



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 13, 2018 – 09:39 PM EDT

PDB ID : 4P5K
Title : Structural Basis of Chronic Beryllium Disease: Bridging the Gap Between Allergy and Autoimmunity
Authors : Clayton, G.M.; Wang, Y.; Crawford, F.; Novikov, A.; Wimberly, B.T.; Kieft, J.S.; Falta, M.T.; Bowerman, N.A.; Marrack, P.; Fontenot, A.P.; Dai, S.; Kappler, J.W.
Deposited on : 2014-03-17
Resolution : 2.59 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

<http://wwpdb.org/validation/2016/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)	:	1.9-1692
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

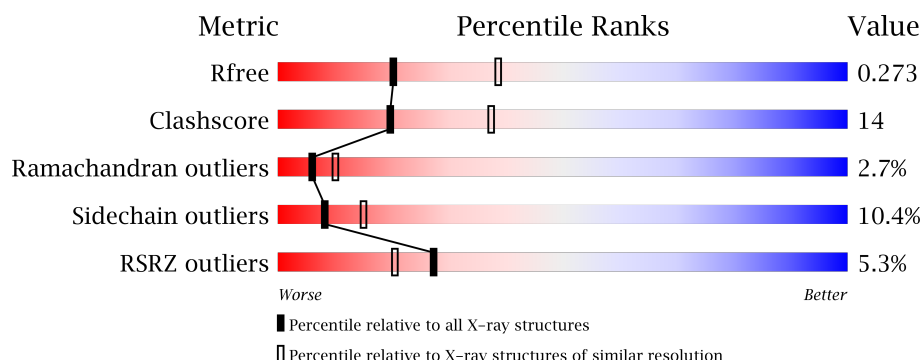
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.59 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	2542 (2.60-2.60)
Clashscore	112137	2895 (2.60-2.60)
Ramachandran outliers	110173	2848 (2.60-2.60)
Sidechain outliers	110143	2848 (2.60-2.60)
RSRZ outliers	101464	2550 (2.60-2.60)

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. 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	183	<div> <div>0.1%</div> <div>67% 28% . .</div> </div>
1	D	183	<div> <div>0.1%</div> <div>72% 23% . .</div> </div>
2	B	212	<div> <div>9%</div> <div>56% 31% 6% 7%</div> </div>
2	E	212	<div> <div>8%</div> <div>58% 29% 6% 7%</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6218 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 HLA class II histocompatibility antigen, DP alpha 1 chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	181	Total	C	N	O	S	0	0	0
			1482	958	240	279	5			
1	D	181	Total	C	N	O	S	0	0	0
			1482	958	240	279	5			

- Molecule 2 is a protein called RAS peptide,HLA class II histocompatibility antigen, DP beta 1 chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	198	Total	C	N	O	S	0	0	0
			1627	1020	287	312	8			
2	E	198	Total	C	N	O	S	0	0	0
			1627	1020	287	312	8			

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-13	GLY	-	linker	PDB ?
B	-12	GLY	-	linker	PDB ?
B	-11	SER	-	linker	PDB ?
B	-10	LEU	-	linker	PDB ?
B	-9	VAL	-	linker	PDB ?
B	-8	PRO	-	linker	PDB ?
B	-7	ARG	-	linker	PDB ?
B	-6	GLY	-	linker	PDB ?
B	-5	SER	-	linker	PDB ?
B	-4	GLY	-	linker	PDB ?
B	-3	GLY	-	linker	PDB ?
B	-2	GLY	-	linker	PDB ?
B	-1	GLY	-	linker	PDB ?
B	3	SER	THR	VARIANT	UNP Q5EP54
E	-13	GLY	-	linker	PDB ?
E	-12	GLY	-	linker	PDB ?

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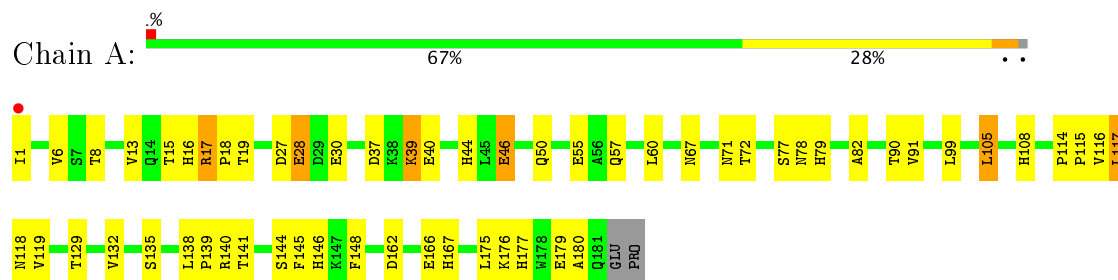
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Chain	Residue	Modelled	Actual	Comment	Reference
E	-11	SER	-	linker	PDB ?
E	-10	LEU	-	linker	PDB ?
E	-9	VAL	-	linker	PDB ?
E	-8	PRO	-	linker	PDB ?
E	-7	ARG	-	linker	PDB ?
E	-6	GLY	-	linker	PDB ?
E	-5	SER	-	linker	PDB ?
E	-4	GLY	-	linker	PDB ?
E	-3	GLY	-	linker	PDB ?
E	-2	GLY	-	linker	PDB ?
E	-1	GLY	-	linker	PDB ?
E	3	SER	THR	VARIANT	UNP Q5EP54

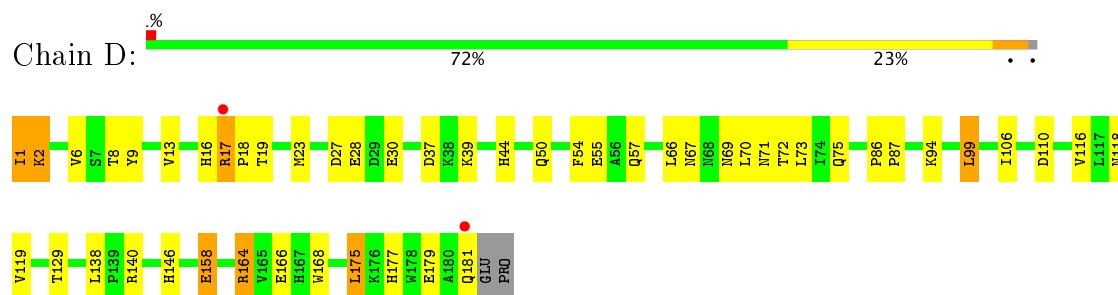
3 Residue-property plots [i](#)

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.

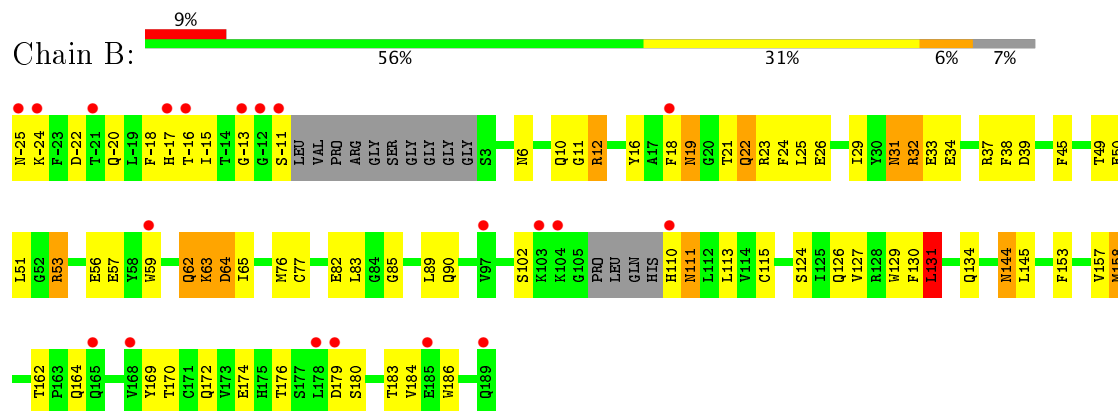
- Molecule 1: HLA class II histocompatibility antigen, DP alpha 1 chain



- Molecule 1: HLA class II histocompatibility antigen, DP alpha 1 chain

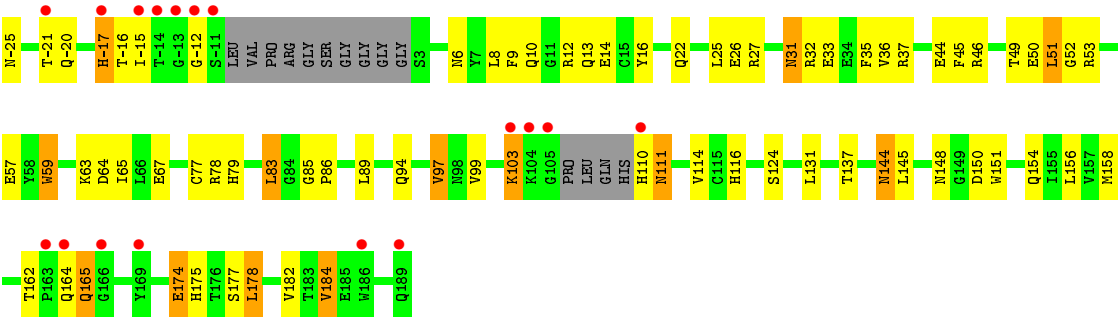


- Molecule 2: RAS peptide,HLA class II histocompatibility antigen, DP beta 1 chain



- Molecule 2: RAS peptide,HLA class II histocompatibility antigen, DP beta 1 chain





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	182.78Å 68.74Å 71.13Å 90.00° 105.72° 90.00°	Depositor
Resolution (Å)	44.62 – 2.59 44.62 – 2.59	Depositor EDS
% Data completeness (in resolution range)	91.2 (44.62-2.59) 91.2 (44.62-2.59)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.26 (at 2.58Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.221 , 0.281 0.219 , 0.273	Depositor DCC
R_{free} test set	1243 reflections (5.39%)	DCC
Wilson B-factor (Å ²)	49.9	Xtriage
Anisotropy	0.146	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 34.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6218	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 4.43% 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	0.76	1/1530 (0.1%)	0.78	0/2088
1	D	0.74	0/1530	0.80	3/2088 (0.1%)
2	B	0.71	1/1665 (0.1%)	0.78	1/2253 (0.0%)
2	E	0.73	1/1665 (0.1%)	0.83	1/2253 (0.0%)
All	All	0.73	3/6390 (0.0%)	0.80	5/8682 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	186	TRP	CD2-CE2	-5.41	1.34	1.41
1	A	55	GLU	CG-CD	5.24	1.59	1.51
2	E	174	GLU	CG-CD	5.00	1.59	1.51

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	-17	HIS	CB-CA-C	-6.21	97.98	110.40
1	D	73	LEU	CA-CB-CG	-6.03	101.43	115.30
1	D	99	LEU	CA-CB-CG	5.93	128.95	115.30
1	D	175	LEU	CA-CB-CG	5.57	128.10	115.30
2	B	131	LEU	CA-CB-CG	5.07	126.95	115.30

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	1482	0	1385	37	0
1	D	1482	0	1385	38	1
2	B	1627	0	1535	51	1
2	E	1627	0	1535	60	0
All	All	6218	0	5840	172	1

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 (172) 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)
2:E:22:GLN:HG2	2:E:77:CYS:HB3	1.40	1.02
2:B:12:ARG:HH11	2:B:12:ARG:CG	1.73	0.99
1:D:54:PHE:HE1	2:E:-21:THR:OG1	1.45	0.98
2:E:144:ASN:HD22	2:E:145:LEU:H	1.16	0.92
1:D:17:ARG:HB2	1:D:18:PRO:CD	2.00	0.91
2:E:22:GLN:HG2	2:E:77:CYS:CB	2.04	0.87
2:B:49:THR:HG22	2:B:51:LEU:H	1.39	0.86
2:B:126:GLN:HB2	2:B:174:GLU:OE1	1.76	0.85
2:B:12:ARG:HH11	2:B:12:ARG:HG3	1.45	0.82
1:D:54:PHE:CE1	2:E:-21:THR:OG1	2.31	0.81
2:B:110:HIS:O	2:B:111:ASN:HB2	1.83	0.78
2:B:12:ARG:HH11	2:B:12:ARG:HG2	1.46	0.78
2:E:116:HIS:HE1	2:E:154:GLN:OE1	1.66	0.78
1:A:30:GLU:OE2	1:A:44:HIS:HD2	1.65	0.78
2:E:-17:HIS:HE1	2:E:45:PHE:CZ	2.04	0.76
1:D:17:ARG:HB2	1:D:18:PRO:HD2	1.69	0.75
2:B:-15:ILE:HA	2:B:59:TRP:CH2	2.23	0.73
2:E:22:GLN:CG	2:E:77:CYS:HB3	2.18	0.73
2:E:-21:THR:HG22	2:E:-20:GLN:H	1.55	0.72
1:D:17:ARG:HG3	1:D:116:VAL:HG13	1.72	0.71
1:D:17:ARG:CB	1:D:18:PRO:CD	2.67	0.71
2:B:-15:ILE:HA	2:B:59:TRP:HH2	1.56	0.70
1:A:17:ARG:CB	1:A:18:PRO:CD	2.69	0.70
2:E:164:GLN:O	2:E:165:GLN:HB2	1.93	0.69
1:D:30:GLU:OE2	1:D:44:HIS:HD2	1.76	0.69
1:A:17:ARG:HB3	1:A:18:PRO:HD3	1.75	0.68
1:A:17:ARG:HG3	1:A:116:VAL:HG13	1.76	0.68
2:E:49:THR:HG22	2:E:51:LEU:H	1.58	0.67
2:B:11:GLY:O	2:B:12:ARG:HD2	1.95	0.67
2:E:175:HIS:HB3	2:E:178:LEU:HD12	1.77	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:71:ASN:O	1:D:75:GLN:HG2	1.95	0.67
1:D:16:HIS:CD2	1:D:17:ARG:HE	2.14	0.66
2:E:14:GLU:OE1	2:E:27:ARG:HD2	1.95	0.65
2:B:10:GLN:HB2	2:B:29:ILE:HB	1.78	0.65
1:D:17:ARG:HB2	1:D:18:PRO:HD3	1.80	0.64
2:E:164:GLN:O	2:E:165:GLN:CB	2.45	0.63
2:B:85:GLY:HA2	2:B:89:LEU:HD13	1.80	0.63
2:E:94:GLN:NE2	2:E:177:SER:O	2.31	0.63
2:B:12:ARG:NH1	2:B:12:ARG:CG	2.46	0.62
2:E:8:LEU:HD21	2:E:10:GLN:CG	2.30	0.62
1:D:6:VAL:HG12	1:D:8:THR:CG2	2.30	0.61
1:D:17:ARG:CB	1:D:18:PRO:HD3	2.30	0.60
2:E:165:GLN:OE1	2:E:165:GLN:HA	2.02	0.59
2:E:116:HIS:CE1	2:E:154:GLN:OE1	2.53	0.59
1:A:17:ARG:HB3	1:A:18:PRO:CD	2.31	0.59
1:D:37:ASP:O	1:D:39:LYS:HD2	2.02	0.59
2:E:116:HIS:CE1	2:E:156:LEU:HD11	2.37	0.59
2:E:49:THR:HG22	2:E:50:GLU:N	2.18	0.59
1:A:72:THR:HG23	2:B:-13:GLY:HA3	1.86	0.58
1:D:16:HIS:HD2	1:D:17:ARG:HE	1.50	0.58
1:A:108:HIS:NE2	1:A:146:HIS:HD2	2.02	0.58
1:A:46:GLU:HG3	1:A:50:GLN:HE22	1.69	0.58
2:B:22:GLN:CG	2:B:77:CYS:HB3	2.34	0.58
2:B:12:ARG:NH1	2:B:12:ARG:HG2	2.16	0.58
1:A:116:VAL:O	1:A:167:HIS:HD2	1.86	0.57
2:E:8:LEU:HD21	2:E:10:GLN:HG3	1.84	0.57
1:D:118:ASN:HB2	1:D:166:GLU:HB2	1.85	0.57
1:D:13:VAL:HG12	1:D:70:LEU:HD22	1.86	0.57
1:D:23:MET:HE1	1:D:138:LEU:HA	1.87	0.56
2:B:172:GLN:HB2	2:B:183:THR:HG22	1.86	0.56
1:A:17:ARG:CG	1:A:116:VAL:HG13	2.35	0.56
2:E:103:LYS:HB3	2:E:111:ASN:HA	1.88	0.56
2:B:-17:HIS:HE1	2:B:45:PHE:HZ	1.54	0.55
2:E:144:ASN:HD22	2:E:145:LEU:N	1.95	0.55
1:D:72:THR:HG23	2:E:-12:GLY:H	1.71	0.55
2:E:110:HIS:O	2:E:111:ASN:HB2	2.07	0.55
2:E:33:GLU:OE2	2:E:49:THR:HG21	2.07	0.55
1:D:119:VAL:HA	1:D:164:ARG:O	2.06	0.55
1:D:1:ILE:O	1:D:2:LYS:HB2	2.07	0.55
1:A:17:ARG:HB2	1:A:18:PRO:HD2	1.87	0.55
1:D:13:VAL:HG11	1:D:67:ASN:OD1	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:82:ALA:HB2	2:B:31:ASN:HB3	1.89	0.54
2:E:-15:ILE:HA	2:E:59:TRP:CH2	2.43	0.54
2:B:113:LEU:HD23	2:B:129:TRP:CH2	2.43	0.54
2:B:179:ASP:OD1	2:B:179:ASP:N	2.40	0.54
1:A:78:ASN:O	1:A:79:HIS:HB2	2.08	0.53
2:B:23:ARG:NH2	2:B:39:ASP:OD2	2.41	0.53
1:D:177:HIS:CE1	1:D:179:GLU:OE1	2.61	0.53
2:B:22:GLN:HG3	2:B:77:CYS:HB3	1.90	0.53
2:E:94:GLN:HA	2:E:177:SER:OG	2.08	0.53
2:B:-13:GLY:C	2:B:-11:SER:H	2.11	0.52
2:B:53:ARG:O	2:B:57:GLU:HG2	2.09	0.52
2:E:148:ASN:HB2	2:E:150:ASP:OD1	2.09	0.52
2:E:150:ASP:O	2:E:151:TRP:HB2	2.08	0.52
2:E:64:ASP:O	2:E:67:GLU:HG2	2.11	0.51
2:B:131:LEU:HD23	2:B:169:TYR:CE2	2.46	0.51
1:A:162:ASP:HA	1:A:176:LYS:O	2.10	0.51
1:A:16:HIS:H	2:B:6:ASN:HD21	1.59	0.51
1:A:139:PRO:HB2	2:B:12:ARG:NH2	2.26	0.50
1:D:16:HIS:HB3	1:D:17:ARG:HG2	1.92	0.50
2:B:50:GLU:OE2	2:B:53:ARG:NH2	2.45	0.50
1:D:69:ASN:HD21	2:E:-15:ILE:H	1.60	0.50
2:E:35:PHE:HD1	2:E:52:GLY:HA2	1.76	0.50
1:A:17:ARG:NE	1:A:17:ARG:H	2.09	0.50
2:E:36:VAL:HG22	2:E:37:ARG:H	1.77	0.50
2:E:-21:THR:HG22	2:E:-20:GLN:N	2.24	0.49
1:D:110:ASP:OD1	1:D:146:HIS:HB3	2.12	0.49
2:B:59:TRP:O	2:B:62:GLN:HB3	2.12	0.49
2:E:-17:HIS:CE1	2:E:45:PHE:CZ	2.94	0.49
2:E:35:PHE:O	2:E:52:GLY:HA3	2.13	0.48
1:D:69:ASN:HD21	2:E:-16:THR:HA	1.79	0.48
2:B:-22:ASP:HB3	2:B:76:MET:HG3	1.96	0.47
2:E:22:GLN:HG2	2:E:77:CYS:HB2	1.94	0.47
1:D:30:GLU:OE2	1:D:44:HIS:CD2	2.62	0.47
2:B:144:ASN:HD22	2:B:145:LEU:H	1.61	0.47
2:B:18:PHE:O	2:B:19:ASN:C	2.53	0.47
2:E:12:ARG:O	2:E:26:GLU:HA	2.14	0.47
1:A:27:ASP:O	1:A:28:GLU:HB2	2.15	0.47
1:A:17:ARG:HB2	1:A:116:VAL:HG22	1.97	0.47
2:B:130:PHE:O	2:B:169:TYR:HA	2.14	0.47
2:E:8:LEU:C	2:E:8:LEU:HD23	2.35	0.47
2:B:16:TYR:CE1	2:B:25:LEU:HD13	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:115:PRO:HG3	1:A:145:PHE:CE1	2.51	0.46
2:B:31:ASN:O	2:B:32:ARG:HB2	2.15	0.46
2:E:178:LEU:HD22	2:E:182:VAL:HG23	1.98	0.46
1:A:91:VAL:HG12	1:A:105:LEU:HD22	1.98	0.46
2:E:99:VAL:HA	2:E:114:VAL:O	2.16	0.46
2:B:33:GLU:OE2	2:B:49:THR:HG21	2.15	0.46
1:D:6:VAL:HG12	1:D:8:THR:HG23	1.96	0.46
2:E:53:ARG:O	2:E:57:GLU:HG3	2.15	0.46
2:B:62:GLN:O	2:B:64:ASP:N	2.50	0.45
2:E:44:GLU:OE1	2:E:46:ARG:NH1	2.48	0.45
2:B:38:PHE:HB2	2:B:45:PHE:CE1	2.52	0.45
1:D:94:LYS:HB2	1:D:106:ILE:HD11	1.99	0.45
1:A:37:ASP:O	1:A:39:LYS:HD2	2.17	0.45
2:E:79:HIS:O	2:E:83:LEU:HB2	2.16	0.45
1:A:27:ASP:O	1:A:28:GLU:CB	2.65	0.44
2:B:24:PHE:CE2	2:B:26:GLU:HB2	2.51	0.44
1:D:86:PRO:HA	1:D:87:PRO:HD3	1.73	0.44
1:A:46:GLU:CG	1:A:50:GLN:HE22	2.29	0.44
2:B:12:ARG:NH1	2:B:12:ARG:HG3	2.19	0.44
1:A:117:LEU:HD13	1:A:119:VAL:HG23	2.00	0.44
1:A:140:ARG:HD3	1:A:144:SER:OG	2.17	0.44
2:B:45:PHE:O	2:B:56:GLU:HG3	2.18	0.44
2:E:-15:ILE:HA	2:E:59:TRP:HH2	1.83	0.44
2:B:89:LEU:H	2:B:89:LEU:CD1	2.32	0.43
1:A:135:SER:HB3	1:A:148:PHE:HB2	2.01	0.43
2:B:59:TRP:HD1	2:B:65:ILE:HG21	1.81	0.43
2:E:36:VAL:HG22	2:E:37:ARG:N	2.33	0.43
2:E:8:LEU:HD21	2:E:10:GLN:HG2	1.99	0.43
2:B:127:VAL:HB	2:B:157:VAL:HG21	2.00	0.43
2:E:97:VAL:HG22	2:E:184:VAL:CG1	2.49	0.43
1:A:139:PRO:HB2	2:B:12:ARG:HH22	1.83	0.43
2:E:97:VAL:HG22	2:E:184:VAL:HG11	2.01	0.43
1:A:114:PRO:HB2	1:A:115:PRO:CD	2.49	0.42
2:E:16:TYR:CE1	2:E:25:LEU:HD13	2.53	0.42
2:E:178:LEU:HD22	2:E:182:VAL:CG2	2.49	0.42
1:A:57:GLN:HA	1:A:60:LEU:HD12	2.01	0.42
2:B:38:PHE:C	2:B:38:PHE:CD2	2.91	0.42
2:E:175:HIS:CB	2:E:178:LEU:HD12	2.48	0.42
1:A:77:SER:O	1:A:78:ASN:HB2	2.18	0.42
1:D:17:ARG:HG3	1:D:116:VAL:CG1	2.45	0.42
1:A:118:ASN:HB2	1:A:166:GLU:HB2	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:67:ASN:O	1:A:71:ASN:OD1	2.37	0.42
2:B:115:CYS:HB2	2:B:129:TRP:CZ2	2.53	0.42
1:A:117:LEU:HD13	1:A:119:VAL:CG2	2.50	0.42
2:B:49:THR:CG2	2:B:50:GLU:N	2.83	0.42
1:D:69:ASN:ND2	2:E:-15:ILE:H	2.18	0.42
2:B:145:LEU:HD21	2:B:153:PHE:CD2	2.55	0.42
2:E:-17:HIS:CE1	2:E:45:PHE:HZ	2.35	0.42
1:D:177:HIS:HE1	1:D:179:GLU:OE1	2.00	0.41
1:D:66:LEU:HD22	2:E:9:PHE:CD2	2.55	0.41
2:E:175:HIS:HB3	2:E:178:LEU:CD1	2.47	0.41
1:D:16:HIS:HB2	2:E:6:ASN:ND2	2.36	0.41
1:D:158:GLU:H	1:D:158:GLU:HG2	1.74	0.41
1:D:13:VAL:HG11	1:D:67:ASN:CG	2.41	0.41
2:B:113:LEU:O	2:B:158:MET:HA	2.21	0.41
1:A:177:HIS:HE1	1:A:179:GLU:OE1	2.04	0.40
2:B:34:GLU:OE1	2:B:37:ARG:HB2	2.21	0.40
1:A:138:LEU:HB2	1:A:146:HIS:CE1	2.57	0.40
1:A:6:VAL:HG12	1:A:8:THR:CG2	2.52	0.40
1:D:9:TYR:HB2	2:E:13:GLN:HB2	2.04	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:-18:PHE:O	1:D:57:GLN:NE2[1_545]	2.12	0.08

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	179/183 (98%)	165 (92%)	12 (7%)	2 (1%)	17 35

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	179/183 (98%)	166 (93%)	8 (4%)	5 (3%)	6	9
2	B	192/212 (91%)	171 (89%)	15 (8%)	6 (3%)	5	8
2	E	192/212 (91%)	172 (90%)	13 (7%)	7 (4%)	4	5
All	All	742/790 (94%)	674 (91%)	48 (6%)	20 (3%)	6	10

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	17	ARG
2	B	19	ASN
2	B	62	GLN
2	B	63	LYS
2	B	111	ASN
1	D	17	ARG
2	E	31	ASN
2	E	111	ASN
2	E	165	GLN
1	A	180	ALA
1	D	168	TRP
2	E	103	LYS
2	B	31	ASN
1	D	27	ASP
1	D	99	LEU
2	B	102	SER
1	D	2	LYS
2	E	86	PRO
2	E	32	ARG
2	E	85	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	162/164 (99%)	146 (90%)	16 (10%)	9	17

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	162/164 (99%)	151 (93%)	11 (7%)	18	37
2	B	179/188 (95%)	154 (86%)	25 (14%)	4	7
2	E	179/188 (95%)	160 (89%)	19 (11%)	8	14
All	All	682/704 (97%)	611 (90%)	71 (10%)	8	15

All (71) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	ILE
1	A	13	VAL
1	A	15	THR
1	A	19	THR
1	A	28	GLU
1	A	39	LYS
1	A	40	GLU
1	A	46	GLU
1	A	90	THR
1	A	99	LEU
1	A	105	LEU
1	A	117	LEU
1	A	129	THR
1	A	132	VAL
1	A	141	THR
1	A	175	LEU
2	B	-25	ASN
2	B	-24	LYS
2	B	-20	GLN
2	B	-16	THR
2	B	12	ARG
2	B	21	THR
2	B	22	GLN
2	B	32	ARG
2	B	53	ARG
2	B	63	LYS
2	B	64	ASP
2	B	82	GLU
2	B	83	LEU
2	B	90	GLN
2	B	124	SER
2	B	131	LEU
2	B	134	GLN

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Mol	Chain	Res	Type
2	B	144	ASN
2	B	158	MET
2	B	162	THR
2	B	164	GLN
2	B	170	THR
2	B	176	THR
2	B	180	SER
2	B	184	VAL
1	D	1	ILE
1	D	19	THR
1	D	28	GLU
1	D	50	GLN
1	D	55	GLU
1	D	129	THR
1	D	140	ARG
1	D	158	GLU
1	D	164	ARG
1	D	175	LEU
1	D	181	GLN
2	E	-25	ASN
2	E	31	ASN
2	E	51	LEU
2	E	59	TRP
2	E	63	LYS
2	E	65	ILE
2	E	78	ARG
2	E	83	LEU
2	E	89	LEU
2	E	97	VAL
2	E	124	SER
2	E	131	LEU
2	E	137	THR
2	E	144	ASN
2	E	158	MET
2	E	162	THR
2	E	174	GLU
2	E	178	LEU
2	E	184	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	44	HIS
1	A	50	GLN
1	A	69	ASN
1	A	146	HIS
1	A	167	HIS
1	A	177	HIS
2	B	-17	HIS
2	B	6	ASN
2	B	79	HIS
2	B	134	GLN
2	B	144	ASN
1	D	16	HIS
1	D	44	HIS
1	D	68	ASN
1	D	69	ASN
1	D	124	ASN
1	D	146	HIS
1	D	177	HIS
2	E	-17	HIS
2	E	6	ASN
2	E	22	GLN
2	E	110	HIS
2	E	116	HIS
2	E	144	ASN

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 ⓘ

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	181/183 (98%)	0.03	1 (0%) 89 88	26, 42, 58, 68	0
1	D	181/183 (98%)	-0.04	2 (1%) 80 77	26, 41, 60, 79	0
2	B	198/212 (93%)	0.61	20 (10%) 8 5	31, 61, 90, 100	0
2	E	198/212 (93%)	0.51	17 (8%) 11 7	30, 50, 78, 102	0
All	All	758/790 (95%)	0.29	40 (5%) 27 20	26, 47, 80, 102	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	E	-11	SER	9.8
2	B	110	HIS	6.3
2	B	-12	GLY	6.2
2	E	104	LYS	5.9
2	E	110	HIS	5.8
2	B	-13	GLY	5.7
2	E	-12	GLY	5.7
2	E	105	GLY	5.6
2	E	-21	THR	4.7
2	E	189	GLN	4.6
2	E	103	LYS	4.5
1	D	17	ARG	4.3
2	B	59	TRP	4.2
2	B	104	LYS	4.2
2	E	166	GLY	4.1
2	B	-21	THR	3.9
2	E	164	GLN	3.5
2	E	186	TRP	3.4
2	B	-11	SER	3.2
2	B	179	ASP	3.1
2	B	-17	HIS	3.0

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Mol	Chain	Res	Type	RSRZ
2	B	103	LYS	2.9
1	D	181	GLN	2.8
2	B	18	PHE	2.8
2	B	178	LEU	2.8
2	B	-16	THR	2.7
1	A	1	ILE	2.5
2	E	-15	ILE	2.4
2	E	169	TYR	2.4
2	E	-17	HIS	2.3
2	B	-24	LYS	2.2
2	E	-13	GLY	2.2
2	B	165	GLN	2.2
2	E	-14	THR	2.2
2	B	185	GLU	2.2
2	E	163	PRO	2.1
2	B	-25	ASN	2.1
2	B	97	VAL	2.1
2	B	189	GLN	2.1
2	B	168	VAL	2.1

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.