# Full wwPDB NMR Structure Validation Report (i)

Feb 19, 2022 – 12:51 PM EST

| | | |
|---|---|---|
| PDB ID | : | 1Q7O |
| Title | : | Determination of f-MLF-OH Peptide Structure with solid-state magic-angle spinning NMR Spectroscopy |
| Authors | : | Rienstra, C.M.; Tucker-Kellogg, L.; Jaroniec, C.P.; Hohwy, M.; Reif, B.; McMahon, M.T.; Tidor, B.; Lozano-Perez, T.; Griffin, R.G. |
| Deposited on | : | 2003-08-19 |

This is a Full wwPDB NMR Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/NMRValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (i)) were used in the production of this report:

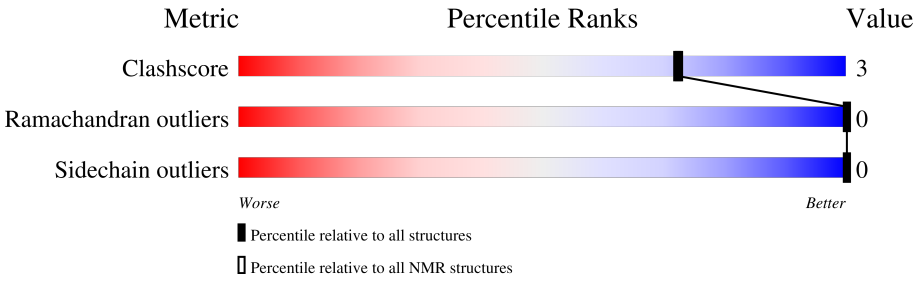
| | | |
|---|---|---|
| MolProbity | : | 4.02b-467 |
| Mogul | : | 1.8.5 (274361), CSD as541be (2020) |
| Percentile statistics | : | 20191225.v01 (using entries in the PDB archive December 25th 2019) |
| RCI | : | v_1n_11_5_13_A (Berjanski et al., 2005) |
| PANAV | : | Wang et al. (2010) |
| ShiftChecker | : | 2.26 |
| Ideal geometry (proteins) | : | Engh & Huber (2001) |
| Ideal geometry (DNA, RNA) | : | Parkinson et al. (1996) |
| Validation Pipeline (wwPDB-VP) | : | 2.26 |

# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:
*SOLID-STATE NMR*

The overall completeness of chemical shifts assignment was not calculated.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.




| Metric | Whole archive (#Entries) | NMR archive (#Entries) |
|---|---|---|
| Clashscore | 158937 | 12864 |
| Ramachandran outliers | 154571 | 11451 |
| Sidechain outliers | 154315 | 11428 |

The table below summarises the geometric issues observed across the polymeric chains and their fit to the experimental data. The red, orange, yellow and green segments indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria. A cyan segment indicates the fraction of residues that are not part of the well-defined cores, and a grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5%

| Mol | Chain | Length | Quality of chain |
|---|---|---|---|
| 1 | A | 3 | 100% |

## 2 Ensemble composition and analysis (i)

This entry contains 20 models.

Cyrange was unable to find well-defined residues.

Error message: Less than two structures selected.

NmrClust was unable to cluster the ensemble.

Error message: Wrapper check: not enough residues in core to run NmrClust

# 3 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 58 atoms, of which 28 are hydrogens and 0 are deuteriums.

- Molecule 1 is a protein called chemotactic peptide.

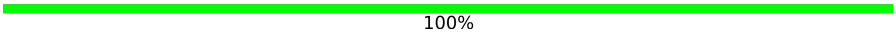
| Mol | Chain | Residues | Atoms | | | | | | Trace |
|---|---|---|---|---|---|---|---|---|---|
| | | | Total | C | H | N | O | S | |
| 1 | A | 3 | 58 | 21 | 28 | 3 | 5 | 1 | 0 |

# 4 Residue-property plots

## 4.1 Average score per residue in the NMR ensemble

These plots are provided for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic is the same as shown in the summary in section 1 of this report. The second graphic shows the sequence where residues are colour-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. Stretches of 2 or more consecutive residues without any outliers are shown as green connectors. Residues which are classified as ill-defined in the NMR ensemble, are shown in cyan with an underline colour-coded according to the previous scheme. Residues which were present in the experimental sample, but not modelled in the final structure are shown in grey.

- Molecule 1: chemotactic peptide

Chain A: 

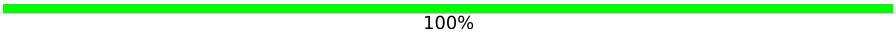

There are no outlier residues in this chain.

## 4.2 Scores per residue for each member of the ensemble

Colouring as in section 4.1 above.

### 4.2.1 Score per residue for model 1

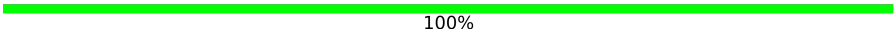
- Molecule 1: chemotactic peptide

Chain A: 


There are no outlier residues in this chain.

### 4.2.2 Score per residue for model 2

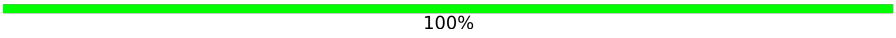
- Molecule 1: chemotactic peptide

Chain A: 


There are no outlier residues in this chain.

### 4.2.3 Score per residue for model 3

- Molecule 1: chemotactic peptide

Chain A: 

There are no outlier residues in this chain.

### 4.2.4 Score per residue for model 4

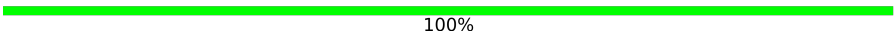
• Molecule 1: chemotactic peptide

Chain A: 



### 4.2.5 Score per residue for model 5

• Molecule 1: chemotactic peptide

Chain A: 


There are no outlier residues in this chain.

### 4.2.6 Score per residue for model 6

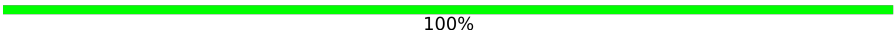
• Molecule 1: chemotactic peptide

Chain A: 



### 4.2.7 Score per residue for model 7

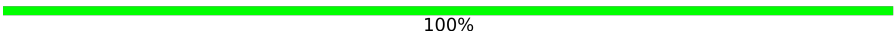
• Molecule 1: chemotactic peptide

Chain A: 


There are no outlier residues in this chain.

### 4.2.8 Score per residue for model 8

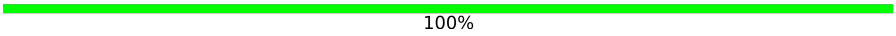
• Molecule 1: chemotactic peptide

Chain A: 


There are no outlier residues in this chain.

### 4.2.9 Score per residue for model 9

• Molecule 1: chemotactic peptide

Chain A: 


There are no outlier residues in this chain.

### 4.2.10 Score per residue for model 10

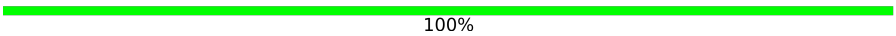
• Molecule 1: chemotactic peptide

Chain A: 



### 4.2.11 Score per residue for model 11

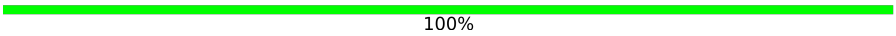
• Molecule 1: chemotactic peptide

Chain A: 


There are no outlier residues in this chain.

### 4.2.12 Score per residue for model 12

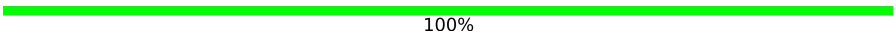
• Molecule 1: chemotactic peptide

Chain A: 


There are no outlier residues in this chain.

### 4.2.13 Score per residue for model 13

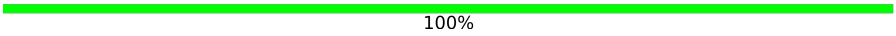
• Molecule 1: chemotactic peptide

Chain A: 


There are no outlier residues in this chain.

#### 4.2.14 Score per residue for model 14

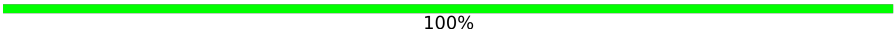
• Molecule 1: chemotactic peptide

Chain A: 


There are no outlier residues in this chain.

#### 4.2.15 Score per residue for model 15

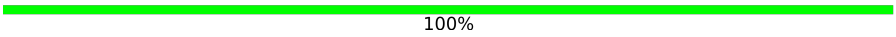
• Molecule 1: chemotactic peptide

Chain A: 


There are no outlier residues in this chain.

#### 4.2.16 Score per residue for model 16

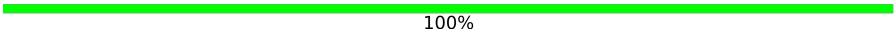
• Molecule 1: chemotactic peptide

Chain A: 


There are no outlier residues in this chain.

#### 4.2.17 Score per residue for model 17

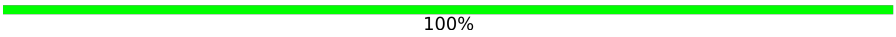
• Molecule 1: chemotactic peptide

Chain A: 


There are no outlier residues in this chain.

#### 4.2.18 Score per residue for model 18

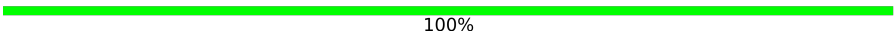
• Molecule 1: chemotactic peptide

Chain A: 


There are no outlier residues in this chain.

#### 4.2.19 Score per residue for model 19

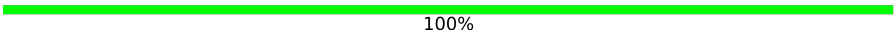
• Molecule 1: chemotactic peptide

Chain A: 

There are no outlier residues in this chain.

### 4.2.20 Score per residue for model 20

• Molecule 1: chemotactic peptide

Chain A: 


There are no outlier residues in this chain.

# 5 Refinement protocol and experimental data overview

The models were refined using the following method: *simulated annealing.*

Of the 20 calculated structures, 20 were deposited, based on the following criterion: *structures with the lowest energy.*

The following table shows the software used for structure solution, optimisation and refinement.

| Software name | Classification | Version |
|---|---|---|
| CNS | refinement | 1.0 |

No chemical shift data was provided. Note: This is a solid-state NMR structure, where hydrogen atoms are typically not assigned a chemical shift value, which may lead to lower completeness of assignment measure.

# 6 Model quality

## 6.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MTY, FME

There are no covalent bond-length or bond-angle outliers.

There are no bond-length outliers.

There are no bond-angle outliers.

There are no chirality outliers.

There are no planarity outliers.

## 6.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in each chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes averaged over the ensemble.

| Mol | Chain | Non-H | H(model) | H(added) | Clashes |
|---|---|---|---|---|---|
| 1 | A | 30 | 28 | 28 | 0±1 |
| All | All | 600 | 560 | 560 | 4 |

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 3.

All unique clashes are listed below, sorted by their clash magnitude.

| Atom-1 | Atom-2 | Clash(Å) | Distance(Å) | Models | |
|---|---|---|---|---|---|
| | | | | Worst | Total |
| 1:A:3:MTY:HD2 | 1:A:3:MTY:C | 0.58 | 2.29 | 4 | 2 |
| 1:A:3:MTY:HD2 | 1:A:3:MTY:O | 0.46 | 2.11 | 4 | 1 |
| 1:A:3:MTY:C | 1:A:3:MTY:CD2 | 0.41 | 2.99 | 4 | 1 |

## 6.3 Torsion angles

### 6.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all NMR entries. The Analysed column shows the number of residues for which the backbone conformation

was analysed and the total number of residues.

| Mol | Chain | Analysed | Favoured | Allowed | Outliers | Percentiles | |
|---|---|---|---|---|---|---|---|
| 1 | A | 1/3 (33%) | 1±0 (100±0%) | 0±0 (0±0%) | 0±0 (0±0%) | 100 | 100 |
| All | All | 20/60 (33%) | 20 (100%) | 0 (0%) | 0 (0%) | 100 | 100 |

There are no Ramachandran outliers.

### 6.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all NMR entries. The Analysed column shows the number of residues for which the sidechain conformation was analysed and the total number of residues.

| Mol | Chain | Analysed | Rotameric | Outliers | Percentiles | |
|---|---|---|---|---|---|---|
| 1 | A | 1/1 (100%) | 1±0 (100±0%) | 0±0 (0±0%) | 100 | 100 |
| All | All | 20/20 (100%) | 20 (100%) | 0 (0%) | 100 | 100 |

There are no protein residues with a non-rotameric sidechain to report.

### 6.3.3 RNA (i)

There are no RNA molecules in this entry.

## 6.4 Non-standard residues in protein, DNA, RNA chains (i)

2 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds for which Mogul statistics could be retrieved, the number of bonds that are observed in the model and the number of bonds that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length is the number of standard deviations the observed value is removed from the expected value. A bond length with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the average root-mean-square of all Z scores of the bond lengths.

| Mol | Type | Chain | Res | Link | Bond lengths | | |
|---|---|---|---|---|---|---|---|
| | | | | | Counts | RMSZ | #Z>2 |
| 1 | FME | A | 1 | 1 | 8,9,10 | 0.36±0.02 | 0±0 (0±0%) |
| 1 | MTY | A | 3 | 1 | 9,12,13 | 0.23±0.04 | 0±0 (0±0%) |

In the following table, the Counts columns list the number of angles for which Mogul statistics could be retrieved, the number of angles that are observed in the model and the number of angles that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond angle is the number of standard deviations the observed value is removed from the expected value. A bond angle with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the average root-mean-square of all Z scores of the bond angles.

| Mol | Type | Chain | Res | Link | Bond angles | | |
|---|---|---|---|---|---|---|---|
| | | | | | Counts | RMSZ | #Z>2 |
| 1 | FME | A | 1 | 1 | 7,9,11 | 0.91±0.05 | 0±0 (0±3%) |
| 1 | MTY | A | 3 | 1 | 10,15,17 | 0.26±0.04 | 0±0 (0±0%) |

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

| Mol | Type | Chain | Res | Link | Chirals | Torsions | Rings |
|---|---|---|---|---|---|---|---|
| 1 | FME | A | 1 | 1 | - | 0±0,7,9,11 | - |
| 1 | MTY | A | 3 | 1 | - | 0±0,4,8,8 | 0±0,1,1,1 |

There are no bond-length outliers.

All unique angle outliers are listed below.

| Mol | Chain | Res | Type | Atoms | Z | Observed($^o$) | Ideal($^o$) | Models | |
|---|---|---|---|---|---|---|---|---|---|
| | | | | | | | | Worst | Total |
| 1 | A | 1 | FME | C-CA-N | 2.01 | 113.36 | 109.73 | 6 | 1 |

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 6.5 Carbohydrates (i)

There are no monosaccharides in this entry.

## 6.6 Ligand geometry (i)

There are no ligands in this entry.

## 6.7 Other polymers (i)

There are no such molecules in this entry.

## 6.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

# 7 Chemical shift validation (i)

No chemical shift data were provided