



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 01:49 am BST

PDB ID : 4B04
Title : Crystal structure of the Catalytic Domain of Human DUSP26 (C152S)
Authors : Won, E.-Y.; Lee, D.Y.; Park, S.G.; Yokoyama, S.; Kim, S.J.; Chi, S.-W.
Deposited on : 2012-06-28
Resolution : 2.21 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

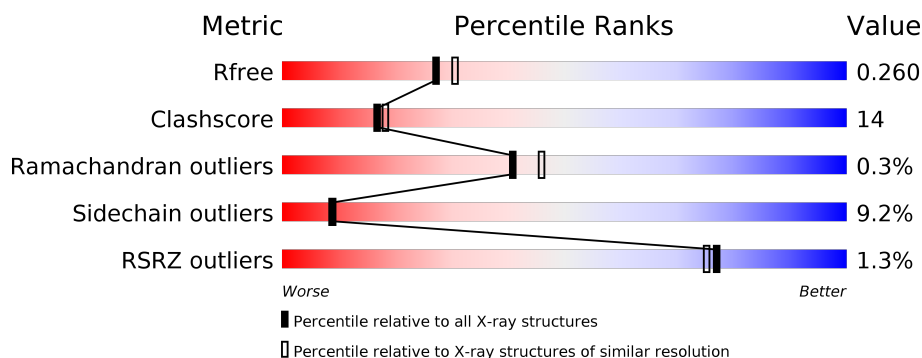
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.21 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	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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	160	<div> <div>71%</div> <div>21%</div> <div>• 5%</div> </div>
1	B	160	<div>4%</div> <div>63%</div> <div>26%</div> <div>6%</div> <div>6%</div>
1	C	160	<div>69%</div> <div>21%</div> <div>• • 6%</div>
1	D	160	<div>%</div> <div>66%</div> <div>22%</div> <div>• 8%</div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 9687 atoms, of which 4821 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 DUAL SPECIFICITY PROTEIN PHOSPHATASE 26.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	152	Total	C	H	N	O	S	0	0	0
			2435	764	1225	232	210	4			
1	B	150	Total	C	H	N	O	S	0	0	0
			2408	756	1212	230	207	3			
1	C	150	Total	C	H	N	O	S	0	0	0
			2384	749	1200	223	208	4			
1	D	147	Total	C	H	N	O	S	0	0	0
			2347	735	1184	223	202	3			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	60	MET	-	expression tag	UNP Q9BV47
A	212	LEU	-	expression tag	UNP Q9BV47
A	213	GLU	-	expression tag	UNP Q9BV47
A	214	HIS	-	expression tag	UNP Q9BV47
A	215	HIS	-	expression tag	UNP Q9BV47
A	216	HIS	-	expression tag	UNP Q9BV47
A	217	HIS	-	expression tag	UNP Q9BV47
A	218	HIS	-	expression tag	UNP Q9BV47
A	219	HIS	-	expression tag	UNP Q9BV47
A	152	SER	CYS	engineered mutation	UNP Q9BV47
B	60	MET	-	expression tag	UNP Q9BV47
B	212	LEU	-	expression tag	UNP Q9BV47
B	213	GLU	-	expression tag	UNP Q9BV47
B	214	HIS	-	expression tag	UNP Q9BV47
B	215	HIS	-	expression tag	UNP Q9BV47
B	216	HIS	-	expression tag	UNP Q9BV47
B	217	HIS	-	expression tag	UNP Q9BV47
B	218	HIS	-	expression tag	UNP Q9BV47
B	219	HIS	-	expression tag	UNP Q9BV47
B	152	SER	CYS	engineered mutation	UNP Q9BV47
C	60	MET	-	expression tag	UNP Q9BV47

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Chain	Residue	Modelled	Actual	Comment	Reference
C	212	LEU	-	expression tag	UNP Q9BV47
C	213	GLU	-	expression tag	UNP Q9BV47
C	214	HIS	-	expression tag	UNP Q9BV47
C	215	HIS	-	expression tag	UNP Q9BV47
C	216	HIS	-	expression tag	UNP Q9BV47
C	217	HIS	-	expression tag	UNP Q9BV47
C	218	HIS	-	expression tag	UNP Q9BV47
C	219	HIS	-	expression tag	UNP Q9BV47
C	152	SER	CYS	engineered mutation	UNP Q9BV47
D	60	MET	-	expression tag	UNP Q9BV47
D	212	LEU	-	expression tag	UNP Q9BV47
D	213	GLU	-	expression tag	UNP Q9BV47
D	214	HIS	-	expression tag	UNP Q9BV47
D	215	HIS	-	expression tag	UNP Q9BV47
D	216	HIS	-	expression tag	UNP Q9BV47
D	217	HIS	-	expression tag	UNP Q9BV47
D	218	HIS	-	expression tag	UNP Q9BV47
D	219	HIS	-	expression tag	UNP Q9BV47
D	152	SER	CYS	engineered mutation	UNP Q9BV47

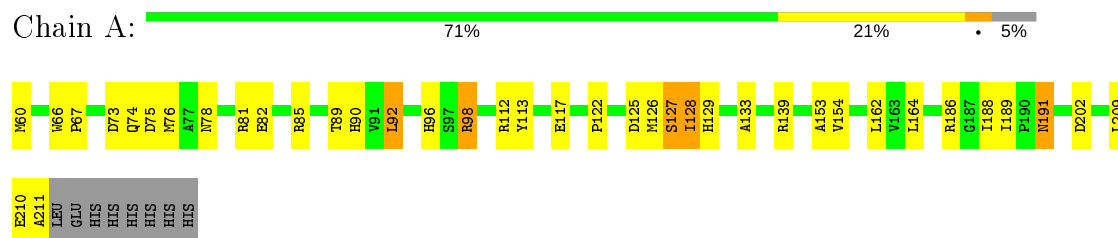
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	44	Total O 44 44	0	0
2	B	22	Total O 22 22	0	0
2	C	27	Total O 27 27	0	0
2	D	20	Total O 20 20	0	0

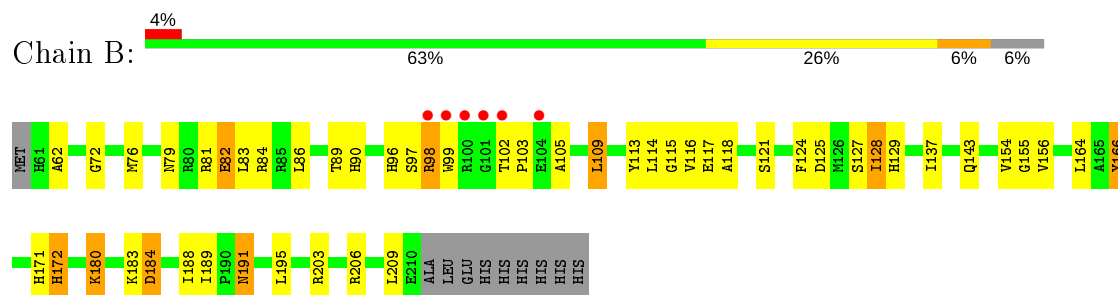
3 Residue-property plots [i](#)

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.

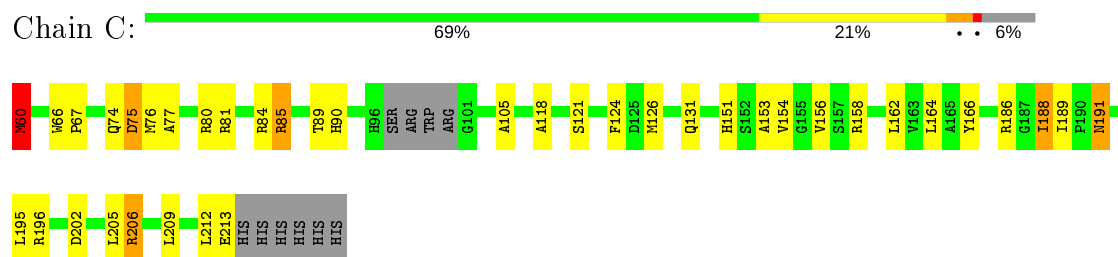
• Molecule 1: DUAL SPECIFICITY PROTEIN PHOSPHATASE 26



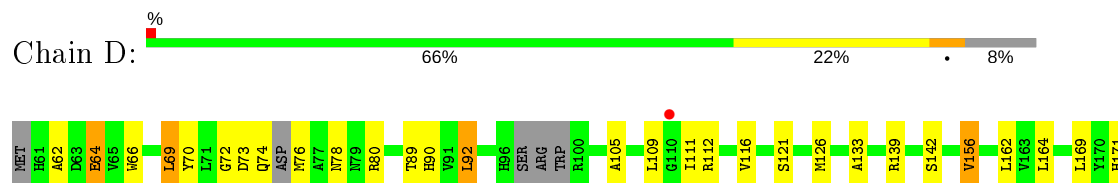
• Molecule 1: DUAL SPECIFICITY PROTEIN PHOSPHATASE 26

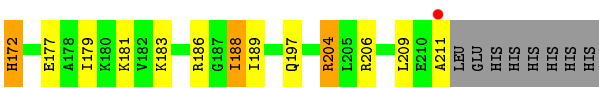


• Molecule 1: DUAL SPECIFICITY PROTEIN PHOSPHATASE 26



• Molecule 1: DUAL SPECIFICITY PROTEIN PHOSPHATASE 26





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	82.60Å 82.78Å 91.72Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.12 – 2.21 40.12 – 2.21	Depositor EDS
% Data completeness (in resolution range)	98.0 (40.12-2.21) 98.0 (40.12-2.21)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.38 (at 2.20Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.200 , 0.266 0.194 , 0.260	Depositor DCC
R_{free} test set	1580 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	28.0	Xtriage
Anisotropy	0.599	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 49.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.008 for k,h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9687	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 49.60 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 7.2858e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² 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	1.08	0/1237	0.97	2/1670 (0.1%)
1	B	0.95	1/1223 (0.1%)	0.93	2/1653 (0.1%)
1	C	1.00	1/1208 (0.1%)	0.93	3/1631 (0.2%)
1	D	0.98	1/1186 (0.1%)	0.96	2/1598 (0.1%)
All	All	1.00	3/4854 (0.1%)	0.95	9/6552 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	166	TYR	CE2-CZ	5.83	1.46	1.38
1	D	177	GLU	CG-CD	5.64	1.60	1.51
1	B	166	TYR	CG-CD1	5.38	1.46	1.39

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	156	VAL	CB-CA-C	-6.89	98.31	111.40
1	D	186	ARG	NE-CZ-NH2	-6.44	117.08	120.30
1	A	202	ASP	CB-CG-OD1	6.11	123.80	118.30
1	A	73	ASP	CB-CG-OD1	5.59	123.33	118.30
1	C	60	MET	CB-CG-SD	-5.53	95.81	112.40
1	C	206	ARG	NE-CZ-NH2	-5.51	117.55	120.30
1	B	184	ASP	CB-CG-OD1	5.50	123.25	118.30
1	C	206	ARG	NE-CZ-NH1	5.49	123.04	120.30
1	B	203	ARG	NE-CZ-NH2	-5.47	117.56	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	1210	1225	1215	44	0
1	B	1196	1212	1201	39	0
1	C	1184	1200	1190	36	0
1	D	1163	1184	1172	34	0
2	A	44	0	0	2	0
2	B	22	0	0	1	0
2	C	27	0	0	1	0
2	D	20	0	0	1	0
All	All	4866	4821	4778	134	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 14.

All (134) 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:B:82:GLU:OE1	1:B:82:GLU:HA	1.69	0.89
1:D:126:MET:HE2	1:D:162:LEU:HD21	1.62	0.80
1:A:74:GLN:HG2	1:A:153:ALA:HA	1.65	0.78
1:B:128:ILE:HG23	1:B:129:HIS:CD2	2.21	0.76
1:A:126:MET:HE1	1:A:129:HIS:CD2	2.21	0.76
1:A:85:ARG:HD2	1:C:85:ARG:HD2	1.66	0.76
1:A:74:GLN:CG	1:A:153:ALA:HA	2.16	0.75
1:C:205:LEU:HD23	2:D:2013:HOH:O	1.89	0.72
1:A:85:ARG:HD2	1:C:85:ARG:CD	2.19	0.72
1:B:76:MET:HE3	1:B:86:LEU:HD12	1.72	0.71
1:D:89:THR:OG1	1:D:90:HIS:HD2	1.73	0.71
1:A:139:ARG:NH2	2:A:2025:HOH:O	2.26	0.67
1:D:126:MET:CE	1:D:162:LEU:HD21	2.24	0.67
1:D:64:GLU:HG3	1:D:70:TYR:CZ	2.28	0.67
1:A:210:GLU:O	1:A:211:ALA:HB2	1.98	0.64
1:A:89:THR:OG1	1:A:90:HIS:HD2	1.81	0.63
1:B:128:ILE:CG2	1:B:129:HIS:CD2	2.81	0.62
1:D:126:MET:HE2	1:D:162:LEU:CD2	2.30	0.62
1:D:204:ARG:HG2	1:D:209:LEU:HD22	1.83	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:191:ASN:HD22	1:A:191:ASN:C	2.03	0.61
1:C:89:THR:OG1	1:C:90:HIS:HD2	1.84	0.60
1:D:62:ALA:HA	1:D:72:GLY:HA2	1.83	0.60
1:C:191:ASN:HD22	1:C:191:ASN:C	2.04	0.60
1:A:122:PRO:HA	1:B:191:ASN:HD21	1.67	0.60
1:C:60:MET:HA	1:C:75:ASP:HB2	1.83	0.60
1:A:126:MET:CE	1:A:129:HIS:CD2	2.84	0.60
1:A:76:MET:SD	1:A:82:GLU:HG2	2.43	0.59
1:A:125:ASP:OD1	1:A:127:SER:HB2	2.04	0.58
1:A:66:TRP:CD1	1:A:67:PRO:HD2	2.38	0.58
1:A:85:ARG:HH11	1:C:85:ARG:HH11	1.49	0.58
1:C:60:MET:HB2	1:C:76:MET:CA	2.33	0.58
1:D:89:THR:OG1	1:D:90:HIS:CD2	2.54	0.58
1:A:210:GLU:O	1:A:211:ALA:CB	2.52	0.58
1:A:189:ILE:HD11	1:B:164:LEU:CD1	2.34	0.57
1:B:76:MET:CE	1:B:86:LEU:HD12	2.34	0.57
1:C:188:ILE:O	1:C:189:ILE:HD13	2.03	0.57
1:A:188:ILE:HG13	1:A:189:ILE:CD1	2.35	0.57
1:B:125:ASP:OD1	1:B:127:SER:CB	2.53	0.57
1:D:64:GLU:HG3	1:D:70:TYR:CE2	2.39	0.57
1:A:98:ARG:HG3	1:A:117:GLU:OE2	2.05	0.56
1:C:126:MET:H	1:D:197:GLN:HE22	1.54	0.56
1:B:98:ARG:HD2	1:B:117:GLU:OE2	2.06	0.56
1:C:74:GLN:HE21	1:C:153:ALA:HA	1.70	0.55
1:C:188:ILE:HD13	1:C:189:ILE:HG12	1.87	0.55
1:C:153:ALA:HB3	1:C:158:ARG:HH22	1.72	0.55
1:A:126:MET:CE	1:A:129:HIS:HB2	2.36	0.55
1:C:202:ASP:OD2	1:C:206:ARG:NH2	2.40	0.55
1:C:188:ILE:HD12	1:C:188:ILE:H	1.73	0.54
1:C:209:LEU:O	1:C:213:GLU:HG3	2.07	0.54
1:D:66:TRP:CZ3	1:D:181:LYS:HE3	2.42	0.54
1:A:96:HIS:HD2	1:A:113:TYR:OH	1.91	0.54
1:D:109:LEU:HB3	1:D:111:ILE:HG13	1.89	0.54
1:B:125:ASP:OD1	1:B:127:SER:HB3	2.07	0.54
1:D:139:ARG:O	1:D:142:SER:HB2	2.07	0.54
1:C:60:MET:HB2	1:C:76:MET:N	2.23	0.54
1:B:118:ALA:HA	1:B:124:PHE:CZ	2.42	0.53
1:C:126:MET:HE2	1:C:162:LEU:HD11	1.91	0.53
1:A:98:ARG:HB2	1:A:117:GLU:OE2	2.08	0.53
1:B:62:ALA:HA	1:B:72:GLY:HA2	1.90	0.52
1:C:66:TRP:CD1	1:C:67:PRO:HD2	2.45	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:189:ILE:HD11	1:B:164:LEU:HD11	1.91	0.52
1:B:105:ALA:O	1:B:109:LEU:HB2	2.10	0.51
1:A:74:GLN:HG3	1:A:153:ALA:HA	1.91	0.51
1:A:66:TRP:CG	1:A:67:PRO:HD2	2.46	0.51
1:A:126:MET:HE1	1:A:129:HIS:HD2	1.74	0.51
1:B:102:THR:HG23	1:B:103:PRO:HD2	1.93	0.51
1:A:188:ILE:HG13	1:A:189:ILE:HD12	1.93	0.50
1:B:76:MET:CE	1:B:86:LEU:CD1	2.89	0.50
1:B:188:ILE:HG13	1:B:189:ILE:HD12	1.94	0.50
1:C:153:ALA:HB3	1:C:158:ARG:NH2	2.27	0.49
1:D:66:TRP:HB3	1:D:69:LEU:HB3	1.95	0.49
1:A:117:GLU:HG3	2:A:2016:HOH:O	2.13	0.49
1:A:164:LEU:CD1	1:B:189:ILE:HD11	2.42	0.49
1:D:179:ILE:HG23	1:D:183:LYS:HE3	1.95	0.49
1:C:89:THR:OG1	1:C:90:HIS:CD2	2.65	0.49
1:A:60:MET:HG2	1:A:75:ASP:HB3	1.95	0.48
1:B:96:HIS:HD2	1:B:113:TYR:OH	1.96	0.48
1:C:131:GLN:OE1	1:D:211:ALA:O	2.31	0.48
1:B:180:LYS:HE3	1:B:184:ASP:OD2	2.14	0.48
1:A:126:MET:CE	1:A:129:HIS:HD2	2.25	0.47
1:A:128:ILE:HA	1:A:128:ILE:HD12	1.64	0.47
1:A:126:MET:HE2	1:A:162:LEU:HD21	1.96	0.47
1:C:74:GLN:NE2	2:C:2002:HOH:O	2.39	0.47
1:A:78:ASN:O	1:A:78:ASN:CG	2.52	0.47
1:C:126:MET:CE	1:C:162:LEU:HD21	2.45	0.46
1:C:189:ILE:HD11	1:D:164:LEU:CD1	2.45	0.46
1:D:73:ASP:O	1:D:76:MET:HB3	2.15	0.46
1:C:80:ARG:NH1	1:C:105:ALA:HB2	2.31	0.46
1:D:179:ILE:CG2	1:D:183:LYS:HE3	2.46	0.46
1:B:79:ASN:OD1	1:B:81:ARG:HG2	2.16	0.46
1:B:89:THR:OG1	1:B:90:HIS:HD2	1.99	0.46
1:A:85:ARG:HD2	1:C:85:ARG:HD3	1.93	0.46
1:B:114:LEU:HD12	1:B:115:GLY:H	1.81	0.45
1:A:164:LEU:HD11	1:B:189:ILE:CD1	2.46	0.45
1:B:116:VAL:HG13	1:B:129:HIS:ND1	2.32	0.45
1:A:164:LEU:CD1	1:B:189:ILE:CD1	2.94	0.45
1:A:90:HIS:HA	1:A:112:ARG:O	2.17	0.45
1:C:60:MET:CA	1:C:75:ASP:HB2	2.46	0.45
1:D:74:GLN:HE21	1:D:78:ASN:ND2	2.15	0.45
1:C:189:ILE:HD11	1:D:164:LEU:HD12	1.98	0.44
1:D:204:ARG:HG2	1:D:209:LEU:CD2	2.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:118:ALA:HA	1:C:124:PHE:CZ	2.52	0.44
1:D:171:HIS:C	1:D:172:HIS:CG	2.90	0.44
1:D:92:LEU:HD21	1:D:133:ALA:HB1	1.99	0.44
1:B:89:THR:C	1:B:90:HIS:CD2	2.91	0.44
1:D:162:LEU:HD23	1:D:162:LEU:HA	1.83	0.44
1:D:126:MET:HE2	1:D:162:LEU:CG	2.49	0.43
1:B:83:LEU:HD23	1:B:83:LEU:HA	1.81	0.43
1:D:126:MET:HG3	1:D:126:MET:O	2.18	0.43
1:A:92:LEU:HD21	1:A:133:ALA:HB1	2.01	0.42
1:A:186:ARG:O	1:B:183:LYS:HE2	2.19	0.42
1:B:191:ASN:C	1:B:191:ASN:HD22	2.22	0.42
1:D:89:THR:HB	1:D:112:ARG:NH2	2.33	0.42
1:A:191:ASN:ND2	1:A:191:ASN:C	2.72	0.42
1:C:212:LEU:HD23	1:C:212:LEU:HA	1.79	0.42
1:A:98:ARG:CB	1:A:117:GLU:OE2	2.67	0.42
1:D:188:ILE:O	1:D:189:ILE:HG13	2.19	0.42
1:C:60:MET:HB2	1:C:76:MET:HA	1.99	0.42
1:B:97:SER:O	1:B:98:ARG:C	2.58	0.42
1:B:171:HIS:C	1:B:172:HIS:CG	2.93	0.42
1:C:188:ILE:C	1:C:189:ILE:HD13	2.41	0.41
1:B:137:ILE:HG22	1:B:166:TYR:CE1	2.56	0.41
1:C:188:ILE:HD11	1:D:179:ILE:HG23	2.03	0.41
1:C:164:LEU:HD12	1:D:189:ILE:CD1	2.51	0.41
1:B:128:ILE:HD12	1:B:128:ILE:O	2.21	0.41
1:A:164:LEU:HD12	1:B:189:ILE:HD11	2.02	0.40
1:B:195:LEU:HA	1:B:195:LEU:HD12	1.99	0.40
1:A:164:LEU:HD11	1:B:189:ILE:HD11	2.02	0.40
1:B:89:THR:HB	2:B:2003:HOH:O	2.21	0.40
1:B:90:HIS:CD2	1:B:90:HIS:N	2.89	0.40
1:C:77:ALA:HB2	1:C:151:HIS:CD2	2.57	0.40
1:D:126:MET:HE1	1:D:162:LEU:HD11	2.03	0.40
1:D:169:LEU:HD23	1:D:169:LEU:HA	1.92	0.40
1:D:80:ARG:HD3	1:D:105:ALA:CB	2.52	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	150/160 (94%)	148 (99%)	2 (1%)	0	100	100
1	B	148/160 (92%)	135 (91%)	11 (7%)	2 (1%)	11	8
1	C	146/160 (91%)	141 (97%)	5 (3%)	0	100	100
1	D	141/160 (88%)	134 (95%)	7 (5%)	0	100	100
All	All	585/640 (91%)	558 (95%)	25 (4%)	2 (0%)	41	46

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	99	TRP
1	B	155	GLY

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	124/132 (94%)	116 (94%)	8 (6%)	17	19
1	B	123/132 (93%)	109 (89%)	14 (11%)	5	5
1	C	122/132 (92%)	109 (89%)	13 (11%)	6	6
1	D	119/132 (90%)	109 (92%)	10 (8%)	11	11
All	All	488/528 (92%)	443 (91%)	45 (9%)	9	9

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	81	ARG
1	A	92	LEU
1	A	98	ARG
1	A	127	SER
1	A	128	ILE
1	A	154	VAL
1	A	191	ASN
1	A	209	LEU
1	B	82	GLU
1	B	84	ARG
1	B	98	ARG
1	B	109	LEU
1	B	121	SER
1	B	128	ILE
1	B	143	GLN
1	B	154	VAL
1	B	156	VAL
1	B	172	HIS
1	B	180	LYS
1	B	191	ASN
1	B	206	ARG
1	B	209	LEU
1	C	60	MET
1	C	75	ASP
1	C	81	ARG
1	C	84	ARG
1	C	85	ARG
1	C	121	SER
1	C	154	VAL
1	C	156	VAL
1	C	186	ARG
1	C	188	ILE
1	C	191	ASN
1	C	195	LEU
1	C	196	ARG
1	D	64	GLU
1	D	69	LEU
1	D	92	LEU
1	D	116	VAL
1	D	121	SER
1	D	156	VAL
1	D	172	HIS
1	D	188	ILE

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Mol	Chain	Res	Type
1	D	204	ARG
1	D	206	ARG

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	74	GLN
1	A	90	HIS
1	A	96	HIS
1	A	191	ASN
1	B	74	GLN
1	B	90	HIS
1	B	93	ASN
1	B	96	HIS
1	B	143	GLN
1	B	185	HIS
1	B	191	ASN
1	C	74	GLN
1	C	90	HIS
1	C	96	HIS
1	C	191	ASN
1	C	207	GLN
1	D	74	GLN
1	D	78	ASN
1	D	90	HIS
1	D	197	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

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, 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	152/160 (95%)	-0.16	0 100 100	16, 29, 55, 71	0
1	B	150/160 (93%)	0.10	6 (4%) 38 36	18, 37, 74, 101	0
1	C	150/160 (93%)	-0.06	0 100 100	17, 35, 62, 82	0
1	D	147/160 (91%)	0.03	2 (1%) 75 73	22, 42, 61, 83	0
All	All	599/640 (93%)	-0.02	8 (1%) 77 75	16, 36, 65, 101	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	104	GLU	4.4
1	B	99	TRP	4.3
1	B	100	ARG	3.9
1	B	101	GLY	3.1
1	D	211	ALA	2.7
1	D	110	GLY	2.6
1	B	98	ARG	2.0
1	B	102	THR	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.