



Full wwPDB X-ray Structure Validation Report ⓘ

May 14, 2020 – 05:27 am BST

PDB ID : 1VK0
Title : X-ray Structure of Gene Product from Arabidopsis Thaliana At5g06450
Authors : Wesenberg, G.E.; Smith, D.W.; Phillips Jr., G.N.; Johnson, K.A.; Bingman, C.A.; Center for Eukaryotic Structural Genomics (CESG)
Deposited on : 2004-04-12
Resolution : 2.10 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
Mogul : 1.8.5 (274361), CSD as541be (2020)
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

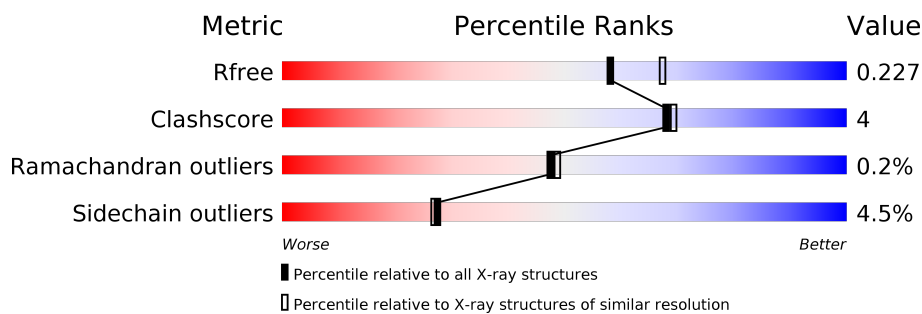
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)

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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	206	86% 10% . .
1	B	206	85% 10% . .
1	C	206	81% 13% . .
1	D	206	86% 8% . .
1	E	206	85% 11% . .
1	F	206	84% 11% . .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10215 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 hypothetical protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	200	Total	C	N	O	S	Se	0	0	0
			1591	1020	266	303	1	1			
1	B	200	Total	C	N	O	S	Se	0	1	0
			1601	1029	267	303	1	1			
1	C	200	Total	C	N	O	S	Se	0	2	0
			1605	1028	272	303	1	1			
1	D	200	Total	C	N	O	S	Se	0	0	0
			1591	1020	266	303	1	1			
1	E	200	Total	C	N	O	S	Se	0	0	0
			1591	1020	266	303	1	1			
1	F	200	Total	C	N	O	S	Se	0	0	0
			1591	1020	266	303	1	1			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	SER	-	CLONING ARTIFACT	UNP Q9FNG3
A	11	MSE	MET	MODIFIED RESIDUE	UNP Q9FNG3
B	1	SER	-	CLONING ARTIFACT	UNP Q9FNG3
B	11	MSE	MET	MODIFIED RESIDUE	UNP Q9FNG3
C	1	SER	-	CLONING ARTIFACT	UNP Q9FNG3
C	11	MSE	MET	MODIFIED RESIDUE	UNP Q9FNG3
D	1	SER	-	CLONING ARTIFACT	UNP Q9FNG3
D	11	MSE	MET	MODIFIED RESIDUE	UNP Q9FNG3
E	1	SER	-	CLONING ARTIFACT	UNP Q9FNG3
E	11	MSE	MET	MODIFIED RESIDUE	UNP Q9FNG3
F	1	SER	-	CLONING ARTIFACT	UNP Q9FNG3
F	11	MSE	MET	MODIFIED RESIDUE	UNP Q9FNG3

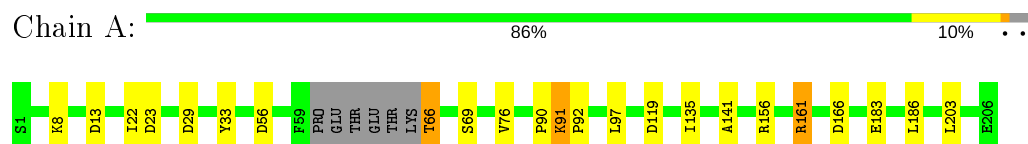
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	118	Total 118	O 118	0	0
2	B	124	Total 124	O 124	0	0
2	C	95	Total 95	O 95	0	0
2	D	96	Total 96	O 96	0	0
2	E	103	Total 103	O 103	0	0
2	F	109	Total 109	O 109	0	0

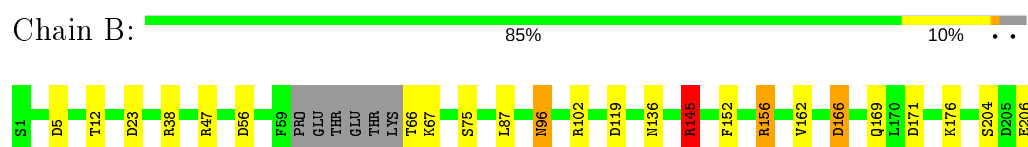
3 Residue-property plots [i](#)

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. 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. 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.

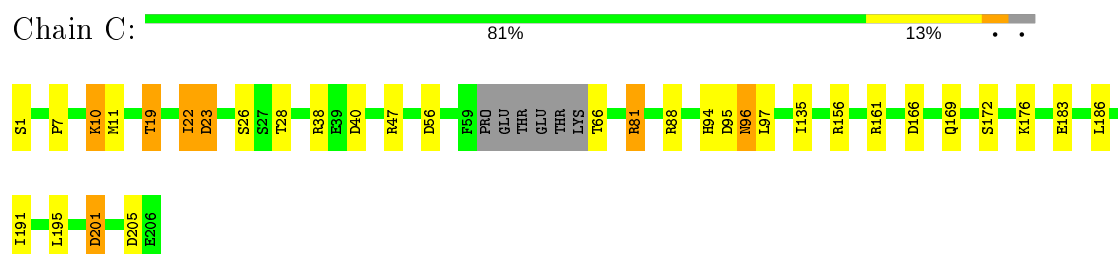
- Molecule 1: hypothetical protein



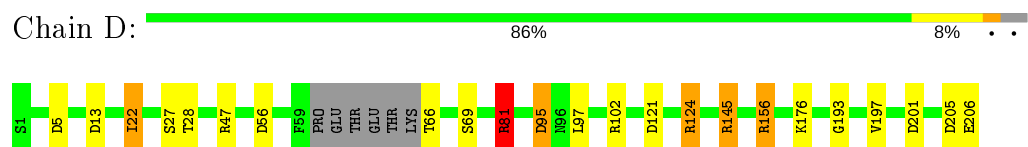
- Molecule 1: hypothetical protein



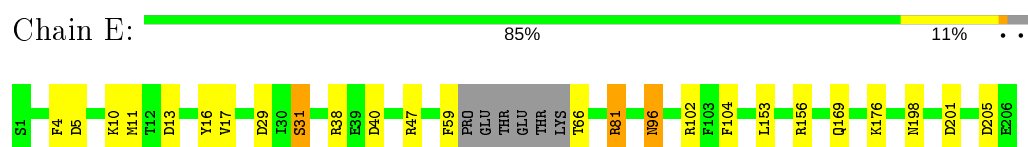
- Molecule 1: hypothetical protein



- Molecule 1: hypothetical protein



- Molecule 1: hypothetical protein



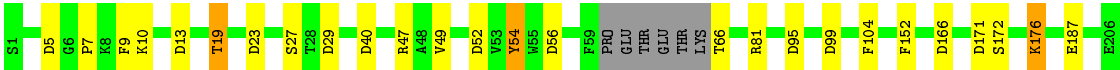
- Molecule 1: hypothetical protein

Chain F:

84%

11%

••



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	120.83Å 120.83Å 185.22Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	34.92 – 2.10 34.92 – 2.10	Depositor EDS
% Data completeness (in resolution range)	100.0 (34.92-2.10) 100.0 (34.92-2.10)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.02 (at 2.10Å)	Xtriage
Refinement program	REFMAC refmac _5.1.24	Depositor
R, R_{free}	0.180 , 0.233 0.180 , 0.227	Depositor DCC
R_{free} test set	4592 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	28.9	Xtriage
Anisotropy	0.053	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 48.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.017 for -h,-k,l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	10215	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.88% 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

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	1.05	0/1623	0.99	6/2196 (0.3%)
1	B	1.07	0/1639	1.06	11/2219 (0.5%)
1	C	1.08	1/1645 (0.1%)	1.11	13/2224 (0.6%)
1	D	1.03	0/1623	1.01	10/2196 (0.5%)
1	E	0.96	0/1623	0.94	6/2196 (0.3%)
1	F	0.98	2/1623 (0.1%)	1.03	13/2196 (0.6%)
All	All	1.03	3/9776 (0.0%)	1.02	59/13227 (0.4%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	11	MSE	SE-CE	-9.00	1.42	1.95
1	F	54	TYR	CD2-CE2	6.76	1.49	1.39
1	F	49	VAL	CB-CG1	5.06	1.63	1.52

All (59) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	161	ARG	NE-CZ-NH1	-9.11	115.74	120.30
1	C	47	ARG	NE-CZ-NH2	8.97	124.78	120.30
1	B	56	ASP	CB-CG-OD2	8.32	125.79	118.30
1	C	81[A]	ARG	NE-CZ-NH2	-8.27	116.17	120.30
1	C	81[B]	ARG	NE-CZ-NH2	-8.27	116.17	120.30
1	C	81[A]	ARG	NE-CZ-NH1	8.14	124.37	120.30
1	C	81[B]	ARG	NE-CZ-NH1	8.14	124.37	120.30
1	B	145	ARG	NE-CZ-NH2	-7.82	116.39	120.30
1	A	166	ASP	CB-CG-OD2	7.77	125.30	118.30
1	F	5	ASP	CB-CG-OD2	7.68	125.21	118.30
1	B	102	ARG	NE-CZ-NH2	7.35	123.97	120.30
1	F	171	ASP	CB-CG-OD2	7.31	124.88	118.30
1	A	29	ASP	CB-CG-OD2	7.30	124.87	118.30
1	D	5	ASP	CB-CG-OD2	7.25	124.82	118.30
1	B	166	ASP	CB-CG-OD2	7.18	124.77	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	47	ARG	NE-CZ-NH1	-7.18	116.71	120.30
1	D	13	ASP	CB-CG-OD2	7.14	124.73	118.30
1	A	56	ASP	CB-CG-OD2	7.06	124.66	118.30
1	C	56	ASP	CB-CG-OD2	7.03	124.62	118.30
1	D	124	ARG	NE-CZ-NH1	7.00	123.80	120.30
1	F	29	ASP	CB-CG-OD2	6.92	124.52	118.30
1	F	23	ASP	CB-CG-OD2	6.83	124.45	118.30
1	B	38	ARG	NE-CZ-NH1	-6.72	116.94	120.30
1	E	40	ASP	CB-CG-OD2	6.51	124.16	118.30
1	F	40	ASP	CB-CG-OD2	6.43	124.09	118.30
1	F	99	ASP	CB-CG-OD2	6.43	124.09	118.30
1	D	205	ASP	CB-CG-OD2	6.39	124.06	118.30
1	C	166	ASP	CB-CG-OD2	6.36	124.03	118.30
1	E	205	ASP	CB-CG-OD2	6.31	123.98	118.30
1	F	47	ARG	NE-CZ-NH2	6.22	123.41	120.30
1	A	13	ASP	CB-CG-OD2	6.06	123.76	118.30
1	F	166	ASP	CB-CG-OD2	6.01	123.71	118.30
1	E	102	ARG	NE-CZ-NH2	6.01	123.31	120.30
1	B	23	ASP	CB-CG-OD2	6.00	123.70	118.30
1	A	23	ASP	CB-CG-OD2	5.97	123.67	118.30
1	C	23	ASP	CB-CG-OD2	5.96	123.66	118.30
1	E	5	ASP	CB-CG-OD2	5.87	123.58	118.30
1	C	201	ASP	CB-CG-OD2	5.82	123.54	118.30
1	D	56	ASP	CB-CG-OD2	5.82	123.54	118.30
1	F	95	ASP	CB-CG-OD2	5.81	123.53	118.30
1	B	102	ARG	NE-CZ-NH1	-5.77	117.42	120.30
1	D	47	ARG	NE-CZ-NH1	-5.73	117.43	120.30
1	F	52	ASP	CB-CG-OD2	5.71	123.44	118.30
1	D	145	ARG	NE-CZ-NH2	5.60	123.10	120.30
1	E	13	ASP	CB-CG-OD2	5.48	123.23	118.30
1	F	13	ASP	CB-CG-OD2	5.47	123.22	118.30
1	D	121	ASP	CB-CG-OD2	5.43	123.19	118.30
1	C	95	ASP	CB-CG-OD2	5.43	123.19	118.30
1	C	40	ASP	CB-CG-OD2	5.39	123.15	118.30
1	B	145	ARG	NE-CZ-NH1	5.39	123.00	120.30
1	F	56	ASP	CB-CG-OD2	5.29	123.06	118.30
1	D	102	ARG	NE-CZ-NH2	5.29	122.94	120.30
1	B	145	ARG	CG-CD-NE	-5.28	100.72	111.80
1	B	119	ASP	CB-CG-OD1	5.13	122.92	118.30
1	A	119	ASP	CB-CG-OD1	5.11	122.90	118.30
1	F	81	ARG	NE-CZ-NH2	-5.11	117.75	120.30
1	D	95	ASP	CB-CG-OD2	5.09	122.89	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	201	ASP	CB-CG-OD2	5.02	122.82	118.30
1	B	5	ASP	CB-CG-OD2	5.01	122.81	118.30

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	1591	0	1575	11	0
1	B	1601	0	1581	11	0
1	C	1605	0	1593	25	0
1	D	1591	0	1575	10	0
1	E	1591	0	1575	16	0
1	F	1591	0	1575	5	0
2	A	118	0	0	0	0
2	B	124	0	0	0	0
2	C	95	0	0	1	0
2	D	96	0	0	0	0
2	E	103	0	0	1	0
2	F	109	0	0	0	0
All	All	10215	0	9474	73	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 4.

All (73) 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:C:19:THR:HG21	1:C:186:LEU:HD21	1.34	1.05
1:C:22:ILE:HD11	1:C:28:THR:HG21	1.39	1.05
1:E:81:ARG:HH12	1:E:198:ASN:HD21	1.11	0.95
1:C:19:THR:CG2	1:C:186:LEU:HD21	2.04	0.87
1:D:22:ILE:CD1	1:D:28:THR:HG21	2.07	0.85
1:B:96:ASN:H	1:B:96:ASN:HD22	1.27	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:10:LYS:O	1:C:10:LYS:HD2	1.80	0.81
1:E:81:ARG:NH1	1:E:198:ASN:HD21	1.85	0.72
1:D:22:ILE:HD11	1:D:28:THR:HG21	1.68	0.72
1:A:91:LYS:HB3	1:A:92:PRO:HD3	1.72	0.71
1:E:81:ARG:HH12	1:E:198:ASN:ND2	1.88	0.70
1:C:22:ILE:CD1	1:C:28:THR:HG21	2.19	0.69
1:F:7:PRO:O	1:F:19:THR:HB	1.93	0.68
1:C:7:PRO:O	1:C:19:THR:HB	1.94	0.68
1:E:11:MSE:HE3	1:E:17:VAL:CG2	2.25	0.65
1:B:96:ASN:N	1:B:96:ASN:HD22	1.96	0.64
1:B:145:ARG:NH2	1:B:162:VAL:O	2.30	0.63
1:E:29:ASP:OD2	1:E:31:SER:HB2	1.97	0.63
1:A:66:THR:HG23	1:A:69:SER:HB3	1.82	0.60
1:D:145:ARG:NH1	1:D:206:GLU:OE2	2.36	0.58
1:E:96:ASN:N	1:E:96:ASN:HD22	2.04	0.55
1:D:81:ARG:NH2	1:D:201:ASP:OD1	2.39	0.55
1:E:11:MSE:HE3	1:E:17:VAL:HG21	1.89	0.54
1:B:145:ARG:NH1	1:B:206:GLU:OE2	2.41	0.53
1:C:19:THR:HG21	1:C:186:LEU:CD2	2.23	0.53
1:E:96:ASN:HD22	1:E:96:ASN:H	1.54	0.53
1:A:141:ALA:HB2	1:A:203:LEU:HD11	1.92	0.51
1:C:10:LYS:N	1:C:10:LYS:HD2	2.26	0.51
1:D:145:ARG:HH11	1:D:206:GLU:CD	2.14	0.51
1:C:22:ILE:HD11	1:C:28:THR:CG2	2.27	0.49
1:C:10:LYS:CD	1:C:10:LYS:O	2.57	0.49
1:E:169:GLN:HA	1:E:169:GLN:HE21	1.78	0.48
1:C:96:ASN:H	1:C:96:ASN:HD22	1.60	0.48
1:F:9:PHE:CD1	1:F:187:GLU:HG3	2.49	0.47
1:D:66:THR:HB	1:D:69:SER:HB3	1.96	0.47
1:E:169:GLN:HA	1:E:169:GLN:NE2	2.30	0.46
1:C:22:ILE:CD1	1:C:28:THR:CG2	2.92	0.46
1:A:22:ILE:HD11	1:A:33:TYR:CD2	2.50	0.46
1:A:156:ARG:HD2	1:B:47:ARG:HD2	1.98	0.46
1:C:169:GLN:HA	1:C:169:GLN:HE21	1.80	0.45
1:C:172:SER:O	1:C:176:LYS:HG2	2.17	0.45
1:D:97:LEU:HD23	1:D:97:LEU:HA	1.85	0.45
1:C:10:LYS:H	1:C:10:LYS:HD2	1.81	0.44
1:C:26:SER:HB3	1:C:94:HIS:CD2	2.52	0.44
1:C:97:LEU:HA	1:C:97:LEU:HD23	1.83	0.44
1:F:54:TYR:N	1:F:54:TYR:CD1	2.85	0.44
1:D:145:ARG:HB3	1:D:145:ARG:HE	1.56	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:75:SER:HA	1:B:87:LEU:O	2.18	0.43
1:F:172:SER:O	1:F:176:LYS:HG3	2.18	0.43
1:C:10:LYS:H	1:C:10:LYS:NZ	2.17	0.43
1:D:156:ARG:HD2	1:E:47:ARG:HD2	1.99	0.43
1:C:191:ILE:O	1:C:195:LEU:HG	2.19	0.43
1:C:23:ASP:OD2	1:C:88:ARG:NH2	2.52	0.43
1:A:91:LYS:HB3	1:A:92:PRO:CD	2.45	0.43
1:E:10:LYS:HB2	1:E:16:TYR:CE1	2.53	0.43
1:A:135:ILE:HD13	1:F:152:PHE:CD1	2.54	0.43
1:D:193:GLY:O	1:D:197:VAL:HG23	2.19	0.42
1:E:4:PHE:N	1:E:4:PHE:CD2	2.87	0.42
1:E:47:ARG:HH11	1:E:47:ARG:HD3	1.72	0.42
1:A:90:PRO:HD2	1:A:97:LEU:HD11	2.02	0.42
1:A:156:ARG:HD2	1:B:47:ARG:CD	2.50	0.42
1:B:169:GLN:OE1	1:B:169:GLN:HA	2.20	0.41
1:A:91:LYS:HD3	1:A:91:LYS:HA	1.83	0.41
1:C:81[A]:ARG:NH1	1:C:201:ASP:OD2	2.54	0.41
1:C:96:ASN:N	1:C:96:ASN:HD22	2.17	0.41
1:B:96:ASN:ND2	1:B:96:ASN:N	2.66	0.41
1:C:10:LYS:CD	1:C:10:LYS:H	2.34	0.41
1:E:153:LEU:HA	1:E:153:LEU:HD23	1.82	0.41
1:E:81:ARG:HB3	2:E:299:HOH:O	2.21	0.41
1:B:156:ARG:NH1	1:B:171:ASP:OD1	2.45	0.40
1:B:152:PHE:CE2	1:C:135:ILE:HD13	2.56	0.40
1:A:161:ARG:HA	1:A:161:ARG:HD2	1.93	0.40
1:C:38:ARG:HD3	2:C:301:HOH:O	2.22	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	196/206 (95%)	194 (99%)	2 (1%)	0	100	100
1	B	197/206 (96%)	194 (98%)	2 (1%)	1 (0%)	29	26
1	C	198/206 (96%)	192 (97%)	6 (3%)	0	100	100
1	D	196/206 (95%)	192 (98%)	3 (2%)	1 (0%)	29	26
1	E	196/206 (95%)	194 (99%)	2 (1%)	0	100	100
1	F	196/206 (95%)	194 (99%)	2 (1%)	0	100	100
All	All	1179/1236 (95%)	1160 (98%)	17 (1%)	2 (0%)	47	49

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	12	THR
1	D	81	ARG

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	173/178 (97%)	166 (96%)	7 (4%)	31	32
1	B	174/178 (98%)	165 (95%)	9 (5%)	23	21
1	C	175/178 (98%)	166 (95%)	9 (5%)	24	22
1	D	173/178 (97%)	166 (96%)	7 (4%)	31	32
1	E	173/178 (97%)	164 (95%)	9 (5%)	23	21
1	F	173/178 (97%)	167 (96%)	6 (4%)	36	38
All	All	1041/1068 (98%)	994 (96%)	47 (4%)	27	27

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	LYS
1	A	66	THR
1	A	76	VAL

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Mol	Chain	Res	Type
1	A	91	LYS
1	A	161	ARG
1	A	183	GLU
1	A	186	LEU
1	B	66	THR
1	B	67	LYS
1	B	96	ASN
1	B	136	ASN
1	B	145	ARG
1	B	156	ARG
1	B	166	ASP
1	B	176	LYS
1	B	204	SER
1	C	1	SER
1	C	10	LYS
1	C	19	THR
1	C	22	ILE
1	C	66	THR
1	C	96	ASN
1	C	156	ARG
1	C	183	GLU
1	C	205	ASP
1	D	22	ILE
1	D	27	SER
1	D	81	ARG
1	D	95	ASP
1	D	124	ARG
1	D	156	ARG
1	D	176	LYS
1	E	31	SER
1	E	38	ARG
1	E	59	PHE
1	E	66	THR
1	E	81	ARG
1	E	96	ASN
1	E	104	PHE
1	E	156	ARG
1	E	176	LYS
1	F	10	LYS
1	F	19	THR
1	F	27	SER
1	F	66	THR

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Mol	Chain	Res	Type
1	F	104	PHE
1	F	176	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	133	ASN
1	B	46	ASN
1	B	96	ASN
1	B	133	ASN
1	B	136	ASN
1	C	94	HIS
1	C	96	ASN
1	C	133	ASN
1	C	169	GLN
1	C	198	ASN
1	D	126	ASN
1	E	96	ASN
1	E	169	GLN
1	E	198	ASN
1	F	44	ASN
1	F	133	ASN

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

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](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.