



Full wwPDB X-ray Structure Validation Report i

Feb 28, 2014 – 10:17 PM GMT

PDB ID : 4F9P
Title : Crystal Structure of the Human BTN3A1 Ectodomain in Complex with the 103.2 Single Chain Antibody
Authors : Palakodeti, A.; Sandstrom, A.; Sundaresan, L.; Harly, C.; Nedellec, S.; Olive, D.; Scotet, E.; Bonneville, M.; Adams, E.J.
Deposited on : 2012-05-19
Resolution : 3.52 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

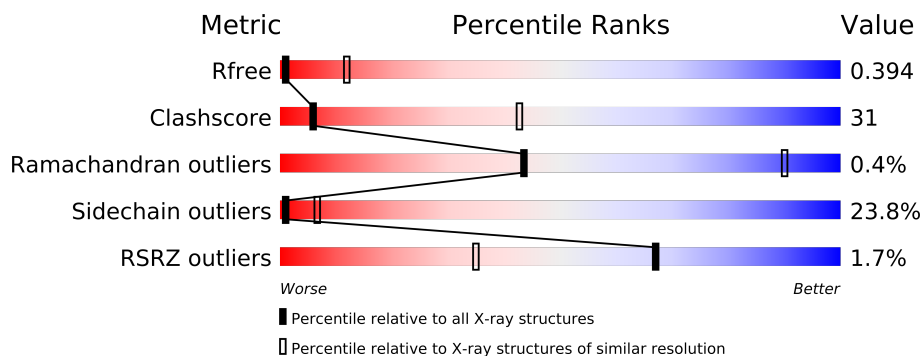
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 3.52 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}	66092	1256 (3.74-3.30)
Clashscore	79885	1020 (3.68-3.36)
Ramachandran outliers	78287	1082 (3.70-3.34)
Sidechain outliers	78261	1082 (3.70-3.34)
RSRZ outliers	66119	1256 (3.74-3.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	220	
1	B	220	
2	C	254	
2	D	254	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5552 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Butyrophilin subfamily 3 member A1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	207	Total	C	N	O	S	0	0	0
			1258	782	227	245	4			
1	B	208	Total	C	N	O	S	0	0	0
			1253	769	225	253	6			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	ALA	-	EXPRESSION TAG	UNP O00481
A	-1	ASP	-	EXPRESSION TAG	UNP O00481
A	0	LEU	-	EXPRESSION TAG	UNP O00481
B	-2	ALA	-	EXPRESSION TAG	UNP O00481
B	-1	ASP	-	EXPRESSION TAG	UNP O00481
B	0	LEU	-	EXPRESSION TAG	UNP O00481

- Molecule 2 is a protein called 103.2 anti-BTN3A1 antibody fragment.

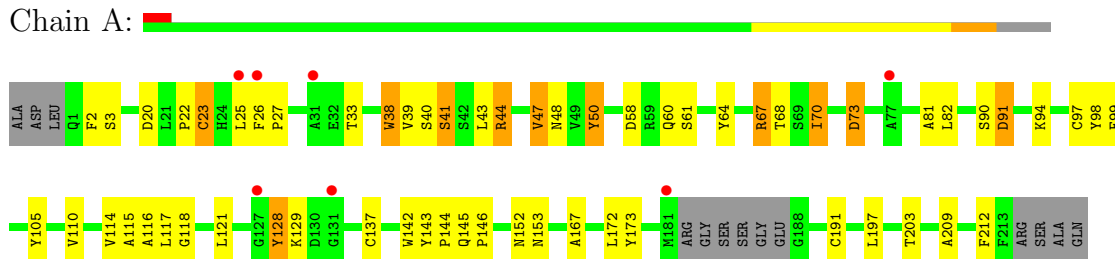
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	224	Total	C	N	O	S	0	0	0
			1550	982	254	310	4			
2	C	220	Total	C	N	O	S	0	0	0
			1491	934	251	301	5			

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

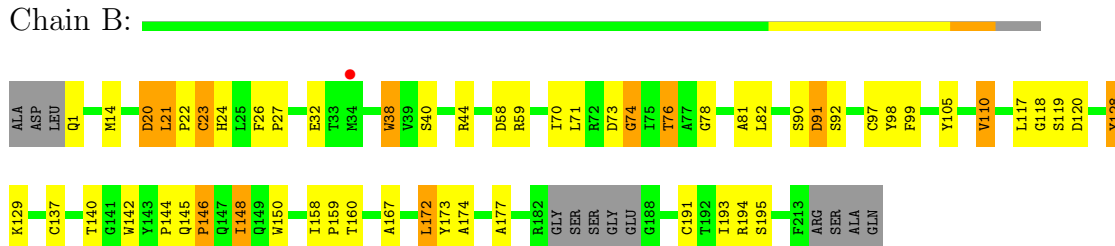
- Molecule 1: Butyrophilin subfamily 3 member A1

Chain A:



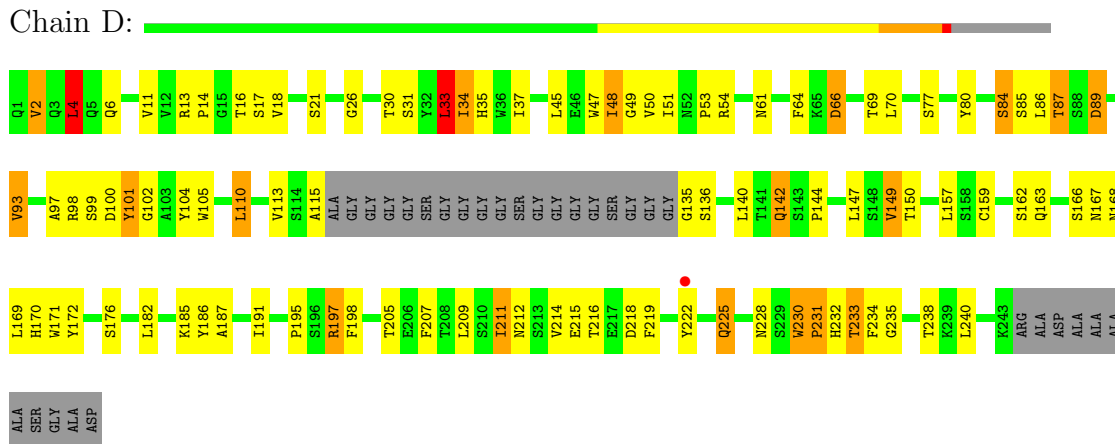
- Molecule 1: Butyrophilin subfamily 3 member A1

Chain B:



- Molecule 2: 103.2 anti-BTN3A1 antibody fragment

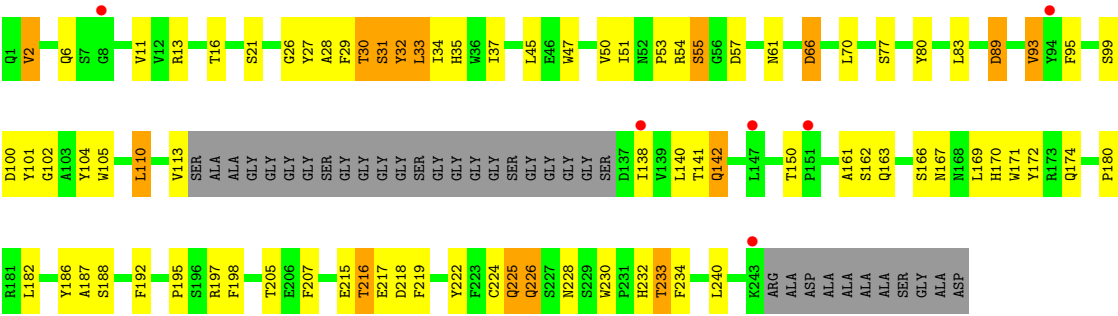
Chain D:



- Molecule 2: 103.2 anti-BTN3A1 antibody fragment

Chain C:





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	88.36Å 117.68Å 121.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.54 – 3.52 41.54 – 3.52	Depositor EDS
% Data completeness (in resolution range)	83.3 (41.54-3.52) 83.3 (41.54-3.52)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.65 (at 3.48Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.2_869)	Depositor
R, R_{free}	0.373 , 0.408 0.365 , 0.394	Depositor DCC
R_{free} test set	668 reflections (4.91%)	DCC
Wilson B-factor (Å ²)	113.5	Xtriage
Anisotropy	0.036	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 116.8	EDS
Estimated twinning fraction	0.021 for -h,l,k	Xtriage
L-test for twinning	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 13609 reflections	Xtriage
F_o, F_c correlation	0.77	EDS
Total number of atoms	5552	wwPDB-VP
Average B, all atoms (Å ²)	126.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.34% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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	0.48	0/1284	0.66	0/1775
1	B	0.45	0/1275	0.70	2/1762 (0.1%)
2	C	0.56	0/1529	0.69	2/2105 (0.1%)
2	D	0.54	0/1589	0.67	3/2187 (0.1%)
All	All	0.51	0/5677	0.68	7/7829 (0.1%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	197	ARG	NE-CZ-NH1	-5.99	117.31	120.30
2	C	102	GLY	N-CA-C	-5.75	98.73	113.10
2	C	33	LEU	CA-CB-CG	5.24	127.35	115.30
1	B	20	ASP	CB-CG-OD2	5.21	122.99	118.30
2	D	33	LEU	CA-CB-CG	5.06	126.94	115.30
2	D	4	LEU	CA-CB-CG	5.06	126.93	115.30
1	B	71	LEU	N-CA-C	-5.01	97.47	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1258	0	880	66	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1253	0	883	45	0
2	C	1491	0	1179	87	0
2	D	1550	0	1309	113	0
All	All	5552	0	4251	304	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 31.

All (304) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:33:THR:HB	2:C:31:SER:O	1.40	1.17
1:A:39:VAL:CG2	1:A:44:ARG:HG3	1.76	1.16
2:C:51:ILE:O	2:C:53:PRO:HD3	1.50	1.12
2:D:48:ILE:N	2:D:48:ILE:HD13	1.57	1.11
1:A:44:ARG:O	1:A:44:ARG:HG2	1.49	1.07
2:D:34:ILE:O	2:D:50:VAL:HG13	1.62	1.00
2:C:47:TRP:HZ2	2:C:50:VAL:HG23	1.26	0.98
2:D:135:GLY:O	2:D:136:SER:OG	1.80	0.98
2:D:48:ILE:HG22	2:D:49:GLY:H	1.30	0.95
1:A:68:THR:HB	1:A:82:LEU:HD11	1.48	0.94
1:A:39:VAL:HG21	1:A:44:ARG:HG3	1.44	0.93
1:B:73:ASP:HA	2:D:186:TYR:CZ	2.03	0.92
2:D:48:ILE:N	2:D:48:ILE:CD1	2.30	0.92
1:A:47:VAL:CG2	1:A:68:THR:HG21	2.04	0.88
1:A:39:VAL:HG22	1:A:44:ARG:HG3	1.53	0.87
1:A:38:TRP:CD1	1:A:82:LEU:HD22	2.11	0.86
1:B:32:GLU:OE1	2:D:33:LEU:HD22	1.75	0.86
2:D:37:ILE:HD11	2:D:105:TRP:HZ3	1.41	0.85
2:C:35:HIS:HB3	2:C:47:TRP:HE1	1.42	0.84
1:A:115:ALA:CB	1:A:144:PRO:HD3	2.06	0.84
2:C:216:THR:O	2:C:219:PHE:CE1	2.30	0.84
2:C:101:TYR:HA	2:C:170:HIS:CE1	2.13	0.83
2:D:47:TRP:C	2:D:48:ILE:HD13	1.99	0.83
2:C:169:LEU:HD23	2:C:187:ALA:HB2	1.59	0.82
2:D:185:LYS:O	2:D:186:TYR:HB2	1.80	0.81
1:B:22:PRO:HA	1:B:81:ALA:HA	1.59	0.81
1:B:148:ILE:O	1:B:148:ILE:HG13	1.81	0.80
2:D:47:TRP:CE2	2:D:48:ILE:O	2.34	0.80
2:C:140:LEU:HD23	2:C:161:ALA:HA	1.64	0.79
2:C:66:ASP:OD1	2:C:66:ASP:N	2.16	0.78
2:C:47:TRP:CZ2	2:C:50:VAL:HG23	2.15	0.77
2:D:230:TRP:HD1	2:D:230:TRP:H	1.29	0.77

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:C:216:THR:O	2:C:219:PHE:CD1	2.38	0.76
1:B:146:PRO:HB2	1:B:194:ARG:O	1.86	0.76
2:D:48:ILE:HG22	2:D:49:GLY:N	2.00	0.76
2:D:37:ILE:HD11	2:D:105:TRP:CZ3	2.20	0.76
1:A:118:GLY:HA3	1:A:142:TRP:NE1	2.01	0.76
2:C:225:GLN:NE2	2:C:232:HIS:HB3	2.02	0.74
2:D:47:TRP:CD2	2:D:48:ILE:O	2.40	0.74
2:C:225:GLN:HG3	2:C:234:PHE:CE1	2.23	0.74
1:B:148:ILE:HD11	1:B:177:ALA:HB3	1.67	0.74
2:D:66:ASP:OD1	2:D:66:ASP:N	2.19	0.73
1:A:39:VAL:HG21	1:A:44:ARG:CG	2.18	0.73
2:C:219:PHE:HA	2:C:240:LEU:HD23	1.70	0.72
2:C:51:ILE:HG23	2:C:51:ILE:O	1.90	0.71
1:A:23:CYS:SG	1:A:97:CYS:CB	2.78	0.71
1:A:38:TRP:NE1	1:A:82:LEU:HB2	2.05	0.71
1:B:118:GLY:HA3	1:B:142:TRP:NE1	2.05	0.71
2:C:142:GLN:NE2	2:C:222:TYR:O	2.24	0.70
1:A:115:ALA:HB1	1:A:144:PRO:HD3	1.72	0.70
2:C:30:THR:HG22	2:C:54:ARG:CB	2.22	0.70
1:B:73:ASP:HA	2:D:186:TYR:CE1	2.25	0.70
2:D:16:THR:O	2:D:86:LEU:HD12	1.89	0.70
2:D:101:TYR:HD1	2:D:101:TYR:N	1.90	0.69
1:A:44:ARG:CG	1:A:44:ARG:O	2.30	0.69
2:D:230:TRP:HB2	2:D:231:PRO:HA	1.74	0.69
2:D:35:HIS:HA	2:D:50:VAL:HG22	1.75	0.69
1:A:2:PHE:CG	1:A:27:PRO:HD2	2.26	0.69
2:D:230:TRP:HA	2:D:231:PRO:C	2.12	0.69
1:B:118:GLY:HA3	1:B:142:TRP:CD1	2.28	0.69
2:D:211:ILE:HD11	2:D:214:VAL:HG22	1.75	0.69
2:C:13:ARG:O	2:C:16:THR:HG22	1.93	0.69
1:B:24:HIS:HA	1:B:78:GLY:O	1.93	0.68
1:A:47:VAL:HG21	1:A:68:THR:HG21	1.73	0.68
1:B:98:TYR:O	1:B:99:PHE:HD1	1.76	0.68
2:D:230:TRP:CD1	2:D:230:TRP:N	2.62	0.68
2:D:169:LEU:HA	2:D:225:GLN:O	1.94	0.68
2:D:13:ARG:O	2:D:16:THR:HB	1.93	0.68
2:C:216:THR:O	2:C:219:PHE:HE1	1.76	0.67
2:D:219:PHE:HA	2:D:240:LEU:HD23	1.77	0.66
2:D:89:ASP:OD1	2:D:89:ASP:N	2.28	0.66
2:D:61:ASN:OD1	2:D:231:PRO:HB3	1.94	0.66
2:C:89:ASP:OD1	2:C:89:ASP:N	2.28	0.66
1:A:38:TRP:HD1	1:A:82:LEU:HD22	1.58	0.65

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:144:PRO:HG3	1:A:197:LEU:HD23	1.77	0.65
2:D:2:VAL:HG23	2:D:26:GLY:HA3	1.79	0.65
1:B:98:TYR:C	1:B:99:PHE:HD1	2.00	0.65
1:B:150:TRP:HD1	1:B:160:THR:HG22	1.62	0.65
2:D:102:GLY:O	2:D:105:TRP:NE1	2.30	0.65
1:A:38:TRP:HE1	1:A:82:LEU:HB2	1.62	0.65
2:D:101:TYR:CD1	2:D:101:TYR:N	2.64	0.64
2:D:182:LEU:HG	2:D:191:ILE:HG13	1.79	0.64
2:D:86:LEU:O	2:D:87:THR:HG22	1.98	0.64
1:A:115:ALA:HB2	1:A:144:PRO:HD3	1.79	0.64
1:B:145:GLN:HB2	1:B:173:TYR:CE2	2.33	0.64
2:C:169:LEU:HD23	2:C:187:ALA:CB	2.28	0.63
2:C:169:LEU:HA	2:C:225:GLN:O	1.97	0.63
2:C:217:GLU:CG	2:C:217:GLU:O	2.47	0.62
2:D:225:GLN:NE2	2:D:232:HIS:HB3	2.15	0.62
2:D:230:TRP:HB3	2:D:232:HIS:HD2	1.65	0.62
1:B:118:GLY:HA3	1:B:142:TRP:CE2	2.35	0.61
2:D:185:LYS:CB	2:D:191:ILE:HD11	2.31	0.61
1:B:21:LEU:O	1:B:82:LEU:N	2.28	0.61
2:C:197:ARG:HH12	2:C:215:GLU:CG	2.13	0.61
2:D:157:LEU:HD22	2:D:238:THR:HG21	1.82	0.61
2:D:30:THR:HA	2:D:53:PRO:HB2	1.82	0.61
2:D:197:ARG:HH12	2:D:215:GLU:CG	2.14	0.61
1:B:32:GLU:OE1	2:D:33:LEU:CD2	2.46	0.61
1:A:61:SER:OG	1:A:64:TYR:HD2	1.84	0.61
2:C:217:GLU:O	2:C:217:GLU:HG2	2.01	0.61
2:D:37:ILE:HD12	2:D:45:LEU:HD11	1.83	0.60
2:C:169:LEU:HD22	2:C:207:PHE:CD2	2.37	0.60
1:B:144:PRO:O	1:B:146:PRO:HD3	2.01	0.60
2:C:2:VAL:HG23	2:C:26:GLY:HA3	1.83	0.60
2:C:34:ILE:O	2:C:50:VAL:HG22	2.02	0.60
1:B:21:LEU:CD1	1:B:82:LEU:CB	2.80	0.60
2:C:172:TYR:O	2:C:222:TYR:HA	2.02	0.59
2:D:84:SER:O	2:D:85:SER:HB3	2.00	0.59
2:D:172:TYR:HD1	2:D:182:LEU:HA	1.67	0.59
1:B:140:THR:HG22	1:B:174:ALA:CB	2.32	0.59
1:A:23:CYS:HG	1:A:97:CYS:CB	2.14	0.59
1:A:47:VAL:HG23	1:A:68:THR:HG21	1.84	0.59
2:D:135:GLY:C	2:D:136:SER:HG	1.93	0.59
2:D:197:ARG:HH21	2:C:192:PHE:HE1	1.50	0.58
2:C:174:GLN:OE1	2:C:180:PRO:HG3	2.03	0.58
2:C:172:TYR:HD1	2:C:182:LEU:HA	1.68	0.58

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:128:TYR:HD1	1:A:129:LYS:N	2.02	0.58
2:D:142:GLN:HG2	2:D:238:THR:OG1	2.03	0.58
2:D:162:SER:OG	2:D:163:GLN:N	2.36	0.58
1:A:40:SER:O	1:A:44:ARG:HB2	2.04	0.58
1:B:20:ASP:HA	1:B:82:LEU:O	2.02	0.58
2:D:172:TYR:HE1	2:D:182:LEU:HD13	1.69	0.58
1:B:140:THR:HG22	1:B:174:ALA:HB1	1.86	0.57
1:A:50:TYR:HB2	1:A:70:ILE:HD13	1.86	0.57
1:A:73:ASP:OD1	1:A:73:ASP:N	2.36	0.57
2:C:37:ILE:HD12	2:C:45:LEU:HD11	1.87	0.57
2:C:197:ARG:HH12	2:C:215:GLU:HG2	1.68	0.57
2:C:138:ILE:HD11	2:C:226:GLN:HG3	1.87	0.56
2:D:225:GLN:HG3	2:D:234:PHE:CE1	2.40	0.56
2:C:215:GLU:HB2	2:C:218:ASP:CG	2.26	0.56
1:A:118:GLY:HA3	1:A:142:TRP:CD1	2.41	0.56
2:C:215:GLU:HB2	2:C:218:ASP:OD1	2.05	0.56
2:D:225:GLN:HG2	2:D:233:THR:O	2.05	0.56
2:D:230:TRP:HB2	2:D:231:PRO:CA	2.35	0.56
1:A:50:TYR:HB2	1:A:70:ILE:CD1	2.36	0.55
2:C:50:VAL:CG1	2:C:51:ILE:N	2.69	0.55
1:A:67:ARG:NH2	1:A:91:ASP:OD2	2.39	0.55
1:A:2:PHE:CD1	1:A:27:PRO:HD2	2.41	0.55
2:D:144:PRO:HD2	2:D:147:LEU:HD11	1.88	0.55
2:D:14:PRO:CG	2:D:115:ALA:HA	2.37	0.55
1:B:23:CYS:HG	1:B:97:CYS:CB	2.20	0.55
2:C:51:ILE:CG2	2:C:51:ILE:O	2.55	0.54
2:D:167:ASN:OD1	2:D:186:TYR:HD1	1.91	0.54
2:D:216:THR:HA	2:D:219:PHE:HE2	1.73	0.54
1:A:2:PHE:CG	1:A:3:SER:N	2.75	0.54
2:C:225:GLN:HG2	2:C:233:THR:O	2.06	0.54
2:D:18:VAL:HG23	2:D:86:LEU:HD11	1.89	0.54
1:A:116:ALA:O	1:A:143:TYR:O	2.25	0.54
2:D:197:ARG:HH22	2:D:215:GLU:HG3	1.72	0.54
2:C:31:SER:HB2	2:C:32:TYR:CD1	2.43	0.54
2:C:162:SER:OG	2:C:163:GLN:N	2.38	0.54
2:C:138:ILE:HD11	2:C:226:GLN:NE2	2.23	0.54
1:A:145:GLN:CB	1:A:173:TYR:CE2	2.91	0.54
2:C:35:HIS:HB3	2:C:47:TRP:NE1	2.20	0.53
2:D:186:TYR:O	2:D:187:ALA:HB3	2.08	0.53
2:D:13:ARG:O	2:D:16:THR:CB	2.56	0.53
1:B:22:PRO:HB3	1:B:81:ALA:HB2	1.91	0.53
2:D:230:TRP:CA	2:D:231:PRO:C	2.78	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:D:231:PRO:HG2	2:D:231:PRO:O	2.08	0.52
2:D:51:ILE:O	2:D:53:PRO:HD3	2.09	0.52
1:B:167:ALA:HA	1:B:172:LEU:O	2.08	0.52
2:C:226:GLN:HG2	2:C:228:ASN:H	1.73	0.52
2:C:195:PRO:HG2	2:C:198:PHE:CE2	2.44	0.52
1:A:91:ASP:N	1:A:91:ASP:OD1	2.42	0.52
1:B:128:TYR:HD1	1:B:129:LYS:N	2.08	0.52
1:B:73:ASP:OD1	1:B:74:GLY:N	2.42	0.52
2:C:218:ASP:N	2:C:218:ASP:OD1	2.41	0.52
2:C:47:TRP:HZ2	2:C:50:VAL:CG2	2.09	0.52
2:D:48:ILE:CG2	2:D:49:GLY:N	2.67	0.52
2:D:195:PRO:HG2	2:D:198:PHE:CE2	2.44	0.52
1:B:91:ASP:OD1	1:B:91:ASP:N	2.42	0.52
2:C:53:PRO:O	2:C:54:ARG:CB	2.58	0.51
2:C:31:SER:HB2	2:C:32:TYR:HD1	1.76	0.51
1:B:22:PRO:CA	1:B:81:ALA:HA	2.34	0.51
2:D:197:ARG:HH22	2:D:215:GLU:CG	2.24	0.50
2:D:50:VAL:CG1	2:D:51:ILE:N	2.74	0.50
1:A:50:TYR:CD1	1:A:50:TYR:C	2.84	0.50
2:C:171:TRP:CZ3	2:C:224:CYS:HB3	2.46	0.50
2:C:169:LEU:CD2	2:C:187:ALA:HB2	2.34	0.50
1:A:47:VAL:O	1:A:60:GLN:HG2	2.12	0.50
2:C:216:THR:O	2:C:219:PHE:HD1	1.92	0.50
2:C:225:GLN:HG3	2:C:234:PHE:CD1	2.46	0.50
2:D:86:LEU:C	2:D:87:THR:CG2	2.80	0.50
2:C:29:PHE:O	2:C:30:THR:C	2.48	0.50
2:D:48:ILE:CG2	2:D:49:GLY:H	2.04	0.50
1:A:114:VAL:O	1:A:143:TYR:CE2	2.65	0.49
2:D:135:GLY:O	2:D:136:SER:CB	2.58	0.49
2:D:172:TYR:CE1	2:D:182:LEU:HD13	2.47	0.49
2:C:28:ALA:O	2:C:31:SER:OG	2.30	0.49
2:C:138:ILE:CD1	2:C:226:GLN:HG3	2.42	0.49
2:C:225:GLN:HE22	2:C:232:HIS:HB3	1.72	0.49
1:A:167:ALA:HA	1:A:172:LEU:O	2.13	0.49
2:D:50:VAL:HG12	2:D:51:ILE:N	2.28	0.49
1:A:26:PHE:HA	1:A:27:PRO:C	2.33	0.49
2:D:47:TRP:HZ2	2:D:50:VAL:HG23	1.77	0.48
2:C:174:GLN:CD	2:C:180:PRO:HG3	2.33	0.48
2:D:231:PRO:O	2:D:231:PRO:CG	2.61	0.48
2:C:197:ARG:HG3	2:C:198:PHE:N	2.28	0.48
1:B:26:PHE:HA	1:B:27:PRO:C	2.33	0.48
1:A:33:THR:HG21	2:C:31:SER:HA	1.95	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:C:105:TRP:CD1	2:C:105:TRP:N	2.82	0.48
2:D:215:GLU:HB2	2:D:218:ASP:CG	2.34	0.48
2:C:47:TRP:O	2:C:61:ASN:ND2	2.47	0.48
1:A:118:GLY:HA3	1:A:142:TRP:CE2	2.49	0.48
2:C:172:TYR:CD1	2:C:182:LEU:HA	2.49	0.48
2:D:197:ARG:NH1	2:D:218:ASP:OD2	2.47	0.48
1:A:23:CYS:CB	1:A:97:CYS:HG	2.26	0.47
1:A:145:GLN:CB	1:A:173:TYR:HE2	2.26	0.47
1:A:115:ALA:HB1	1:A:144:PRO:CD	2.41	0.47
1:A:43:LEU:N	1:A:44:ARG:HA	2.29	0.47
1:A:38:TRP:CE3	1:A:97:CYS:HB2	2.50	0.47
2:C:32:TYR:N	2:C:32:TYR:CD1	2.83	0.47
1:B:145:GLN:H	1:B:173:TYR:HE2	1.62	0.47
1:A:152:ASN:CG	1:A:153:ASN:H	2.18	0.47
2:C:55:SER:O	2:C:55:SER:OG	2.30	0.47
1:A:209:ALA:HB3	1:A:212:PHE:HD2	1.80	0.47
2:D:169:LEU:HD22	2:D:207:PHE:CD2	2.50	0.46
1:B:150:TRP:HB2	1:B:158:ILE:HB	1.96	0.46
1:A:61:SER:HG	1:A:64:TYR:HD2	1.62	0.46
2:C:104:TYR:C	2:C:105:TRP:CD1	2.88	0.46
2:D:197:ARG:HD3	2:D:212:ASN:O	2.15	0.46
2:C:226:GLN:NE2	2:C:232:HIS:HA	2.31	0.46
1:B:150:TRP:CD1	1:B:160:THR:HG22	2.48	0.46
2:D:195:PRO:HG2	2:D:198:PHE:HE2	1.81	0.46
1:B:44:ARG:O	1:B:44:ARG:HG3	2.15	0.46
2:D:142:GLN:OE1	2:D:235:GLY:HA3	2.16	0.45
1:B:38:TRP:CE3	1:B:97:CYS:HB2	2.51	0.45
1:A:143:TYR:CG	1:A:144:PRO:HA	2.52	0.45
2:D:149:VAL:HG11	2:D:214:VAL:HG21	1.98	0.45
2:D:172:TYR:CD1	2:D:182:LEU:HA	2.48	0.45
1:B:145:GLN:HB2	1:B:173:TYR:CD2	2.52	0.45
2:D:4:LEU:HD21	2:D:104:TYR:O	2.16	0.45
2:D:34:ILE:O	2:D:50:VAL:CG1	2.48	0.45
2:D:225:GLN:HE21	2:D:232:HIS:HB3	1.79	0.45
1:B:21:LEU:HG	1:B:82:LEU:CB	2.47	0.45
1:B:98:TYR:C	1:B:99:PHE:CD1	2.85	0.45
2:D:144:PRO:O	2:D:238:THR:HG23	2.16	0.45
2:D:4:LEU:HD21	2:D:104:TYR:C	2.36	0.45
1:B:74:GLY:O	1:B:76:THR:N	2.48	0.45
1:A:142:TRP:O	1:A:173:TYR:HB2	2.16	0.45
1:A:20:ASP:HA	1:A:82:LEU:O	2.16	0.44
1:A:114:VAL:O	1:A:143:TYR:CD2	2.70	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:D:230:TRP:HB3	2:D:232:HIS:CD2	2.47	0.44
1:A:2:PHE:CZ	1:A:26:PHE:O	2.70	0.44
1:B:119:SER:N	1:B:140:THR:O	2.38	0.44
2:C:195:PRO:HG2	2:C:198:PHE:HE2	1.81	0.44
2:D:218:ASP:O	2:D:222:TYR:OH	2.21	0.44
1:B:40:SER:O	1:B:44:ARG:HB2	2.16	0.44
2:D:93:VAL:HG13	2:D:110:LEU:HD12	2.00	0.44
2:D:98:ARG:HA	2:D:99:SER:HA	1.88	0.44
1:A:48:ASN:CG	1:A:60:GLN:HE21	2.21	0.44
2:C:226:GLN:OE1	2:C:233:THR:N	2.49	0.44
1:A:98:TYR:O	1:A:99:PHE:CD1	2.71	0.44
2:D:48:ILE:HD12	2:D:64:PHE:CE2	2.53	0.44
2:D:97:ALA:HB1	2:D:102:GLY:HA2	1.98	0.44
2:D:157:LEU:HD12	2:D:209:LEU:HD23	2.00	0.44
1:B:21:LEU:HD12	1:B:82:LEU:CB	2.48	0.43
2:C:32:TYR:N	2:C:32:TYR:HD1	2.16	0.43
2:C:93:VAL:HG13	2:C:110:LEU:HD12	2.00	0.43
2:C:225:GLN:HE21	2:C:232:HIS:HB3	1.83	0.43
2:C:197:ARG:HH22	2:C:215:GLU:CG	2.32	0.43
1:A:121:LEU:HD12	1:A:203:THR:C	2.39	0.43
1:B:158:ILE:CG2	1:B:159:PRO:HD2	2.49	0.43
2:D:230:TRP:HA	2:D:231:PRO:O	2.18	0.43
2:D:215:GLU:HB2	2:D:218:ASP:OD2	2.18	0.43
2:C:21:SER:HA	2:C:80:TYR:HD1	1.83	0.43
1:A:41:SER:HB2	1:A:94:LYS:O	2.19	0.43
2:D:21:SER:HA	2:D:80:TYR:HD1	1.83	0.43
2:D:53:PRO:O	2:D:54:ARG:CB	2.65	0.43
2:D:140:LEU:O	2:D:235:GLY:HA2	2.19	0.43
2:D:101:TYR:O	2:D:182:LEU:HD22	2.19	0.43
2:C:216:THR:C	2:C:219:PHE:HE1	2.22	0.42
2:D:166:SER:OG	2:D:167:ASN:N	2.52	0.42
2:C:216:THR:HA	2:C:219:PHE:HE1	1.83	0.42
1:B:21:LEU:HB2	1:B:110:VAL:HG11	2.00	0.42
2:D:197:ARG:HG3	2:D:198:PHE:N	2.34	0.42
2:D:14:PRO:HG2	2:D:115:ALA:HA	2.01	0.42
1:A:38:TRP:HE1	1:A:82:LEU:HD13	1.85	0.42
2:D:159:CYS:HB2	2:D:171:TRP:CH2	2.55	0.42
2:C:166:SER:OG	2:C:167:ASN:N	2.52	0.42
1:A:39:VAL:HG22	1:A:40:SER:O	2.20	0.42
2:C:186:TYR:C	2:C:188:SER:H	2.23	0.42
1:A:58:ASP:C	1:A:60:GLN:N	2.72	0.41
2:C:169:LEU:HB3	2:C:187:ALA:HB2	2.01	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:98:TYR:O	1:A:99:PHE:HD1	2.03	0.41
2:D:197:ARG:HH12	2:D:215:GLU:HG2	1.85	0.41
2:D:100:ASP:O	2:D:170:HIS:NE2	2.54	0.41
2:C:99:SER:O	2:C:100:ASP:CB	2.68	0.41
2:C:104:TYR:C	2:C:105:TRP:HD1	2.23	0.41
2:D:48:ILE:HD12	2:D:64:PHE:CD2	2.55	0.41
2:C:140:LEU:CD2	2:C:161:ALA:HB2	2.51	0.41
2:D:14:PRO:HG3	2:D:115:ALA:HA	2.02	0.41
1:A:144:PRO:HG3	1:A:197:LEU:CD2	2.50	0.41
2:D:197:ARG:HH12	2:D:215:GLU:HB2	1.85	0.41
2:D:168:ASN:ND2	2:D:228:ASN:OD1	2.54	0.41
2:C:186:TYR:O	2:C:188:SER:N	2.53	0.41
2:D:86:LEU:O	2:D:87:THR:CG2	2.67	0.40
2:D:216:THR:HA	2:D:219:PHE:CE2	2.55	0.40
2:C:37:ILE:HG13	2:C:95:PHE:HB2	2.03	0.40
1:A:22:PRO:HA	1:A:81:ALA:HA	2.02	0.40
2:C:197:ARG:NH1	2:C:215:GLU:HG2	2.36	0.40
2:D:47:TRP:C	2:D:48:ILE:CD1	2.78	0.40
2:C:169:LEU:HD22	2:C:207:PHE:CG	2.56	0.40
1:B:128:TYR:CD1	1:B:129:LYS:N	2.88	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	203/220 (92%)	195 (96%)	7 (3%)	1 (0%)	38	87
1	B	204/220 (93%)	194 (95%)	8 (4%)	2 (1%)	22	78
2	C	216/254 (85%)	203 (94%)	13 (6%)	0	100	100
2	D	220/254 (87%)	206 (94%)	14 (6%)	0	100	100
All	All	843/948 (89%)	798 (95%)	42 (5%)	3 (0%)	43	90

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	146	PRO
1	A	146	PRO
1	B	74	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	69/178 (39%)	51 (74%)	18 (26%)	1	5
1	B	71/178 (40%)	48 (68%)	23 (32%)	0	3
2	C	127/201 (63%)	100 (79%)	27 (21%)	1	9
2	D	144/201 (72%)	114 (79%)	30 (21%)	2	10
All	All	411/758 (54%)	313 (76%)	98 (24%)	1	6

All (98) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	23	CYS
1	A	25	LEU
1	A	38	TRP
1	A	41	SER
1	A	44	ARG
1	A	47	VAL
1	A	50	TYR
1	A	67	ARG
1	A	70	ILE
1	A	73	ASP
1	A	90	SER
1	A	91	ASP
1	A	105	TYR
1	A	110	VAL
1	A	117	LEU
1	A	128	TYR
1	A	137	CYS
1	A	191	CYS
1	B	1	GLN
1	B	14	MET

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Mol	Chain	Res	Type
1	B	21	LEU
1	B	23	CYS
1	B	38	TRP
1	B	58	ASP
1	B	59	ARG
1	B	70	ILE
1	B	76	THR
1	B	90	SER
1	B	91	ASP
1	B	92	SER
1	B	105	TYR
1	B	110	VAL
1	B	117	LEU
1	B	120	ASP
1	B	128	TYR
1	B	137	CYS
1	B	148	ILE
1	B	172	LEU
1	B	191	CYS
1	B	193	ILE
1	B	195	SER
2	D	2	VAL
2	D	4	LEU
2	D	6	GLN
2	D	11	VAL
2	D	17	SER
2	D	31	SER
2	D	33	LEU
2	D	34	ILE
2	D	48	ILE
2	D	66	ASP
2	D	69	THR
2	D	70	LEU
2	D	77	SER
2	D	84	SER
2	D	87	THR
2	D	89	ASP
2	D	93	VAL
2	D	101	TYR
2	D	110	LEU
2	D	113	VAL
2	D	142	GLN

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Mol	Chain	Res	Type
2	D	149	VAL
2	D	150	THR
2	D	176	SER
2	D	205	THR
2	D	211	ILE
2	D	225	GLN
2	D	230	TRP
2	D	231	PRO
2	D	233	THR
2	C	2	VAL
2	C	6	GLN
2	C	11	VAL
2	C	27	TYR
2	C	30	THR
2	C	31	SER
2	C	32	TYR
2	C	33	LEU
2	C	55	SER
2	C	57	ASP
2	C	66	ASP
2	C	70	LEU
2	C	77	SER
2	C	83	LEU
2	C	89	ASP
2	C	93	VAL
2	C	110	LEU
2	C	113	VAL
2	C	141	THR
2	C	142	GLN
2	C	150	THR
2	C	205	THR
2	C	216	THR
2	C	225	GLN
2	C	226	GLN
2	C	230	TRP
2	C	233	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	60	GLN
2	D	59	HIS

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Mol	Chain	Res	Type
2	D	225	GLN
2	D	232	HIS
2	C	225	GLN

5.3.3 RNA ⓘ

There are no RNA chains 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	207/220 (94%)	0.25	7 (3%)	43 20	44, 118, 244, 500	0
1	B	208/220 (94%)	0.17	1 (0%)	88 65	56, 124, 192, 409	0
2	C	220/254 (86%)	0.26	6 (2%)	52 25	51, 134, 208, 257	0
2	D	224/254 (88%)	0.12	1 (0%)	90 71	46, 112, 180, 236	0
All	All	859/948 (90%)	0.20	15 (1%)	67 35	44, 122, 207, 500	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	31	ALA	5.3
1	A	127	GLY	5.3
2	C	147	LEU	4.0
2	C	94	TYR	3.9
1	A	77	ALA	2.7
2	C	138	ILE	2.5
1	A	26	PHE	2.4
1	A	181	MET	2.4
1	B	34	MET	2.3
2	C	151	PRO	2.2
1	A	25	LEU	2.2
2	D	222	TYR	2.1
2	C	243	LYS	2.1
2	C	8	GLY	2.0
1	A	131	GLY	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.