



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 03:24 PM UTC

PDB ID : 6M3P / pdb_00006m3p
Title : Crystal structure of AnkG/beta2-spectrin complex
Authors : Li, J.; Chen, K.; Zhu, R.; Zhang, M.
Deposited on : 2020-03-04
Resolution : 3.31 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	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

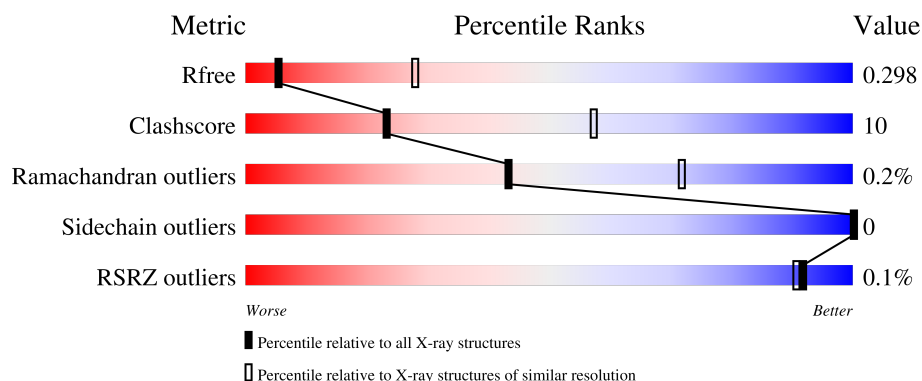
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.31 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}	180053	1153 (3.34-3.30)
Clashscore	190562	1193 (3.34-3.30)
Ramachandran outliers	187476	1172 (3.34-3.30)
Sidechain outliers	187428	1171 (3.34-3.30)
RSRZ outliers	180081	1153 (3.34-3.30)

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 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 $\leq 5\%$. 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	C	491	 65% 21% 14%
1	E	491	 66% 21% 13%
2	A	320	 79% 15% 6%
2	B	320	 79% 15% 6%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10198 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 Ankyrin-3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	E	426	Total	C	N	O	S	0	0	0
			2972	1893	502	557	20			
1	C	422	Total	C	N	O	S	0	0	0
			2946	1875	498	553	20			

- Molecule 2 is a protein called Spectrin beta chain, non-erythrocytic 1.

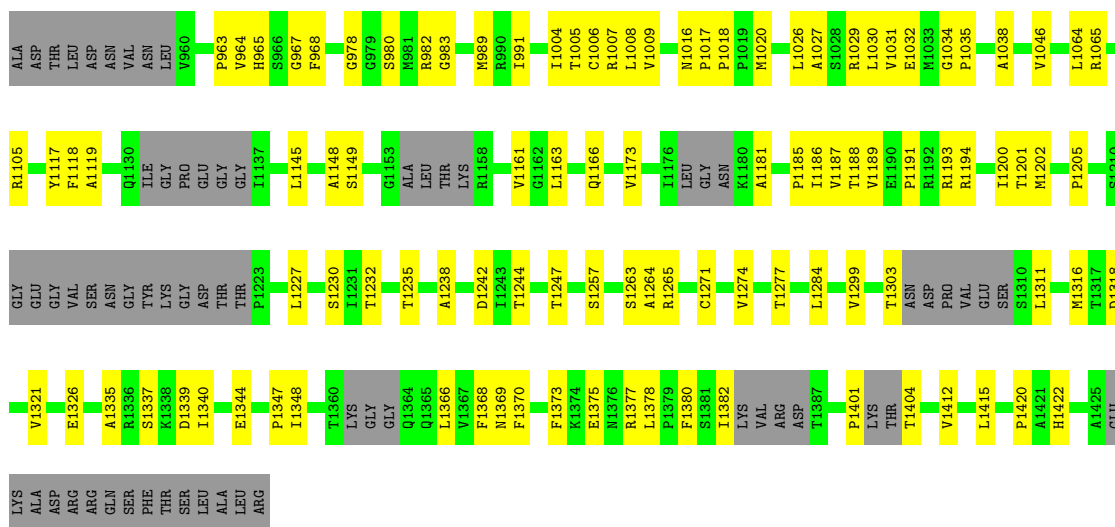
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	301	Total	C	N	O	S	0	0	0
			2140	1326	375	434	5			
2	B	301	Total	C	N	O	S	0	0	0
			2140	1326	375	434	5			

3 Residue-property plots

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.

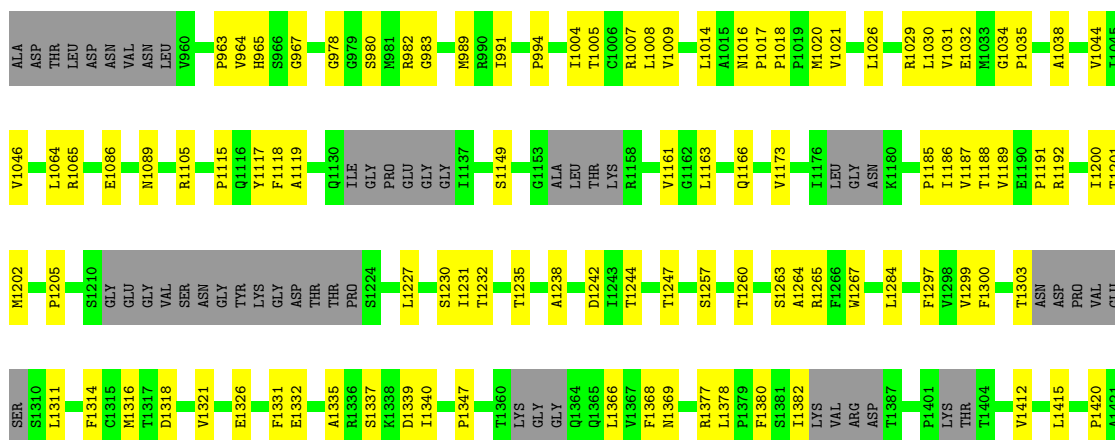
• Molecule 1: Ankyrin-3

Chain E: 



• Molecule 1: Ankyrin-3

Chain C: 



H1422	ALA
	LYS
	LYS
	ALA
	GLU
	LYS
	ALA
	ASP
	ARG
	GLN
	SER
	PHE
	THR
	SER
	LEU
	ALA
	LEU
	ARG

- Molecule 2: Spectrin beta chain, non-erythrocytic 1

Chain A:

79%

15%

6%

ALA	HIS	LYS	ALA	GLN	LYS	TYR	PHE	ASP	A11	N25	E28	GLU	LYS	ALA	LYS	D33	M40	A55	H73	P74	E75	R78	K85	L89	Y90	R100	L104	D105	E106	L110	N114	I124	R127	S133	T157	G161	D162	E163	R164	V165	D166
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T167	D173	I176	N177	H180	A183	I186	W189	W197	R207	L211	L217	F220	Y221	I227	F228	I231	Q232	E241	LEU	GLY	ARG	ASP	GLN	R247	Q269	V270	Q274	L280	D287	D291	T292	Q293	E296	L300	T320
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- Molecule 2: Spectrin beta chain, non-erythrocytic 1

Chain B:

%

79%

15%

6%

ALA	HIS	LYS	ALA	GLN	TYR	PHE	ASP	A11	N25	N26	S27	E28	GLU	LYS	ALA	LYS	D33	E34	M40	A55	E56	T57	L61	V69	S72	H73	P74	E75	S76	K85	L89	Y90	L96	R100	L104	L110	N114	R115	I124	L136	G161
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H162	E163	R164	V165	D166	T167	D173	I176	N177	H180	A183	I186	W189	W197	L200	L217	F220	Y221	I227	I231	E241	LEU	GLY	ARG	ASP	GLN	R247	Q269	V270	L280	D287	D291	G312	R313	R316	T320
------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	------	-----	-----	-----	-----	-----	------	------	------	------	------	------	------	------	------	------

4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	66.28Å 66.31Å 160.93Å 99.47° 99.78° 100.87°	Depositor
Resolution (Å)	44.99 – 3.31 44.99 – 3.31	Depositor EDS
% Data completeness (in resolution range)	95.2 (44.99-3.31) 95.1 (44.99-3.31)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.35 (at 3.32Å)	Xtriage
Refinement program	PHENIX v1.15.2	Depositor
R, R_{free}	0.255 , 0.295 0.254 , 0.298	Depositor DCC
R_{free} test set	1733 reflections (4.49%)	wwPDB-VP
Wilson B-factor (Å ²)	98.7	Xtriage
Anisotropy	0.459	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 148.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.387 for k,h,-h-k-l 0.026 for -k,-h,-l 0.025 for -h,-k,h+k+l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10198	wwPDB-VP
Average B, all atoms (Å ²)	107.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.56% 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 [i](#)

5.1 Standard geometry [i](#)

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	C	0.25	0/3006	0.51	0/4126
1	E	0.25	0/3033	0.50	0/4164
2	A	0.23	0/2172	0.52	0/2970
2	B	0.23	0/2172	0.49	0/2970
All	All	0.24	0/10383	0.51	0/14230

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	C	2946	0	2680	64	0
1	E	2972	0	2708	70	0
2	A	2140	0	1817	34	0
2	B	2140	0	1817	34	0
All	All	10198	0	9022	198	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 10.

All (198) 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:A:217:LEU:HA	2:A:280:LEU:HD21	1.58	0.85
1:C:1161:VAL:HG23	1:C:1191:PRO:HG2	1.57	0.85
2:B:217:LEU:HA	2:B:280:LEU:HD21	1.58	0.85
1:C:1035:PRO:HB2	1:C:1038:ALA:HB2	1.60	0.84
1:E:1035:PRO:HB2	1:E:1038:ALA:HB2	1.58	0.84
1:E:1326:GLU:OE2	1:E:1377:ARG:NH1	2.10	0.84
1:E:1161:VAL:HG23	1:E:1191:PRO:HG2	1.59	0.83
1:C:1326:GLU:OE2	1:C:1377:ARG:NH1	2.12	0.81
1:E:1188:THR:HG23	1:E:1265:ARG:HG3	1.63	0.80
1:C:1008:LEU:HD22	1:C:1029:ARG:HB2	1.65	0.77
1:C:1173:VAL:HG22	1:C:1284:LEU:HD11	1.67	0.77
1:E:1008:LEU:HD22	1:E:1029:ARG:HB2	1.67	0.76
1:E:1173:VAL:HG22	1:E:1284:LEU:HD11	1.68	0.76
1:C:1149:SER:HB3	1:C:1201:THR:HB	1.72	0.72
1:C:1188:THR:HG23	1:C:1265:ARG:HG3	1.71	0.70
2:A:100:ARG:HH21	2:A:100:ARG:HG3	1.59	0.66
1:E:1149:SER:HB3	1:E:1201:THR:HB	1.78	0.64
1:C:1185:PRO:HD2	1:C:1205:PRO:HG2	1.81	0.63
1:E:1311:LEU:HD12	1:E:1382:ILE:HD12	1.82	0.62
1:E:1189:VAL:HG21	1:E:1263:SER:HB2	1.81	0.62
2:B:176:ILE:HG12	2:B:183:ALA:HB1	1.82	0.61
1:E:983:GLY:N	1:E:989:MET:O	2.28	0.61
1:C:1303:THR:HG21	1:C:1422:HIS:H	1.64	0.61
1:E:1271:CYS:SG	1:E:1277:THR:HG22	2.42	0.60
1:C:1311:LEU:HD12	1:C:1382:ILE:HD12	1.84	0.60
1:E:1401:PRO:HG2	1:E:1404:THR:HG22	1.83	0.59
2:B:100:ARG:HH21	2:B:100:ARG:HG3	1.68	0.59
2:B:180:HIS:HB3	2:B:183:ALA:HB2	1.83	0.58
1:E:1299:VAL:HB	1:E:1335:ALA:HB3	1.83	0.58
2:A:180:HIS:HB3	2:A:183:ALA:HB2	1.85	0.58
2:B:173:ASP:O	2:B:177:ASN:HB2	2.02	0.58
2:B:287:ASP:N	2:B:287:ASP:OD1	2.36	0.58
1:E:1230:SER:O	1:E:1264:ALA:HB1	2.04	0.58
2:A:227:ILE:HD11	2:A:269:GLN:HB3	1.84	0.58
1:C:1316:MET:HE1	1:C:1321:VAL:HG22	1.85	0.58
1:E:1347:PRO:HA	1:E:1369:ASN:HA	1.86	0.57
1:C:1200:ILE:HG22	1:C:1260:THR:O	2.04	0.57
2:A:161:GLY:O	2:A:165:VAL:HG23	2.05	0.56
1:C:964:VAL:HG22	1:C:978:GLY:O	2.05	0.56
1:E:1316:MET:HE1	1:E:1321:VAL:HG22	1.88	0.56
1:E:991:ILE:HG12	1:E:1046:VAL:HG13	1.86	0.56
1:C:1299:VAL:HB	1:C:1335:ALA:HB3	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1230:SER:O	1:C:1264:ALA:HB1	2.05	0.55
2:B:11:ALA:HB2	2:B:61:LEU:HD21	1.89	0.55
2:B:40:MET:HG2	2:B:104:LEU:HD22	1.86	0.55
1:E:1016:ASN:O	1:E:1016:ASN:ND2	2.39	0.55
1:C:1232:THR:OG1	1:C:1238:ALA:HA	2.07	0.54
2:B:227:ILE:HD11	2:B:269:GLN:HB3	1.87	0.54
2:A:287:ASP:OD1	2:A:287:ASP:N	2.39	0.54
1:E:1064:LEU:HB2	1:E:1119:ALA:HB3	1.90	0.54
1:E:964:VAL:HG22	1:E:978:GLY:O	2.08	0.54
1:E:1005:THR:H	1:E:1034:GLY:HA3	1.73	0.54
2:B:163:GLU:O	2:B:167:THR:OG1	2.24	0.54
1:E:1185:PRO:HD2	1:E:1205:PRO:HG2	1.90	0.53
2:A:100:ARG:HG3	2:A:100:ARG:NH2	2.22	0.53
1:C:1189:VAL:CG2	1:C:1263:SER:HB2	2.39	0.53
2:A:176:ILE:HG12	2:A:183:ALA:HB1	1.90	0.53
1:E:1193:ARG:HH12	2:B:313:ARG:HA	1.74	0.53
1:C:1368:PHE:CD2	1:C:1378:LEU:HB2	2.44	0.52
1:C:1016:ASN:O	1:C:1016:ASN:ND2	2.39	0.52
1:E:1148:ALA:HB1	1:E:1200:ILE:HD11	1.92	0.52
1:C:1202:MET:O	1:C:1257:SER:HA	2.08	0.52
2:B:25:MET:O	2:B:100:ARG:HD2	2.10	0.52
1:C:1005:THR:H	1:C:1034:GLY:HA3	1.73	0.52
1:C:994:PRO:HD2	1:C:1044:VAL:HG12	1.92	0.51
1:E:1032:GLU:HG3	1:E:1117:TYR:CE1	2.45	0.51
1:E:1340:ILE:HG21	1:E:1412:VAL:O	2.11	0.51
1:E:1375:GLU:OE2	1:E:1377:ARG:NH2	2.43	0.51
1:C:963:PRO:HB3	1:C:978:GLY:HA3	1.92	0.51
1:E:963:PRO:HB3	1:E:978:GLY:HA3	1.92	0.51
1:C:1007:ARG:O	1:C:1031:VAL:HA	2.10	0.51
2:A:293:GLN:NE2	2:A:296:GLU:OE1	2.41	0.51
2:B:73:HIS:HA	2:B:76:SER:HB2	1.91	0.51
1:E:1193:ARG:NH1	2:B:312:GLY:O	2.44	0.51
2:A:163:GLU:O	2:A:167:THR:OG1	2.24	0.51
1:E:1366:LEU:HD21	1:E:1380:PHE:CE1	2.47	0.50
2:A:124:ILE:HD13	2:A:197:TRP:HA	1.93	0.50
1:E:1017:PRO:HB2	1:E:1018:PRO:HD2	1.94	0.49
2:B:55:ALA:HB2	2:B:90:TYR:OH	2.12	0.49
1:C:1064:LEU:HB2	1:C:1119:ALA:HB3	1.94	0.49
1:C:1189:VAL:HG21	1:C:1263:SER:HB2	1.93	0.49
1:E:1189:VAL:CG2	1:E:1263:SER:HB2	2.41	0.49
1:C:1242:ASP:OD1	1:C:1244:THR:OG1	2.28	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:25:MET:O	2:A:100:ARG:HD2	2.12	0.49
2:A:40:MET:HG2	2:A:104:LEU:HD22	1.95	0.49
1:C:983:GLY:N	1:C:989:MET:O	2.32	0.49
2:A:73:HIS:N	2:A:74:PRO:HD2	2.28	0.49
1:C:1017:PRO:HB2	1:C:1018:PRO:HD2	1.95	0.49
2:A:55:ALA:HB2	2:A:90:TYR:OH	2.13	0.48
1:E:1227:LEU:HD23	1:E:1247:THR:HG21	1.96	0.48
2:A:114:ASN:HB2	2:A:189:TRP:CH2	2.48	0.48
1:C:1326:GLU:HB3	1:C:1331:PHE:HB2	1.93	0.48
2:B:114:ASN:HB2	2:B:189:TRP:CH2	2.49	0.48
2:B:124:ILE:HD13	2:B:197:TRP:HA	1.96	0.48
1:E:1020:MET:O	1:E:1235:THR:HG22	2.14	0.48
1:C:1347:PRO:HA	1:C:1369:ASN:HA	1.95	0.48
2:A:221:TYR:OH	2:A:291:ASP:HB3	2.14	0.48
1:E:1005:THR:HB	1:E:1034:GLY:HA2	1.96	0.48
2:A:173:ASP:O	2:A:177:ASN:HB2	2.14	0.47
2:B:85:LYS:HG3	2:B:89:LEU:HD23	1.96	0.47
1:E:1348:ILE:HD11	1:E:1368:PHE:CZ	2.49	0.47
2:A:110:LEU:HB2	2:A:180:HIS:CE1	2.49	0.47
1:E:1232:THR:OG1	1:E:1238:ALA:HA	2.14	0.47
1:E:1007:ARG:O	1:E:1031:VAL:HA	2.16	0.47
1:E:1368:PHE:CD2	1:E:1378:LEU:HB2	2.50	0.46
1:E:1018:PRO:HG2	1:E:1026:LEU:HD21	1.98	0.46
1:C:1161:VAL:HG23	1:C:1191:PRO:CG	2.38	0.46
1:E:1166:GLN:HB2	1:E:1186:ILE:HB	1.97	0.46
1:E:991:ILE:HD11	1:E:1006:CYS:SG	2.55	0.46
2:A:114:ASN:HB2	2:A:189:TRP:CZ3	2.51	0.46
2:B:221:TYR:OH	2:B:291:ASP:HB3	2.16	0.46
2:A:220:PHE:CD1	2:A:220:PHE:C	2.94	0.46
1:E:1181:ALA:HA	1:E:1271:CYS:HB3	1.97	0.45
1:E:1193:ARG:NH1	2:B:313:ARG:HA	2.31	0.45
1:E:1242:ASP:OD1	1:E:1244:THR:OG1	2.35	0.45
2:B:72:SER:C	2:B:74:PRO:HD3	2.41	0.45
1:C:991:ILE:HG12	1:C:1046:VAL:HG13	1.98	0.45
1:C:1163:LEU:HD11	1:C:1187:VAL:HG13	1.98	0.45
2:B:100:ARG:HG3	2:B:100:ARG:NH2	2.31	0.45
1:E:1274:VAL:O	1:E:1277:THR:HG23	2.17	0.45
2:B:114:ASN:HB2	2:B:189:TRP:CZ3	2.51	0.45
1:E:1009:VAL:CG2	1:E:1030:LEU:HB2	2.47	0.45
1:E:1065:ARG:HB3	1:E:1118:PHE:CE1	2.51	0.45
1:E:1202:MET:O	1:E:1257:SER:HA	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:293:GLN:NE2	2:A:293:GLN:HA	2.32	0.45
1:E:1163:LEU:HD11	1:E:1187:VAL:HG13	1.99	0.44
2:A:73:HIS:N	2:A:74:PRO:CD	2.80	0.44
1:C:1014:LEU:HD13	1:C:1117:TYR:CZ	2.53	0.44
1:C:1166:GLN:HB2	1:C:1186:ILE:HB	1.99	0.44
1:C:1231:ILE:HG22	1:C:1264:ALA:HB2	1.99	0.44
1:E:1020:MET:HE3	1:E:1020:MET:HB2	1.82	0.44
1:C:1227:LEU:HD23	1:C:1247:THR:HG21	2.00	0.44
2:B:220:PHE:HD1	2:B:221:TYR:CD1	2.36	0.44
1:C:1009:VAL:CG2	1:C:1030:LEU:HB2	2.47	0.44
1:C:1303:THR:HG23	1:C:1420:PRO:HG2	1.99	0.44
2:B:161:GLY:O	2:B:165:VAL:HG23	2.18	0.44
1:C:1297:PHE:N	1:C:1337:SER:OG	2.43	0.43
1:C:1031:VAL:HG23	1:C:1118:PHE:HB2	2.00	0.43
2:A:75:GLU:O	2:A:78:ARG:N	2.51	0.43
1:C:1004:ILE:HA	1:C:1035:PRO:HD3	1.99	0.43
1:C:1314:PHE:CE1	1:C:1377:ARG:HD3	2.53	0.43
2:A:106:GLU:HB3	2:A:180:HIS:NE2	2.33	0.43
2:A:127:ARG:NH1	2:A:157:THR:OG1	2.51	0.43
1:E:1194:ARG:HA	2:B:316:ARG:O	2.18	0.43
1:E:1027:ALA:O	1:E:1373:PHE:HZ	2.02	0.42
1:C:1005:THR:HB	1:C:1034:GLY:HA2	2.00	0.42
1:E:1284:LEU:HD23	1:E:1284:LEU:HA	1.90	0.42
1:C:1337:SER:CB	1:C:1415:LEU:HD21	2.49	0.42
1:C:1284:LEU:HD23	1:C:1284:LEU:HA	1.84	0.42
1:E:965:HIS:HA	1:E:980:SER:O	2.19	0.42
1:E:1105:ARG:NH2	1:E:1339:ASP:OD2	2.48	0.42
1:E:1337:SER:HB3	1:E:1415:LEU:HD11	2.02	0.42
1:C:1032:GLU:HG3	1:C:1117:TYR:CE1	2.54	0.42
2:B:34:GLU:OE1	2:B:115:ARG:NH1	2.52	0.42
1:E:1303:THR:HG21	1:E:1422:HIS:H	1.84	0.42
2:A:176:ILE:HD11	2:A:186:ILE:HG22	2.01	0.42
2:B:220:PHE:CD1	2:B:220:PHE:C	2.97	0.42
1:E:1008:LEU:HD23	1:E:1008:LEU:HA	1.86	0.42
2:A:231:ILE:HD11	2:A:270:VAL:HG21	2.00	0.42
1:C:1318:ASP:OD1	1:C:1318:ASP:C	2.62	0.42
2:A:85:LYS:HG3	2:A:89:LEU:HD23	2.01	0.41
1:C:1035:PRO:CB	1:C:1038:ALA:HB2	2.42	0.41
1:C:1366:LEU:HD11	1:C:1380:PHE:CZ	2.55	0.41
1:C:967:GLY:HA2	1:C:982:ARG:O	2.20	0.41
1:E:1161:VAL:HG23	1:E:1191:PRO:CG	2.42	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:293:GLN:HA	2:A:293:GLN:HE21	1.85	0.41
2:A:207:ARG:CZ	2:A:211:LEU:HD21	2.51	0.41
2:A:274:GLN:OE1	2:A:300:LEU:HD11	2.21	0.41
1:C:1105:ARG:NH2	1:C:1339:ASP:OD2	2.42	0.41
1:C:1231:ILE:C	1:C:1231:ILE:HD12	2.46	0.41
2:B:231:ILE:HD11	2:B:270:VAL:HG21	2.02	0.41
1:E:1401:PRO:HG2	1:E:1404:THR:CG2	2.47	0.41
2:A:220:PHE:HD1	2:A:221:TYR:CD1	2.39	0.41
1:C:965:HIS:HA	1:C:980:SER:O	2.20	0.41
2:B:176:ILE:HD11	2:B:186:ILE:HG22	2.02	0.41
1:E:1337:SER:CB	1:E:1415:LEU:HD21	2.51	0.41
1:E:967:GLY:HA2	1:E:982:ARG:O	2.21	0.41
2:B:26:MET:CG	2:B:96:LEU:HD11	2.51	0.41
2:B:110:LEU:HB2	2:B:180:HIS:CE1	2.56	0.41
1:C:1021:VAL:CG1	1:C:1192:ARG:HH12	2.34	0.41
1:C:1065:ARG:NH2	1:C:1115:PRO:HB3	2.35	0.41
1:E:1318:ASP:OD1	1:E:1318:ASP:C	2.64	0.40
1:C:1020:MET:O	1:C:1235:THR:HG22	2.22	0.40
1:C:1227:LEU:HD12	1:C:1267:TRP:O	2.21	0.40
2:B:124:ILE:HG23	2:B:200:LEU:HB2	2.03	0.40
1:E:1031:VAL:HG23	1:E:1118:PHE:HB2	2.03	0.40
2:A:228:PHE:O	2:A:232:GLN:HG2	2.21	0.40
1:C:1018:PRO:HG2	1:C:1026:LEU:HD21	2.02	0.40
1:C:1300:PHE:HA	1:C:1332:GLU:O	2.21	0.40
1:E:1311:LEU:O	1:E:1380:PHE:HB2	2.21	0.40
1:C:1340:ILE:HG21	1:C:1412:VAL:O	2.21	0.40
2:B:57:THR:O	2:B:61:LEU:HB2	2.21	0.40
1:E:968:PHE:CZ	1:E:991:ILE:HD12	2.56	0.40
1:E:1344:GLU:HA	1:E:1370:PHE:O	2.21	0.40
1:C:1086:GLU:O	1:C:1089:ASN:HB2	2.21	0.40
1:E:1004:ILE:HA	1:E:1035:PRO:HD3	2.03	0.40
1:E:1145:LEU:O	1:E:1205:PRO:HD3	2.22	0.40
1:C:1020:MET:HE3	1:C:1020:MET:HB2	1.85	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	C	404/491 (82%)	381 (94%)	23 (6%)	0	100	100
1	E	408/491 (83%)	382 (94%)	25 (6%)	1 (0%)	43	72
2	A	295/320 (92%)	287 (97%)	6 (2%)	2 (1%)	18	49
2	B	295/320 (92%)	289 (98%)	6 (2%)	0	100	100
All	All	1402/1622 (86%)	1339 (96%)	60 (4%)	3 (0%)	43	72

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	133	SER
2	A	73	HIS
1	E	1420	PRO

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	C	290/431 (67%)	290 (100%)	0	100	100
1	E	292/431 (68%)	292 (100%)	0	100	100
2	A	185/272 (68%)	185 (100%)	0	100	100
2	B	185/272 (68%)	185 (100%)	0	100	100
All	All	952/1406 (68%)	952 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	1261	ASN
2	A	194	ASN
2	A	281	GLN
1	C	1414	ASN
2	B	222	HIS
2	B	281	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å²)	Q<0.9	
1	C	422/491 (85%)	-1.00	0	100	100	68, 107, 145, 183	0
1	E	426/491 (86%)	-1.08	0	100	100	58, 108, 145, 171	0
2	A	301/320 (94%)	-1.14	0	100	100	65, 104, 151, 187	0
2	B	301/320 (94%)	-1.12	2 (0%)	84	70	64, 105, 147, 201	0
All	All	1450/1622 (89%)	-1.08	2 (0%)	92	91	58, 106, 147, 201	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	69	VAL	2.8
2	B	136	LEU	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 oligosaccharides 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.