



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 08:48 PM GMT

PDB ID : 1M5S  
Title : Formylmethanofuran:tetrahydromethanopterin formyltransferase from Methanosarcina barkeri  
Authors : Mamat, B.; Roth, A.; Grimm, C.; Ermler, U.; Tziatzios, C.; Schubert, D.; Thauer, R.K.; Shima, S.  
Deposited on : 2002-07-10  
Resolution : 1.85 Å(reported)

This is a Full wwPDB X-ray Structure Validation 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  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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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 : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

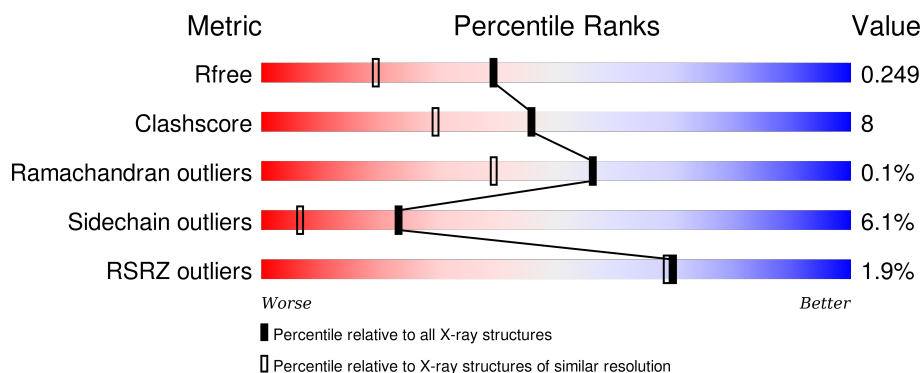
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.85 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))
$R_{free}$	91344	1745 (1.86-1.86)
Clashscore	102246	1898 (1.86-1.86)
Ramachandran outliers	100387	1875 (1.86-1.86)
Sidechain outliers	100360	1875 (1.86-1.86)
RSRZ outliers	91569	1747 (1.86-1.86)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	297	<div> <div>3%</div> <div>81%15%</div> <div>•</div> </div>
1	B	297	<div> <div>2%</div> <div>84%13%</div> <div>•</div> </div>
1	C	297	<div> <div>%</div> <div>81%15%</div> <div>•</div> </div>
1	D	297	<div> <div>2%</div> <div>81%18%</div> <div>••</div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 9642 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 Formylmethanofuran--tetrahydromethanopterin formyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	297	Total	C	N	O	S	5	1	0
			2233	1427	362	434	10			
1	B	297	Total	C	N	O	S	0	0	0
			2227	1423	361	433	10			
1	C	297	Total	C	N	O	S	13	0	0
			2227	1423	361	433	10			
1	D	297	Total	C	N	O	S	6	0	0
			2227	1423	361	433	10			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	115	ASP	PHE	CONFLICT	UNP P55301
B	1115	ASP	PHE	CONFLICT	UNP P55301
C	2115	ASP	PHE	CONFLICT	UNP P55301
D	3115	ASP	PHE	CONFLICT	UNP P55301

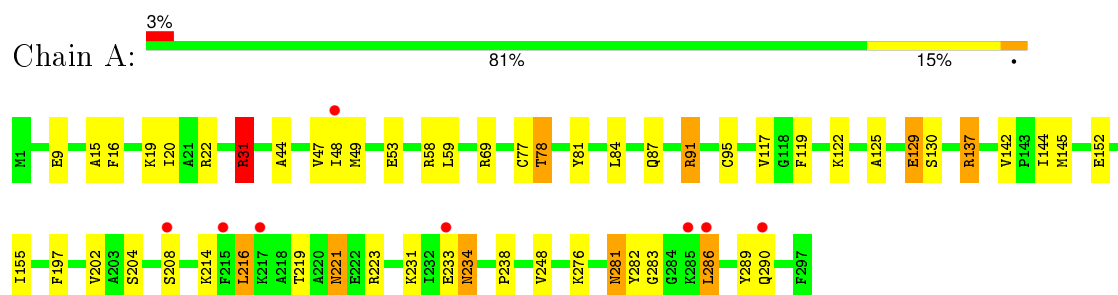
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	179	Total	O	0	0
			179	179		
2	B	168	Total	O	0	0
			168	168		
2	C	194	Total	O	0	0
			194	194		
2	D	187	Total	O	0	0
			187	187		

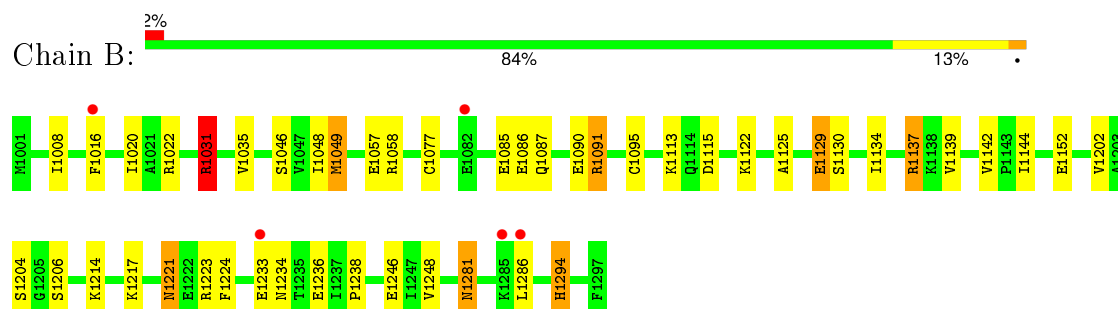
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

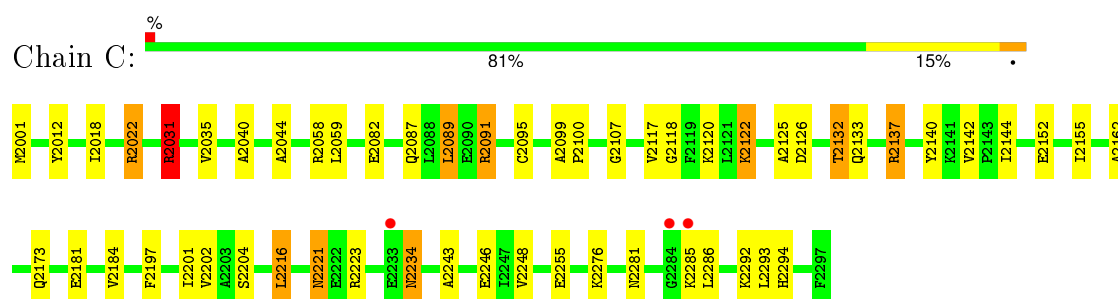
- Molecule 1: Formylmethanofuran--tetrahydromethanopterin formyltransferase



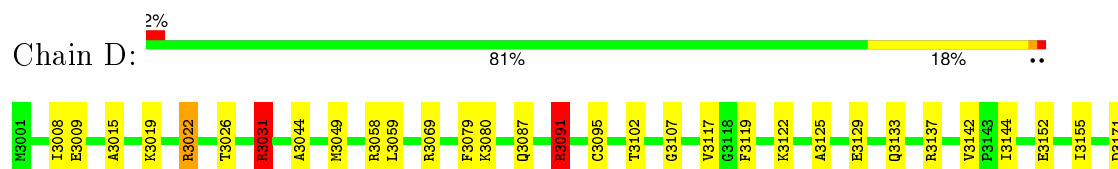
- Molecule 1: Formylmethanofuran--tetrahydromethanopterin formyltransferase

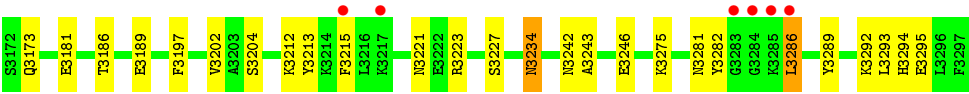


- Molecule 1: Formylmethanofuran--tetrahydromethanopterin formyltransferase



- Molecule 1: Formylmethanofuran--tetrahydromethanopterin formyltransferase





## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	127.10Å 83.80Å 126.40Å 90.00° 108.40° 90.00°	Depositor
Resolution (Å)	19.78 – 1.85 19.78 – 1.85	Depositor EDS
% Data completeness (in resolution range)	91.0 (19.78-1.85) 91.1 (19.78-1.85)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.04	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	7.96 (at 1.85Å)	Xtriage
Refinement program	CNS 0.3	Depositor
R, $R_{free}$	0.197 , 0.248 0.201 , 0.249	Depositor DCC
$R_{free}$ test set	4913 reflections (5.02%)	DCC
Wilson B-factor (Å <sup>2</sup> )	23.6	Xtriage
Anisotropy	0.492	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 51.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 97796 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	9642	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.91% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

## 5 Model quality

### 5.1 Standard geometry

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	0.42	0/2275	0.64	6/3075 (0.2%)
1	B	0.37	0/2266	0.59	4/3064 (0.1%)
1	C	0.38	0/2266	0.61	5/3064 (0.2%)
1	D	0.38	0/2266	0.60	5/3064 (0.2%)
All	All	0.39	0/9073	0.61	20/12267 (0.2%)

There are no bond length outliers.

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	2137	ARG	NE-CZ-NH2	7.90	124.25	120.30
1	A	22	ARG	NE-CZ-NH2	7.77	124.19	120.30
1	D	3069	ARG	NE-CZ-NH2	7.50	124.05	120.30
1	C	2058	ARG	NE-CZ-NH2	7.49	124.05	120.30
1	B	1031	ARG	NE-CZ-NH2	7.47	124.03	120.30
1	C	2091	ARG	NE-CZ-NH2	7.47	124.03	120.30
1	B	1091	ARG	NE-CZ-NH2	7.46	124.03	120.30
1	A	58	ARG	NE-CZ-NH2	7.45	124.03	120.30
1	A	137	ARG	NE-CZ-NH2	7.42	124.01	120.30
1	B	1137	ARG	NE-CZ-NH2	7.42	124.01	120.30
1	A	91	ARG	NE-CZ-NH2	7.40	124.00	120.30
1	A	69	ARG	NE-CZ-NH2	7.40	124.00	120.30
1	D	3091	ARG	NE-CZ-NH2	7.39	124.00	120.30
1	C	2031	ARG	NE-CZ-NH2	7.36	123.98	120.30
1	C	2022	ARG	NE-CZ-NH2	7.35	123.97	120.30
1	D	3058	ARG	NE-CZ-NH2	7.32	123.96	120.30
1	B	1022	ARG	NE-CZ-NH2	7.32	123.96	120.30
1	D	3031	ARG	NE-CZ-NH2	7.30	123.95	120.30
1	D	3022	ARG	NE-CZ-NH2	7.25	123.92	120.30
1	A	31	ARG	NE-CZ-NH2	6.56	123.58	120.30

There are no chirality outliers.

There are no planarity outliers.

## 5.2 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 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	2233	0	2251	35	0
1	B	2227	0	2238	33	0
1	C	2227	0	2238	43	0
1	D	2227	0	2238	35	0
2	A	179	0	0	2	0
2	B	168	0	0	3	0
2	C	194	0	0	0	0
2	D	187	0	0	1	0
All	All	9642	0	8965	137	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 8.

All (137) 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)
1:D:3091:ARG:HH11	1:D:3091:ARG:HG2	1.36	0.89
1:A:125:ALA:HB3	1:A:129:GLU:HG3	1.60	0.82
1:C:2001:MET:HE1	1:C:2294:HIS:HA	1.62	0.82
1:B:1087:GLN:HE21	1:B:1091:ARG:HE	1.37	0.72
1:C:2234:ASN:H	1:C:2234:ASN:HD22	1.40	0.70
1:D:3087:GLN:HE21	1:D:3091:ARG:HE	1.40	0.69
1:C:2181:GLU:CD	1:D:3031:ARG:HH22	1.97	0.68
1:B:1125:ALA:HB3	1:B:1129:GLU:HG3	1.76	0.67
1:A:234:ASN:H	1:A:234:ASN:HD22	1.44	0.66
1:B:1134:ILE:HD12	1:B:1139:VAL:HG21	1.80	0.63
1:D:3022:ARG:HD3	1:D:3107:GLY:O	1.99	0.63
1:B:1125:ALA:HB2	1:B:1142:VAL:HG13	1.81	0.61
1:D:3202:VAL:HB	1:D:3246:GLU:HB3	1.82	0.61
1:B:1016:PHE:HE2	1:B:1286:LEU:HB3	1.64	0.61
1:A:122:LYS:O	1:A:129:GLU:HG2	2.01	0.60
2:A:5096:HOH:O	1:B:1031:ARG:HD2	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:3091:ARG:NH1	1:D:3091:ARG:HG2	2.08	0.60
1:D:3117:VAL:HB	1:D:3155:ILE:HD12	1.82	0.60
1:C:2125:ALA:HB2	1:C:2142:VAL:HG13	1.84	0.60
1:A:221:ASN:C	1:A:221:ASN:HD22	2.05	0.59
1:B:1087:GLN:NE2	1:B:1091:ARG:HE	2.00	0.59
1:C:2089:LEU:HD13	1:C:2117:VAL:HG13	1.85	0.58
1:A:282:TYR:HB3	1:A:286:LEU:HD13	1.87	0.57
1:A:31:ARG:HD3	1:A:31:ARG:C	2.25	0.57
1:D:3087:GLN:NE2	1:D:3091:ARG:HE	2.03	0.57
1:A:119:PHE:O	1:A:122:LYS:HG2	2.05	0.57
1:D:3125:ALA:HB2	1:D:3142:VAL:HG13	1.88	0.56
1:C:2044:ALA:HB2	1:C:2091:ARG:NH2	2.21	0.56
1:B:1091:ARG:O	1:B:1095:CYS:HB2	2.05	0.55
1:A:91:ARG:O	1:A:95:CYS:HB2	2.07	0.55
1:A:216:LEU:HD23	1:A:219:THR:HG22	1.88	0.55
1:C:2087:GLN:NE2	1:C:2091:ARG:HE	2.05	0.54
1:D:3091:ARG:O	1:D:3095:CYS:HB2	2.08	0.54
1:D:3223:ARG:HD2	1:D:3234:ASN:HB3	1.89	0.54
1:C:2234:ASN:ND2	1:C:2234:ASN:H	2.04	0.54
1:A:137:ARG:HD3	1:A:152:GLU:OE1	2.06	0.54
1:A:290:GLN:O	1:A:290:GLN:HG3	2.07	0.54
1:D:3282:TYR:HB3	1:D:3286:LEU:HD13	1.89	0.54
1:A:221:ASN:HD22	1:A:223:ARG:H	1.56	0.54
1:C:2031:ARG:HH22	1:D:3181:GLU:CD	2.12	0.53
1:D:3008:ILE:O	1:D:3294:HIS:HE1	1.91	0.53
1:A:125:ALA:HB2	1:A:142:VAL:HG13	1.89	0.53
1:A:281:ASN:HD22	1:A:283:GLY:H	1.58	0.52
1:C:2117:VAL:HB	1:C:2155:ILE:HD12	1.92	0.52
1:C:2202:VAL:HB	1:C:2246:GLU:HB3	1.91	0.52
1:D:3125:ALA:HB3	1:D:3129:GLU:HB3	1.91	0.52
1:C:2197:PHE:CG	1:C:2202:VAL:HG22	2.46	0.51
1:A:238:PRO:HG2	2:A:5920:HOH:O	2.09	0.51
1:A:221:ASN:ND2	1:A:223:ARG:H	2.09	0.51
1:A:204:SER:HA	1:B:1144:ILE:HB	1.93	0.51
1:C:2132:THR:HG22	1:C:2133:GLN:H	1.75	0.51
1:C:2223:ARG:HD2	1:C:2234:ASN:O	2.11	0.51
1:A:144:ILE:HB	1:B:1204:SER:HA	1.93	0.51
1:C:2040:ALA:O	1:C:2091:ARG:HD3	2.11	0.51
1:B:1236:GLU:O	1:B:1238:PRO:HD3	2.11	0.50
1:C:2133:GLN:HA	1:C:2137:ARG:O	2.11	0.50
1:A:202:VAL:HG21	1:A:248:VAL:HG23	1.92	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1057:GLU:HG2	1:B:1058:ARG:HG2	1.94	0.49
1:C:2087:GLN:HE21	1:C:2091:ARG:HE	1.60	0.49
1:C:2087:GLN:O	1:C:2091:ARG:HG3	2.12	0.49
1:B:1223:ARG:HD2	1:B:1234:ASN:ND2	2.28	0.48
1:C:2144:ILE:HB	1:D:3204:SER:HA	1.95	0.48
1:B:1008:ILE:O	1:B:1294:HIS:HE1	1.96	0.48
1:B:1046:SER:OG	1:B:1048:ILE:HG12	2.14	0.48
1:B:1202:VAL:HB	1:B:1246:GLU:HB3	1.95	0.48
1:D:3212:LYS:HE2	2:D:5849:HOH:O	2.12	0.48
1:C:2118:GLY:HA3	1:C:2152:GLU:O	2.14	0.48
1:C:2181:GLU:CD	1:D:3031:ARG:NH2	2.66	0.48
1:B:1221:ASN:HD21	1:B:1223:ARG:HG3	1.79	0.48
1:C:2234:ASN:N	1:C:2234:ASN:HD22	2.01	0.47
1:B:1221:ASN:HD22	1:B:1223:ARG:H	1.62	0.47
1:C:2031:ARG:NH1	1:C:2031:ARG:HG3	2.29	0.47
1:D:3091:ARG:NH1	1:D:3091:ARG:CG	2.77	0.47
1:B:1137:ARG:HD3	1:B:1152:GLU:CD	2.35	0.47
1:A:197:PHE:CG	1:A:202:VAL:HG22	2.50	0.47
1:A:78:THR:HG21	1:A:84:LEU:N	2.29	0.47
1:D:3044:ALA:HB2	1:D:3091:ARG:NH2	2.30	0.47
1:D:3173:GLN:HA	1:D:3243:ALA:CB	2.45	0.47
1:C:2221:ASN:HD21	1:C:2223:ARG:HG3	1.79	0.47
1:A:15:ALA:HB3	1:A:289:TYR:HB2	1.95	0.47
1:D:3119:PHE:O	1:D:3122:LYS:HG2	2.15	0.46
1:C:2031:ARG:HG3	1:C:2031:ARG:HH11	1.80	0.46
1:B:1134:ILE:O	1:B:1137:ARG:HB2	2.16	0.46
1:D:3197:PHE:CG	1:D:3202:VAL:HG22	2.50	0.46
1:C:2202:VAL:HG21	1:C:2248:VAL:HG23	1.98	0.46
1:A:117:VAL:HB	1:A:155:ILE:HD12	1.98	0.45
1:C:2223:ARG:HD3	1:C:2234:ASN:OD1	2.17	0.45
1:B:1031:ARG:O	1:B:1035:VAL:HG23	2.16	0.45
1:A:31:ARG:CD	1:A:31:ARG:C	2.85	0.45
1:A:9:GLU:HG3	1:A:276:LYS:HD3	1.97	0.45
1:A:20:ILE:HG12	1:A:77:CYS:SG	2.57	0.45
1:C:2126:ASP:HA	1:D:3213:TYR:CD2	2.51	0.45
1:D:3234:ASN:HD22	1:D:3234:ASN:HA	1.62	0.44
1:C:2022:ARG:HD3	1:C:2107:GLY:HA3	1.98	0.44
1:D:3171:ASP:HB3	1:D:3275:LYS:HE2	2.00	0.44
1:B:1085:GLU:HG3	1:B:1115:ASP:HB3	1.99	0.44
1:B:1086:GLU:O	1:B:1090:GLU:HG3	2.18	0.44
1:A:16:PHE:N	1:A:16:PHE:CD1	2.86	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:31:ARG:HD3	1:A:31:ARG:O	2.18	0.44
1:B:1221:ASN:C	1:B:1221:ASN:HD22	2.21	0.44
1:B:1020:ILE:HG12	1:B:1077:CYS:SG	2.58	0.44
1:B:1214:LYS:HE2	2:B:5666:HOH:O	2.18	0.43
1:C:2173:GLN:HA	1:C:2243:ALA:CB	2.48	0.43
1:C:2091:ARG:O	1:C:2095:CYS:HB2	2.19	0.43
1:A:78:THR:HG22	1:A:84:LEU:HB2	2.01	0.43
1:A:47:VAL:HG22	1:A:53:GLU:OE2	2.19	0.43
1:C:2012:TYR:HB3	1:C:2292:LYS:HG2	2.01	0.43
1:C:2204:SER:HA	1:D:3144:ILE:HB	2.01	0.43
1:C:2122:LYS:HD2	1:C:2140:TYR:CD2	2.54	0.42
1:C:2234:ASN:N	1:C:2234:ASN:ND2	2.67	0.42
1:B:1202:VAL:HG21	1:B:1248:VAL:HG23	2.00	0.42
1:D:3019:LYS:HG2	1:D:3079:PHE:O	2.19	0.42
1:A:44:ALA:HB2	1:A:91:ARG:NH2	2.35	0.42
1:A:145:MET:HA	1:B:1224:PHE:O	2.19	0.42
1:B:1281:ASN:C	1:B:1281:ASN:HD22	2.23	0.42
1:D:3137:ARG:HD3	1:D:3152:GLU:OE1	2.20	0.42
1:C:2031:ARG:O	1:C:2035:VAL:HG23	2.20	0.42
1:A:19:LYS:O	1:A:77:CYS:HA	2.20	0.42
1:D:3026:THR:O	1:D:3102:THR:HA	2.20	0.42
1:C:2255:GLU:CD	1:C:2255:GLU:H	2.22	0.42
1:A:234:ASN:N	1:A:234:ASN:HD22	2.13	0.41
1:B:1031:ARG:NH2	2:B:5244:HOH:O	2.51	0.41
1:C:2044:ALA:HB2	1:C:2091:ARG:HH22	1.85	0.41
1:D:3292:LYS:HD2	1:D:3295:GLU:OE2	2.20	0.41
1:D:3171:ASP:O	1:D:3242:ASN:HB2	2.21	0.41
1:C:2216:LEU:HA	1:C:2216:LEU:HD12	1.96	0.41
1:C:2031:ARG:CD	1:C:2031:ARG:C	2.89	0.41
1:D:3213:TYR:HB3	1:D:3215:PHE:CE1	2.56	0.41
1:C:2184:VAL:HG21	1:C:2201:ILE:HG13	2.03	0.41
1:B:1137:ARG:HD3	1:B:1152:GLU:OE1	2.20	0.41
1:B:1223:ARG:NH1	2:B:6074:HOH:O	2.53	0.41
1:C:2099:ALA:HA	1:C:2100:PRO:HD3	1.90	0.41
1:C:2018:ILE:HG21	1:C:2162:ALA:HB2	2.02	0.41
1:B:1048:ILE:C	1:B:1049:MET:HG2	2.40	0.40
1:D:3186:THR:O	1:D:3189:GLU:HB2	2.21	0.40
1:A:87:GLN:O	1:A:91:ARG:HB2	2.22	0.40
1:D:3015:ALA:HB3	1:D:3289:TYR:HB2	2.02	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 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 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	296/297 (100%)	286 (97%)	9 (3%)	1 (0%)	46	29
1	B	295/297 (99%)	289 (98%)	6 (2%)	0	100	100
1	C	295/297 (99%)	288 (98%)	7 (2%)	0	100	100
1	D	295/297 (99%)	286 (97%)	9 (3%)	0	100	100
All	All	1181/1188 (99%)	1149 (97%)	31 (3%)	1 (0%)	56	39

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	49	MET

### 5.3.2 Protein sidechains

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	229/228 (100%)	213 (93%)	16 (7%)	19	5
1	B	228/228 (100%)	216 (95%)	12 (5%)	28	10
1	C	228/228 (100%)	213 (93%)	15 (7%)	21	6
1	D	228/228 (100%)	215 (94%)	13 (6%)	25	9
All	All	913/912 (100%)	857 (94%)	56 (6%)	23	6

All (56) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	31	ARG
1	A	48	ILE
1	A	59	LEU
1	A	78	THR
1	A	81	TYR
1	A	129	GLU
1	A	130	SER
1	A	208	SER
1	A	214	LYS
1	A	216	LEU
1	A	221	ASN
1	A	231	LYS
1	A	233	GLU
1	A	234	ASN
1	A	281	ASN
1	A	286	LEU
1	B	1031	ARG
1	B	1049	MET
1	B	1113	LYS
1	B	1122	LYS
1	B	1129	GLU
1	B	1130	SER
1	B	1206	SER
1	B	1217	LYS
1	B	1221	ASN
1	B	1233	GLU
1	B	1281	ASN
1	B	1294	HIS
1	C	2031	ARG
1	C	2059	LEU
1	C	2082	GLU
1	C	2089	LEU
1	C	2120	LYS
1	C	2122	LYS
1	C	2132	THR
1	C	2216	LEU
1	C	2221	ASN
1	C	2234	ASN
1	C	2276	LYS
1	C	2281	ASN
1	C	2285	LYS
1	C	2286	LEU
1	C	2293	LEU

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Mol	Chain	Res	Type
1	D	3009	GLU
1	D	3031	ARG
1	D	3049	MET
1	D	3059	LEU
1	D	3080	LYS
1	D	3091	ARG
1	D	3133	GLN
1	D	3221	ASN
1	D	3227	SER
1	D	3234	ASN
1	D	3281	ASN
1	D	3286	LEU
1	D	3293	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	221	ASN
1	A	234	ASN
1	A	281	ASN
1	B	1087	GLN
1	B	1221	ASN
1	B	1242	ASN
1	B	1281	ASN
1	B	1294	HIS
1	C	2087	GLN
1	C	2221	ASN
1	C	2234	ASN
1	C	2281	ASN
1	C	2290	GLN
1	D	3087	GLN
1	D	3133	GLN
1	D	3221	ASN
1	D	3234	ASN
1	D	3242	ASN
1	D	3281	ASN
1	D	3294	HIS

### 5.3.3 RNA ⓘ

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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

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

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

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	297/297 (100%)	-0.12	8 (2%) 58 55	15, 29, 57, 73	2 (0%)
1	B	297/297 (100%)	-0.12	5 (1%) 73 72	15, 31, 58, 76	0
1	C	297/297 (100%)	-0.17	3 (1%) 84 84	15, 29, 53, 84	5 (1%)
1	D	297/297 (100%)	-0.27	6 (2%) 68 67	14, 28, 52, 76	2 (0%)
All	All	1188/1188 (100%)	-0.17	22 (1%) 70 69	14, 29, 56, 84	9 (0%)

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	3285	LYS	4.3
1	B	1286	LEU	4.0
1	D	3284	GLY	3.9
1	C	2285	LYS	3.7
1	B	1285	LYS	3.5
1	C	2284	GLY	3.4
1	D	3286	LEU	3.3
1	A	215	PHE	3.3
1	A	233	GLU	3.2
1	C	2233	GLU	3.2
1	A	286	LEU	3.2
1	D	3283	GLY	3.1
1	B	1233	GLU	3.0
1	B	1016	PHE	2.8
1	A	48	ILE	2.5
1	A	208	SER	2.5
1	A	285	LYS	2.5
1	A	290	GLN	2.4
1	A	217	LYS	2.3
1	D	3217	LYS	2.1
1	D	3215	PHE	2.1
1	B	1082	GLU	2.0



## 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 carbohydrates in this entry.

## 6.4 Ligands [i](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.