

Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG060917\  
 Data File : BG027362.D  
 Acq On : 10 Jun 2017 00:43  
 Operator : SJ/MA  
 Sample : I3563-03  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 OK-04-060817

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:\HPCHEM1\BNA\_G\METHODS\8270-BG060517.M  
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

| peak<br># | R.T.<br>min | first<br>scan | max<br>scan | last<br>scan | PK<br>TY | peak<br>height | corr.<br>area | corr.<br>% max. | % of<br>total |
|-----------|-------------|---------------|-------------|--------------|----------|----------------|---------------|-----------------|---------------|
| 1         | 4.552       | 142           | 147         | 158          | rVB      | 148620         | 224048        | 1.89%           | 0.485%        |
| 2         | 4.669       | 160           | 167         | 175          | rBV      | 95083          | 154006        | 1.30%           | 0.333%        |
| 3         | 5.128       | 238           | 245         | 254          | rBV      | 896293         | 1463112       | 12.32%          | 3.165%        |
| 4         | 5.221       | 254           | 261         | 300          | rBV      | 4067685        | 10133010      | 85.35%          | 21.918%       |
| 5         | 6.144       | 408           | 418         | 428          | rBV      | 914424         | 1577932       | 13.29%          | 3.413%        |
| 6         | 7.102       | 575           | 581         | 588          | rBV      | 277861         | 434157        | 3.66%           | 0.939%        |
| 7         | 7.648       | 663           | 674         | 680          | rBV      | 6274904        | 11872064      | 100.00%         | 25.680%       |
| 8         | 8.089       | 741           | 749         | 760          | rVB      | 976449         | 1789535       | 15.07%          | 3.871%        |
| 9         | 8.553       | 822           | 828         | 836          | rVB      | 173364         | 320996        | 2.70%           | 0.694%        |
| 10        | 8.882       | 877           | 884         | 896          | rVB      | 654227         | 1260977       | 10.62%          | 2.728%        |
| 11        | 9.716       | 1018          | 1026        | 1038         | rVB      | 593970         | 1137164       | 9.58%           | 2.460%        |
| 12        | 10.086      | 1082          | 1089        | 1101         | rBV      | 1459294        | 2372552       | 19.98%          | 5.132%        |
| 13        | 11.373      | 1295          | 1308        | 1314         | rBV      | 263929         | 531131        | 4.47%           | 1.149%        |
| 14        | 12.566      | 1505          | 1511        | 1550         | rBV2     | 655093         | 1883821       | 15.87%          | 4.075%        |
| 15        | 13.759      | 1707          | 1714        | 1722         | rBV      | 1663610        | 2666776       | 22.46%          | 5.768%        |
| 16        | 14.064      | 1760          | 1766        | 1789         | rBV      | 426850         | 1151134       | 9.70%           | 2.490%        |
| 17        | 15.133      | 1942          | 1948        | 1956         | rBV2     | 147878         | 273799        | 2.31%           | 0.592%        |
| 18        | 15.462      | 1994          | 2004        | 2020         | rVB      | 109085         | 257164        | 2.17%           | 0.556%        |
| 19        | 16.620      | 2193          | 2201        | 2209         | rBV      | 1268708        | 2090826       | 17.61%          | 4.523%        |
| 20        | 17.883      | 2409          | 2416        | 2428         | rBV2     | 405929         | 721328        | 6.08%           | 1.560%        |
| 21        | 20.462      | 2848          | 2855        | 2864         | rBV2     | 2448148        | 3459887       | 29.14%          | 7.484%        |
| 22        | 22.272      | 3155          | 3163        | 3175         | rVB3     | 204168         | 456213        | 3.84%           | 0.987%        |

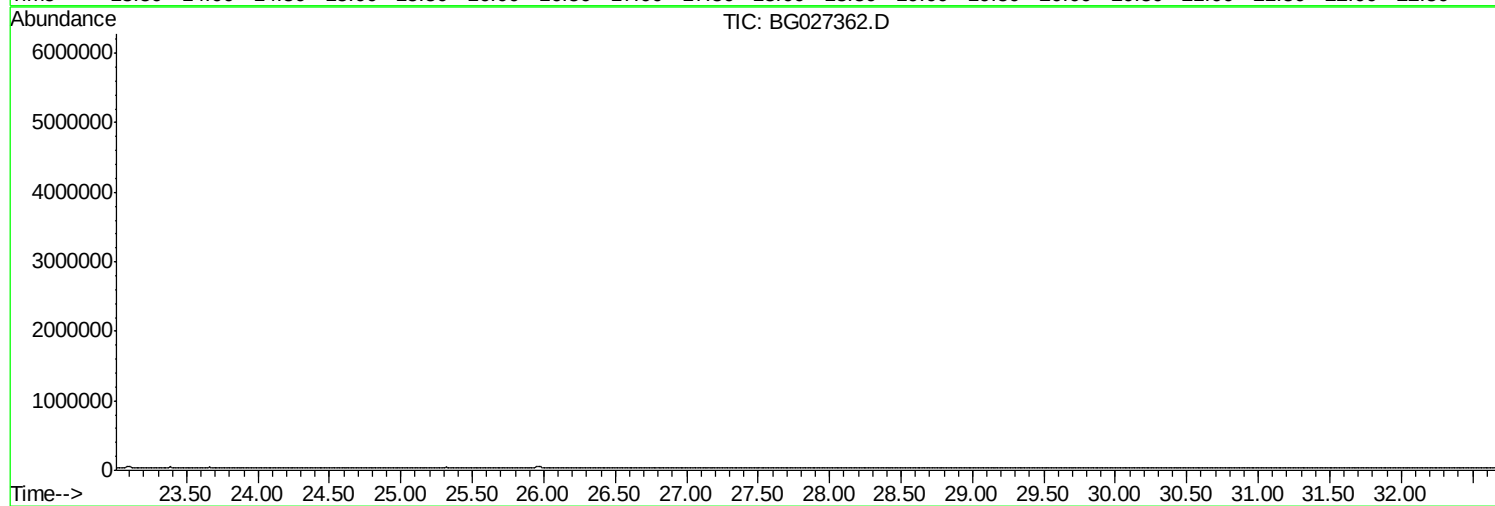
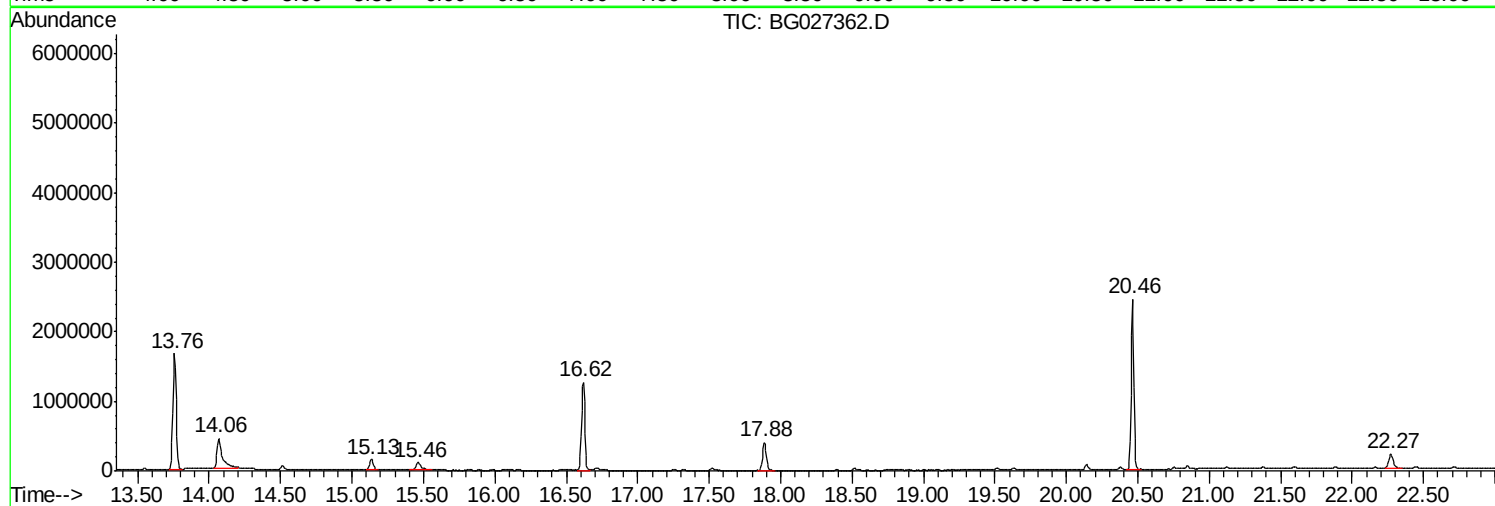
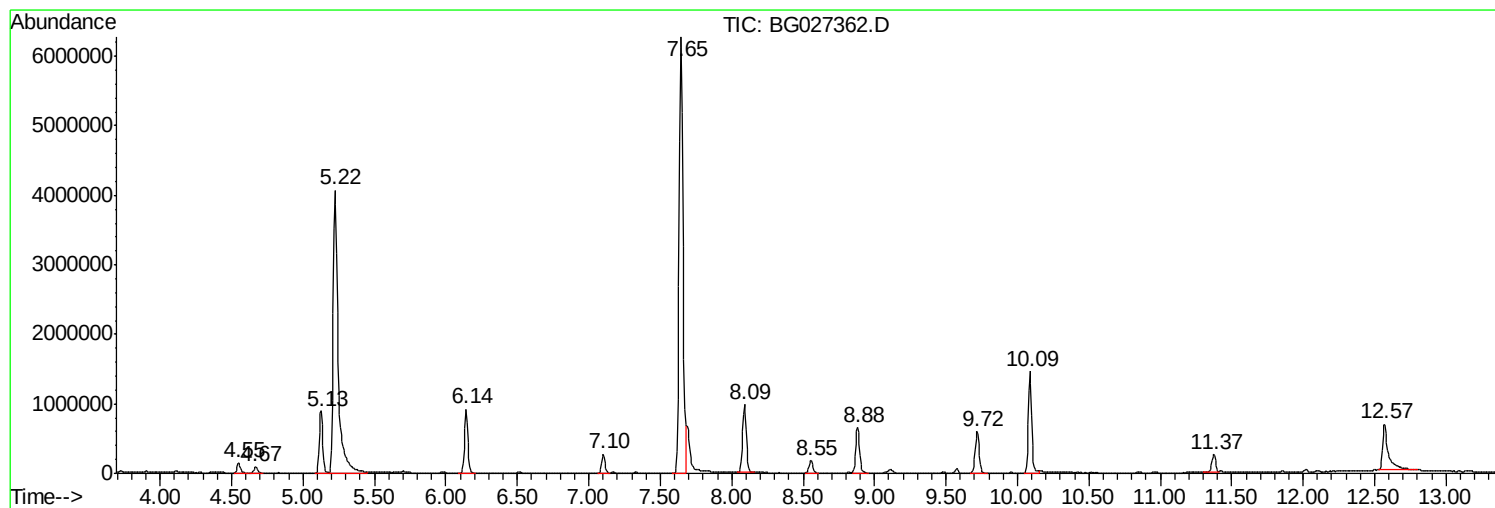
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ClientSampleId :  
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Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG060517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P



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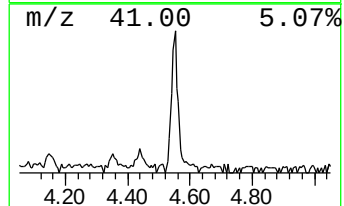
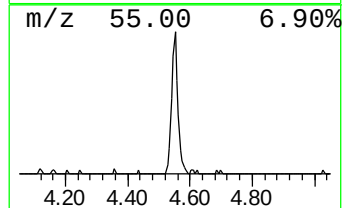
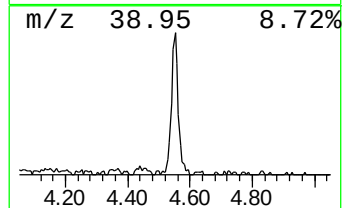
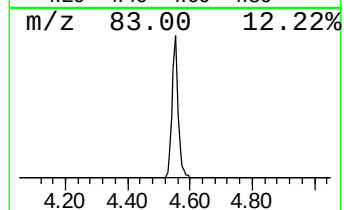
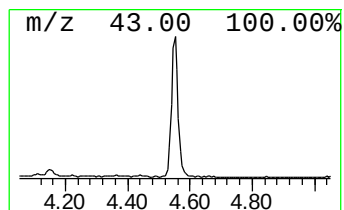
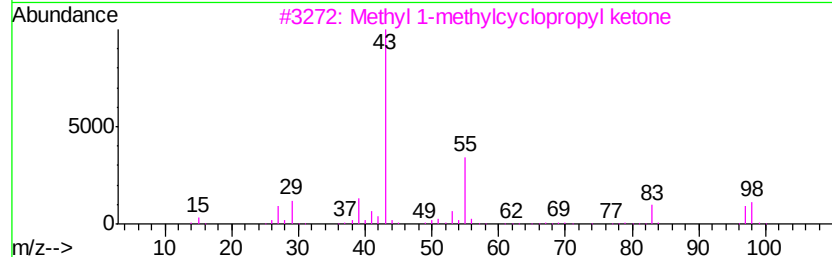
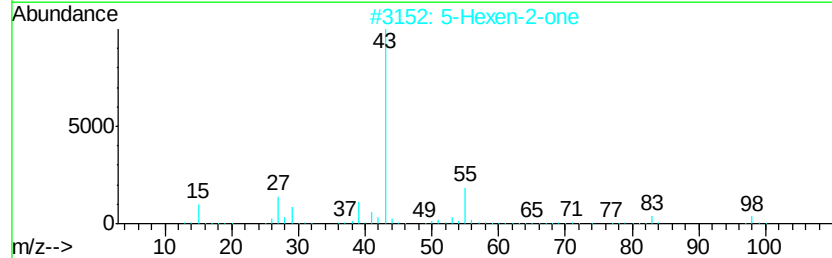
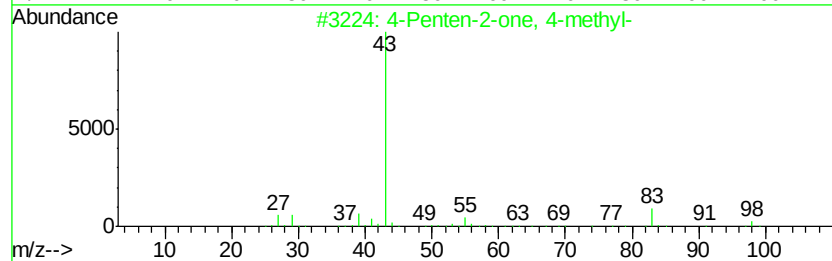
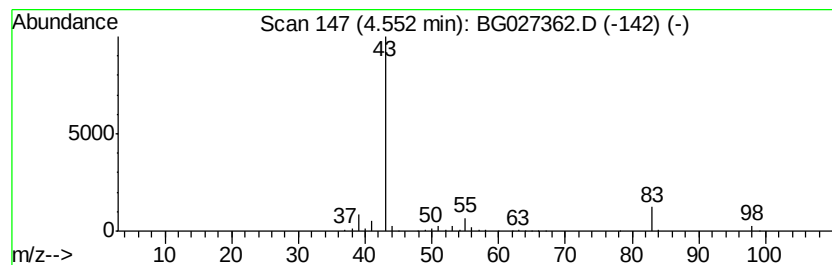
Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG060517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 1 4-Penten-2-one, 4-methyl- Concentration Rank 10

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 4.55 | 13.96 ng | 224048 | 1,4-Dichlorobenzene-d4 | 8.56 |

| Hit# | of | Tentative ID                      | MW  | MolForm | CAS#        | Qual |
|------|----|-----------------------------------|-----|---------|-------------|------|
| 1    | 5  | 4-Penten-2-one, 4-methyl-         | 98  | C6H10O  | 003744-02-3 | 72   |
| 2    |    | 5-Hexen-2-one                     | 98  | C6H10O  | 000109-49-9 | 49   |
| 3    |    | Methyl 1-methylcyclopropyl ketone | 98  | C6H10O  | 001567-75-5 | 40   |
| 4    |    | 2-Vinylethyl acetate              | 114 | C6H10O2 | 001576-84-7 | 25   |
| 5    |    | Isoxazole, 5-methyl-              | 83  | C4H5NO  | 005765-44-6 | 9    |



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Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG060517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

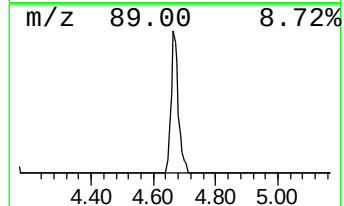
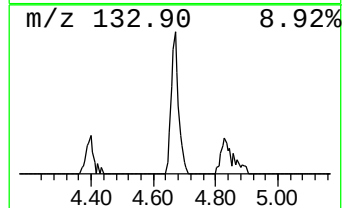
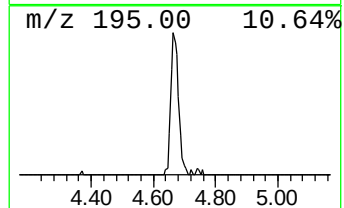
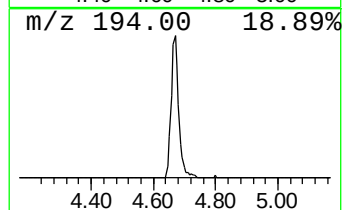
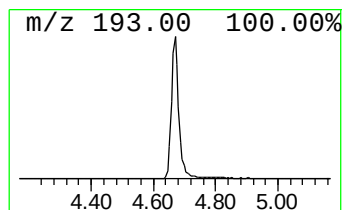
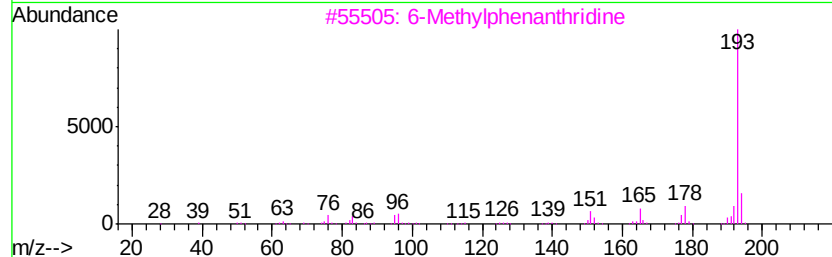
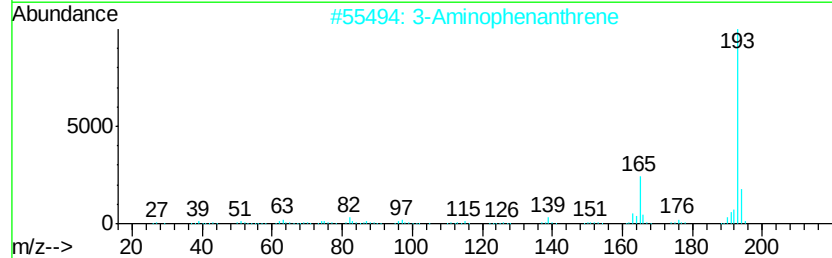
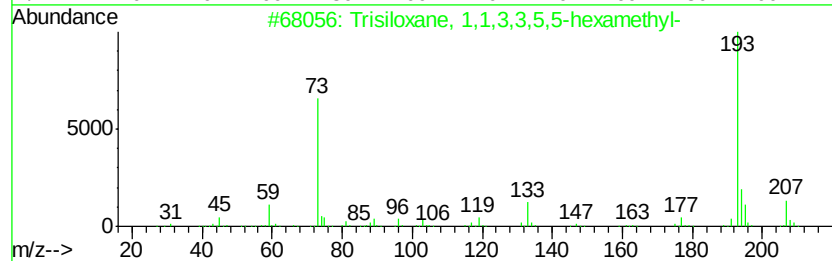
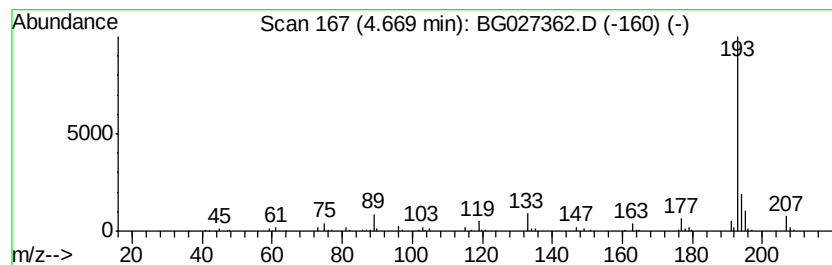
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 2 Trisiloxane, 1,1,3,3,5,5-he... Concentration Rank 11

| R.T. | EstConc | Area   | Relative to ISTD       | R.T. |
|------|---------|--------|------------------------|------|
| 4.67 | 9.60 ng | 154006 | 1,4-Dichlorobenzene-d4 | 8.56 |

| Hit# | of | 5 | Tentative ID                        | MW  | MolForm    | CAS#        | Qual |
|------|----|---|-------------------------------------|-----|------------|-------------|------|
| 1    |    |   | Trisiloxane, 1,1,3,3,5,5-hexamet... | 208 | C6H2002Si3 | 001189-93-1 | 80   |
| 2    |    |   | 3-Aminophenanthrene                 | 193 | C14H11N    | 001892-54-2 | 64   |
| 3    |    |   | 6-Methylphenanthridine              | 193 | C14H11N    | 003955-65-5 | 64   |
| 4    |    |   | 3-Phenylindole                      | 193 | C14H11N    | 001504-16-1 | 59   |
| 5    |    |   | 1-Anthracenamine                    | 193 | C14H11N    | 000610-49-1 | 59   |



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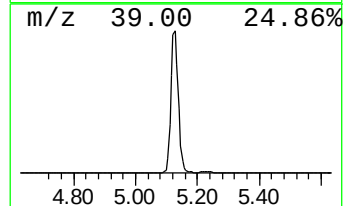
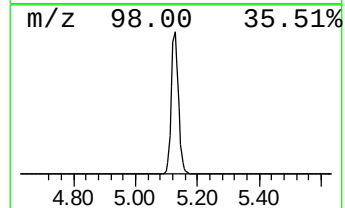
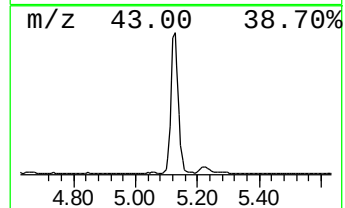
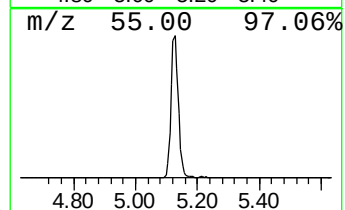
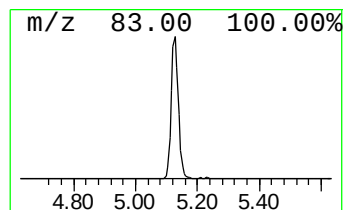
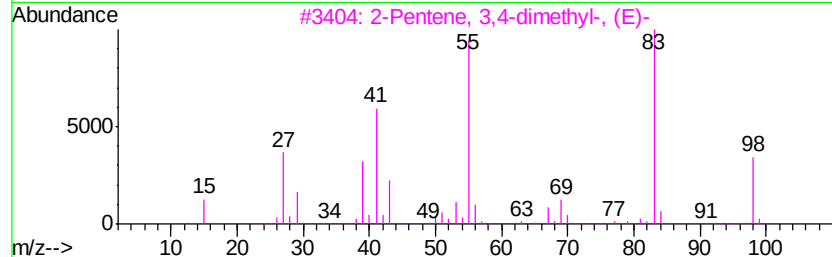
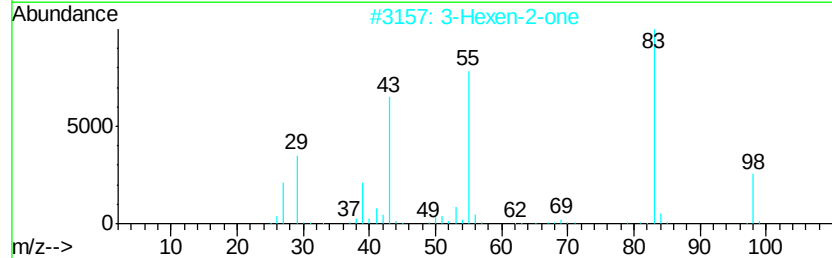
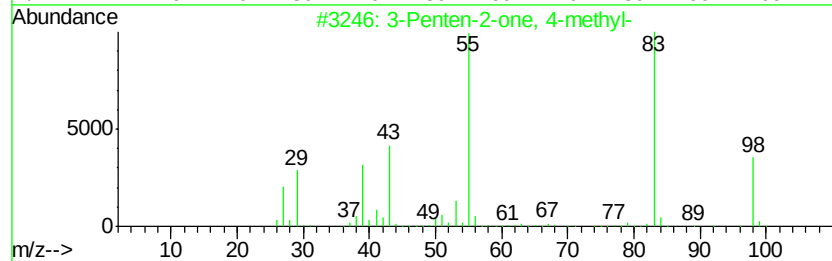
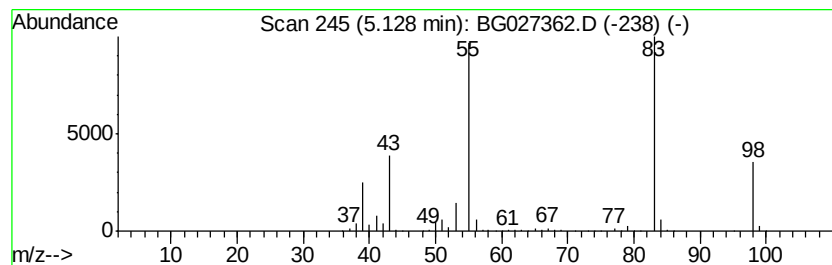
Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG060517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 3 3-Penten-2-one, 4-methyl- Concentration Rank 3

| R.T. | EstConc  | Area    | Relative to ISTD       | R.T. |
|------|----------|---------|------------------------|------|
| 5.13 | 91.16 ng | 1463110 | 1,4-Dichlorobenzene-d4 | 8.56 |

| Hit# | of | Tentative ID                   | MW | MolForm | CAS#        | Qual |
|------|----|--------------------------------|----|---------|-------------|------|
| 1    | 5  | 3-Penten-2-one, 4-methyl-      | 98 | C6H10O  | 000141-79-7 | 91   |
| 2    |    | 3-Hexen-2-one                  | 98 | C6H10O  | 000763-93-9 | 91   |
| 3    |    | 2-Pentene, 3,4-dimethyl-, (E)- | 98 | C7H14   | 004914-92-5 | 90   |
| 4    |    | 2-Pentene, 2,4-dimethyl-       | 98 | C7H14   | 000625-65-0 | 87   |
| 5    |    | 2-Pentene, 3,4-dimethyl-, (Z)- | 98 | C7H14   | 004914-91-4 | 86   |



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Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG060517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

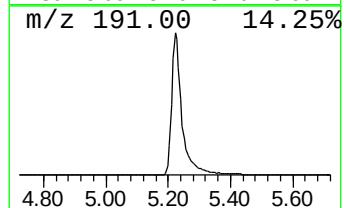
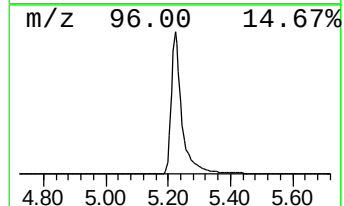
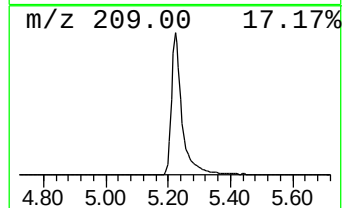
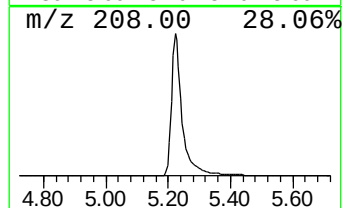
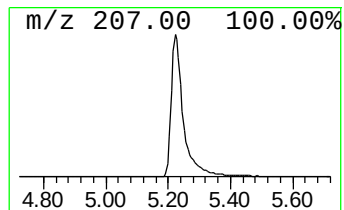
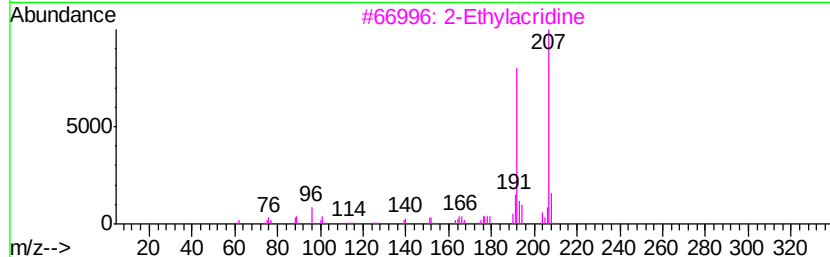
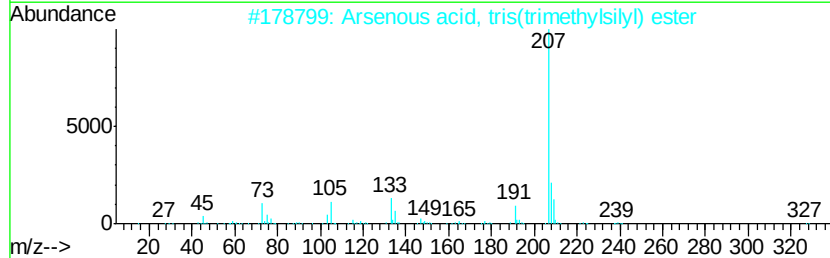
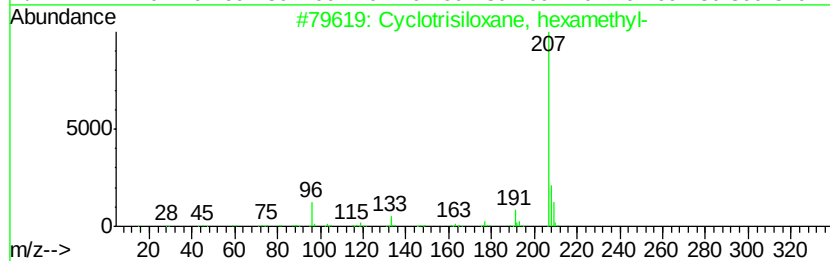
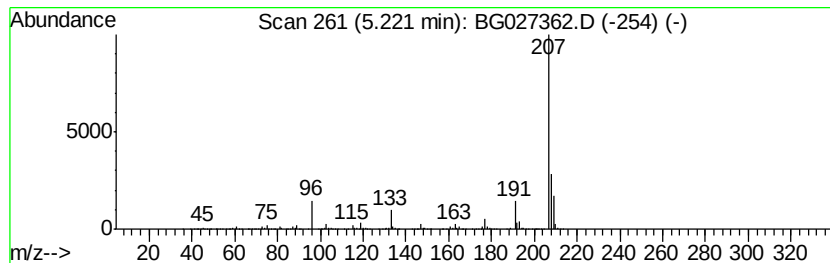
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 4 Cyclotrisiloxane, hexamethyl- Concentration Rank 1

| R.T. | EstConc   | Area     | Relative to ISTD       | R.T. |
|------|-----------|----------|------------------------|------|
| 5.22 | 631.35 ng | 10133000 | 1,4-Dichlorobenzene-d4 | 8.56 |

| Hit# of | 5 | Tentative ID                        | MW  | MolForm      | CAS#         | Qual |
|---------|---|-------------------------------------|-----|--------------|--------------|------|
| 1       |   | Cyclotrisiloxane, hexamethyl-       | 222 | C6H18O3Si3   | 000541-05-9  | 91   |
| 2       |   | Arsenous acid, tris(trimethylsil... | 342 | C9H27AsO3Si3 | 055429-29-3  | 78   |
| 3       |   | 2-Ethylacridine                     | 207 | C15H13N      | 055751-83-2  | 38   |
| 4       |   | 2-p-Nitrophenyl-oxadiazol-1,3,4-... | 207 | C8H5N3O4     | 1000147-64-6 | 9    |
| 5       |   | 1,3,5-Triazine, 2-chloro-4,6-bis... | 207 | C5H6ClN3S2   | 004407-40-3  | 9    |



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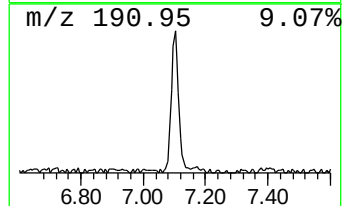
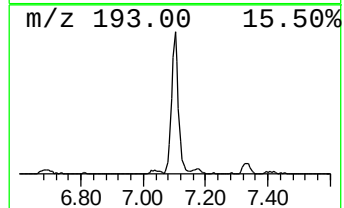
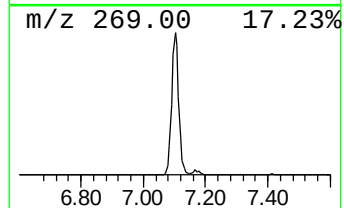
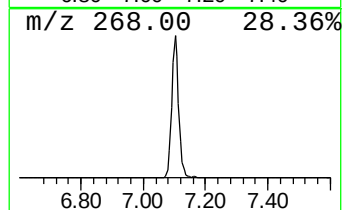
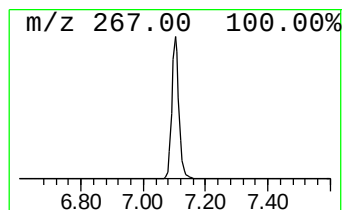
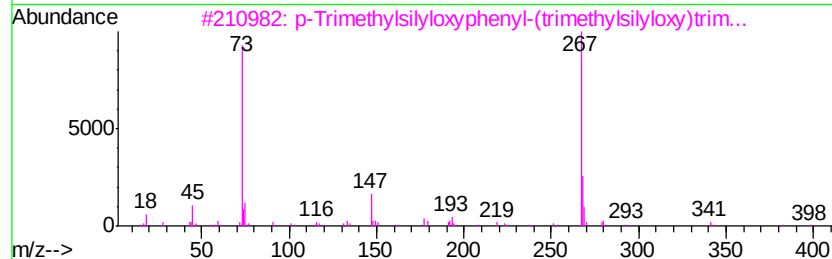
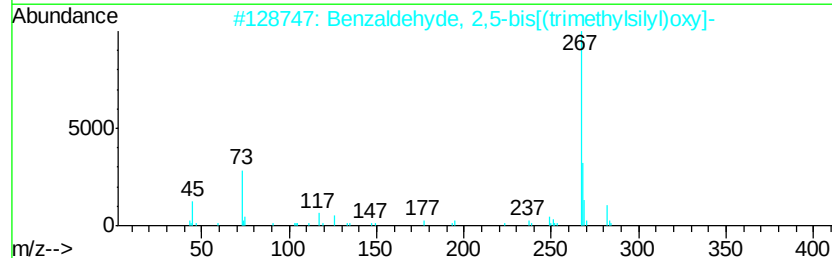
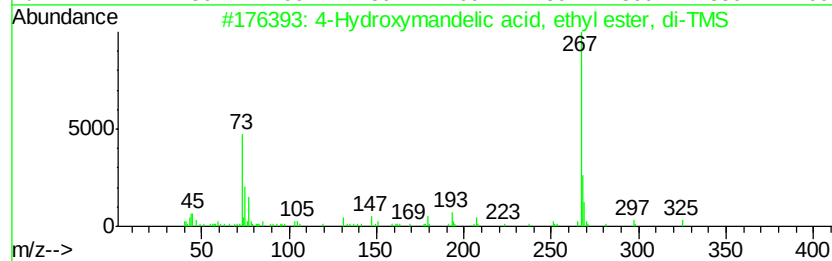
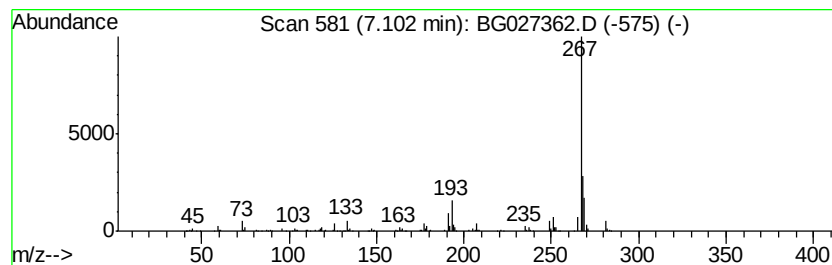
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TIC Integration Parameters: LSCINT.P

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Peak Number 5 4-Hydroxymandelic acid, eth... Concentration Rank 8

| R.T. | EstConc  | Area   | Relative to ISTD       | R.T. |
|------|----------|--------|------------------------|------|
| 7.10 | 27.05 ng | 434157 | 1,4-Dichlorobenzene-d4 | 8.56 |

| Hit# of | 5 | Tentative ID                        | MW  | MolForm      | CAS#         | Qual |
|---------|---|-------------------------------------|-----|--------------|--------------|------|
| 1       |   | 4-Hydroxymandelic acid, ethyl es... | 340 | C16H28O4Si2  | 1000071-53-3 | 64   |
| 2       |   | Benzaldehyde, 2,5-bis[(trimethyl... | 282 | C13H22O3Si2  | 056114-69-3  | 59   |
| 3       |   | p-Trimethylsilyloxyphenyl-(trime... | 398 | C18H34O4Si3  | 1000079-05-1 | 59   |
| 4       |   | Benzeneethanamine, N-butyl-.beta... | 353 | C18H35N02Si2 | 040629-66-1  | 59   |
| 5       |   | 3-Hydroxymandelic acid, ethyl es... | 340 | C16H28O4Si2  | 1000071-88-9 | 59   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG060917\  
Data File : BG027362.D  
Acq On : 10 Jun 2017 00:43  
Operator : SJ/MA  
Sample : I3563-03  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
OK-04-060817

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG060517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

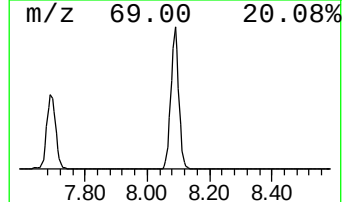
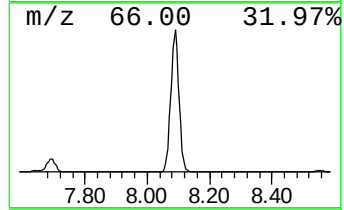
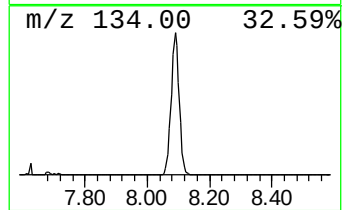
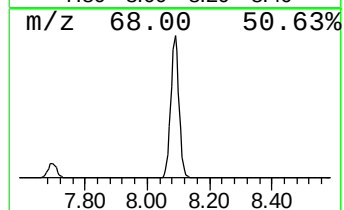
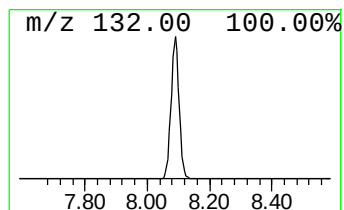
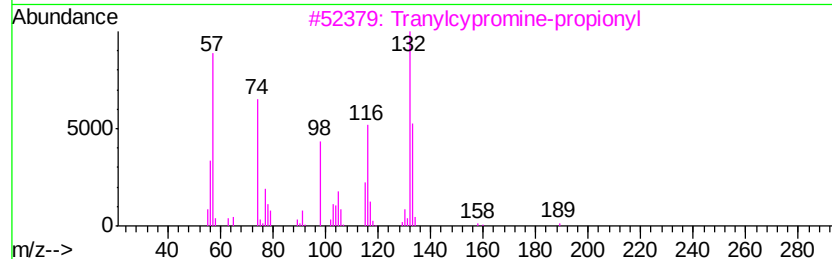
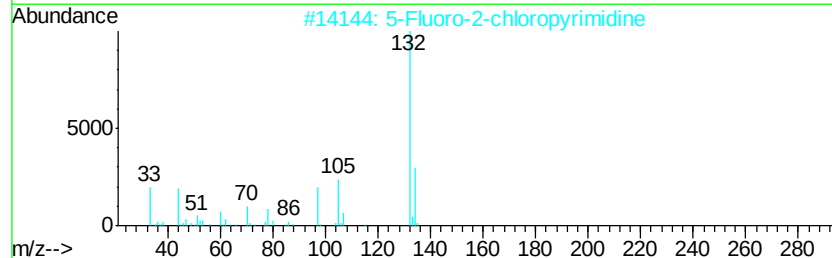
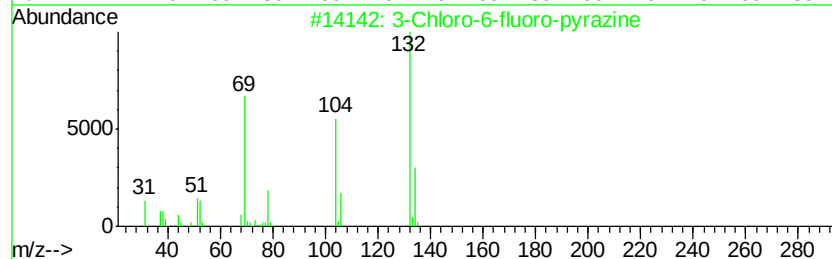
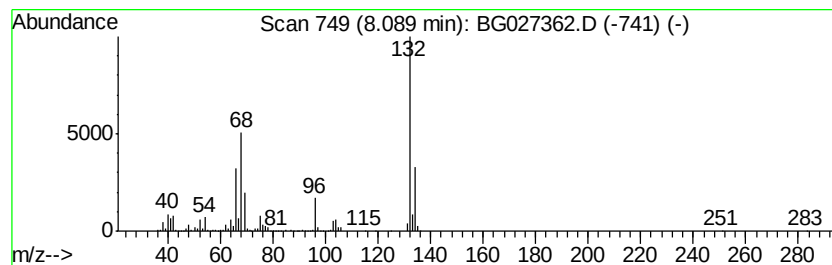
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 6 unknown8.09 Concentration Rank 2

| R.T. | EstConc   | Area    | Relative to ISTD       | R.T. |
|------|-----------|---------|------------------------|------|
| 8.09 | 111.50 ng | 1789540 | 1,4-Dichlorobenzene-d4 | 8.56 |

| Hit# of 5 | Tentative ID                     | MW  | MolForm   | CAS#         | Qual |
|-----------|----------------------------------|-----|-----------|--------------|------|
| 1         | 3-Chloro-6-fluoro-pyrazine       | 132 | C4H2ClFN2 | 1000146-10-7 | 27   |
| 2         | 5-Fluoro-2-chloropyrimidine      | 132 | C4H2ClFN2 | 062802-42-0  | 12   |
| 3         | Tranylcypromine-propionyl        | 189 | C12H15NO  | 1000123-86-3 | 12   |
| 4         | (5-Methyl-2-pyridyl)acetonitrile | 132 | C8H8N2    | 1000241-93-9 | 10   |
| 5         | 2-Cyclohexen-1-one               | 96  | C6H8O     | 000930-68-7  | 9    |





Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG060917\  
Data File : BG027362.D  
Acq On : 10 Jun 2017 00:43  
Operator : SJ/MA  
Sample : I3563-03  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
OK-04-060817

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG060517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

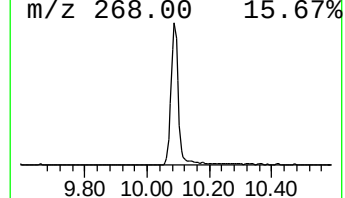
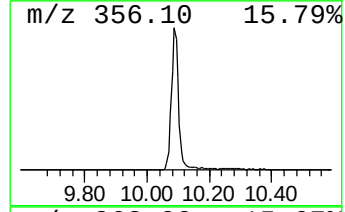
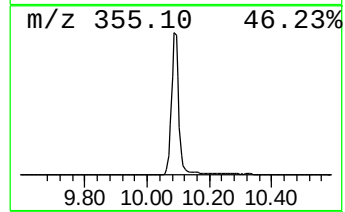
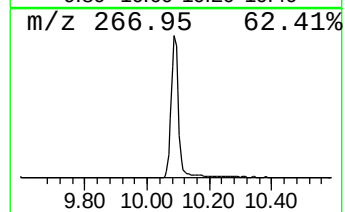
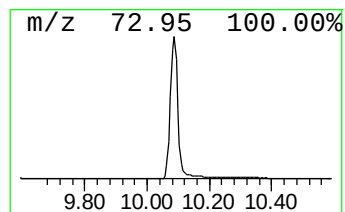
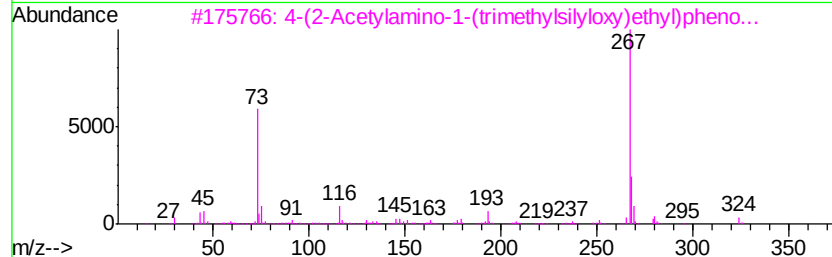
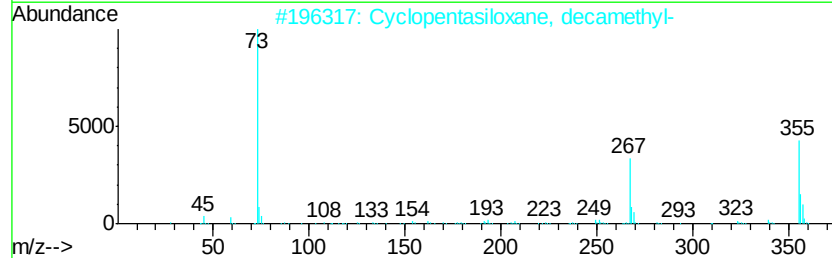
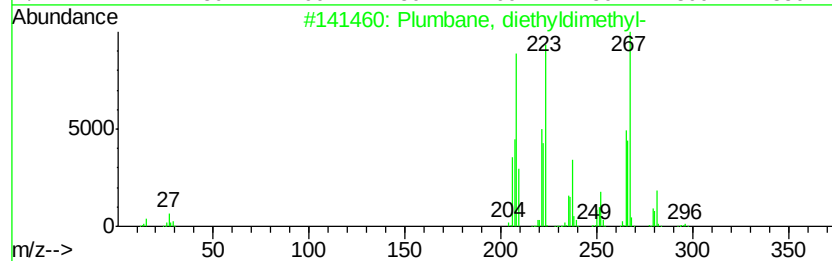
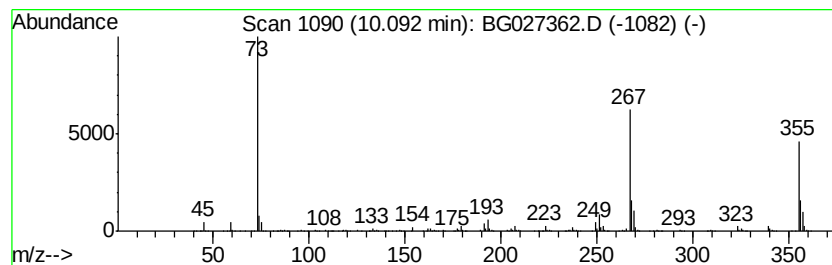
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 8 Plumbane, diethyldimethyl- Concentration Rank 4

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 10.09 | 89.34 ng | 2372550 | Naphthalene-d8   | 11.37 |

| Hit# | of | Tentative ID                         | MW  | MolForm        | CAS#         | Qual |
|------|----|--------------------------------------|-----|----------------|--------------|------|
| 1    | 5  | Plumbane, diethyldimethyl-           | 296 | C6H16Pb        | 001762-27-2  | 78   |
| 2    |    | Cyclopentasiloxane, decamethyl-      | 370 | C10H30O5Si5    | 000541-02-6  | 74   |
| 3    |    | 4-(2-Acetylamino-1-(trimethylsil...  | 339 | C16H29N03Si2   | 1000373-43-5 | 43   |
| 4    |    | 3-Amino-2-phenazinol ditms           | 355 | C18H25N3O5Si2  | 1000332-15-5 | 38   |
| 5    |    | N-(Trifluoroacetyl)-O,0',0''-tris... | 495 | C20H36F3N04Si3 | 054135-51-2  | 37   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG060917\  
Data File : BG027362.D  
Acq On : 10 Jun 2017 00:43  
Operator : SJ/MA  
Sample : I3563-03  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
OK-04-060817

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG060517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

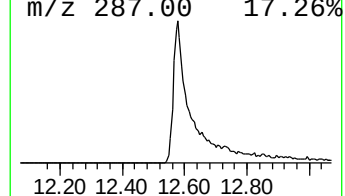
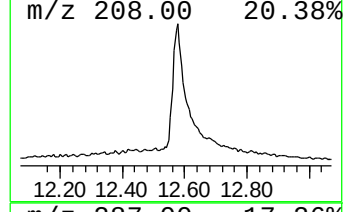
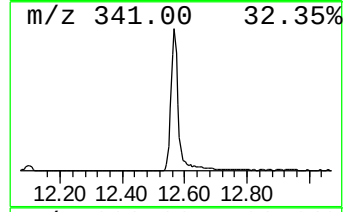
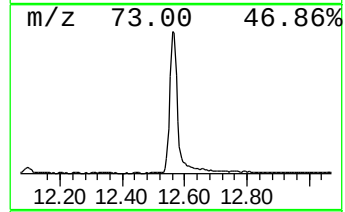
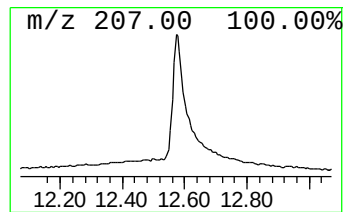
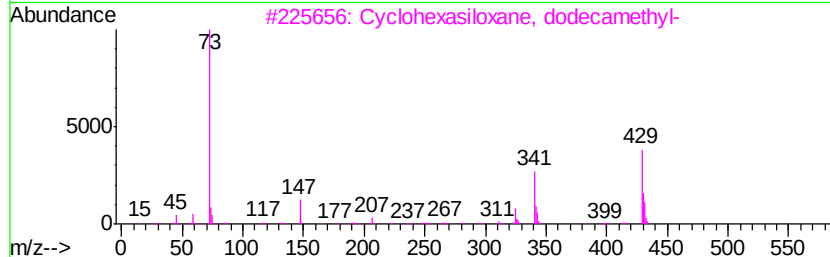
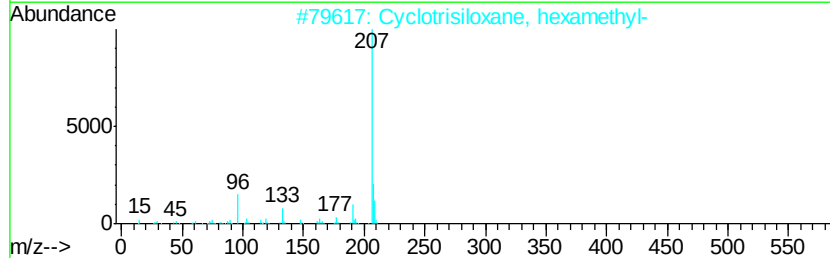
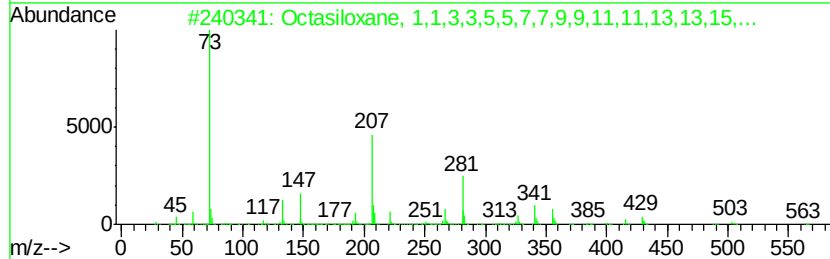
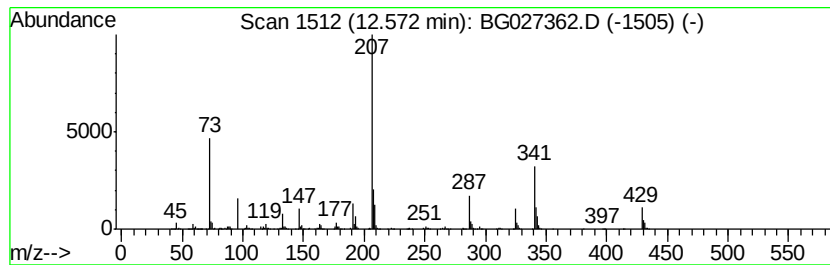
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 9 unknown12.57 Concentration Rank 7

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 12.57 | 70.94 ng | 1883820 | Naphthalene-d8   | 11.37 |

| Hit# | of | Tentative ID                        | MW  | MolForm      | CAS#        | Qual |
|------|----|-------------------------------------|-----|--------------|-------------|------|
| 1    | 5  | Octasiloxane, 1,1,3,3,5,5,7,7,9,... | 578 | C16H5007Si8  | 019095-24-0 | 45   |
| 2    |    | Cyclotrisiloxane, hexamethyl-       | 222 | C6H1803Si3   | 000541-05-9 | 43   |
| 3    |    | Cyclohexasiloxane, dodecamethyl-    | 444 | C12H3606Si6  | 000540-97-6 | 41   |
| 4    |    | Arsenous acid, tris(trimethylsil... | 342 | C9H27As03Si3 | 055429-29-3 | 38   |
| 5    |    | Methyltris(trimethylsiloxyl)silane  | 310 | C10H3003Si4  | 017928-28-8 | 38   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG060917\  
Data File : BG027362.D  
Acq On : 10 Jun 2017 00:43  
Operator : SJ/MA  
Sample : I3563-03  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
OK-04-060817

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG060517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

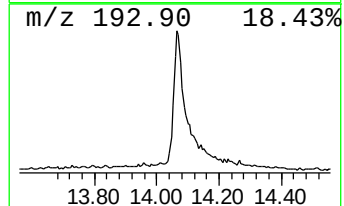
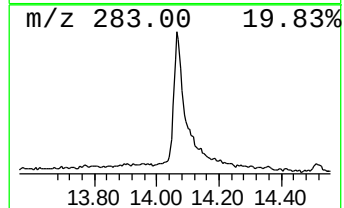
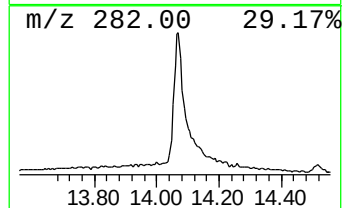
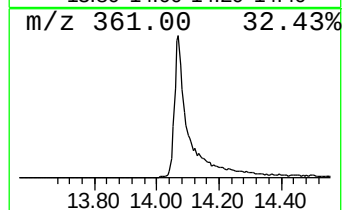
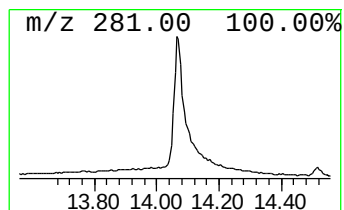
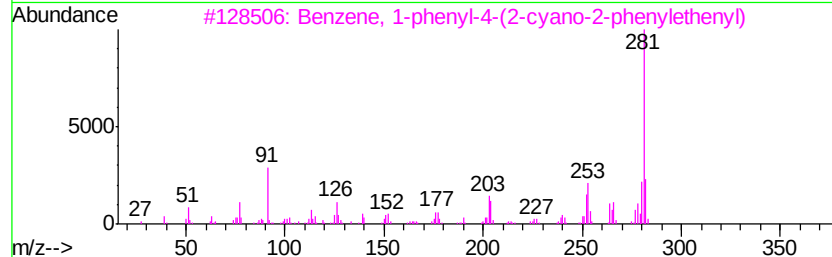
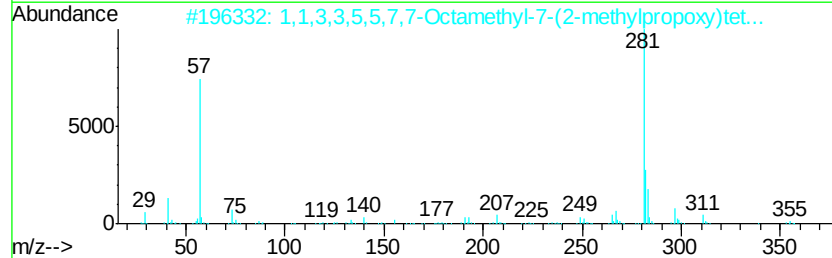
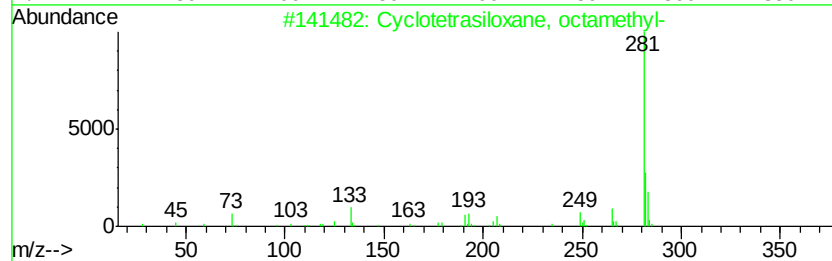
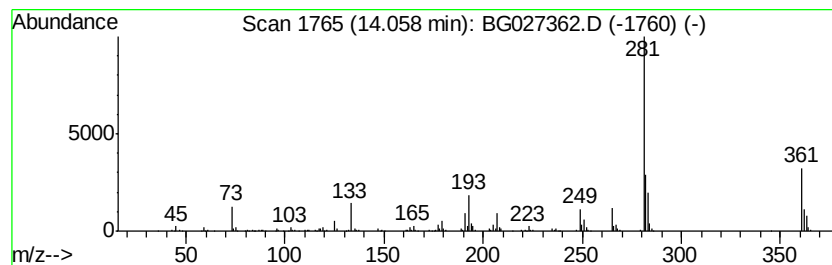
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 10 Cyclotetrasiloxane, octamet... Concentration Rank 5

| R.T.  | EstConc  | Area    | Relative to ISTD | R.T.  |
|-------|----------|---------|------------------|-------|
| 14.06 | 84.09 ng | 1151130 | Acenaphthene-d10 | 15.13 |

| Hit# | of | Tentative ID                        | MW  | MolForm     | CAS#         | Qual |
|------|----|-------------------------------------|-----|-------------|--------------|------|
| 1    | 5  | Cyclotetrasiloxane, octamethyl-     | 296 | C8H24O4Si4  | 000556-67-2  | 76   |
| 2    |    | 1,1,3,3,5,5,7,7-Octamethyl-7-(2-... | 370 | C12H34O5Si4 | 1000364-61-2 | 50   |
| 3    |    | Benzene, 1-phenyl-4-(2-cyano-2-p... | 281 | C21H15N     | 027869-56-3  | 50   |
| 4    |    | 5-(p-Aminophenyl)-4-(p-tolyl)-2-... | 281 | C16H15N3S   | 094460-46-5  | 47   |
| 5    |    | 4H-1,2,4-Triazole-3-thiol, 4-all... | 281 | C16H15N3S   | 031803-13-1  | 46   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG060917\  
Data File : BG027362.D  
Acq On : 10 Jun 2017 00:43  
Operator : SJ/MA  
Sample : I3563-03  
Misc :  
ALS Vial : 15 Sample Multiplier: 1

Instrument :  
BNA\_G  
ClientSampleId :  
OK-04-060817

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG060517.M  
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

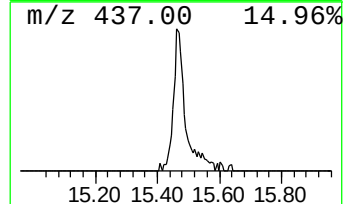
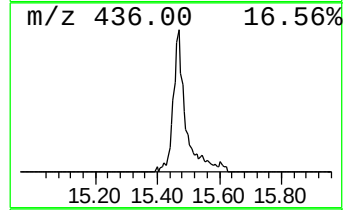
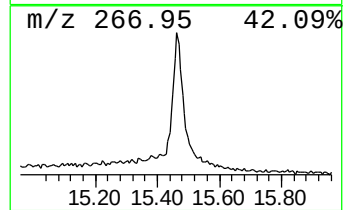
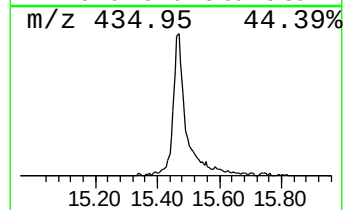
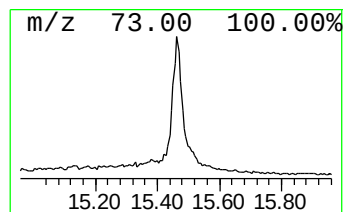
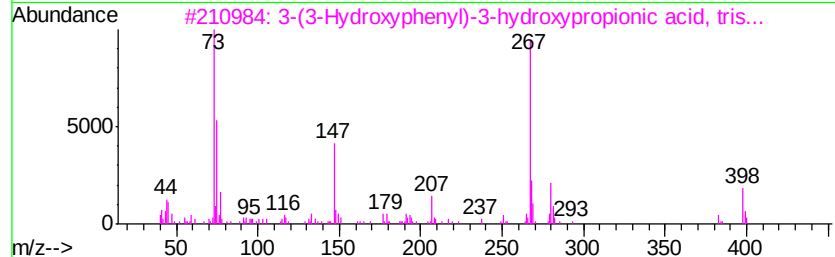
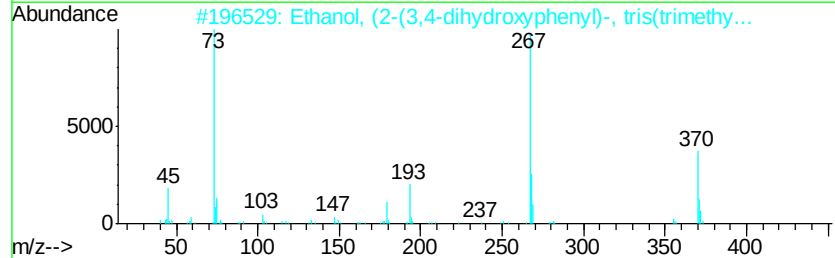
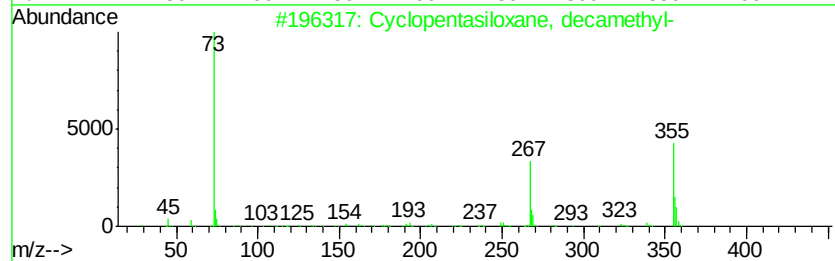
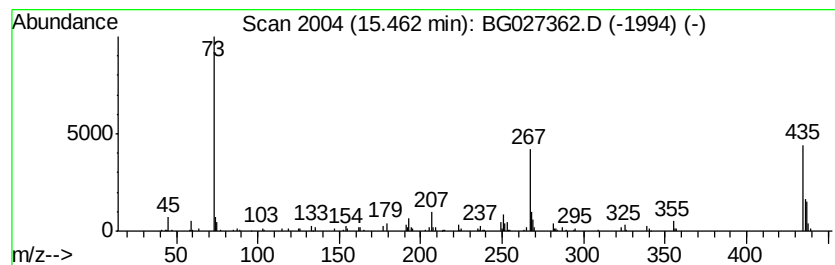
TIC Library : C:\DATABASE\NIST11.L  
TIC Integration Parameters: LSCINT.P

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Peak Number 11 unknown15.46 Concentration Rank 9

| R.T.  | EstConc  | Area   | Relative to ISTD | R.T.  |
|-------|----------|--------|------------------|-------|
| 15.46 | 18.78 ng | 257164 | Acenaphthene-d10 | 15.13 |

| Hit# | of | Tentative ID                         | MW  | MolForm     | CAS#         | Qual |
|------|----|--------------------------------------|-----|-------------|--------------|------|
| 1    | 5  | Cyclopentasiloxane, decamethyl-      | 370 | C10H30O5Si5 | 000541-02-6  | 38   |
| 2    |    | Ethanol, (2-(3,4-dihydroxyphenyl)... | 370 | C17H34O3Si3 | 068595-80-2  | 35   |
| 3    |    | 3-(3-Hydroxyphenyl)-3-hydroxypro...  | 398 | C18H34O4Si3 | 1000071-69-7 | 32   |
| 4    |    | Benzoic acid, 2-[(trimethylsilyl)... | 282 | C13H22O3Si2 | 003789-85-3  | 25   |
| 5    |    | 11H-Dibenzo[b,e][1,4]diazepin-11...  | 267 | C16H17N3O   | 013450-73-2  | 18   |



Data Path : Z:\HPCHEM1\BNA\_G\DATA\BG060917\  
 Data File : BG027362.D  
 Acq On : 10 Jun 2017 00:43  
 Operator : SJ/MA  
 Sample : I3563-03  
 Misc :  
 ALS Vial : 15 Sample Multiplier: 1

Instrument :  
 BNA\_G  
 ClientSampleId :  
 OK-04-060817

Quant Method : Z:\HPCHEM1\BNA\_G\METHODS\8270-BG060517.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L  
 TIC Integration Parameters: LSCINT.P

| TIC Top Hit name     | RT    | EstConc | Units | Response | --Internal Standard-- |       |        |      |
|----------------------|-------|---------|-------|----------|-----------------------|-------|--------|------|
|                      |       |         |       |          | #                     | RT    | Resp   | Conc |
| 4-Penten-2-one, 4... | 4.55  | 14.0    | ng    | 224048   | 1                     | 8.56  | 320996 | 20.0 |
| Trisiloxane, 1,1,... | 4.67  | 9.6     | ng    | 154006   | 1                     | 8.56  | 320996 | 20.0 |
| 3-Penten-2-one, 4... | 5.13  | 91.2    | ng    | 1463110  | 1                     | 8.56  | 320996 | 20.0 |
| Cyclotrisiloxane,... | 5.22  | 631.4   | ng    | 10133000 | 1                     | 8.56  | 320996 | 20.0 |
| 4-Hydroxymandelic... | 7.10  | 27.1    | ng    | 434157   | 1                     | 8.56  | 320996 | 20.0 |
| unknown8.09          | 8.09  | 111.5   | ng    | 1789540  | 1                     | 8.56  | 320996 | 20.0 |
| Plumbane, diethyl... | 10.09 | 89.3    | ng    | 2372550  | 2                     | 11.37 | 531131 | 20.0 |
| unknown12.57         | 12.57 | 70.9    | ng    | 1883820  | 2                     | 11.37 | 531131 | 20.0 |
| Cyclotetrasiloxan... | 14.06 | 84.1    | ng    | 1151130  | 3                     | 15.13 | 273799 | 20.0 |
| unknown15.46         | 15.46 | 18.8    | ng    | 257164   | 3                     | 15.13 | 273799 | 20.0 |