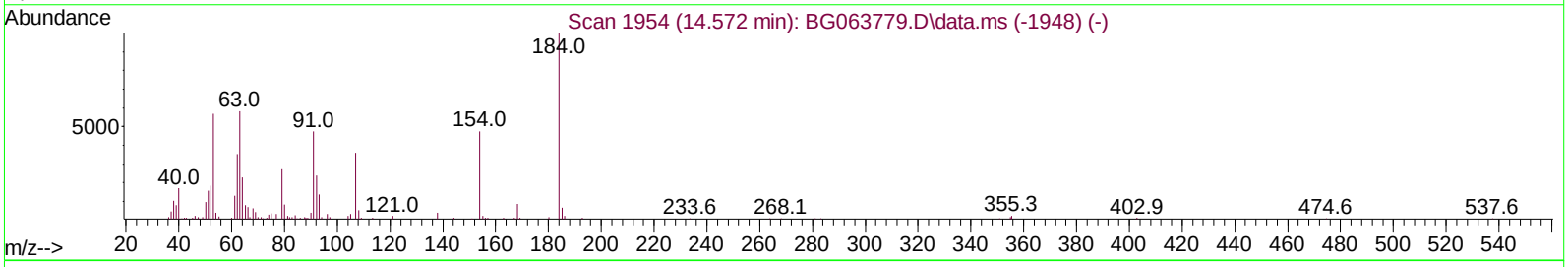
**Instrument :**
BNA_G
**ClientSampleId :**
E2AR6MS

**Manual IntegrationsAPPROVED**

Reviewed By :Yogesh Patel 12/24/2024
Supervised By :mohammad ahmed 12/26/2024




TIC: BG063788.D\data.ms

**(53) 2,4-Dinitrophenol**

**14.571min (-0.009) 34.75 ng/ul m**

**response 82736**

| Ion | Exp% | Act% |
|---|---|---|
| 184.00 | 100.00 | 100.00 |
| 63.00 | 65.10 | 76.30 |
| 154.00 | 59.70 | 59.31 |
| 0.00 | 0.00 | 0.00 |

Quantitation Report (Qedit)

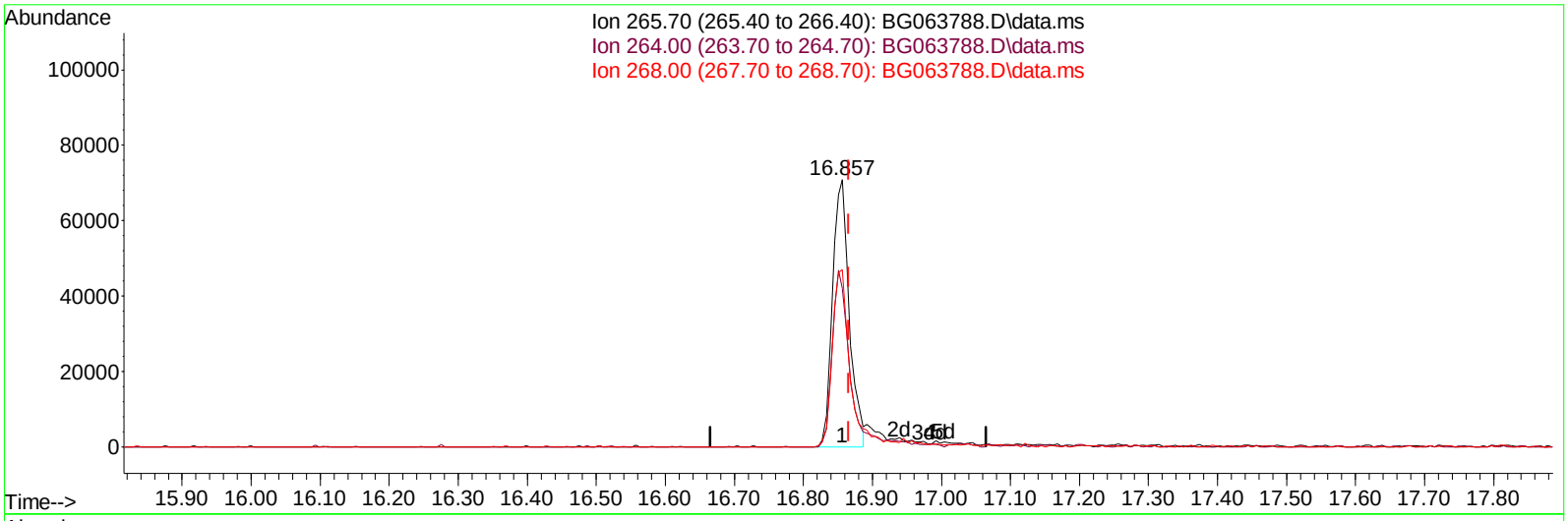
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122324\
Data File : BG063788.D
Acq On : 23 Dec 2024 17:37
Operator : RC/JU
Sample : P5302-04MS
Misc :
ALS Vial : 11 Sample Multiplier: 1

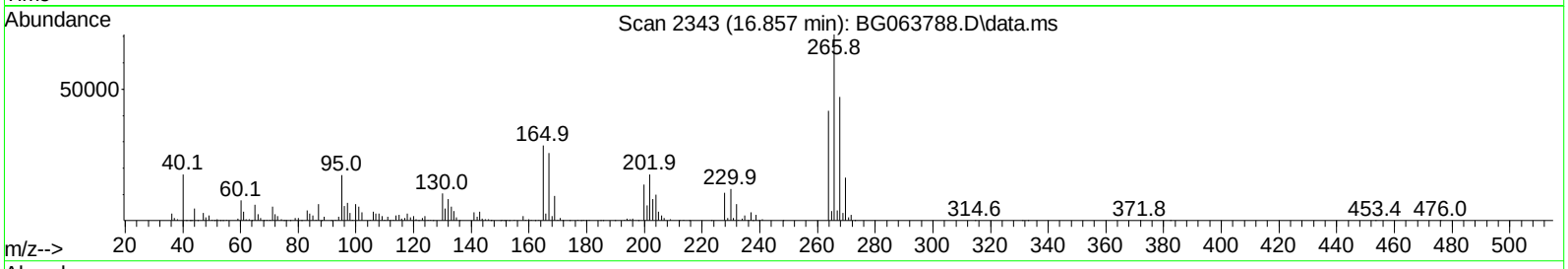
Quant Time: Dec 23 23:53:50 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121124.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Dec 11 23:59:23 2024
Response via : Initial Calibration

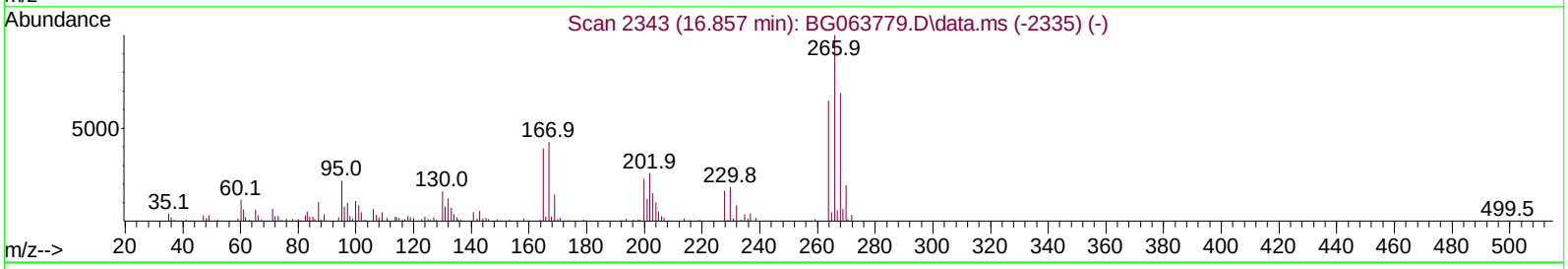
**Instrument :**
BNA_G
**ClientSampleId :**
E2AR6MS

**Manual IntegrationsAPPROVED**

Reviewed By :Yogesh Patel 12/24/2024
Supervised By :mohammad ahmed 12/26/2024




TIC: BG063788.D\data.ms

**(71) Pentachlorophenol (C)**

**16.857min (-0.009) 26.58 ng/ul**

**response 119894**

| Ion | Exp% | Act% |
|---|---|---|
| 265.70 | 100.00 | 100.00 |
| 264.00 | 61.60 | 59.21 |
| 268.00 | 63.10 | 66.48 |
| 0.00 | 0.00 | 0.00 |

Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122324\
Data File : BG063788.D
Acq On : 23 Dec 2024 17:37
Operator : RC/JU
Sample : P5302-04MS
Misc :
ALS Vial : 11 Sample Multiplier: 1

Quant Time: Dec 23 23:53:50 2024
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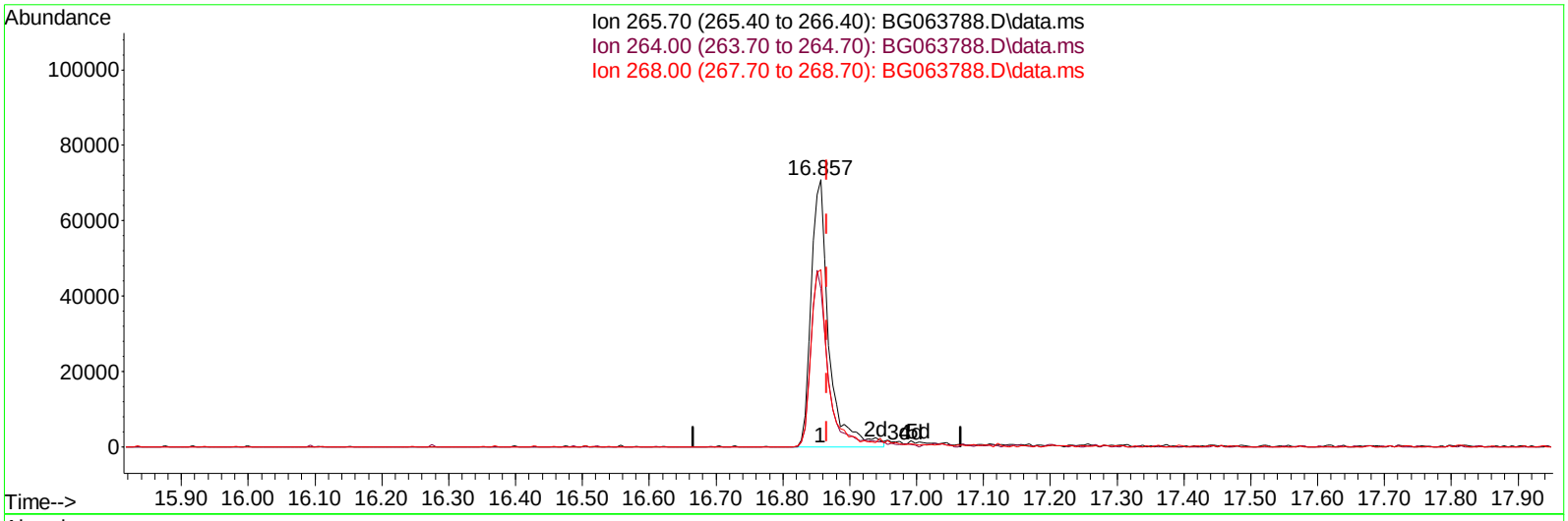
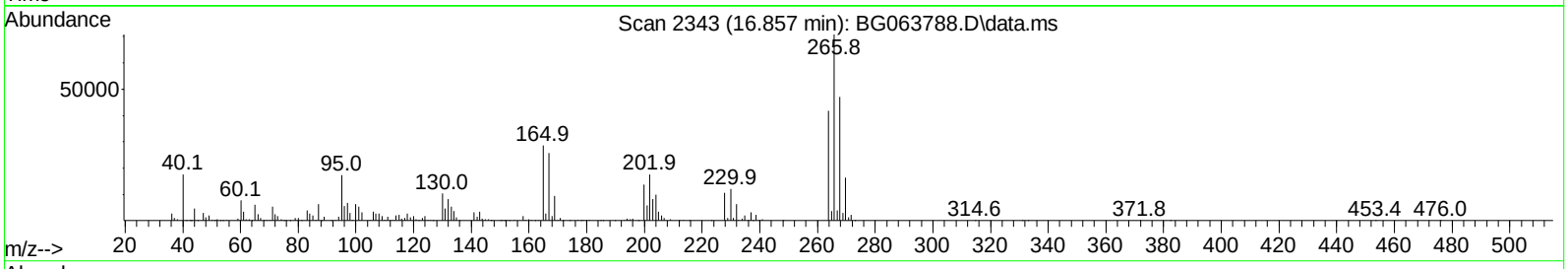
Quant Title : SVOA CALIBRATION
QLast Update : Wed Dec 11 23:59:23 2024
Response via : Initial Calibration

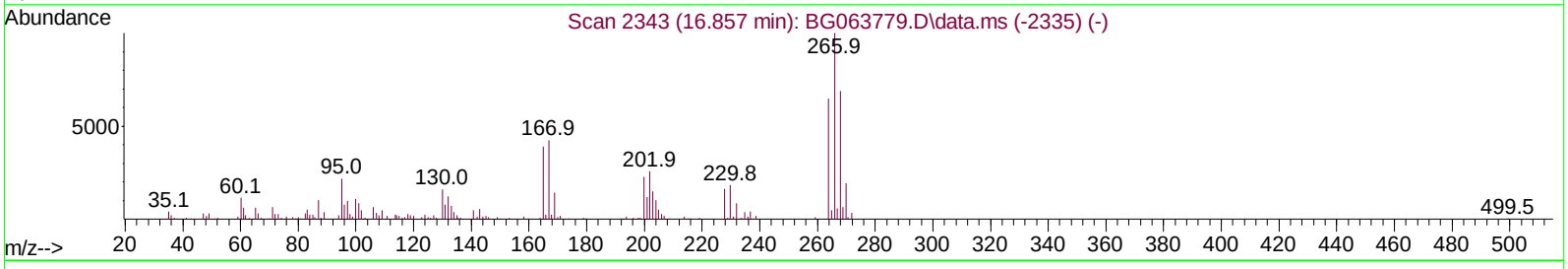
**Instrument :**
BNA_G
**ClientSampleId :**
E2AR6MS

**Manual IntegrationsAPPROVED**

Reviewed By :Yogesh Patel 12/24/2024
Supervised By :mohammad ahmed 12/26/2024




TIC: BG063788.D\data.ms

**(71) Pentachlorophenol (C)**

**16.857min (-0.009) 29.19 ng/ul m**

**response 131663**

| Ion | Exp% | Act% |
|---|---|---|
| 265.70 | 100.00 | 100.00 |
| 264.00 | 61.60 | 59.21 |
| 268.00 | 63.10 | 66.48 |
| 0.00 | 0.00 | 0.00 |

Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122324\
Data File : BG063788.D
Acq On : 23 Dec 2024 17:37
Operator : RC/JU
Sample : P5302-04MS
Misc :
ALS Vial : 11 Sample Multiplier: 1

Quant Time: Dec 23 23:58:17 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121124.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Dec 11 23:59:23 2024
Response via : Initial Calibration

**Instrument :**
BNA_G
**ClientSampleId :**
E2AR6MS

**Manual IntegrationsAPPROVED**

Reviewed By :Yogesh Patel 12/24/2024
Supervised By :mohammad ahmed 12/26/2024

| Compound | R.T. | QIon | Response | Conc | Units | Dev(Min) |
|---|---|---|---|---|---|---|
| Internal Standards | | | | | | |
| 1) 1,4-Dichlorobenzene-d4 | 7.797 | 152 | 108173 | 20.000 | ng/ul | -0.02 |
| 20) Naphthalene-d8 | 10.588 | 136 | 445066 | 20.000 | ng/ul | -0.02 |
| 38) Acenaphthene-d10 | 14.436 | 164 | 367863 | 20.000 | ng/ul | -0.02 |
| 64) Phenanthrene-d10 | 17.192 | 188 | 936976 | 20.000 | ng/ul | -0.02 |
| 79) Chrysene-d12 | 21.440 | 240 | 964530 | 20.000 | ng/ul | -0.02 |
| 88) Perylene-d12 | 24.448 | 264 | 1031283 | 20.000 | ng/ul | -0.03 |
| System Monitoring Compounds | | | | | | |
| 3) 1,4-Dioxane-d8 | 3.261 | 96 | 11524m | 4.175 | ng/uL | -0.02 |
| 4) Pyridine-d5 | 3.666 | 84 | 164236 | 18.861 | ng/ul | -0.02 |
| 7) Phenol-d5 | 6.986 | 99 | 216788 | 21.760 | ng/ul | 0.00 |
| 9) Bis-(2-Chloroethyl)eth... | 7.127 | 67 | 142771 | 20.749 | ng/ul | -0.02 |
| 11) 2-Chlorophenol-d4 | 7.338 | 132 | 134491 | 20.852 | ng/ul | 0.00 |
| 15) 4-Methylphenol-d8 | 8.519 | 113 | 164692 | 21.292 | ng/ul | 0.00 |
| 21) Nitrobenzene-d5 | 8.948 | 128 | 74250 | 23.365 | ng/ul | -0.02 |
| 24) 2-Nitrophenol-d4 | 9.671 | 143 | 83623 | 23.475 | ng/ul | -0.02 |
| 28) 2,4-Dichlorophenol-d3 | 10.217 | 165 | 170348 | 24.700 | ng/ul | -0.02 |
| 31) 4-Chloroaniline-d4 | 10.723 | 131 | 194893 | 22.256 | ng/ul | -0.02 |
| 46) Dimethylphthalate-d6 | 13.843 | 166 | 609754 | 24.360 | ng/ul | -0.02 |
| 49) Acenaphthylene-d8 | 14.131 | 160 | 661766 | 23.192 | ng/ul | -0.02 |
| 54) 4-Nitrophenol-d4 | 14.671 | 143 | 86868 | 24.493 | ng/ul | 0.00 |
| 60) Fluorene-d10 | 15.435 | 176 | 521390 | 23.559 | ng/ul | -0.02 |
| 65) 4,6-Dinitro-2-methylph... | 15.570 | 200 | 96633 | 19.864 | ng/ul | 0.00 |
| 73) Anthracene-d10 | 17.292 | 188 | 951463 | 23.619 | ng/ul | -0.02 |
| 81) Pyrene-d10 | 19.589 | 212 | 1025190 | 20.425 | ng/ul | -0.02 |
| 92) Benzo(a)pyrene-d12 | 24.242 | 264 | 933835 | 19.282 | ng/ul | -0.03 |
| Target Compounds | | | | | | Qvalue |
| 2) 1,4-Dioxane | 3.296 | 88 | 35423 | 10.843 | ng/uL | 95 |
| 5) Pyridine | 3.684 | 79 | 254115 | 27.245 | ng/ul | 94 |
| 6) Benzaldehyde | 6.945 | 77 | 25745 | 4.213 | ng/ul | 93 |
| 8) Phenol | 7.009 | 94 | 325998 | 30.703 | ng/ul | 99 |
| 10) Bis(2-Chloroethyl)ether | 7.221 | 93 | 230876 | 28.420 | ng/ul | 93 |
| 12) 2-Chlorophenol | 7.368 | 128 | 198389 | 30.387 | ng/ul | 98 |
| 13) 2-Methylphenol | 8.249 | 108 | 228465 | 29.771 | ng/ul | 97 |
| 14) 2,2'-oxybis(1-Chloropr... | 8.314 | 45 | 343187 | 28.456 | ng/ul | 98 |
| 16) Acetophenone | 8.619 | 105 | 375312 | 28.607 | ng/ul | 95 |
| 17) N-Nitroso-di-n-propyla... | 8.602 | 70 | 222929 | 28.478 | ng/ul | 96 |
| 18) 4-Methylphenol | 8.578 | 108 | 254638 | 30.318 | ng/ul | 96 |
| 19) Hexachloroethane | 8.872 | 117 | 82941 | 25.855 | ng/ul | 92 |
| 22) Nitrobenzene | 8.995 | 77 | 312622 | 29.460 | ng/ul | 98 |
| 23) Isophorone | 9.518 | 82 | 620452 | 30.945 | ng/ul | 98 |
| 25) 2-Nitrophenol | 9.706 | 139 | 115492 | 31.548 | ng/ul | 86 |
| 26) 2,4-Dimethylphenol | 9.777 | 107 | 261883 | 31.525 | ng/ul | 95 |
| 27) Bis(2-Chloroethoxy)met... | 10.000 | 93 | 342339 | 31.567 | ng/ul | 96 |
| 29) 2,4-Dichlorophenol | 10.247 | 162 | 227713 | 33.229 | ng/ul | 95 |
| 30) Naphthalene | 10.635 | 128 | 692419 | 31.117 | ng/ul | 99 |
| 32) 4-Chloroaniline | 10.752 | 127 | 171372 | 20.569 | ng/ul | 98 |
| 33) Hexachlorobutadiene | 10.923 | 225 | 151483 | 27.209 | ng/ul | 99 |
| 34) Caprolactam | 11.498 | 113 | 89092m | 37.610 | ng/ul | |
| 35) 4-Chloro-3-methylphenol | 11.892 | 107 | 275882 | 34.291 | ng/ul | 95 |

Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122324\
Data File : BG063788.D
Acq On : 23 Dec 2024 17:37
Operator : RC/JU
Sample : P5302-04MS
Misc :
ALS Vial : 11 Sample Multiplier: 1

Quant Time: Dec 23 23:58:17 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121124.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Dec 11 23:59:23 2024
Response via : Initial Calibration

**Instrument :**
BNA_G
**ClientSampleId :**
E2AR6MS

**Manual IntegrationsAPPROVED**

Reviewed By :Yogesh Patel 12/24/2024
Supervised By :mohammad ahmed 12/26/2024

| Compound | R.T. | QIon | Response | Conc Units | Dev(Min) |
|---|---|---|---|---|---|
| 36) 2-Methylnaphthalene | 12.250 | 142 | 515083 | 32.208 ng/ul | 97 |
| 37) 1-Methylnaphthalene | 12.474 | 142 | 519328 | 31.880 ng/ul | 98 |
| 39) 1,2,4,5-Tetrachloroben... | 12.626 | 216 | 312997 | 26.578 ng/ul | 99 |
| 40) Hexachlorocyclopentadiene | 12.603 | 237 | 72062 | 15.479 ng/ul | 88 |
| 41) 2,4,6-Trichlorophenol | 12.873 | 196 | 196855 | 29.576 ng/ul | 88 |
| 42) 2,4,5-Trichlorophenol | 12.955 | 196 | 214219 | 30.840 ng/ul | 99 |
| 43) 1,1'-Biphenyl | 13.267 | 154 | 731473 | 30.151 ng/ul | 96 |
| 44) 2-Chloronaphthalene | 13.308 | 162 | 581607 | 29.855 ng/ul | 98 |
| 45) 2-Nitroaniline | 13.519 | 65 | 215590 | 33.038 ng/ul | 96 |
| 47) Dimethylphthalate | 13.896 | 163 | 798570 | 31.832 ng/ul | 98 |
| 48) 2,6-Dinitrotoluene | 14.019 | 165 | 165294 | 35.478 ng/ul# | 91 |
| 50) Acenaphthylene | 14.160 | 152 | 974230 | 31.842 ng/ul | 98 |
| 51) 3-Nitroaniline | 14.348 | 138 | 152375 | 35.373 ng/ul | 98 |
| 52) Acenaphthene | 14.501 | 153 | 677404 | 31.237 ng/ul | 96 |
| 53) 2,4-Dinitrophenol | 14.571 | 184 | 82736m | 34.748 ng/ul | |
| 55) 4-Nitrophenol | 14.683 | 109 | 105476 | 28.641 ng/ul | 96 |
| 56) Dibenzofuran | 14.841 | 168 | 980180 | 32.060 ng/ul | 97 |
| 57) 2,4-Dinitrotoluene | 14.812 | 165 | 248308 | 35.867 ng/ul# | 100 |
| 58) 2,3,4,6-Tetrachlorophenol | 15.076 | 232 | 196161 | 32.761 ng/ul | 94 |
| 59) Diethylphthalate | 15.259 | 149 | 822283 | 33.906 ng/ul | 96 |
| 61) Fluorene | 15.494 | 166 | 803428 | 32.910 ng/ul | 100 |
| 62) 4-Chlorophenyl-phenyle... | 15.482 | 204 | 434649 | 31.958 ng/ul | 96 |
| 63) 4-Nitroaniline | 15.517 | 138 | 179851 | 41.704 ng/ul | 94 |
| 66) 4,6-Dinitro-2-methylph... | 15.588 | 198 | 163897 | 32.702 ng/ul | 90 |
| 67) N-Nitrosodiphenylamine | 15.699 | 169 | 721817 | 31.813 ng/ul | 98 |
| 68) 4-Bromophenyl-phenylether | 16.381 | 248 | 297686 | 31.174 ng/ul | 97 |
| 69) Hexachlorobenzene | 16.504 | 284 | 325333 | 30.409 ng/ul | 96 |
| 70) Atrazine | 16.657 | 200 | 303243 | 31.617 ng/ul | 96 |
| 71) Pentachlorophenol | 16.857 | 266 | 131663m | 29.188 ng/ul | |
| 72) Phenanthrene | 17.233 | 178 | 1491990 | 33.043 ng/ul | 99 |
| 74) Anthracene | 17.327 | 178 | 1493209 | 32.674 ng/ul | 99 |
| 75) 1,2,3,4-Tetrachloroben... | 13.237 | 216 | 326277 | 26.099 ng/ul | 97 |
| 76) Pentachlorobenzene | 14.765 | 250 | 346567 | 27.983 ng/ul | 99 |
| 77) Carbazole | 17.597 | 167 | 1307279 | 35.338 ng/ul | 99 |
| 78) Di-n-butylphthalate | 18.155 | 149 | 1547385 | 34.636 ng/ul | 98 |
| 80) Fluoranthene | 19.254 | 202 | 1837273 | 32.206 ng/ul | 98 |
| 82) Pyrene | 19.618 | 202 | 1939078 | 31.948 ng/ul# | 94 |
| 83) Butylbenzylphthalate | 20.511 | 149 | 645938 | 34.669 ng/ul | 98 |
| 84) 3,3'-Dichlorobenzidine | 21.340 | 252 | 465096 | 26.712 ng/ul | 97 |
| 85) Benzo(a)anthracene | 21.422 | 228 | 1866317 | 30.875 ng/ul | 99 |
| 86) Bis(2-ethylhexyl)phtha... | 21.334 | 149 | 965069 | 35.608 ng/ul | 98 |
| 87) Chrysene | 21.481 | 228 | 1799704 | 31.238 ng/ul | 98 |
| 89) Di-n-octyl phthalate | 22.462 | 149 | 1643422 | 36.921 ng/ul | 100 |
| 90) Benzo(b)fluoranthene | 23.502 | 252 | 1870729 | 31.916 ng/ul | 99 |
| 91) Benzo(k)fluoranthene | 23.561 | 252 | 1856031 | 31.745 ng/ul | 100 |
| 93) Benzo(a)pyrene | 24.313 | 252 | 1723051 | 31.319 ng/ul | 100 |
| 94) Indeno(1,2,3-cd)pyrene | 27.850 | 276 | 2149529 | 31.993 ng/ul | 98 |
| 95) Dibenzo(a,h)anthracene | 27.897 | 278 | 1727593 | 32.135 ng/ul | 96 |
| 96) Benzo(g,h,i)perylene | 28.919 | 276 | 1665743 | 31.243 ng/ul | 99 |

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

**Instrument :**
BNA_G
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E2AR6MS

**Manual IntegrationsAPPROVED**

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Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122324\
Data File : BG063788.D
Acq On : 23 Dec 2024 17:37
Operator : RC/JU
Sample : P5302-04MS
Misc :
ALS Vial : 11 Sample Multiplier: 1

Quant Time: Dec 23 23:58:17 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121124.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Dec 11 23:59:23 2024
Response via : Initial Calibration

**Instrument :**
BNA_G
**ClientSampleId :**
E2AR6MS

**Manual IntegrationsAPPROVED**

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