



Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 08:23 am BST

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
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)	:	2.11

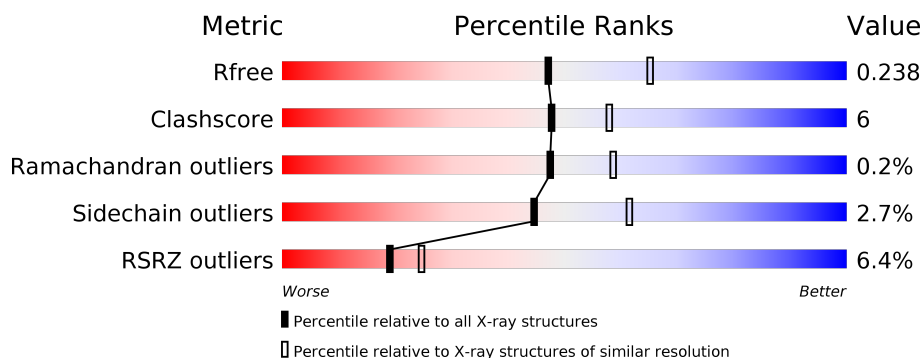
1 Overall quality at a glance

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}	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (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 ≥ 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	319	<div> <div>6%</div> <div> <div></div> <div>85%</div> <div>14%</div> </div> <div></div> </div>
1	B	319	<div> <div>8%</div> <div> <div></div> <div>86%</div> <div>13%</div> </div> <div></div> </div>
1	C	319	<div> <div>9%</div> <div> <div></div> <div>87%</div> <div>12%</div> </div> <div></div> </div>
1	D	319	<div> <div>3%</div> <div> <div></div> <div>83%</div> <div>15%</div> </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

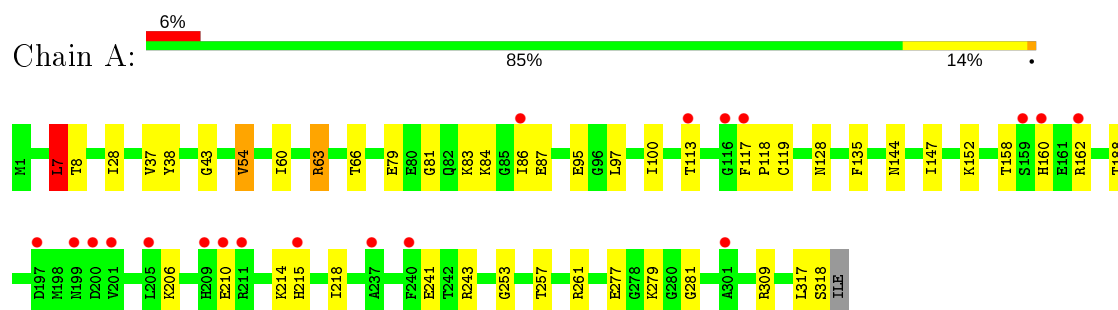
- Molecule 2 is water.

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		

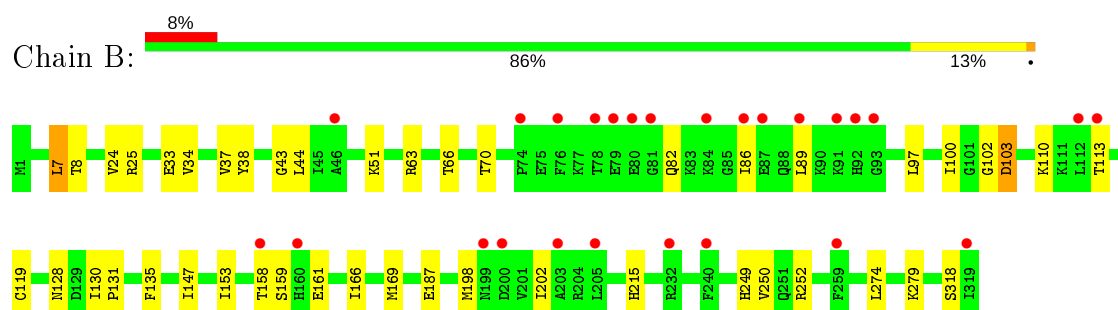
3 Residue-property plots [i](#)

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.

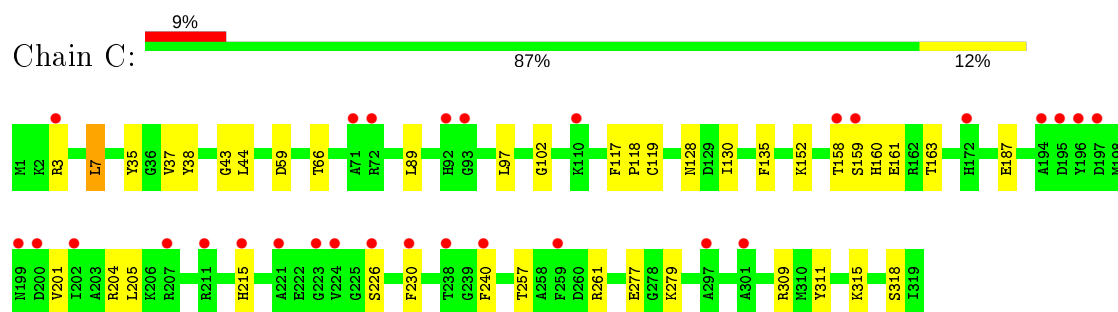
- Molecule 1: 6-phosphofructokinase



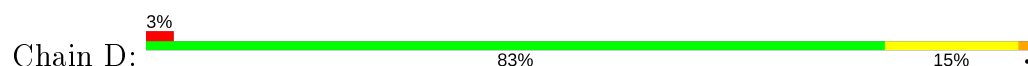
- Molecule 1: 6-phosphofructokinase

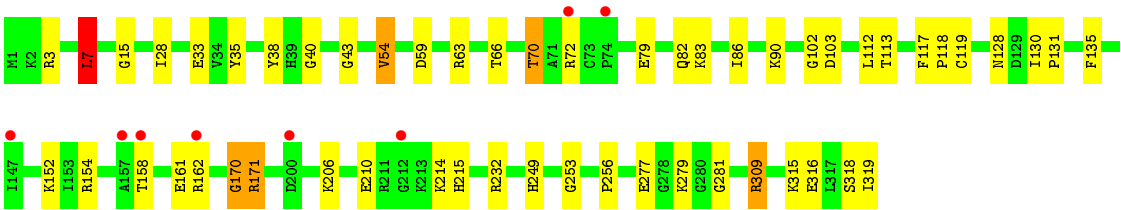


- Molecule 1: 6-phosphofructokinase



- Molecule 1: 6-phosphofructokinase





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	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
$\langle I/\sigma(I) \rangle$ ¹	3.20 (at 2.29Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.201 , 0.254 0.187 , 0.238	Depositor DCC
R_{free} test set	3204 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	43.6	Xtriage
Anisotropy	0.244	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 45.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	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 ²)	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.*

¹Intensities estimated from amplitudes.

²Theoretical 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%)	41	50
1	D	318/319 (100%)	310 (98%)	6 (2%)	2 (1%)	25	31
All	All	1275/1276 (100%)	1241 (97%)	31 (2%)	3 (0%)	47	58

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%)	38	53
1	B	243/246 (99%)	237 (98%)	6 (2%)	47	65
1	C	242/246 (98%)	238 (98%)	4 (2%)	60	76
1	D	242/246 (98%)	234 (97%)	8 (3%)	38	53
All	All	968/984 (98%)	942 (97%)	26 (3%)	44	61

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, 95th 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 ²)	Q<0.9
1	A	318/319 (99%)	0.23	19 (5%) 21 28	24, 45, 76, 95	0
1	B	319/319 (100%)	0.42	26 (8%) 11 15	26, 45, 81, 102	0
1	C	319/319 (100%)	0.51	29 (9%) 9 12	27, 50, 84, 100	0
1	D	319/319 (100%)	0.19	8 (2%) 57 64	25, 41, 67, 88	0
All	All	1275/1276 (99%)	0.34	82 (6%) 19 25	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	3.9
1	C	202	ILE	3.8
1	B	113	THR	3.7
1	B	80	GLU	3.7
1	C	72	ARG	3.7
1	D	72	ARG	3.7
1	B	86	ILE	3.5
1	B	74	PRO	3.4
1	A	209	HIS	3.3
1	A	301	ALA	3.3
1	C	200	ASP	3.2
1	A	205	LEU	3.1
1	D	74	PRO	3.1
1	D	158	THR	3.1
1	C	238	THR	3.0
1	D	200	ASP	3.0

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Mol	Chain	Res	Type	RSRZ
1	C	71	ALA	3.0
1	A	160	HIS	3.0
1	B	78	THR	3.0
1	A	215	HIS	3.0
1	A	86	ILE	2.9
1	A	199	ASN	2.9
1	B	240	PHE	2.8
1	C	195	ASP	2.8
1	B	232	ARG	2.8
1	B	158	THR	2.7
1	B	319	ILE	2.7
1	D	157	ALA	2.7
1	A	116	GLY	2.7
1	B	92	HIS	2.7
1	C	223	GLY	2.7
1	C	159	SER	2.6
1	A	211	ARG	2.6
1	D	162	ARG	2.6
1	A	197	ASP	2.6
1	C	158	THR	2.5
1	C	196	TYR	2.5
1	C	301	ALA	2.5
1	B	87	GLU	2.5
1	B	76	PHE	2.5
1	B	203	ALA	2.5
1	B	89	LEU	2.4
1	A	117	PHE	2.4
1	C	297	ALA	2.4
1	A	201	VAL	2.4
1	C	240	PHE	2.4
1	A	237	ALA	2.4
1	A	200	ASP	2.4
1	C	197	ASP	2.4
1	C	207	ARG	2.4
1	C	211	ARG	2.4
1	C	92	HIS	2.4
1	C	3	ARG	2.3
1	B	112	LEU	2.3
1	B	91	LYS	2.3
1	C	221	ALA	2.3
1	C	224	VAL	2.3
1	D	147	ILE	2.3

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Mol	Chain	Res	Type	RSRZ
1	D	212	GLY	2.2
1	A	159	SER	2.2
1	B	81	GLY	2.2
1	A	162	ARG	2.2
1	A	240	PHE	2.1
1	C	110	LYS	2.1
1	A	210	GLU	2.1
1	B	205	LEU	2.1
1	C	172	HIS	2.1
1	B	200	ASP	2.1
1	C	93	GLY	2.1
1	C	194	ALA	2.1
1	C	199	ASN	2.1
1	B	46	ALA	2.1
1	B	84	LYS	2.0
1	B	93	GLY	2.0
1	B	199	ASN	2.0
1	C	215	HIS	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.