



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 16, 2018 – 09:57 AM EST

PDB ID : 4I36  
Title : Crystal Structure of the Bacillus stearothermophilus Phosphofructokinase Mutant D12A  
Authors : Mosser, R.; Reddy, M.; Bruning, J.B.; Sacchettini, J.C.; Reinhart, G.D.  
Deposited on : 2012-11-25  
Resolution : 2.30 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  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20030736  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
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) : rb-20030736

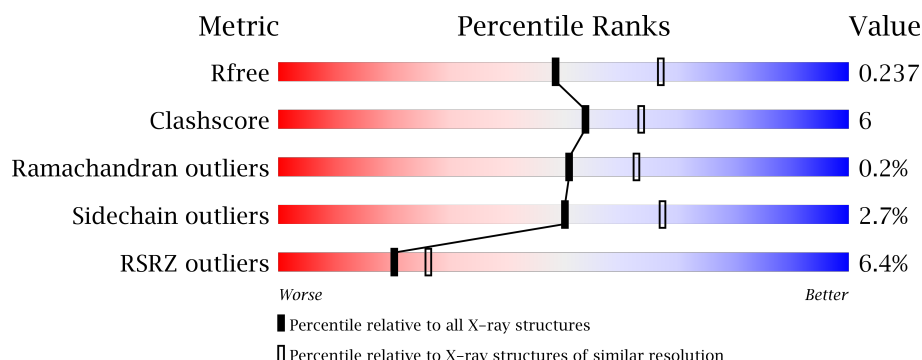
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 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))
$R_{free}$	100719	4130 (2.30-2.30)
Clashscore	112137	4751 (2.30-2.30)
Ramachandran outliers	110173	4705 (2.30-2.30)
Sidechain outliers	110143	4704 (2.30-2.30)
RSRZ outliers	101464	4156 (2.30-2.30)

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	319	<div> <div>6%</div> <div>85%</div> <div>14%</div> <div>.</div> </div>
1	B	319	<div> <div>8%</div> <div>86%</div> <div>13%</div> <div>.</div> </div>
1	C	319	<div> <div>9%</div> <div>87%</div> <div>12%</div> <div>.</div> </div>
1	D	319	<div> <div>3%</div> <div>83%</div> <div>15%</div> <div>.</div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 9846 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 6-phosphofructokinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	318	Total	C	N	O	S	0	2	0
			2370	1479	430	452	9			
1	B	319	Total	C	N	O	S	0	3	0
			2388	1495	427	457	9			
1	C	319	Total	C	N	O	S	0	2	0
			2380	1493	427	451	9			
1	D	319	Total	C	N	O	S	0	1	0
			2379	1487	431	452	9			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	12	ALA	ASP	ENGINEERED MUTATION	UNP P00512
B	12	ALA	ASP	ENGINEERED MUTATION	UNP P00512
C	12	ALA	ASP	ENGINEERED MUTATION	UNP P00512
D	12	ALA	ASP	ENGINEERED MUTATION	UNP P00512

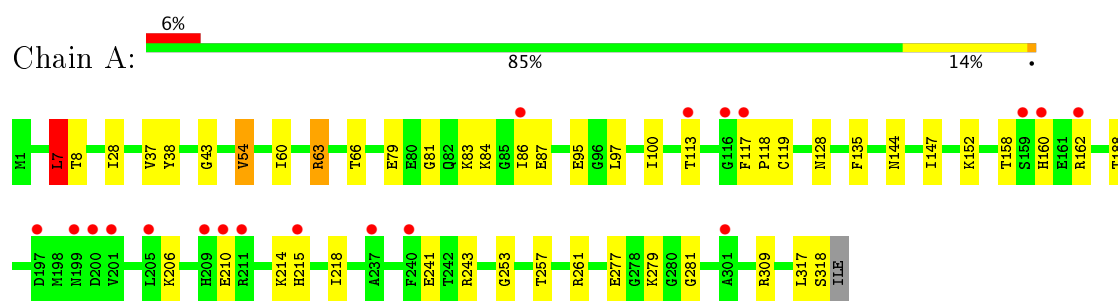
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	76	Total	O	0	0
			76	76		
2	B	72	Total	O	0	0
			72	72		
2	C	75	Total	O	0	0
			75	75		
2	D	106	Total	O	0	0
			106	106		

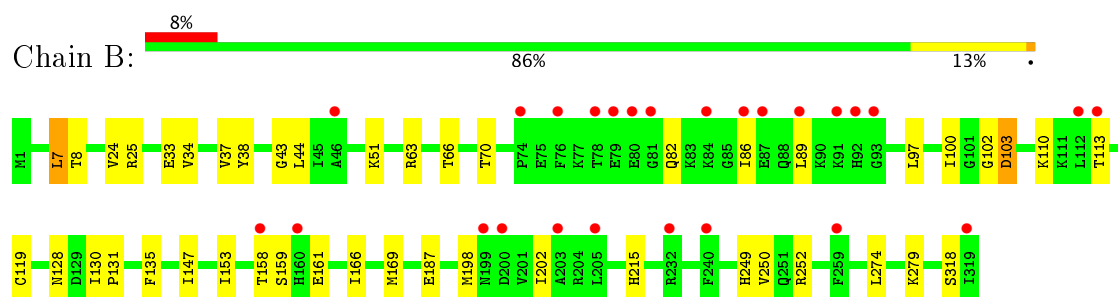
### 3 Residue-property plots [i](#)

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

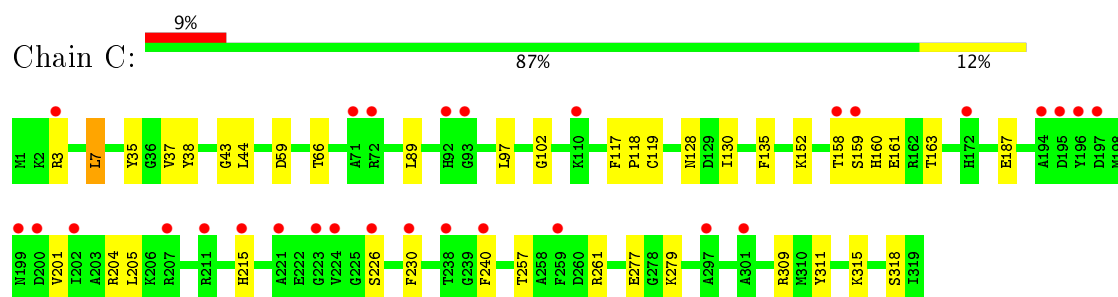
- Molecule 1: 6-phosphofructokinase



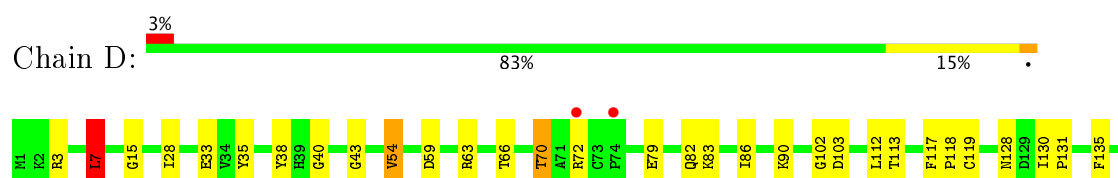
- Molecule 1: 6-phosphofructokinase

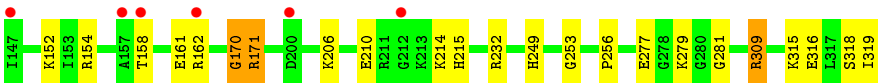


- Molecule 1: 6-phosphofructokinase



- Molecule 1: 6-phosphofructokinase





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	96.05Å 112.66Å 129.73Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.89 – 2.30 38.89 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.9 (38.89-2.30) 99.9 (38.89-2.30)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	0.07	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.20 (at 2.29Å)	Xtriage
Refinement program	PHENIX	Depositor
R, $R_{free}$	0.201 , 0.254 0.185 , 0.237	Depositor DCC
$R_{free}$ test set	3205 reflections (5.08%)	DCC
Wilson B-factor (Å <sup>2</sup> )	43.6	Xtriage
Anisotropy	0.244	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 45.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	9846	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.93% 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](#)

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	0.28	0/2412	0.45	1/3258 (0.0%)
1	B	0.27	0/2433	0.45	0/3284
1	C	0.26	0/2422	0.45	0/3267
1	D	0.29	0/2417	0.48	1/3262 (0.0%)
All	All	0.28	0/9684	0.46	2/13071 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	7	LEU	CA-CB-CG	6.13	129.40	115.30
1	A	7	LEU	CA-CB-CG	5.58	128.12	115.30

There are no chirality outliers.

There are no planarity outliers.

### 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	2370	0	2358	39	0
1	B	2388	0	2380	28	0
1	C	2380	0	2388	26	0
1	D	2379	0	2377	36	0
2	A	76	0	0	1	0
2	B	72	0	0	2	0
2	C	75	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	106	0	0	4	0
All	All	9846	0	9503	116	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 6.

All (116) 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:28:ILE:HD13	1:D:54:VAL:HG13	1.57	0.86
1:D:102:GLY:HA2	1:D:130:ILE:HD11	1.61	0.81
1:C:3:ARG:HD2	1:C:35:TYR:CE1	2.20	0.77
1:D:3:ARG:HG2	1:D:33:GLU:HB2	1.65	0.77
1:D:277:GLU:HG3	1:D:279:LYS:HD3	1.68	0.74
1:D:79:GLU:O	1:D:83:LYS:HG2	1.86	0.74
1:D:38:TYR:O	1:D:43:GLY:HA3	1.89	0.71
1:B:97:LEU:HB3	1:B:119[B]:CYS:SG	2.30	0.71
1:C:38:TYR:O	1:C:43:GLY:HA3	1.90	0.71
1:A:188:THR:HB	1:A:218:ILE:HD12	1.76	0.67
1:D:113:THR:HG21	1:D:281:GLY:N	2.12	0.66
1:B:7:LEU:HB3	1:B:37:VAL:HB	1.79	0.65
1:D:3:ARG:HD3	1:D:35:TYR:CE1	2.32	0.65
1:B:187:GLU:HA	1:B:318:SER:HB2	1.79	0.65
1:A:38:TYR:O	1:A:43:GLY:HA3	1.98	0.64
1:A:28:ILE:HD13	1:A:54:VAL:HG13	1.79	0.63
1:A:60:ILE:HA	1:A:63:ARG:HG3	1.81	0.62
1:C:226:SER:O	1:C:230:PHE:HD2	1.82	0.61
1:A:113:THR:HG21	1:A:281:GLY:H	1.66	0.61
1:C:7:LEU:HB3	1:C:37:VAL:HB	1.83	0.60
1:C:102:GLY:HA2	1:C:130:ILE:HD11	1.82	0.60
1:B:158:THR:HG22	1:B:159:SER:H	1.66	0.59
1:C:59:ASP:HB2	1:D:214:LYS:HE3	1.85	0.59
1:D:315:LYS:HD2	2:D:492:HOH:O	2.02	0.59
1:A:113:THR:HG21	1:A:281:GLY:N	2.18	0.59
1:A:97:LEU:HB3	1:A:119[A]:CYS:SG	2.43	0.58
1:B:102:GLY:HA2	1:B:130:ILE:HD11	1.85	0.58
1:B:38:TYR:O	1:B:43:GLY:HA3	2.02	0.58
1:D:206:LYS:O	1:D:210:GLU:HG2	2.03	0.58
1:B:44:LEU:HD11	1:B:89:LEU:HG	1.86	0.58
1:B:128:ASN:OD1	1:B:135:PHE:HA	2.04	0.57
1:C:158:THR:O	1:C:160:HIS:N	2.37	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:81:GLY:HA2	1:A:84:LYS:HE3	1.86	0.56
1:B:158:THR:HG22	1:B:159:SER:N	2.21	0.56
1:B:33:GLU:OE2	1:B:51:LYS:HE3	2.06	0.56
1:C:44:LEU:HD11	1:C:89:LEU:HG	1.88	0.56
1:A:158:THR:HG22	1:D:70:THR:HG23	1.88	0.54
1:D:7:LEU:C	1:D:7:LEU:HD23	2.28	0.54
1:D:82:GLN:O	1:D:86:ILE:HG13	2.08	0.54
1:D:7:LEU:HG	1:D:40:GLY:HA2	1.90	0.53
1:D:113:THR:HG21	1:D:281:GLY:H	1.71	0.53
1:A:206:LYS:O	1:A:210:GLU:HG2	2.08	0.53
1:A:160:HIS:HD2	2:D:500:HOH:O	1.91	0.53
1:D:112:LEU:HD13	1:D:119[B]:CYS:SG	2.49	0.53
1:C:158:THR:O	1:C:161:GLU:N	2.42	0.53
1:A:210:GLU:OE1	1:A:210:GLU:HA	2.10	0.52
1:B:250:VAL:HG23	1:C:152:LYS:HB2	1.92	0.52
1:A:160:HIS:CD2	2:D:500:HOH:O	2.63	0.51
1:C:97:LEU:HB3	1:C:119[A]:CYS:SG	2.50	0.51
1:D:154:ARG:O	1:D:215:HIS:CE1	2.64	0.51
1:A:7:LEU:HB3	1:A:37:VAL:HB	1.92	0.51
1:A:162:ARG:O	1:A:162:ARG:HG2	2.11	0.51
1:C:89:LEU:HD13	1:C:117:PHE:CE1	2.47	0.50
1:B:110:LYS:O	1:B:113:THR:HB	2.13	0.49
1:D:158:THR:HB	1:D:161:GLU:HB2	1.94	0.49
1:A:113:THR:HG22	1:A:119[A]:CYS:H	1.78	0.49
1:B:250:VAL:HG22	1:C:152:LYS:HD3	1.95	0.49
1:A:128:ASN:OD1	1:A:135:PHE:HA	2.13	0.49
1:A:113:THR:HG22	1:A:119[B]:CYS:H	1.79	0.48
1:D:232:ARG:NH2	2:D:481:HOH:O	2.47	0.48
1:D:128:ASN:OD1	1:D:135:PHE:HA	2.14	0.48
1:A:113:THR:CG2	1:A:119[A]:CYS:H	2.27	0.48
1:C:7:LEU:CD1	1:C:97:LEU:HD21	2.43	0.48
1:A:160:HIS:O	1:A:241:GLU:HB2	2.13	0.48
1:A:113:THR:HA	1:A:117:PHE:O	2.14	0.47
1:A:113:THR:CG2	1:A:119[B]:CYS:H	2.27	0.47
1:D:113:THR:HG21	1:D:281:GLY:CA	2.44	0.47
1:D:15:GLY:N	1:D:256:PRO:HG3	2.30	0.47
1:C:311:TYR:CE2	1:C:315:LYS:HE3	2.50	0.47
1:C:3:ARG:HD2	1:C:35:TYR:CD1	2.50	0.47
1:A:188:THR:HB	1:A:218:ILE:CD1	2.41	0.46
1:D:318:SER:O	1:D:319:ILE:C	2.53	0.46
1:C:187:GLU:HA	1:C:318:SER:HB2	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:442:HOH:O	1:D:158:THR:HG23	2.16	0.46
1:D:309:ARG:NH2	1:D:316:GLU:OE2	2.48	0.46
1:B:63:ARG:NH1	1:D:63:ARG:HD2	2.30	0.45
1:C:277:GLU:OE2	1:C:279:LYS:HE2	2.16	0.45
1:A:8:THR:HA	1:A:100:ILE:O	2.16	0.45
1:C:257:THR:O	1:C:261:ARG:HG3	2.16	0.45
1:B:169:MET:HE1	1:B:249:HIS:HA	1.99	0.45
1:B:250:VAL:HG22	1:C:152:LYS:CD	2.47	0.44
1:C:163:THR:HG21	1:C:205:LEU:HD11	1.99	0.44
1:A:317:LEU:O	1:B:25:ARG:HD2	2.18	0.44
1:B:198:MET:O	1:B:202:ILE:HG12	2.17	0.44
1:C:201:VAL:O	1:C:204:ARG:HB2	2.17	0.44
1:A:162:ARG:HG3	1:A:214:LYS:O	2.17	0.43
1:A:261:ARG:CZ	1:B:147:ILE:HD12	2.49	0.43
1:B:130:ILE:HA	1:B:131:PRO:HD3	1.89	0.43
1:D:130:ILE:HA	1:D:131:PRO:HD3	1.85	0.43
1:A:277:GLU:OE1	1:A:279:LYS:HE3	2.17	0.43
1:B:8:THR:HA	1:B:100:ILE:O	2.18	0.43
1:D:117:PHE:HA	1:D:118:PRO:HD3	1.88	0.43
1:A:113:THR:HG22	1:A:118:PRO:HA	2.01	0.43
1:A:160:HIS:ND1	1:D:72:ARG:NH2	2.67	0.43
1:A:144:ASN:HA	1:A:147:ILE:HG12	2.01	0.42
1:C:205:LEU:HD22	1:C:240:PHE:CD2	2.54	0.42
1:A:152:LYS:O	1:D:253:GLY:HA3	2.20	0.42
1:B:153:ILE:HD12	1:B:166:ILE:HD11	2.00	0.42
1:C:226:SER:O	1:C:230:PHE:CD2	2.68	0.42
1:B:252:ARG:NH1	2:B:461:HOH:O	2.34	0.42
1:A:95:GLU:O	1:A:118:PRO:HD2	2.20	0.42
1:D:90:LYS:HB2	1:D:90:LYS:HE3	1.71	0.42
1:A:243:ARG:HD2	1:D:249:HIS:CE1	2.55	0.42
1:B:103:ASP:N	1:B:103:ASP:OD1	2.52	0.42
1:B:7:LEU:C	1:B:7:LEU:CD2	2.88	0.42
1:B:82:GLN:O	1:B:86:ILE:HG13	2.20	0.42
1:B:24:VAL:HA	1:B:34:VAL:HG21	2.00	0.42
1:A:83:LYS:O	1:A:87:GLU:HG2	2.20	0.41
1:A:160:HIS:ND1	1:D:72:ARG:CZ	2.84	0.41
1:D:170:GLY:O	1:D:171:ARG:CB	2.68	0.41
1:A:257:THR:O	1:A:261:ARG:HG3	2.20	0.41
1:C:117:PHE:HA	1:C:118:PRO:HD3	1.79	0.41
1:A:253:GLY:HA3	1:D:152:LYS:O	2.21	0.41
1:A:318:SER:HB3	2:B:467:HOH:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:128:ASN:OD1	1:C:135:PHE:HA	2.21	0.40
1:B:274:LEU:HD22	1:B:279:LYS:HD2	2.03	0.40

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	318/319 (100%)	311 (98%)	7 (2%)	0	100	100
1	B	320/319 (100%)	312 (98%)	8 (2%)	0	100	100
1	C	319/319 (100%)	308 (97%)	10 (3%)	1 (0%)	44	55
1	D	318/319 (100%)	310 (98%)	6 (2%)	2 (1%)	28	34
All	All	1275/1276 (100%)	1241 (97%)	31 (2%)	3 (0%)	51	63

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	159	SER
1	D	171	ARG
1	D	170	GLY

### 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	241/246 (98%)	233 (97%)	8 (3%)	43	59
1	B	243/246 (99%)	237 (98%)	6 (2%)	53	70
1	C	242/246 (98%)	238 (98%)	4 (2%)	66	81
1	D	242/246 (98%)	234 (97%)	8 (3%)	43	59
All	All	968/984 (98%)	942 (97%)	26 (3%)	50	67

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	LEU
1	A	54	VAL
1	A	63	ARG
1	A	66	THR
1	A	79	GLU
1	A	86	ILE
1	A	215	HIS
1	A	309	ARG
1	B	7	LEU
1	B	66	THR
1	B	70	THR
1	B	103	ASP
1	B	161	GLU
1	B	215	HIS
1	C	7	LEU
1	C	66	THR
1	C	215	HIS
1	C	309	ARG
1	D	7	LEU
1	D	54	VAL
1	D	59	ASP
1	D	66	THR
1	D	70	THR
1	D	103	ASP
1	D	162	ARG
1	D	309	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	209	HIS

### 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 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	318/319 (99%)	0.23	19 (5%)	23 29	24, 45, 76, 95	0
1	B	319/319 (100%)	0.42	26 (8%)	12 17	26, 45, 81, 102	0
1	C	319/319 (100%)	0.51	29 (9%)	10 14	27, 50, 84, 100	0
1	D	319/319 (100%)	0.19	8 (2%)	58 65	25, 41, 67, 88	0
All	All	1275/1276 (99%)	0.34	82 (6%)	20 26	24, 45, 79, 102	0

All (82) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	259[A]	PHE	6.5
1	B	160	HIS	4.5
1	B	259[A]	PHE	4.4
1	B	79	GLU	4.1
1	C	230	PHE	4.0
1	A	113	THR	4.0
1	C	202	ILE	3.8
1	B	113	THR	3.8
1	C	72	ARG	3.7
1	B	80	GLU	3.7
1	D	72	ARG	3.7
1	B	86	ILE	3.6
1	B	74	PRO	3.4
1	A	301	ALA	3.4
1	A	209	HIS	3.4
1	C	200	ASP	3.2
1	D	158	THR	3.1
1	D	74	PRO	3.1
1	A	205	LEU	3.1
1	C	238	THR	3.0
1	C	71	ALA	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	215	HIS	3.0
1	A	160	HIS	3.0
1	D	200	ASP	3.0
1	A	199	ASN	3.0
1	B	78	THR	3.0
1	A	86	ILE	2.9
1	C	195	ASP	2.9
1	B	232	ARG	2.8
1	B	240	PHE	2.8
1	B	319	ILE	2.8
1	B	158	THR	2.7
1	A	116	GLY	2.7
1	D	157	ALA	2.6
1	B	92	HIS	2.6
1	C	159	SER	2.6
1	C	223	GLY	2.6
1	A	211	ARG	2.6
1	D	162	ARG	2.6
1	C	301	ALA	2.6
1	C	158	THR	2.5
1	A	197	ASP	2.5
1	C	196	TYR	2.5
1	B	203	ALA	2.5
1	B	87	GLU	2.5
1	C	297	ALA	2.4
1	B	89	LEU	2.4
1	A	117	PHE	2.4
1	A	201	VAL	2.4
1	A	237	ALA	2.4
1	B	76	PHE	2.4
1	C	3	ARG	2.4
1	C	197	ASP	2.4
1	C	207	ARG	2.3
1	C	211	ARG	2.3
1	A	200	ASP	2.3
1	C	240	PHE	2.3
1	C	92	HIS	2.3
1	B	91	LYS	2.3
1	C	221	ALA	2.3
1	B	112	LEU	2.3
1	D	212	GLY	2.3
1	D	147	ILE	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	159	SER	2.3
1	C	224	VAL	2.2
1	B	81	GLY	2.2
1	A	162	ARG	2.2
1	A	210	GLU	2.1
1	A	240	PHE	2.1
1	C	110	LYS	2.1
1	B	200	ASP	2.1
1	B	205	LEU	2.1
1	C	93	GLY	2.1
1	C	194	ALA	2.1
1	C	172	HIS	2.1
1	B	84	LYS	2.1
1	B	93	GLY	2.1
1	C	215	HIS	2.0
1	B	199	ASN	2.0
1	C	199	ASN	2.0
1	B	46	ALA	2.0
1	C	226	SER	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.