



# wwPDB X-ray Structure Validation Summary Report

Jan 30, 2024 – 04:29 PM EST

PDB ID : 1FZ6  
Title : METHANE MONOOXYGENASE HYDROXYLASE, FORM II SOAKED IN  
1 M METHANOL  
Authors : Whittington, D.A.; Sazinsky, M.H.; Lippard, S.J.  
Deposited on : 2000-10-03  
Resolution : 2.05 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the  symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.36  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.36

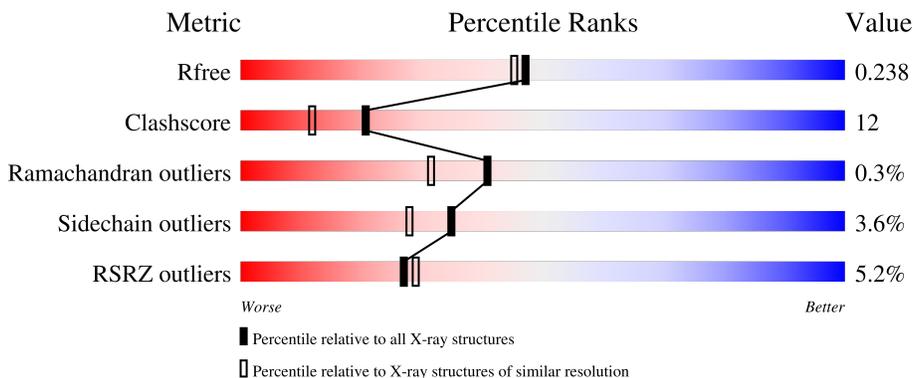
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.05 Å.

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)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	1692 (2.04-2.04)
Clashscore	141614	1773 (2.04-2.04)
Ramachandran outliers	138981	1752 (2.04-2.04)
Sidechain outliers	138945	1752 (2.04-2.04)
RSRZ outliers	127900	1672 (2.04-2.04)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. 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  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	527	 2% 72% 23% . .
1	B	527	 4% 69% 26% . .
2	C	389	 % 82% 16% .
2	D	389	 9% 70% 28% .
3	E	170	 3% 78% 15% . . .

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Mol	Chain	Length	Quality of chain
3	F	170	 <p>A horizontal bar chart representing the quality of chain. The bar is divided into three segments: a red segment on the left labeled '21%', a green segment in the middle labeled '68%', and a yellow segment on the right labeled '28%'. There are two small black dots at the far right end of the bar.</p>

## 2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 18586 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called METHANE MONOOXYGENASE COMPONENT A, ALPHA CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	510	4177	2673	719	767	18	40	0	0
1	B	510	4173	2671	718	766	18	36	0	0

- Molecule 2 is a protein called METHANE MONOOXYGENASE COMPONENT A, BETA CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	C	388	3193	2054	551	580	8	30	0	0
2	D	388	3193	2054	551	580	8	42	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	370	ARG	ALA	conflict	UNP P18798
D	370	ARG	ALA	conflict	UNP P18798

- Molecule 3 is a protein called METHANE MONOOXYGENASE COMPONENT A, GAMMA CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	E	166	1368	867	246	250	5	4	0	0
3	F	166	1366	866	245	250	5	8	0	0

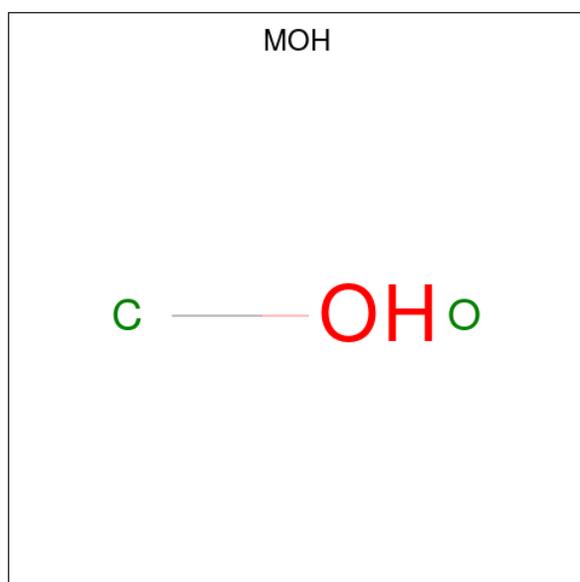
- Molecule 4 is FE (III) ION (three-letter code: FE) (formula: Fe).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	2	Total Fe 2 2	0	0
4	B	2	Total Fe 2 2	0	0

- Molecule 5 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total Ca 1 1	0	0
5	B	1	Total Ca 1 1	0	0
5	C	1	Total Ca 1 1	0	0

- Molecule 6 is METHANOL (three-letter code: MOH) (formula: CH<sub>4</sub>O).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	B	1	Total C O 2 1 1	0	0

- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	259	Total O 259 259	0	0

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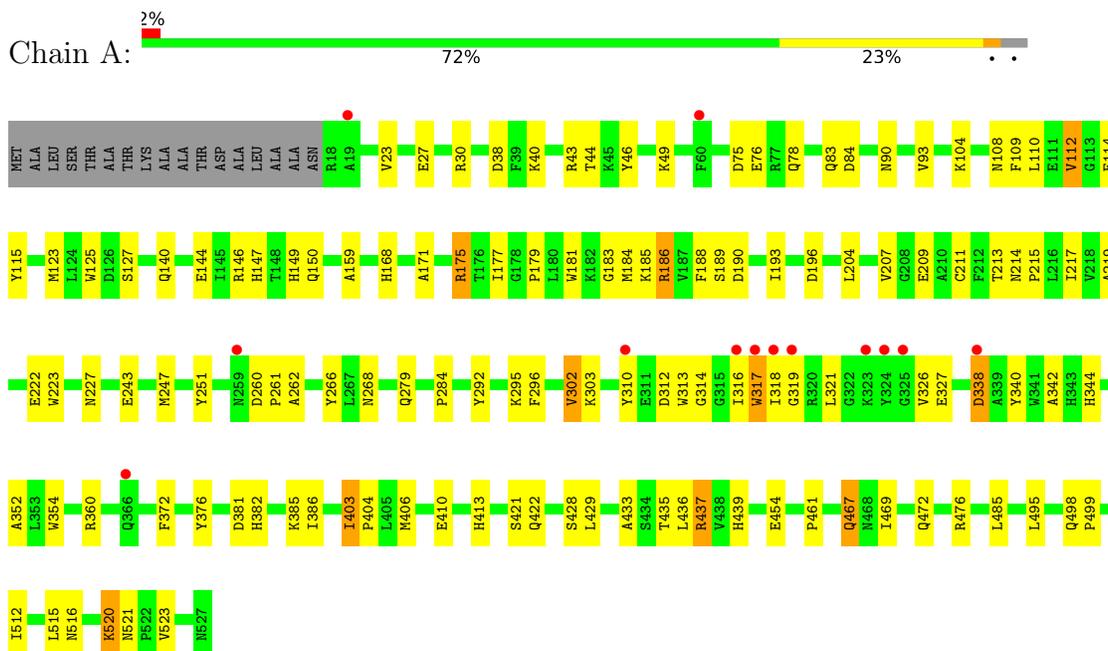
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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	B	239	Total O 239 239	0	0
7	C	286	Total O 286 286	0	0
7	D	144	Total O 144 144	0	0
7	E	125	Total O 125 125	0	0
7	F	54	Total O 54 54	0	0

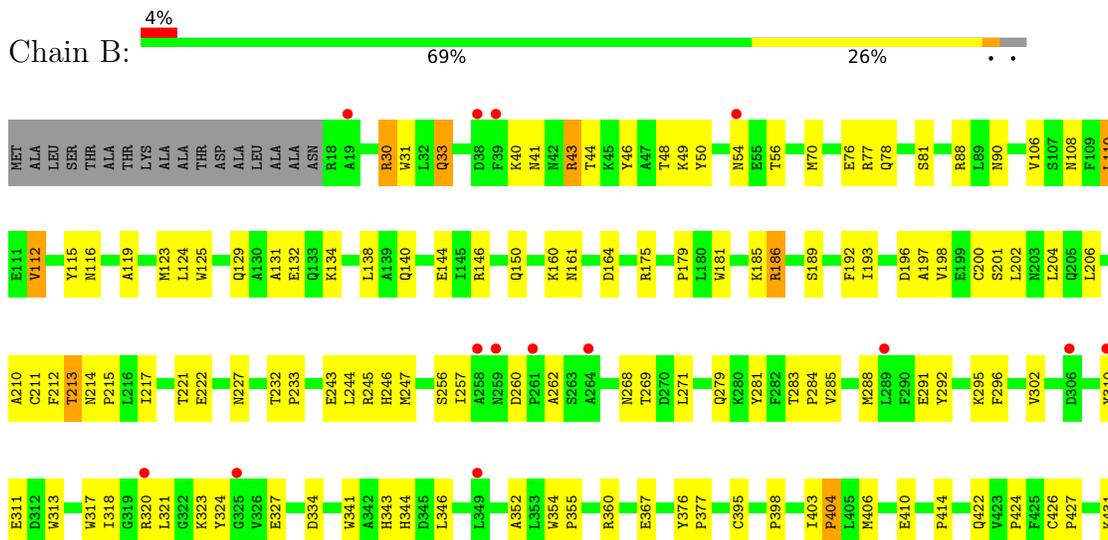
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-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. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

#### • Molecule 1: METHANE MONOOXYGENASE COMPONENT A, ALPHA CHAIN

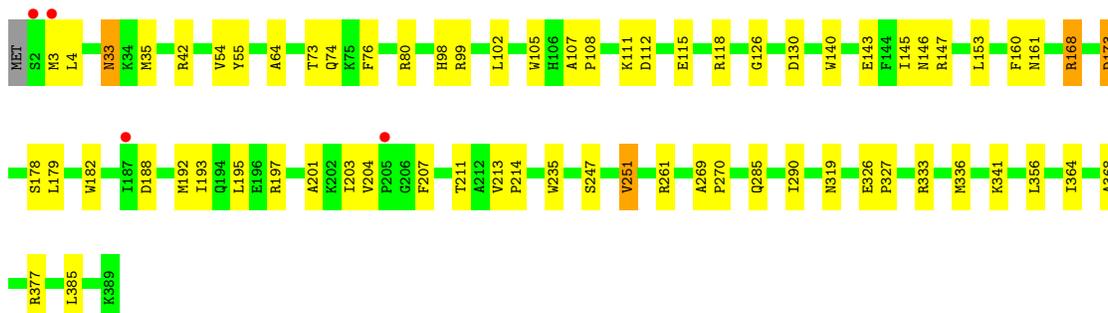
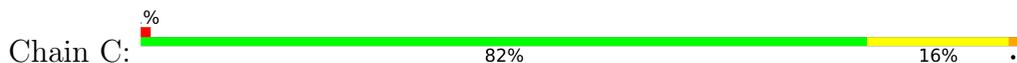


#### • Molecule 1: METHANE MONOOXYGENASE COMPONENT A, ALPHA CHAIN

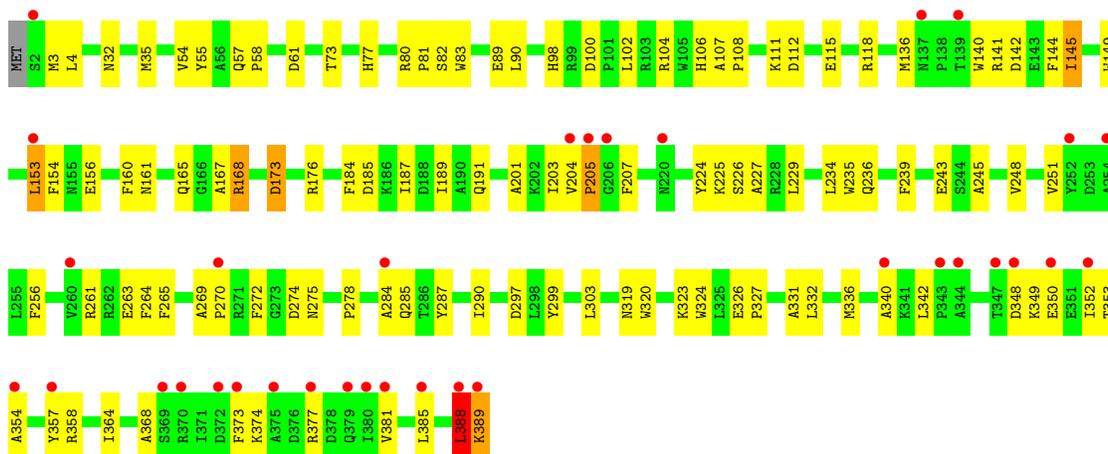




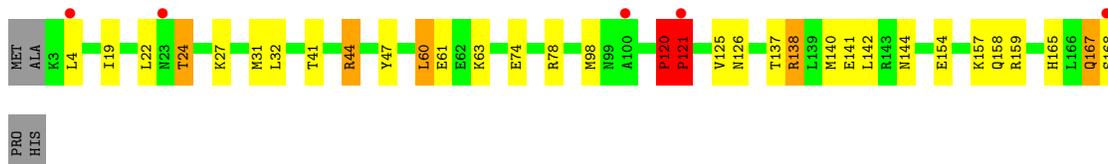
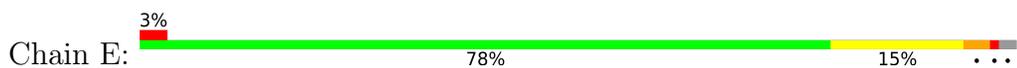
- Molecule 2: METHANE MONOOXYGENASE COMPONENT A, BETA CHAIN



- Molecule 2: METHANE MONOOXYGENASE COMPONENT A, BETA CHAIN

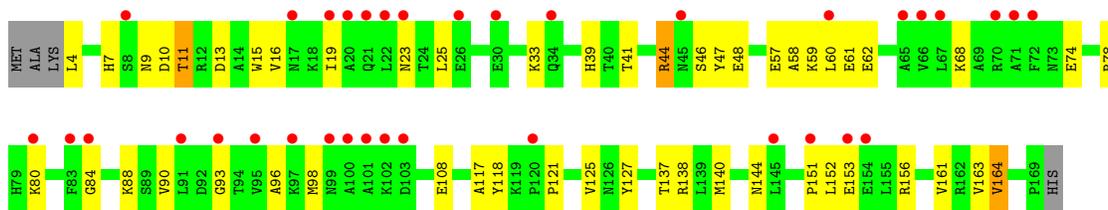


- Molecule 3: METHANE MONOOXYGENASE COMPONENT A, GAMMA CHAIN



- Molecule 3: METHANE MONOOXYGENASE COMPONENT A, GAMMA CHAIN





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	70.67Å 171.29Å 221.34Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.90 – 2.05 29.87 – 2.05	Depositor EDS
% Data completeness (in resolution range)	89.0 (29.90-2.05) 89.1 (29.87-2.05)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.72 (at 2.04Å)	Xtrriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.218 , 0.244 0.211 , 0.238	Depositor DCC
$R_{free}$ test set	5330 reflections (3.26%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	28.2	Xtrriage
Anisotropy	0.249	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 49.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	18586	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	35.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.25% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality i

### 5.1 Standard geometry i

Bond lengths and bond angles in the following residue types are not validated in this section: MOH, CA, FE

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	A	0.39	1/4302 (0.0%)	0.61	1/5842 (0.0%)
1	B	0.40	0/4298	0.60	0/5837
2	C	0.40	0/3289	0.60	0/4464
2	D	1.03	9/3289 (0.3%)	0.85	11/4464 (0.2%)
3	E	0.82	7/1396 (0.5%)	1.08	7/1880 (0.4%)
3	F	0.34	0/1395	0.57	1/1881 (0.1%)
All	All	0.60	17/17969 (0.1%)	0.70	20/24368 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
3	E	0	1

The worst 5 of 17 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	389	LYS	CB-CG	35.74	2.49	1.52
2	D	389	LYS	CA-C	32.22	2.36	1.52
3	E	167	GLN	C-N	-18.30	0.92	1.34
2	D	389	LYS	CA-CB	14.82	1.86	1.53
2	D	388	LEU	CA-C	13.22	1.87	1.52

The worst 5 of 20 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	120	PRO	C-N-CD	-27.86	59.30	120.60
2	D	389	LYS	N-CA-CB	-20.98	72.84	110.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	389	LYS	N-CA-C	16.70	156.10	111.00
2	D	389	LYS	CA-C-O	15.21	152.04	120.10
3	E	121	PRO	CA-N-CD	-14.31	91.47	111.50

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	E	120	PRO	Mainchain

## 5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the 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 within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4177	0	3975	99	0
1	B	4173	0	3969	125	0
2	C	3193	0	3042	56	0
2	D	3193	0	3042	107	0
3	E	1368	0	1362	33	0
3	F	1366	0	1357	44	0
4	A	2	0	0	0	0
4	B	2	0	0	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
5	C	1	0	0	0	0
6	B	2	0	0	0	0
7	A	259	0	0	6	0
7	B	239	0	0	6	0
7	C	286	0	0	3	0
7	D	144	0	0	0	0
7	E	125	0	0	2	0
7	F	54	0	0	0	0
All	All	18586	0	16747	415	0

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 12.

The worst 5 of 415 close contacts within the same asymmetric unit are listed below, sorted by

their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:389:LYS:CG	2:D:389:LYS:CD	1.75	1.59
2:D:389:LYS:CB	2:D:389:LYS:CA	1.86	1.50
3:E:167:GLN:C	3:E:167:GLN:CA	1.80	1.49
2:D:388:LEU:C	2:D:388:LEU:CA	1.87	1.42
2:D:389:LYS:CB	2:D:389:LYS:N	2.07	1.17

There are no symmetry-related clashes.

## 5.3 Torsion angles [\(i\)](#)

### 5.3.1 Protein backbone [\(i\)](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

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	508/527 (96%)	487 (96%)	21 (4%)	0	100	100
1	B	508/527 (96%)	479 (94%)	28 (6%)	1 (0%)	47	39
2	C	386/389 (99%)	372 (96%)	12 (3%)	2 (0%)	29	18
2	D	386/389 (99%)	362 (94%)	21 (5%)	3 (1%)	19	10
3	E	164/170 (96%)	159 (97%)	4 (2%)	1 (1%)	25	15
3	F	164/170 (96%)	153 (93%)	11 (7%)	0	100	100
All	All	2116/2172 (97%)	2012 (95%)	97 (5%)	7 (0%)	41	31

5 of 7 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	E	121	PRO
1	B	40	LYS
2	D	388	LEU
2	C	64	ALA
2	D	251	VAL

### 5.3.2 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 X-ray entries followed by that with respect to entries of similar resolution.

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	432/442 (98%)	416 (96%)	16 (4%)	34	27
1	B	431/442 (98%)	412 (96%)	19 (4%)	28	21
2	C	322/323 (100%)	312 (97%)	10 (3%)	40	33
2	D	322/323 (100%)	314 (98%)	8 (2%)	47	40
3	E	144/147 (98%)	138 (96%)	6 (4%)	30	22
3	F	144/147 (98%)	139 (96%)	5 (4%)	36	29
All	All	1795/1824 (98%)	1731 (96%)	64 (4%)	35	28

5 of 64 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	E	121	PRO
3	F	11	THR
1	B	186	ARG
1	B	125	TRP
3	F	23	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 61 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	227	ASN
3	E	165	HIS
1	B	344	HIS
3	E	144	ASN
3	F	144	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 7 are monoatomic - leaving 1 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or 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 length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
6	MOH	B	9001	4	1,1,1	0.46	0	-		

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
3	E	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	E	167:GLN	C	168:SER	N	0.92

## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	510/527 (96%)	0.16	13 (2%) 57 61	21, 32, 49, 63	10 (1%)
1	B	510/527 (96%)	0.15	19 (3%) 41 45	20, 33, 50, 62	9 (1%)
2	C	388/389 (99%)	-0.16	4 (1%) 82 84	18, 25, 39, 54	7 (1%)
2	D	388/389 (99%)	0.59	34 (8%) 10 10	21, 40, 61, 79	10 (2%)
3	E	166/170 (97%)	0.05	5 (3%) 50 54	21, 30, 44, 59	1 (0%)
3	F	166/170 (97%)	1.12	35 (21%) 1 0	37, 49, 67, 75	2 (1%)
All	All	2128/2172 (97%)	0.24	110 (5%) 27 29	18, 33, 55, 79	39 (1%)

The worst 5 of 110 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	389	LYS	9.2
2	D	205	PRO	5.9
3	F	100	ALA	5.4
2	C	2	SER	5.1
3	F	101	ALA	4.8

### 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

### 6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
5	CA	C	5007	1/1	0.49	0.09	77,77,77,77	0
6	MOH	B	9001	2/2	0.90	0.20	43,43,43,45	0
5	CA	B	5005	1/1	0.94	0.10	68,68,68,68	0
5	CA	A	5006	1/1	0.97	0.13	45,45,45,45	0
4	FE	A	5002	1/1	0.98	0.03	42,42,42,42	0
4	FE	B	5003	1/1	0.99	0.04	30,30,30,30	0
4	FE	B	5004	1/1	0.99	0.02	41,41,41,41	0
4	FE	A	5001	1/1	0.99	0.02	30,30,30,30	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.