



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 19, 2023 – 07:24 PM JST

PDB ID : 6M3P
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 : 4.02b-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

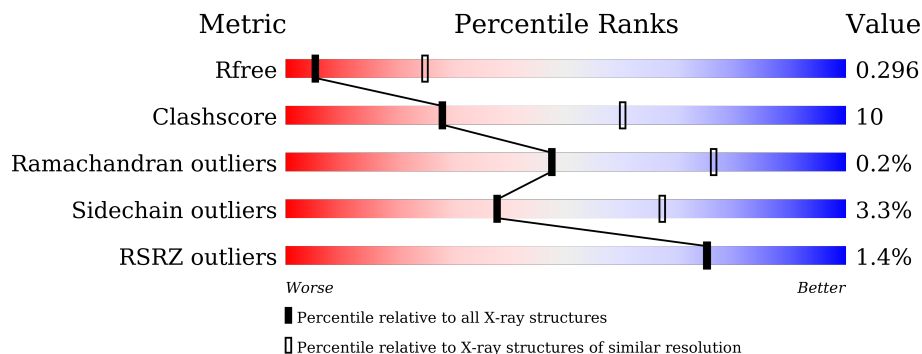
1 Overall quality at a glance

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}	130704	1089 (3.36-3.28)
Clashscore	141614	1137 (3.36-3.28)
Ramachandran outliers	138981	1115 (3.36-3.28)
Sidechain outliers	138945	1114 (3.36-3.28)
RSRZ outliers	127900	1059 (3.36-3.28)

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	 2% 65% 21% 14%
1	E	491	 2% 65% 21% 13%
2	A	320	 76% 17% 6%
2	B	320	 77% 17% 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
			Total	C	N	O	S			
1	E	426	2972	1893	502	557	20	0	0	0
1	C	422	2946	1875	498	553	20	0	0	0

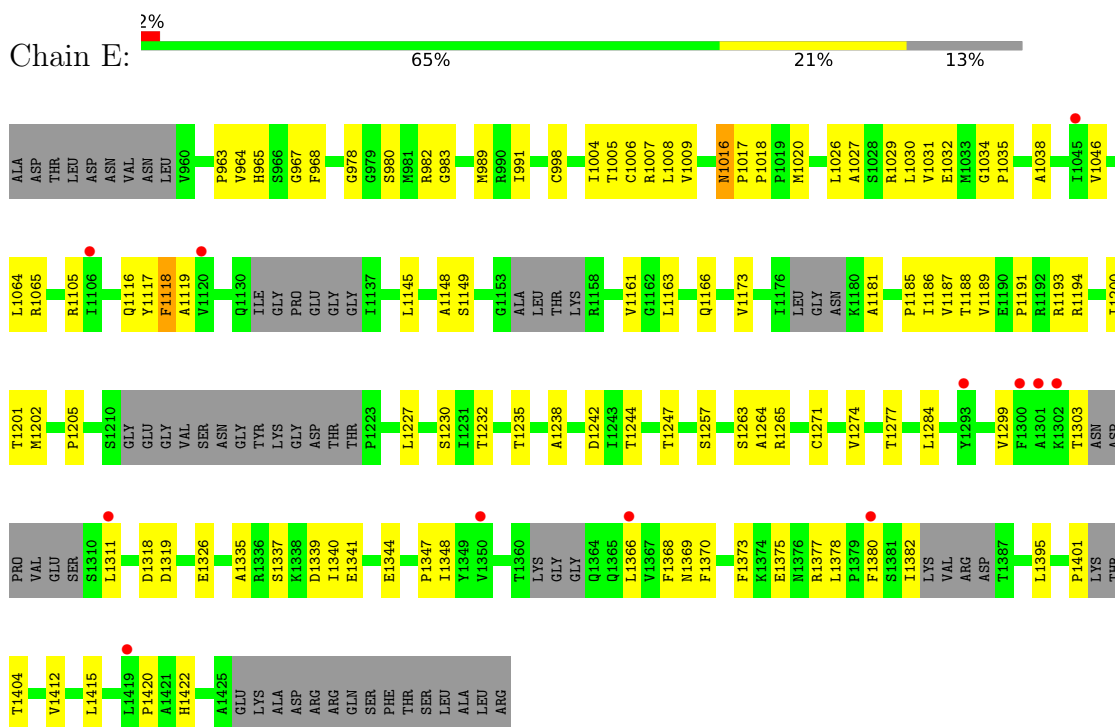
- Molecule 2 is a protein called Spectrin beta chain, non-erythrocytic 1.

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

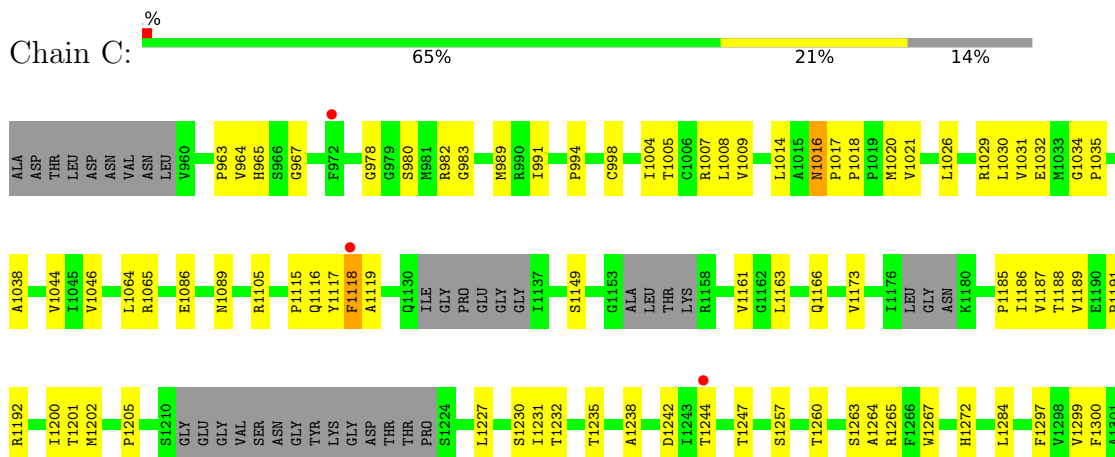
3 Residue-property plots [i](#)

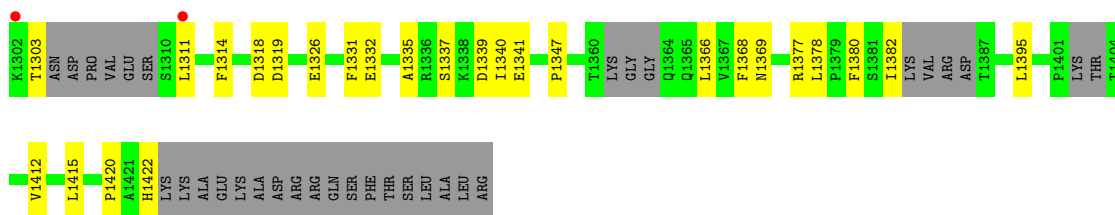
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Ankyrin-3

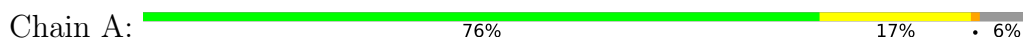


- Molecule 1: Ankyrin-3

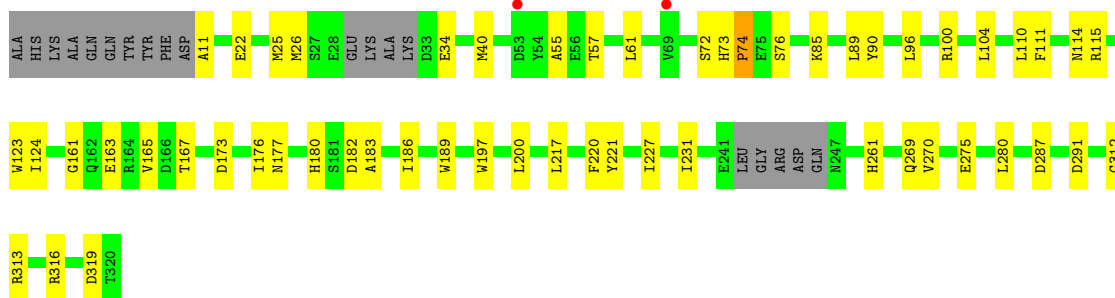
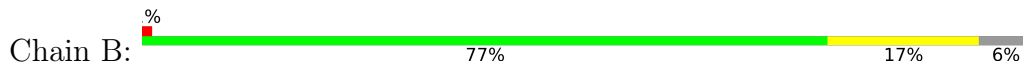




• Molecule 2: Spectrin beta chain, non-erythrocytic 1



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



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 0.255 , 0.296	Depositor DCC
R_{free} test set	1723 reflections (4.69%)	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)	0.29 , 86.7	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	0.92	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	0.31	0/3006	0.55	1/4126 (0.0%)
1	E	0.32	0/3033	0.55	1/4164 (0.0%)
2	A	0.29	0/2172	0.55	1/2970 (0.0%)
2	B	0.29	0/2172	0.50	0/2970
All	All	0.31	0/10383	0.54	3/14230 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	127	ARG	NE-CZ-NH1	8.27	124.43	120.30
1	C	1318	ASP	CB-CG-OD1	5.26	123.03	118.30
1	E	1318	ASP	CB-CG-OD1	5.13	122.92	118.30

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	60	0
1	E	2972	0	2708	67	0
2	A	2140	0	1817	34	0
2	B	2140	0	1817	33	0
All	All	10198	0	9022	190	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 (190) 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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:964:VAL:HG22	1:C:978:GLY:O	2.05	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
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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
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
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:1202:MET:O	1:E:1257:SER:HA	2.17	0.45
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: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
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
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
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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
2:A:220:PHE:CD1	2:A:220:PHE:C	2.94	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: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
2:B:72:SER:C	2:B:74:PRO:HD3	2.41	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

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	C	404/491 (82%)	381 (94%)	23 (6%)	0	100	100
1	E	408/491 (83%)	382 (94%)	25 (6%)	1 (0%)	47	76
2	A	295/320 (92%)	287 (97%)	6 (2%)	2 (1%)	22	55
2	B	295/320 (92%)	289 (98%)	6 (2%)	0	100	100
All	All	1402/1622 (86%)	1339 (96%)	60 (4%)	3 (0%)	47	76

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 [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	C	290/431 (67%)	282 (97%)	8 (3%)	43	71
1	E	292/431 (68%)	285 (98%)	7 (2%)	49	74
2	A	185/272 (68%)	177 (96%)	8 (4%)	29	61
2	B	185/272 (68%)	177 (96%)	8 (4%)	29	61
All	All	952/1406 (68%)	921 (97%)	31 (3%)	38	67

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	998	CYS
1	E	1016	ASN
1	E	1116	GLN
1	E	1118	PHE
1	E	1319	ASP

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Mol	Chain	Res	Type
1	E	1341	GLU
1	E	1395	LEU
2	A	22	GLU
2	A	111	PHE
2	A	153	PHE
2	A	182	ASP
2	A	202	GLU
2	A	222	HIS
2	A	261	HIS
2	A	275	GLU
1	C	998	CYS
1	C	1016	ASN
1	C	1116	GLN
1	C	1118	PHE
1	C	1272	HIS
1	C	1319	ASP
1	C	1341	GLU
1	C	1395	LEU
2	B	22	GLU
2	B	74	PRO
2	B	111	PHE
2	B	123	TRP
2	B	182	ASP
2	B	261	HIS
2	B	275	GLU
2	B	319	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 [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

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	C	422/491 (85%)	-0.10	5 (1%) 79 79	68, 107, 145, 183	0
1	E	426/491 (86%)	-0.09	12 (2%) 53 51	58, 108, 145, 171	0
2	A	301/320 (94%)	-0.35	1 (0%) 94 95	65, 104, 151, 187	0
2	B	301/320 (94%)	-0.34	2 (0%) 87 89	64, 105, 147, 201	0
All	All	1450/1622 (89%)	-0.20	20 (1%) 75 75	58, 106, 147, 201	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	1311	LEU	4.3
2	B	69	VAL	4.0
1	C	972	PHE	3.1
1	E	1419	LEU	2.9
2	A	53	ASP	2.8
2	B	53	ASP	2.7
1	E	1301	ALA	2.6
1	E	1366	LEU	2.5
1	E	1311	LEU	2.5
1	E	1120	VAL	2.3
1	E	1300	PHE	2.3
1	C	1302	LYS	2.3
1	E	1045	ILE	2.2
1	E	1302	LYS	2.2
1	C	1244	THR	2.2
1	E	1106	ILE	2.1
1	E	1293	TYR	2.1
1	C	1118	PHE	2.1
1	E	1380	PHE	2.0
1	E	1350	VAL	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.