Full wwPDB X-ray Structure Validation Report ⓘ

Oct 30, 2023 – 01:16 PM JST

PDB ID	:	5B02
Title	:	Structure of the prenyltransferase MoeN5 with a fusion protein tag of Sso7d
Authors	:	Ko, T.-P.; Zhang, L.; Chen, C.-C.; Guo, R.-T.; Oldfield, E.O.
Deposited on	:	2015-10-27
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 types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

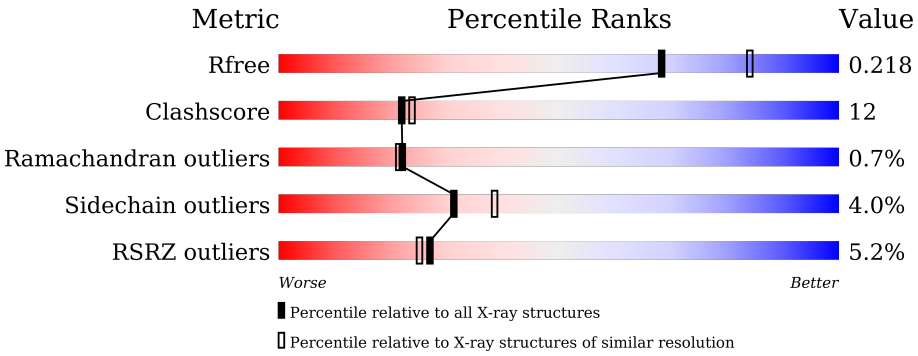
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation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

1 Overall quality at a glance (i)

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	5912 (2.24-2.20)
Clashscore	141614	6646 (2.24-2.20)
Ramachandran outliers	138981	6543 (2.24-2.20)
Sidechain outliers	138945	6544 (2.24-2.20)
RSRZ outliers	127900	5797 (2.24-2.20)

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 of 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 <=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	343	2% 63% 11% 23%
1	B	343	4% 80% 15%
1	C	343	% 59% 15% 24%
1	D	343	11% 75% 20%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 9942 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 MoeN5,DNA-binding protein 7d.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	264	Total	C	N	O	S	0	0	0
			2013	1243	374	385	11			
1	B	332	Total	C	N	O	S	0	0	0
			2533	1571	463	485	14			
1	C	262	Total	C	N	O	S	0	0	0
			1998	1234	372	381	11			
1	D	333	Total	C	N	O	S	0	0	0
			2544	1579	466	485	14			

There are 76 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-13	MET	-	expression tag	UNP A0A010
A	-12	ALA	-	expression tag	UNP A0A010
A	-11	HIS	-	expression tag	UNP A0A010
A	-10	HIS	-	expression tag	UNP A0A010
A	-9	HIS	-	expression tag	UNP A0A010
A	-8	HIS	-	expression tag	UNP A0A010
A	-7	HIS	-	expression tag	UNP A0A010
A	-6	HIS	-	expression tag	UNP A0A010
A	-5	VAL	-	expression tag	UNP A0A010
A	-4	ASP	-	expression tag	UNP A0A010
A	-3	ASP	-	expression tag	UNP A0A010
A	-2	ASP	-	expression tag	UNP A0A010
A	-1	ASP	-	expression tag	UNP A0A010
A	0	LYS	-	expression tag	UNP A0A010
A	261	ALA	-	linker	UNP A0A010
A	262	GLY	-	linker	UNP A0A010
A	263	ALA	-	linker	UNP A0A010
A	264	GLY	-	linker	UNP A0A010
A	265	ALA	-	linker	UNP A0A010
B	-13	MET	-	expression tag	UNP A0A010
B	-12	ALA	-	expression tag	UNP A0A010

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-11	HIS	-	expression tag	UNP A0A010
B	-10	HIS	-	expression tag	UNP A0A010
B	-9	HIS	-	expression tag	UNP A0A010
B	-8	HIS	-	expression tag	UNP A0A010
B	-7	HIS	-	expression tag	UNP A0A010
B	-6	HIS	-	expression tag	UNP A0A010
B	-5	VAL	-	expression tag	UNP A0A010
B	-4	ASP	-	expression tag	UNP A0A010
B	-3	ASP	-	expression tag	UNP A0A010
B	-2	ASP	-	expression tag	UNP A0A010
B	-1	ASP	-	expression tag	UNP A0A010
B	0	LYS	-	expression tag	UNP A0A010
B	261	ALA	-	linker	UNP A0A010
B	262	GLY	-	linker	UNP A0A010
B	263	ALA	-	linker	UNP A0A010
B	264	GLY	-	linker	UNP A0A010
B	265	ALA	-	linker	UNP A0A010
C	-13	MET	-	expression tag	UNP A0A010
C	-12	ALA	-	expression tag	UNP A0A010
C	-11	HIS	-	expression tag	UNP A0A010
C	-10	HIS	-	expression tag	UNP A0A010
C	-9	HIS	-	expression tag	UNP A0A010
C	-8	HIS	-	expression tag	UNP A0A010
C	-7	HIS	-	expression tag	UNP A0A010
C	-6	HIS	-	expression tag	UNP A0A010
C	-5	VAL	-	expression tag	UNP A0A010
C	-4	ASP	-	expression tag	UNP A0A010
C	-3	ASP	-	expression tag	UNP A0A010
C	-2	ASP	-	expression tag	UNP A0A010
C	-1	ASP	-	expression tag	UNP A0A010
C	0	LYS	-	expression tag	UNP A0A010
C	261	ALA	-	linker	UNP A0A010
C	262	GLY	-	linker	UNP A0A010
C	263	ALA	-	linker	UNP A0A010
C	264	GLY	-	linker	UNP A0A010
C	265	ALA	-	linker	UNP A0A010
D	-13	MET	-	expression tag	UNP A0A010
D	-12	ALA	-	expression tag	UNP A0A010
D	-11	HIS	-	expression tag	UNP A0A010
D	-10	HIS	-	expression tag	UNP A0A010
D	-9	HIS	-	expression tag	UNP A0A010
D	-8	HIS	-	expression tag	UNP A0A010

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-7	HIS	-	expression tag	UNP A0A010
D	-6	HIS	-	expression tag	UNP A0A010
D	-5	VAL	-	expression tag	UNP A0A010
D	-4	ASP	-	expression tag	UNP A0A010
D	-3	ASP	-	expression tag	UNP A0A010
D	-2	ASP	-	expression tag	UNP A0A010
D	-1	ASP	-	expression tag	UNP A0A010
D	0	LYS	-	expression tag	UNP A0A010
D	261	ALA	-	linker	UNP A0A010
D	262	GLY	-	linker	UNP A0A010
D	263	ALA	-	linker	UNP A0A010
D	264	GLY	-	linker	UNP A0A010
D	265	ALA	-	linker	UNP A0A010

- Molecule 2 is water.

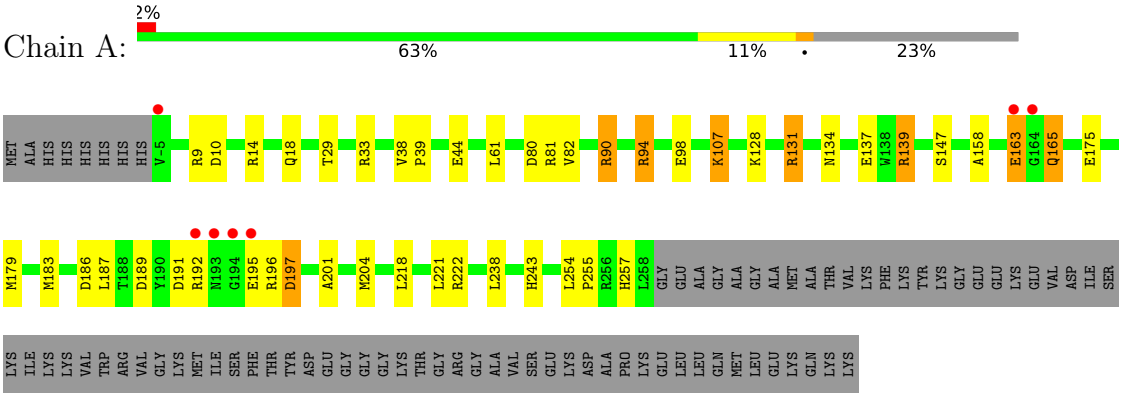
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	229	Total O 229 229	0	0
2	B	235	Total O 235 235	0	0
2	C	176	Total O 176 176	0	0
2	D	214	Total O 214 214	0	0

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.

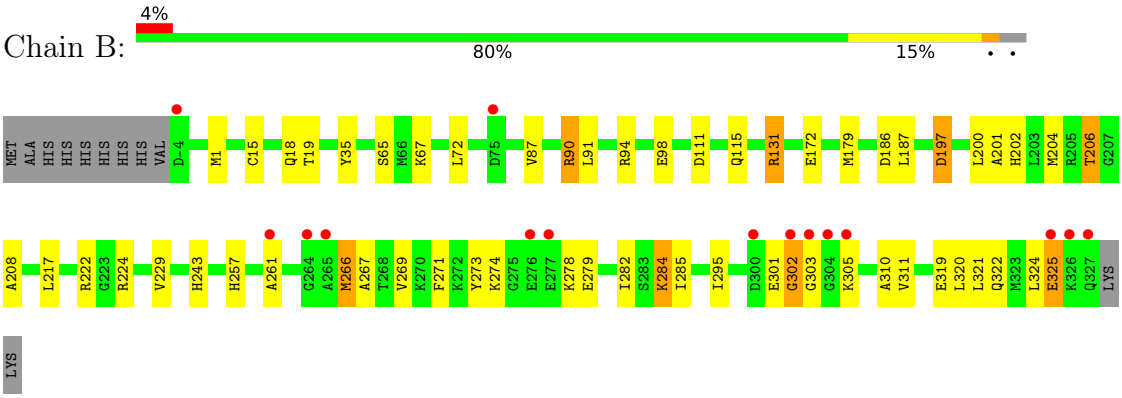
• Molecule 1: MoeN5,DNA-binding protein 7d

Chain A:



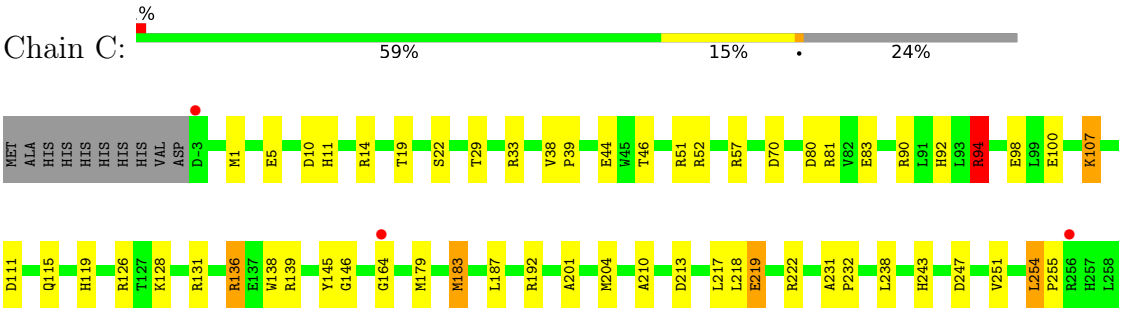
• Molecule 1: MoeN5,DNA-binding protein 7d

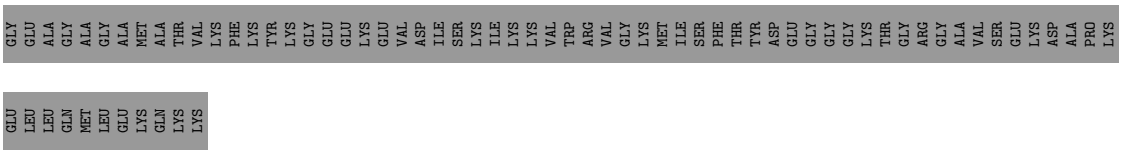
Chain B:



• Molecule 1: MoeN5,DNA-binding protein 7d

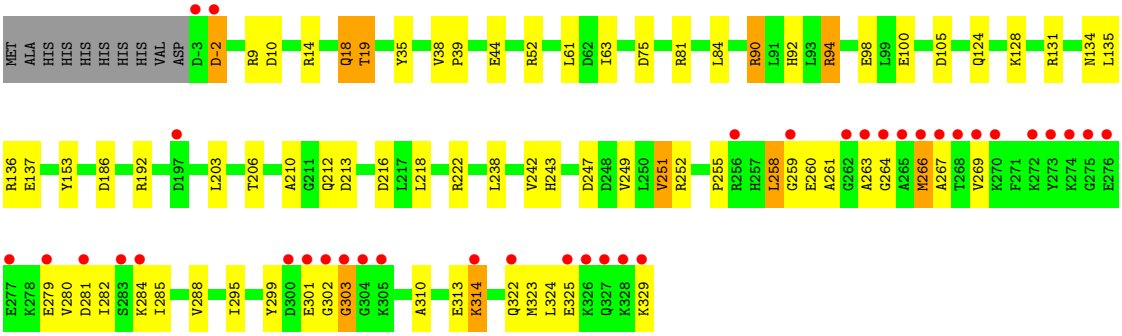
Chain C:





• Molecule 1: MoeN5,DNA-binding protein 7d

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 2 2 2	Depositor
Cell constants a, b, c, α, β, γ	137.88Å 217.44Å 104.64Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.00 – 2.21 24.98 – 2.21	Depositor EDS
% Data completeness (in resolution range)	96.1 (25.00-2.21) 96.0 (24.98-2.21)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.62 (at 2.22Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.176 , 0.218 0.176 , 0.218	Depositor DCC
R_{free} test set	3842 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å^2)	25.6	Xtriage
Anisotropy	0.222	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 43.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	9942	wwPDB-VP
Average B, all atoms (Å^2)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.57% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 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	0.78	0/2045	0.84	2/2780 (0.1%)
1	B	0.80	2/2572 (0.1%)	0.86	2/3478 (0.1%)
1	C	0.86	2/2030 (0.1%)	0.92	4/2759 (0.1%)
1	D	0.72	0/2583	0.83	5/3489 (0.1%)
All	All	0.79	4/9230 (0.0%)	0.86	13/12506 (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
1	D	0	1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	219	GLU	CG-CD	6.77	1.62	1.51
1	B	15	CYS	CB-SG	-6.36	1.71	1.82
1	C	183	MET	SD-CE	5.59	2.09	1.77
1	B	172	GLU	CG-CD	5.36	1.59	1.51

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	94	ARG	NE-CZ-NH2	-9.68	115.46	120.30
1	C	94	ARG	NE-CZ-NH1	8.42	124.51	120.30
1	D	52	ARG	NE-CZ-NH1	-7.10	116.75	120.30
1	D	94	ARG	NE-CZ-NH1	7.05	123.83	120.30
1	B	224	ARG	NE-CZ-NH1	6.64	123.62	120.30
1	A	94	ARG	NE-CZ-NH1	5.99	123.29	120.30
1	D	258	LEU	N-CA-C	-5.92	95.03	111.00
1	A	107	LYS	CD-CE-NZ	-5.58	98.87	111.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	72	LEU	CA-CB-CG	5.45	127.84	115.30
1	C	94	ARG	CD-NE-CZ	5.41	131.17	123.60
1	D	-2	ASP	N-CA-C	-5.39	96.43	111.00
1	C	94	ARG	CG-CD-NE	-5.19	100.91	111.80
1	D	9	ARG	NE-CZ-NH2	5.06	122.83	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	153	TYR	Sidechain

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	2013	0	1996	41	0
1	B	2533	0	2532	53	0
1	C	1998	0	1983	61	0
1	D	2544	0	2554	68	0
2	A	229	0	0	4	0
2	B	235	0	0	3	0
2	C	176	0	0	6	0
2	D	214	0	0	1	0
All	All	9942	0	9065	215	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.

All (215) 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:1:MET:CE	1:B:1:MET:SD	2.01	1.48
1:C:183:MET:CE	1:C:183:MET:SD	2.09	1.40
1:D:131:ARG:HG2	1:D:131:ARG:HH11	1.36	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:222:ARG:HH11	1:D:243:HIS:HD2	1.27	0.83
1:D:267:ALA:HB1	1:D:282:ILE:HB	1.61	0.83
1:B:273:TYR:CE2	1:B:274:LYS:HD2	2.16	0.81
1:D:325:GLU:HA	1:D:329:LYS:HB3	1.65	0.79
1:D:10:ASP:HB2	2:D:521:HOH:O	1.85	0.77
1:B:222:ARG:HH11	1:B:243:HIS:HD2	1.33	0.76
1:B:284:LYS:HA	1:B:284:LYS:HE3	1.71	0.73
1:C:22:SER:N	1:C:83:GLU:OE1	2.21	0.73
1:A:61:LEU:HD13	1:A:94:ARG:HG3	1.71	0.72
1:C:145:TYR:HD1	1:C:146:GLY:H	1.37	0.72
1:C:131:ARG:HG2	1:C:131:ARG:HH11	1.55	0.71
1:D:10:ASP:O	1:D:14:ARG:HG3	1.89	0.71
1:D:94:ARG:HD2	1:D:98:GLU:OE2	1.90	0.71
1:A:192:ARG:HD2	1:C:80:ASP:HB2	1.72	0.71
1:C:145:TYR:CD1	1:C:146:GLY:N	2.58	0.70
1:D:325:GLU:HB2	1:D:329:LYS:HD3	1.73	0.70
1:D:247:ASP:O	1:D:251:VAL:HG13	1.92	0.69
1:D:19:THR:HG21	1:D:90:ARG:HG3	1.73	0.69
1:D:251:VAL:O	1:D:255:PRO:HG2	1.93	0.69
1:B:179:MET:CE	1:B:217:LEU:HD21	2.22	0.69
1:A:189:ASP:HB3	1:A:195:GLU:OE1	1.92	0.68
1:D:131:ARG:HG2	1:D:131:ARG:NH1	2.09	0.68
1:D:75:ASP:HB2	1:D:81:ARG:NH1	2.09	0.67
1:C:94:ARG:HD2	1:C:98:GLU:OE2	1.95	0.66
1:B:187:LEU:HD21	1:B:204:MET:CE	2.25	0.66
1:D:61:LEU:HD13	1:D:94:ARG:HG3	1.77	0.66
1:D:19:THR:CG2	1:D:90:ARG:HG3	2.26	0.66
1:A:195:GLU:O	1:A:196:ARG:HD2	1.97	0.65
1:B:295:ILE:HD12	1:B:324:LEU:HD11	1.79	0.65
1:A:94:ARG:HD2	1:A:98:GLU:OE2	1.97	0.65
1:C:183:MET:CE	1:C:183:MET:HB3	2.28	0.64
1:C:187:LEU:HD21	1:C:204:MET:HE1	1.80	0.64
1:C:136:ARG:HH11	1:C:136:ARG:HB3	1.64	0.63
1:B:111:ASP:O	1:B:115:GLN:HG2	1.98	0.62
1:C:201:ALA:HA	1:C:204:MET:HE3	1.80	0.62
1:C:222:ARG:NH1	1:C:243:HIS:HD2	1.98	0.61
1:D:218:LEU:HD11	1:D:249:VAL:CG1	2.30	0.61
1:C:145:TYR:HB2	2:C:532:HOH:O	1.99	0.61
1:C:179:MET:CE	1:C:217:LEU:HD11	2.29	0.61
1:C:1:MET:HE2	2:C:504:HOH:O	2.00	0.61
1:C:247:ASP:O	1:C:251:VAL:HG12	2.00	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:134:ASN:HD21	1:D:136:ARG:NH2	1.99	0.61
1:C:81:ARG:HD3	1:D:81:ARG:NH2	2.16	0.60
1:D:75:ASP:HB2	1:D:81:ARG:HH12	1.64	0.60
1:C:145:TYR:HD1	1:C:146:GLY:N	1.98	0.59
1:A:187:LEU:HD21	1:A:204:MET:CE	2.32	0.59
1:A:38:VAL:HG12	1:A:39:PRO:HD3	1.84	0.59
1:D:222:ARG:NH1	1:D:243:HIS:HD2	1.96	0.59
1:B:200:LEU:HG	1:B:204:MET:HE2	1.85	0.59
1:B:187:LEU:HD21	1:B:204:MET:HE1	1.85	0.58
1:C:183:MET:HB3	1:C:183:MET:HE3	1.84	0.58
1:A:179:MET:HE2	1:A:221:LEU:HD11	1.85	0.58
1:D:218:LEU:HD11	1:D:249:VAL:HG12	1.84	0.58
1:D:281:ASP:HB3	1:D:284:LYS:HG3	1.86	0.58
1:D:255:PRO:O	1:D:258:LEU:O	2.22	0.58
1:C:5:GLU:OE2	1:C:33:ARG:HG2	2.03	0.57
1:C:219:GLU:OE1	1:C:222:ARG:NE	2.38	0.57
1:C:107:LYS:HD3	1:C:107:LYS:O	2.05	0.57
1:D:192:ARG:HH11	1:D:192:ARG:HG2	1.70	0.57
1:B:200:LEU:HG	1:B:204:MET:CE	2.35	0.57
1:A:107:LYS:O	1:A:107:LYS:HD3	2.05	0.56
1:B:301:GLU:HB3	1:B:305:LYS:HD3	1.86	0.56
1:D:124:GLN:HE21	1:D:128:LYS:HE2	1.71	0.56
1:B:35:TYR:CE2	1:B:67:LYS:HG3	2.41	0.55
1:B:301:GLU:OE2	1:B:305:LYS:HE3	2.05	0.55
1:A:201:ALA:HB1	1:A:257:HIS:CE1	2.42	0.54
1:B:301:GLU:CD	1:B:305:LYS:HE3	2.28	0.54
1:C:111:ASP:O	1:C:115:GLN:HG2	2.08	0.54
1:C:219:GLU:CD	1:C:222:ARG:HE	2.09	0.54
1:C:222:ARG:HH11	1:C:243:HIS:CD2	2.26	0.54
1:C:1:MET:C	1:C:1:MET:SD	2.86	0.54
1:C:201:ALA:HA	1:C:204:MET:CE	2.38	0.53
1:B:266:MET:HE2	1:B:267:ALA:N	2.23	0.53
1:B:65:SER:HB2	1:B:91:LEU:HB2	1.91	0.53
1:B:243:HIS:HE1	2:B:559:HOH:O	1.92	0.53
1:C:187:LEU:CD2	1:C:204:MET:HE1	2.38	0.53
1:D:222:ARG:HH11	1:D:243:HIS:CD2	2.17	0.53
1:D:269:VAL:O	1:D:279:GLU:HA	2.08	0.53
1:A:94:ARG:HD3	1:A:98:GLU:HG3	1.91	0.53
1:C:44:GLU:HG2	1:C:238:LEU:HG	1.91	0.53
1:D:18:GLN:O	1:D:18:GLN:HG2	2.08	0.53
1:C:38:VAL:HG12	1:C:39:PRO:HD3	1.90	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:319:GLU:CD	1:B:319:GLU:H	2.13	0.52
1:D:238:LEU:O	1:D:242:VAL:HG23	2.10	0.52
1:D:280:VAL:HG21	1:D:299:TYR:CZ	2.45	0.52
1:D:288:VAL:HG12	1:D:324:LEU:HD13	1.92	0.52
1:C:183:MET:CE	1:C:183:MET:CB	2.88	0.52
1:A:90:ARG:C	1:A:90:ARG:HD3	2.30	0.52
1:A:9:ARG:HD2	2:A:468:HOH:O	2.10	0.52
1:B:187:LEU:HD21	1:B:204:MET:HE3	1.90	0.52
1:B:295:ILE:CD1	1:B:324:LEU:HD11	2.39	0.52
1:D:313:GLU:HG3	1:D:314:LYS:N	2.25	0.51
1:C:210:ALA:HB3	1:C:213:ASP:OD2	2.10	0.51
1:D:281:ASP:HB3	1:D:284:LYS:CG	2.40	0.51
1:D:61:LEU:CD1	1:D:94:ARG:HG3	2.41	0.51
1:D:251:VAL:CG2	1:D:252:ARG:N	2.74	0.51
1:A:187:LEU:HD21	1:A:204:MET:HE1	1.92	0.51
1:D:267:ALA:HB1	1:D:282:ILE:CB	2.38	0.50
1:B:98:GLU:OE1	2:B:401:HOH:O	2.19	0.50
1:A:192:ARG:HD2	1:C:80:ASP:CB	2.41	0.50
1:C:10:ASP:OD2	1:C:14:ARG:NH1	2.44	0.50
1:A:38:VAL:CG1	1:A:39:PRO:HD3	2.42	0.50
1:D:75:ASP:CB	1:D:81:ARG:NH1	2.75	0.50
1:C:11:HIS:CE1	1:C:57:ARG:HD2	2.47	0.49
1:D:325:GLU:CB	1:D:329:LYS:HD3	2.41	0.49
1:B:273:TYR:HB3	1:B:278:LYS:HD2	1.94	0.49
1:B:303:GLY:C	1:B:305:LYS:H	2.14	0.49
1:B:322:GLN:O	1:B:325:GLU:HG3	2.12	0.49
1:B:269:VAL:HG21	1:B:320:LEU:HD22	1.95	0.49
1:D:203:LEU:O	1:D:206:THR:HG22	2.12	0.48
1:D:285:ILE:HG22	1:D:323:MET:HE2	1.95	0.48
1:B:19:THR:O	1:B:19:THR:HG22	2.14	0.48
1:B:201:ALA:HB1	1:B:257:HIS:CE1	2.49	0.48
1:C:179:MET:HE2	1:C:217:LEU:HD11	1.96	0.48
1:A:131:ARG:HG3	1:A:197:ASP:HB3	1.96	0.47
1:C:14:ARG:HG3	2:C:486:HOH:O	2.13	0.47
1:A:222:ARG:NH1	1:A:243:HIS:ND1	2.61	0.47
1:B:285:ILE:CD1	1:B:320:LEU:HD13	2.44	0.47
1:C:139:ARG:HG3	1:C:179:MET:HE1	1.96	0.47
1:C:126:ARG:NH2	2:C:402:HOH:O	2.31	0.47
1:D:252:ARG:O	1:D:255:PRO:HD2	2.14	0.47
1:B:295:ILE:O	1:B:310:ALA:HA	2.15	0.47
1:D:259:GLY:O	1:D:260:GLU:HG2	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:202:HIS:O	1:B:206:THR:HB	2.15	0.47
1:C:100:GLU:OE1	1:D:100:GLU:OE1	2.33	0.47
1:D:90:ARG:C	1:D:90:ARG:HD3	2.36	0.47
1:D:210:ALA:O	1:D:213:ASP:HB2	2.14	0.47
1:B:90:ARG:HD3	1:B:90:ARG:C	2.35	0.46
1:B:179:MET:HE1	1:B:217:LEU:HD11	1.97	0.46
1:C:131:ARG:HH11	1:C:131:ARG:CG	2.25	0.46
1:C:136:ARG:HH11	1:C:136:ARG:CG	2.29	0.46
1:C:119:HIS:CD2	2:C:402:HOH:O	2.69	0.45
1:C:46:THR:HG21	1:C:51:ARG:HG2	1.99	0.45
1:C:138:TRP:CH2	1:C:183:MET:HG2	2.51	0.45
1:A:183:MET:HE2	1:A:218:LEU:HD21	1.98	0.45
1:A:147:SER:O	1:A:175:GLU:HG2	2.17	0.45
1:C:222:ARG:NH1	1:C:243:HIS:CD2	2.80	0.45
1:D:267:ALA:C	1:D:282:ILE:HG12	2.37	0.45
1:A:134:ASN:OD1	1:A:137:GLU:HG3	2.17	0.45
1:B:187:LEU:HD23	1:B:201:ALA:HB2	1.98	0.45
1:D:38:VAL:HG12	1:D:39:PRO:HD3	1.97	0.45
1:D:259:GLY:HA2	1:D:322:GLN:HG3	1.99	0.45
1:D:325:GLU:CA	1:D:329:LYS:HB3	2.42	0.45
1:A:187:LEU:CD2	1:A:204:MET:CE	2.95	0.45
1:D:44:GLU:O	1:D:44:GLU:HG3	2.17	0.44
1:D:295:ILE:O	1:D:310:ALA:HA	2.17	0.44
1:A:139:ARG:HD3	1:A:179:MET:CE	2.48	0.44
1:B:321:LEU:O	1:B:324:LEU:HB2	2.17	0.44
1:A:61:LEU:CD1	1:A:94:ARG:HG3	2.44	0.44
1:A:192:ARG:HA	2:A:494:HOH:O	2.18	0.44
1:C:136:ARG:HB2	2:C:448:HOH:O	2.17	0.44
1:C:183:MET:CE	1:C:218:LEU:HG	2.47	0.44
1:D:218:LEU:HD11	1:D:249:VAL:HG11	1.99	0.44
1:C:136:ARG:HH11	1:C:136:ARG:CB	2.29	0.44
1:A:10:ASP:OD2	1:C:192:ARG:HD2	2.18	0.44
1:A:187:LEU:CD2	1:A:204:MET:HE3	2.48	0.44
1:B:267:ALA:O	1:B:282:ILE:HG12	2.18	0.44
1:A:44:GLU:HG2	1:A:238:LEU:HG	1.99	0.43
1:A:187:LEU:HD22	1:A:204:MET:HE3	1.99	0.43
1:B:302:GLY:H	1:B:305:LYS:HB3	1.82	0.43
1:A:14:ARG:HD2	2:A:543:HOH:O	2.17	0.43
1:C:81:ARG:HD3	1:D:81:ARG:HH22	1.82	0.43
1:D:280:VAL:HG11	1:D:299:TYR:CD2	2.53	0.43
1:C:19:THR:HG22	1:C:19:THR:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:87:VAL:O	1:B:91:LEU:HG	2.18	0.43
1:B:90:ARG:NH1	1:B:94:ARG:HB2	2.34	0.43
1:B:269:VAL:O	1:B:279:GLU:HA	2.18	0.43
1:D:258:LEU:HD12	1:D:258:LEU:HA	1.81	0.43
1:D:301:GLU:O	1:D:303:GLY:N	2.51	0.43
1:B:206:THR:HG22	1:B:208:ALA:H	1.83	0.43
1:B:266:MET:CE	1:B:267:ALA:N	2.82	0.43
1:B:115:GLN:HG3	2:B:581:HOH:O	2.18	0.43
1:A:165:GLN:HE21	1:A:165:GLN:HB2	1.60	0.43
1:C:254:LEU:N	1:C:255:PRO:CD	2.82	0.43
1:B:271:PHE:HB3	1:B:311:VAL:CG1	2.49	0.43
1:A:10:ASP:CG	1:C:192:ARG:HD2	2.39	0.42
1:A:80:ASP:OD1	1:A:82:VAL:HB	2.19	0.42
1:B:131:ARG:HA	1:B:131:ARG:HD2	1.83	0.42
1:C:70:ASP:OD2	1:C:128:LYS:HE3	2.19	0.42
1:A:33:ARG:HG2	1:A:33:ARG:NH1	2.34	0.42
1:A:81:ARG:HG2	1:A:81:ARG:HH11	1.85	0.42
1:D:134:ASN:OD1	1:D:137:GLU:HG3	2.19	0.42
1:D:266:MET:H	1:D:266:MET:HG3	1.64	0.42
1:D:285:ILE:HG22	1:D:323:MET:CE	2.49	0.42
1:C:231:ALA:HA	1:C:232:PRO:HD3	1.85	0.42
1:B:271:PHE:HB3	1:B:311:VAL:HG12	2.01	0.42
1:C:183:MET:CE	1:C:183:MET:CG	2.94	0.42
1:D:131:ARG:NH1	1:D:131:ARG:CG	2.76	0.42
1:A:158:ALA:CB	1:A:165:GLN:HG2	2.49	0.42
1:B:303:GLY:C	1:B:305:LYS:N	2.74	0.42
1:B:131:ARG:HG3	1:B:197:ASP:HB3	2.02	0.41
1:C:131:ARG:HG2	1:C:131:ARG:NH1	2.31	0.41
1:D:35:TYR:HB3	1:D:63:ILE:HG22	2.02	0.41
1:B:285:ILE:O	1:B:285:ILE:HG22	2.19	0.41
1:C:92:HIS:NE2	1:D:92:HIS:NE2	2.63	0.41
1:A:131:ARG:CG	1:A:131:ARG:HH11	2.33	0.41
1:A:195:GLU:C	1:A:196:ARG:HD2	2.41	0.41
1:C:38:VAL:N	1:C:39:PRO:CD	2.84	0.41
1:D:124:GLN:NE2	1:D:128:LYS:HE2	2.34	0.41
1:A:9:ARG:HG2	2:A:443:HOH:O	2.20	0.41
1:B:179:MET:HE2	1:B:217:LEU:HD21	2.01	0.41
1:B:266:MET:HA	1:B:266:MET:HE3	2.01	0.41
1:D:135:LEU:CD2	1:D:213:ASP:HB3	2.51	0.41
1:D:288:VAL:HG12	1:D:324:LEU:CD1	2.50	0.41
1:A:163:GLU:H	1:A:163:GLU:HG2	1.79	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1:MET:CE	1:B:1:MET:HB2	2.52	0.40
1:B:301:GLU:HB3	1:B:305:LYS:CD	2.50	0.40
1:C:131:ARG:CG	1:C:131:ARG:NH1	2.84	0.40
1:A:254:LEU:HB2	1:A:255:PRO:HD3	2.02	0.40
1:D:212:GLN:NE2	1:D:216:ASP:OD1	2.48	0.40
1:D:251:VAL:HG22	1:D:252:ARG:N	2.36	0.40
1:D:84:LEU:HD23	1:D:84:LEU:HA	1.94	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	262/343 (76%)	256 (98%)	6 (2%)	0	100 100
1	B	330/343 (96%)	311 (94%)	17 (5%)	2 (1%)	25 25
1	C	260/343 (76%)	252 (97%)	7 (3%)	1 (0%)	34 37
1	D	331/343 (96%)	310 (94%)	16 (5%)	5 (2%)	10 7
All	All	1183/1372 (86%)	1129 (95%)	46 (4%)	8 (1%)	22 21

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	261	ALA
1	C	164	GLY
1	D	302	GLY
1	B	261	ALA
1	D	263	ALA
1	D	303	GLY
1	B	302	GLY
1	D	264	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	208/270 (77%)	197 (95%)	11 (5%)	22	26
1	B	260/270 (96%)	250 (96%)	10 (4%)	33	41
1	C	206/270 (76%)	199 (97%)	7 (3%)	37	46
1	D	261/270 (97%)	252 (97%)	9 (3%)	37	46
All	All	935/1080 (87%)	898 (96%)	37 (4%)	31	38

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	GLN
1	A	29	THR
1	A	90	ARG
1	A	128	LYS
1	A	131	ARG
1	A	139	ARG
1	A	163	GLU
1	A	165	GLN
1	A	186	ASP
1	A	191	ASP
1	A	197	ASP
1	B	18	GLN
1	B	90	ARG
1	B	131	ARG
1	B	186	ASP
1	B	197	ASP
1	B	206	THR
1	B	229	VAL
1	B	266	MET
1	B	284	LYS
1	B	325	GLU
1	C	29	THR
1	C	52	ARG
1	C	90	ARG

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Mol	Chain	Res	Type
1	C	94	ARG
1	C	107	LYS
1	C	136	ARG
1	C	254	LEU
1	D	-2	ASP
1	D	18	GLN
1	D	19	THR
1	D	90	ARG
1	D	105	ASP
1	D	186	ASP
1	D	251	VAL
1	D	266	MET
1	D	314	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	18	GLN
1	B	243	HIS
1	C	119	HIS
1	C	165	GLN
1	C	243	HIS
1	D	124	GLN
1	D	165	GLN
1	D	243	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 monosaccharides 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

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^{th} 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	264/343 (76%)	-0.47	7 (2%) 54 52	17, 26, 52, 81	0
1	B	332/343 (96%)	-0.30	15 (4%) 33 31	16, 28, 66, 84	0
1	C	262/343 (76%)	-0.51	3 (1%) 80 79	15, 29, 45, 73	0
1	D	333/343 (97%)	0.04	37 (11%) 5 4	15, 30, 100, 120	0
All	All	1191/1372 (86%)	-0.29	62 (5%) 27 25	15, 28, 78, 120	0

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	266	MET	9.5
1	D	264	GLY	9.2
1	D	303	GLY	8.1
1	B	264	GLY	7.8
1	D	265	ALA	6.2
1	D	328	LYS	6.2
1	D	267	ALA	6.0
1	D	326	LYS	6.0
1	D	304	GLY	5.6
1	B	304	GLY	5.1
1	D	305	LYS	4.9
1	D	302	GLY	4.9
1	D	325	GLU	4.7
1	D	268	THR	4.7
1	D	275	GLY	4.5
1	B	326	LYS	4.4
1	B	305	LYS	4.3
1	D	329	LYS	4.2
1	D	283	SER	4.1
1	D	263	ALA	4.1
1	B	325	GLU	4.1

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Mol	Chain	Res	Type	RSRZ
1	B	261	ALA	3.7
1	A	164	GLY	3.6
1	D	276	GLU	3.6
1	C	-3	ASP	3.5
1	D	327	GLN	3.4
1	B	303	GLY	3.3
1	B	265	ALA	3.1
1	D	262	GLY	3.1
1	B	302	GLY	3.0
1	D	281	ASP	3.0
1	D	322	GLN	3.0
1	A	194	GLY	2.9
1	D	274	LYS	2.9
1	A	-5	VAL	2.8
1	D	314	LYS	2.7
1	D	277	GLU	2.7
1	B	276	GLU	2.6
1	D	284	LYS	2.6
1	A	195	GLU	2.6
1	D	300	ASP	2.6
1	A	193	ASN	2.6
1	A	163	GLU	2.5
1	B	327	GLN	2.5
1	D	-2	ASP	2.5
1	D	197	ASP	2.4
1	D	270	LYS	2.3
1	D	259	GLY	2.3
1	D	-3	ASP	2.3
1	D	269	VAL	2.3
1	D	301	GLU	2.2
1	D	272	LYS	2.2
1	C	164	GLY	2.2
1	B	-4	ASP	2.1
1	D	256	ARG	2.1
1	B	300	ASP	2.1
1	B	277	GLU	2.1
1	A	192	ARG	2.1
1	C	256	ARG	2.1
1	B	75	ASP	2.1
1	D	279	GLU	2.1
1	D	273	TYR	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 monosaccharides 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.