



Full wwPDB X-ray Structure Validation Report ⓘ

May 21, 2020 – 07:09 am BST

PDB ID : 5X5Y
Title : A membrane protein complex
Authors : Luo, Q.; Yang, X.; Huang, Y.
Deposited on : 2017-02-18
Resolution : 3.46 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

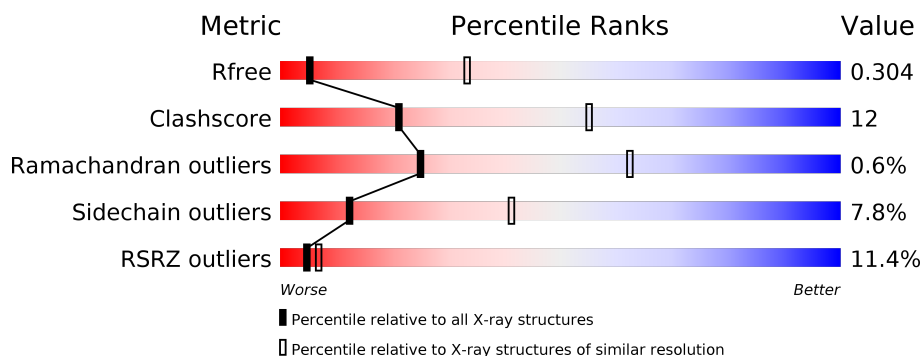
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.46 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	1291 (3.52-3.40)
Clashscore	141614	1372 (3.52-3.40)
Ramachandran outliers	138981	1337 (3.52-3.40)
Sidechain outliers	138945	1338 (3.52-3.40)
RSRZ outliers	127900	1205 (3.52-3.40)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. 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	247	<div> <div>9%</div> <div>64%</div> <div>31%</div> <div>..</div> </div>
1	B	247	<div> <div>8%</div> <div>62%</div> <div>33%</div> <div>..</div> </div>
2	G	355	<div> <div>15%</div> <div>73%</div> <div>23%</div> <div>..</div> </div>
3	F	362	<div> <div>11%</div> <div>63%</div> <div>31%</div> <div>..</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9178 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 Probable ATP-binding component of ABC transporter.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	239	Total	C	N	O	S	0	0	0
			1841	1153	336	345	7			
1	A	239	Total	C	N	O	S	0	0	0
			1841	1153	336	345	7			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	242	HIS	-	expression tag	UNP Q9HVV6
B	243	HIS	-	expression tag	UNP Q9HVV6
B	244	HIS	-	expression tag	UNP Q9HVV6
B	245	HIS	-	expression tag	UNP Q9HVV6
B	246	HIS	-	expression tag	UNP Q9HVV6
B	247	HIS	-	expression tag	UNP Q9HVV6
A	242	HIS	-	expression tag	UNP Q9HVV6
A	243	HIS	-	expression tag	UNP Q9HVV6
A	244	HIS	-	expression tag	UNP Q9HVV6
A	245	HIS	-	expression tag	UNP Q9HVV6
A	246	HIS	-	expression tag	UNP Q9HVV6
A	247	HIS	-	expression tag	UNP Q9HVV6

- Molecule 2 is a protein called Uncharacterized protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	G	351	Total	C	N	O	S	0	0	0
			2734	1782	468	473	11			

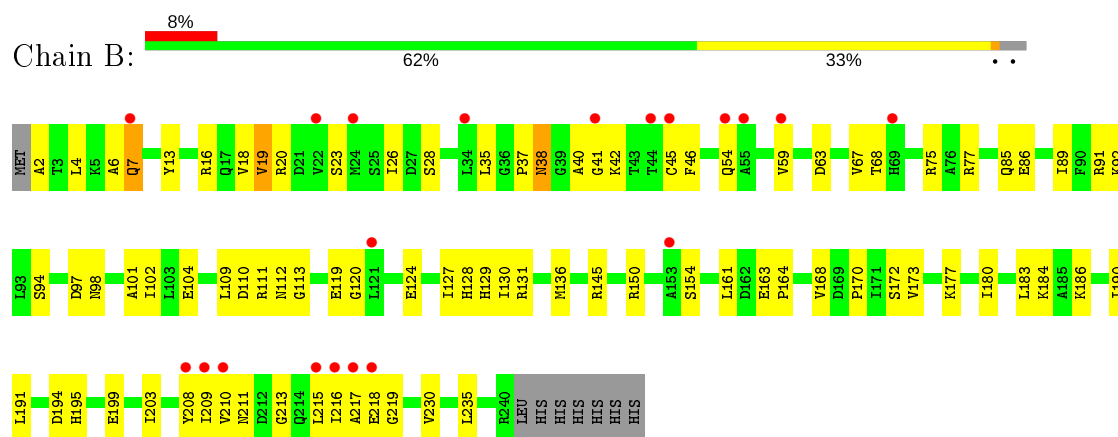
- Molecule 3 is a protein called Uncharacterized protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	F	351	Total	C	N	O	S	0	0	0
			2762	1815	461	475	11			

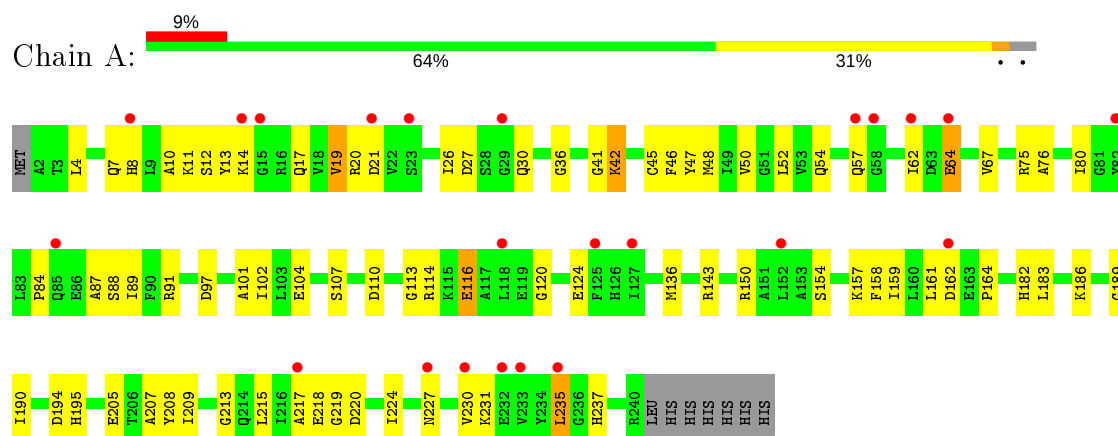
3 Residue-property plots [i](#)

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.

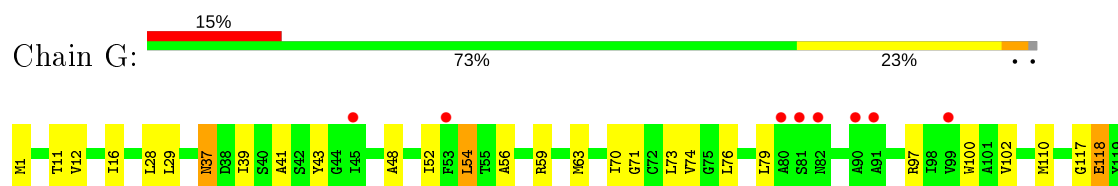
- Molecule 1: Probable ATP-binding component of ABC transporter

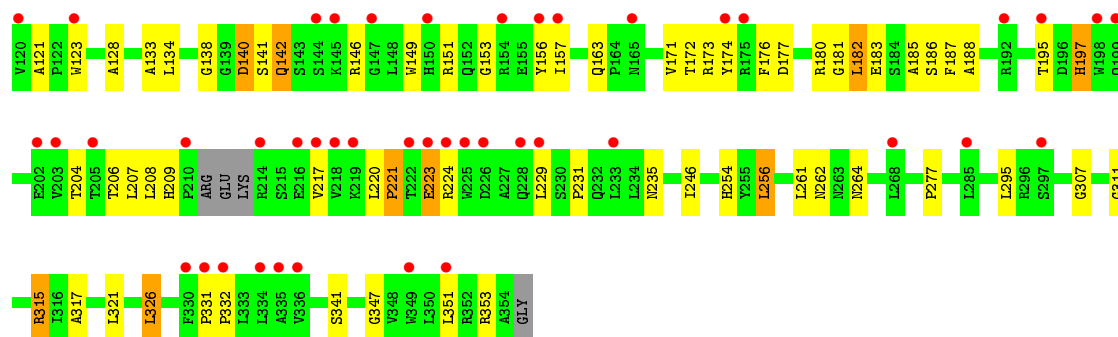


- Molecule 1: Probable ATP-binding component of ABC transporter

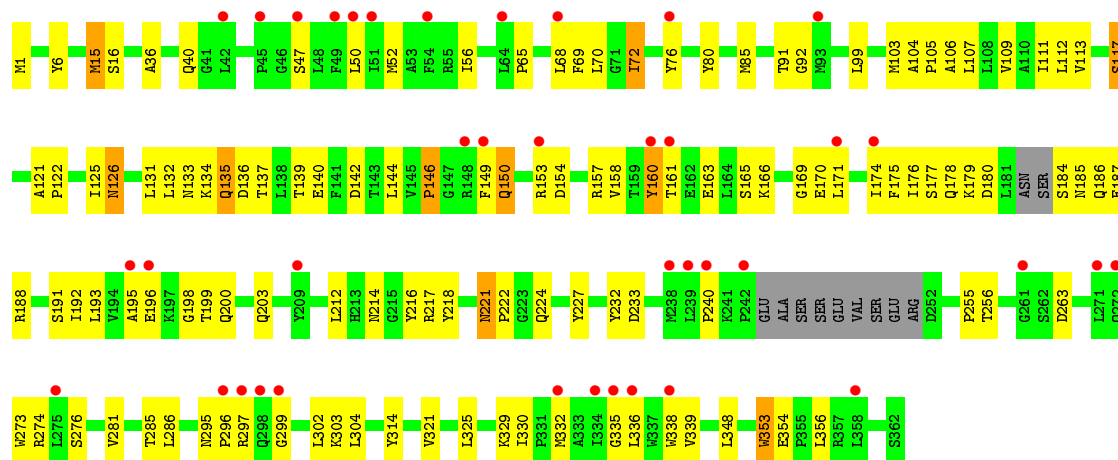


- Molecule 2: Uncharacterized protein





• Molecule 3: Uncharacterized protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	69.95Å 159.87Å 166.79Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.64 – 3.46 45.64 – 3.46	Depositor EDS
% Data completeness (in resolution range)	97.7 (45.64-3.46) 97.7 (45.64-3.46)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.51 (at 3.48Å)	Xtriage
Refinement program	PHENIX 1.11.1_2575	Depositor
R, R_{free}	0.269 , 0.304 0.269 , 0.304	Depositor DCC
R_{free} test set	1195 reflections (4.88%)	wwPDB-VP
Wilson B-factor (Å ²)	135.7	Xtriage
Anisotropy	0.429	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.24 , 73.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	0.039 for -h,l,k	Xtriage
F_o, F_c correlation	0.86	EDS
Total number of atoms	9178	wwPDB-VP
Average B, all atoms (Å ²)	145.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.79% 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	A	0.25	0/1867	0.46	0/2517
1	B	0.24	0/1867	0.45	0/2517
2	G	0.25	0/2796	0.43	0/3807
3	F	0.26	0/2820	0.47	0/3826
All	All	0.25	0/9350	0.45	0/12667

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	A	1841	0	1869	51	0
1	B	1841	0	1869	54	1
2	G	2734	0	2814	58	1
3	F	2762	0	2891	76	0
All	All	9178	0	9443	225	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 12.

All (225) 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)
3:F:132:LEU:O	3:F:136:ASP:HB2	1.70	0.92
2:G:117:GLY:O	2:G:121:ALA:HB3	1.77	0.83
3:F:133:ASN:O	3:F:137:THR:HB	1.83	0.77
2:G:182:LEU:HD11	2:G:185:ALA:HB2	1.66	0.77
3:F:65:PRO:HB3	3:F:281:VAL:HG22	1.66	0.75
3:F:174:ILE:HB	3:F:195:ALA:HB3	1.70	0.74
1:B:102:ILE:HD11	1:B:154:SER:HB2	1.69	0.73
2:G:134:LEU:HA	2:G:138:GLY:HA3	1.72	0.72
2:G:54:LEU:HB3	2:G:128:ALA:HB2	1.73	0.71
2:G:97:ARG:HH21	2:G:100:TRP:HD1	1.38	0.71
1:B:68:THR:O	1:B:75:ARG:NH2	2.23	0.70
3:F:99:LEU:O	3:F:103:MET:HB2	1.92	0.70
3:F:135:GLN:O	3:F:139:THR:HB	1.92	0.69
3:F:135:GLN:O	3:F:139:THR:CB	2.41	0.69
2:G:37:ASN:OD1	2:G:37:ASN:N	2.27	0.68
3:F:47:SER:HB2	3:F:50:LEU:HD12	1.76	0.68
2:G:180:ARG:HH12	3:F:196:GLU:HB2	1.60	0.67
1:A:231:LYS:HA	1:A:235:LEU:HB2	1.75	0.66
3:F:160:TYR:CD1	3:F:175:PHE:HB3	2.30	0.66
2:G:151:ARG:NH1	2:G:153:GLY:O	2.29	0.66
1:B:41:GLY:O	1:B:45:CYS:CB	2.44	0.66
3:F:175:PHE:HA	3:F:193:LEU:O	1.95	0.66
3:F:325:LEU:HD12	3:F:330:ILE:HB	1.77	0.65
1:B:120:GLY:O	1:B:124:GLU:HB2	1.97	0.63
1:B:209:ILE:HG21	1:B:230:VAL:HG13	1.79	0.63
2:G:117:GLY:O	2:G:121:ALA:CB	2.46	0.63
1:A:4:LEU:HD12	1:A:62:ILE:HB	1.80	0.63
3:F:195:ALA:HB2	3:F:212:LEU:HD13	1.81	0.63
1:A:120:GLY:O	1:A:124:GLU:HB2	1.99	0.63
1:B:67:VAL:HG22	1:B:75:ARG:HG2	1.80	0.63
1:A:89:ILE:HD13	1:A:143:ARG:HG2	1.80	0.62
1:B:102:ILE:HG13	3:F:91:THR:HG21	1.80	0.62
1:A:11:LYS:H	1:A:19:VAL:HG13	1.64	0.62
1:B:89:ILE:HA	1:B:150:ARG:HH22	1.65	0.61
1:B:19:VAL:HB	1:B:41:GLY:HA3	1.82	0.61
3:F:169:GLY:O	3:F:200:GLN:N	2.33	0.61
1:B:102:ILE:HG21	1:B:150:ARG:HG3	1.83	0.60
1:A:50:VAL:HG23	1:A:52:LEU:H	1.67	0.60
2:G:207:LEU:HD22	2:G:208:LEU:H	1.67	0.59
3:F:144:LEU:HB3	3:F:146:PRO:HA	1.82	0.59
1:B:209:ILE:HB	1:B:217:ALA:HB3	1.83	0.59
2:G:206:THR:OG1	3:F:222:PRO:O	2.20	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:142:GLN:HE22	3:F:179:LYS:HD2	1.67	0.59
1:A:64:GLU:HG2	1:A:64:GLU:O	2.03	0.59
3:F:113:VAL:O	3:F:117:SER:HB3	2.03	0.58
3:F:122:PRO:O	3:F:126:ASN:HB2	2.02	0.58
3:F:126:ASN:ND2	3:F:256:THR:O	2.36	0.58
1:B:77:ARG:NH2	3:F:92:GLY:O	2.34	0.58
1:A:19:VAL:HA	1:A:213:GLY:HA2	1.83	0.58
1:A:182:HIS:CE1	1:A:186:LYS:HE3	2.39	0.58
3:F:171:LEU:N	3:F:198:GLY:O	2.34	0.57
1:B:2:ALA:N	1:B:28:SER:HG	2.01	0.57
1:B:109:LEU:HD22	1:B:113:GLY:HA3	1.85	0.57
1:B:208:TYR:HB3	1:B:215:LEU:HD11	1.86	0.57
1:B:199:GLU:OE1	1:B:199:GLU:N	2.38	0.56
2:G:220:LEU:HG	2:G:221:PRO:HD2	1.86	0.56
1:A:161:LEU:HB3	1:A:164:PRO:HG3	1.86	0.56
1:B:41:GLY:O	1:B:45:CYS:HB3	2.05	0.56
3:F:221:ASN:N	3:F:221:ASN:OD1	2.40	0.55
1:A:194:ASP:OD1	1:A:195:HIS:N	2.39	0.55
1:B:85:GLN:HB3	1:B:163:GLU:HB2	1.88	0.55
2:G:48:ALA:O	2:G:52:ILE:HG12	2.06	0.55
1:B:41:GLY:O	1:B:45:CYS:HB2	2.06	0.55
3:F:131:LEU:O	3:F:135:GLN:HG2	2.06	0.55
3:F:170:GLU:HA	3:F:199:THR:HA	1.87	0.55
3:F:184:SER:OG	3:F:185:ASN:N	2.39	0.55
1:B:94:SER:O	1:B:98:ASN:ND2	2.38	0.54
3:F:176:ILE:O	3:F:192:ILE:HA	2.07	0.54
2:G:39:ILE:HG22	2:G:41:ALA:H	1.73	0.54
2:G:28:LEU:HD21	2:G:56:ALA:HB2	1.88	0.54
3:F:198:GLY:HA2	3:F:212:LEU:HA	1.89	0.53
1:A:104:GLU:HA	1:A:114:ARG:HD2	1.91	0.53
1:B:6:ALA:O	1:B:23:SER:HA	2.08	0.53
2:G:140:ASP:N	2:G:140:ASP:OD1	2.42	0.53
1:A:91:ARG:HB3	1:A:136:MET:HB3	1.90	0.53
1:A:102:ILE:HD11	1:A:154:SER:HB3	1.90	0.53
1:B:194:ASP:OD1	1:B:195:HIS:N	2.41	0.53
3:F:157:ARG:HB2	3:F:178:GLN:HG2	1.90	0.52
2:G:142:GLN:NE2	3:F:179:LYS:HD2	2.25	0.52
1:A:47:TYR:HA	1:A:50:VAL:HG22	1.92	0.52
3:F:136:ASP:O	3:F:140:GLU:CB	2.57	0.52
3:F:68:LEU:O	3:F:72:ILE:HG12	2.10	0.52
1:A:110:ASP:O	1:A:114:ARG:N	2.37	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:165:SER:OG	3:F:166:LYS:N	2.43	0.52
2:G:73:LEU:HD21	2:G:307:GLY:HA3	1.91	0.52
3:F:165:SER:HB2	3:F:170:GLU:HB2	1.90	0.52
2:G:177:ASP:OD1	2:G:181:GLY:N	2.43	0.52
2:G:311:GLY:O	2:G:315:ARG:HB2	2.10	0.51
1:A:209:ILE:HD11	1:A:224:ILE:HD13	1.93	0.51
1:B:211:ASN:HB3	1:B:216:ILE:HG21	1.93	0.51
3:F:106:ALA:HB2	3:F:285:THR:HG21	1.92	0.51
3:F:135:GLN:O	3:F:139:THR:OG1	2.28	0.51
3:F:286:LEU:HD21	3:F:348:LEU:HD22	1.93	0.51
1:B:183:LEU:HB3	1:B:190:ILE:HD11	1.93	0.50
1:A:52:LEU:HD12	2:G:295:LEU:HD12	1.94	0.50
3:F:335:GLY:O	3:F:338:TRP:HB3	2.11	0.50
1:B:210:VAL:HG22	1:B:215:LEU:HD12	1.92	0.50
2:G:204:THR:HG23	2:G:217:VAL:HG22	1.93	0.50
3:F:121:ALA:HB3	3:F:122:PRO:HD3	1.94	0.50
3:F:157:ARG:HA	3:F:177:SER:O	2.12	0.50
1:A:10:ALA:HA	1:A:19:VAL:HG22	1.94	0.49
2:G:180:ARG:HH12	3:F:196:GLU:CB	2.25	0.49
1:B:35:LEU:HB2	1:B:209:ILE:HD13	1.94	0.49
3:F:80:TYR:CZ	3:F:85:MET:HG2	2.47	0.49
2:G:149:TRP:NE1	3:F:158:VAL:HG21	2.27	0.49
1:B:86:GLU:HG2	3:F:297:ARG:NH1	2.28	0.49
3:F:125:ILE:HG21	3:F:255:PRO:HB2	1.94	0.49
2:G:173:ARG:HB2	2:G:186:SER:OG	2.12	0.49
1:A:87:ALA:HA	1:A:143:ARG:HH21	1.77	0.49
1:A:218:GLU:HG2	1:A:219:GLY:N	2.27	0.49
1:B:16:ARG:HD2	1:B:18:VAL:HG22	1.95	0.48
3:F:70:LEU:HG	3:F:304:LEU:HD22	1.96	0.48
1:A:183:LEU:HD23	1:A:190:ILE:HD11	1.95	0.48
1:A:227:ASN:O	1:A:231:LYS:NZ	2.37	0.48
1:A:46:PHE:O	1:A:50:VAL:HG13	2.14	0.48
1:B:161:LEU:HB3	1:B:164:PRO:HG3	1.94	0.48
2:G:11:THR:HG21	2:G:79:LEU:HD13	1.95	0.48
3:F:136:ASP:O	3:F:140:GLU:HB2	2.13	0.47
3:F:191:SER:HA	3:F:218:TYR:O	2.14	0.47
1:A:182:HIS:HE1	1:A:186:LYS:HE3	1.78	0.47
1:A:208:TYR:HB3	1:A:215:LEU:HD11	1.97	0.47
2:G:141:SER:HA	2:G:187:PHE:HE2	1.78	0.47
2:G:16:ILE:HG12	2:G:71:GLY:HA3	1.96	0.47
2:G:176:PHE:HE1	2:G:182:LEU:HD12	1.80	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:295:ASN:N	3:F:296:PRO:HD2	2.30	0.47
2:G:171:VAL:HB	2:G:188:ALA:HB3	1.96	0.47
2:G:256:LEU:HB3	2:G:261:LEU:HB3	1.96	0.46
2:G:133:ALA:O	2:G:138:GLY:N	2.48	0.46
1:B:4:LEU:HB3	1:B:26:ILE:HG23	1.97	0.46
2:G:73:LEU:CD2	2:G:307:GLY:HA3	2.45	0.46
1:A:205:GLU:O	1:A:220:ASP:HA	2.16	0.46
3:F:214:ASN:HA	3:F:233:ASP:HA	1.97	0.46
1:B:177:LYS:HD3	1:B:203:ILE:HD12	1.97	0.46
1:A:11:LYS:N	1:A:19:VAL:HG13	2.30	0.46
1:A:4:LEU:CD1	1:A:80:ILE:HD11	2.46	0.46
3:F:107:LEU:O	3:F:111:ILE:HG12	2.15	0.45
3:F:353:TRP:O	3:F:356:LEU:HG	2.16	0.45
3:F:52:MET:O	3:F:56:ILE:HG13	2.16	0.45
3:F:158:VAL:O	3:F:176:ILE:HA	2.15	0.45
1:A:113:GLY:HA2	1:A:116:GLU:OE1	2.17	0.45
3:F:354:GLU:N	3:F:354:GLU:OE2	2.50	0.45
1:B:218:GLU:HG2	1:B:219:GLY:N	2.32	0.45
2:G:317:ALA:O	2:G:321:LEU:HB2	2.16	0.45
3:F:171:LEU:HB2	3:F:174:ILE:HD11	1.99	0.45
2:G:264:ASN:HB3	2:G:326:LEU:HG	1.99	0.45
2:G:118:GLU:HG3	2:G:246:ILE:HG13	1.98	0.45
2:G:176:PHE:HD1	2:G:182:LEU:HA	1.82	0.45
1:B:180:ILE:HG23	1:B:190:ILE:HD13	1.99	0.45
3:F:132:LEU:O	3:F:136:ASP:CB	2.53	0.45
3:F:221:ASN:HB2	3:F:224:GLN:HG3	1.98	0.45
2:G:70:ILE:O	2:G:74:VAL:HG23	2.17	0.45
2:G:229:LEU:HD12	2:G:229:LEU:HA	1.90	0.44
1:A:27:ASP:O	1:A:30:GLN:HG2	2.18	0.44
1:A:207:ALA:O	1:A:218:GLU:HA	2.18	0.44
1:B:129:HIS:ND1	1:B:130:ILE:HG23	2.33	0.44
1:A:11:LYS:HB2	1:A:48:MET:SD	2.58	0.44
2:G:197:HIS:CD2	2:G:223:GLU:HG2	2.53	0.44
1:A:158:PHE:CD1	1:A:189:GLY:HA3	2.53	0.43
3:F:161:THR:HG22	3:F:163:GLU:H	1.83	0.43
1:B:92:LYS:HB2	3:F:6:TYR:CE1	2.53	0.43
1:B:173:VAL:O	1:B:177:LYS:HG2	2.18	0.43
3:F:144:LEU:HD11	3:F:150:GLN:HB2	2.01	0.43
1:B:128:HIS:CD2	1:B:131:ARG:HH21	2.36	0.43
3:F:321:VAL:HG11	3:F:339:VAL:HG21	2.01	0.43
1:B:38:ASN:N	1:B:38:ASN:OD1	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:110:MET:HB3	2:G:277:PRO:HG3	1.99	0.43
1:A:76:ALA:HA	1:A:80:ILE:O	2.18	0.43
1:A:97:ASP:O	1:A:101:ALA:HB2	2.18	0.43
3:F:15:MET:HG3	3:F:16:SER:N	2.33	0.43
3:F:273:TRP:HA	3:F:276:SER:HB3	2.01	0.43
1:A:230:VAL:O	1:A:235:LEU:N	2.40	0.43
2:G:142:GLN:OE1	3:F:179:LYS:HB2	2.18	0.43
2:G:151:ARG:HB2	3:F:149:PHE:HZ	1.84	0.43
1:B:91:ARG:HE	1:B:136:MET:HB3	1.84	0.42
1:B:37:PRO:HD2	1:B:210:VAL:O	2.19	0.42
1:B:184:LYS:HE2	1:B:184:LYS:HB3	1.92	0.42
1:A:14:LYS:HA	1:A:14:LYS:HD3	1.71	0.42
1:A:36:GLY:O	1:A:42:LYS:HE2	2.19	0.42
1:B:7:GLN:CG	1:B:59:VAL:HB	2.50	0.42
1:B:7:GLN:H	1:B:7:GLN:HG2	1.67	0.42
2:G:1:MET:HG2	2:G:97:ARG:HD3	2.02	0.42
1:A:41:GLY:O	1:A:45:CYS:HB2	2.20	0.42
1:B:104:GLU:OE2	3:F:1:MET:N	2.45	0.42
1:B:37:PRO:HG2	1:B:40:ALA:HB2	2.01	0.42
2:G:331:PRO:HA	2:G:332:PRO:HD3	1.91	0.42
1:A:67:VAL:HB	1:A:75:ARG:HG2	2.00	0.42
1:A:227:ASN:HB3	1:A:230:VAL:HG22	2.02	0.42
3:F:104:ALA:HB3	3:F:105:PRO:HD3	2.02	0.42
1:B:119:GLU:OE1	1:B:131:ARG:NH2	2.52	0.41
2:G:157:ILE:HA	2:G:172:THR:O	2.19	0.41
2:G:326:LEU:HA	2:G:326:LEU:HD13	1.75	0.41
1:A:88:SER:O	1:A:150:ARG:NH2	2.49	0.41
2:G:12:VAL:HG21	2:G:102:VAL:HG13	2.03	0.41
1:A:208:TYR:HD2	1:A:215:LEU:HD21	1.84	0.41
1:A:97:ASP:O	1:A:101:ALA:CB	2.69	0.41
3:F:221:ASN:O	3:F:227:TYR:HE1	2.04	0.41
2:G:76:LEU:HD11	2:G:102:VAL:HG21	2.02	0.41
1:A:84:PRO:HA	1:A:162:ASP:HB3	2.03	0.41
1:A:4:LEU:HB3	1:A:26:ILE:HG23	2.01	0.41
3:F:36:ALA:O	3:F:40:GLN:HG3	2.20	0.41
3:F:72:ILE:O	3:F:76:TYR:HB2	2.21	0.41
1:B:127:ILE:HG22	1:B:145:ARG:HH21	1.86	0.41
1:B:168:VAL:HG13	1:B:172:SER:HB2	2.01	0.41
3:F:134:LYS:O	3:F:137:THR:HG22	2.21	0.41
3:F:72:ILE:H	3:F:72:ILE:HG12	1.60	0.41
1:A:159:ILE:HB	1:A:190:ILE:HD13	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:18:VAL:O	1:B:213:GLY:HA2	2.21	0.41
1:B:97:ASP:O	1:B:101:ALA:HB2	2.21	0.41
2:G:156:TYR:HB2	2:G:174:TYR:HB2	2.02	0.41
1:B:46:PHE:HE1	1:B:191:LEU:HD22	1.86	0.41
3:F:212:LEU:HD11	3:F:232:TYR:OH	2.21	0.41
1:A:12:SER:HB3	1:A:17:GLN:HE21	1.86	0.40
1:B:170:PRO:O	1:B:173:VAL:HG22	2.22	0.40
2:G:176:PHE:CE1	2:G:182:LEU:HD12	2.56	0.40
2:G:231:PRO:O	2:G:235:ASN:HB2	2.21	0.40
1:A:26:ILE:HD11	1:A:190:ILE:N	2.36	0.40
1:B:7:GLN:HG2	1:B:59:VAL:HB	2.02	0.40
3:F:109:VAL:O	3:F:113:VAL:HG23	2.21	0.40
2:G:12:VAL:HG21	2:G:102:VAL:HG22	2.03	0.40
1:A:209:ILE:HB	1:A:217:ALA:HB3	2.03	0.40
2:G:347:GLY:O	2:G:351:LEU:HG	2.21	0.40
2:G:54:LEU:HD13	2:G:54:LEU:HA	1.91	0.40
2:G:182:LEU:HD23	2:G:207:LEU:HD23	2.04	0.40
1:B:230:VAL:HG12	1:B:235:LEU:HD13	2.03	0.40
2:G:28:LEU:HD12	2:G:59:ARG:HD2	2.03	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)
1:B:112:ASN:ND2	2:G:262:ASN:OD1[2_455]	2.18	0.02

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	237/247 (96%)	233 (98%)	3 (1%)	1 (0%)	34 70

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	237/247 (96%)	231 (98%)	6 (2%)	0	100	100
2	G	347/355 (98%)	328 (94%)	17 (5%)	2 (1%)	25	62
3	F	345/362 (95%)	318 (92%)	23 (7%)	4 (1%)	13	48
All	All	1166/1211 (96%)	1110 (95%)	49 (4%)	7 (1%)	25	62

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	F	153	ARG
3	F	240	PRO
1	A	237	HIS
2	G	221	PRO
2	G	195	THR
3	F	146	PRO
3	F	299	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	196/204 (96%)	182 (93%)	14 (7%)	14	46
1	B	196/204 (96%)	185 (94%)	11 (6%)	21	53
2	G	287/290 (99%)	264 (92%)	23 (8%)	12	40
3	F	292/302 (97%)	264 (90%)	28 (10%)	8	32
All	All	971/1000 (97%)	895 (92%)	76 (8%)	12	41

All (76) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	7	GLN
1	B	13	TYR
1	B	19	VAL
1	B	20	ARG

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Mol	Chain	Res	Type
1	B	38	ASN
1	B	42	LYS
1	B	54	GLN
1	B	63	ASP
1	B	110	ASP
1	B	111	ARG
1	B	186	LYS
1	A	7	GLN
1	A	8	HIS
1	A	13	TYR
1	A	19	VAL
1	A	20	ARG
1	A	21	ASP
1	A	42	LYS
1	A	54	GLN
1	A	57	GLN
1	A	64	GLU
1	A	107	SER
1	A	116	GLU
1	A	157	LYS
1	A	235	LEU
2	G	29	LEU
2	G	37	ASN
2	G	43	TYR
2	G	54	LEU
2	G	63	MET
2	G	118	GLU
2	G	123	TRP
2	G	140	ASP
2	G	142	GLN
2	G	146	ARG
2	G	163	GLN
2	G	182	LEU
2	G	183	GLU
2	G	197	HIS
2	G	209	HIS
2	G	223	GLU
2	G	224	ARG
2	G	254	HIS
2	G	256	LEU
2	G	315	ARG
2	G	326	LEU

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Mol	Chain	Res	Type
2	G	341	SER
2	G	353	ARG
3	F	15	MET
3	F	69	PHE
3	F	72	ILE
3	F	112	LEU
3	F	117	SER
3	F	126	ASN
3	F	135	GLN
3	F	142	ASP
3	F	150	GLN
3	F	154	ASP
3	F	160	TYR
3	F	180	ASP
3	F	186	GLN
3	F	187	GLU
3	F	188	ARG
3	F	203	GLN
3	F	216	TYR
3	F	217	ARG
3	F	221	ASN
3	F	263	ASP
3	F	274	ARG
3	F	302	LEU
3	F	303	LYS
3	F	314	TYR
3	F	329	LYS
3	F	332	MET
3	F	336	LEU
3	F	353	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	182	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [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	A	239/247 (96%)	0.54	23 (9%) 8 10	103, 144, 182, 206	0
1	B	239/247 (96%)	0.55	20 (8%) 11 13	100, 137, 163, 193	0
2	G	351/355 (98%)	0.70	52 (14%) 2 3	83, 143, 190, 220	0
3	F	351/362 (96%)	0.41	39 (11%) 5 7	93, 150, 209, 230	0
All	All	1180/1211 (97%)	0.55	134 (11%) 5 7	83, 143, 194, 230	0

All (134) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	G	228	GLN	9.0
1	B	54	GLN	8.8
2	G	223	GLU	8.3
2	G	226	ASP	8.2
3	F	160	TYR	7.4
2	G	165	ASN	6.7
1	B	215	LEU	6.5
2	G	224	ARG	6.5
3	F	334	ILE	6.1
3	F	298	GLN	5.9
2	G	225	TRP	5.5
2	G	331	PRO	4.9
3	F	161	THR	4.7
3	F	209	TYR	4.7
2	G	219	LYS	4.6
2	G	222	THR	4.6
1	B	55	ALA	4.5
3	F	240	PRO	4.4
3	F	335	GLY	4.4
1	B	216	ILE	4.4
2	G	175	ARG	4.3

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Mol	Chain	Res	Type	RSRZ
3	F	239	LEU	4.3
3	F	174	ILE	4.2
1	A	23	SER	4.2
1	B	24	MET	4.0
1	B	41	GLY	4.0
1	A	15	GLY	4.0
1	B	208	TYR	4.0
3	F	299	GLY	3.9
2	G	91	ALA	3.9
2	G	81	SER	3.9
1	A	233	VAL	3.9
2	G	229	LEU	3.7
2	G	210	PRO	3.6
1	A	232	GLU	3.6
2	G	297	SER	3.5
2	G	156	TYR	3.5
1	B	209	ILE	3.5
2	G	198	TRP	3.4
2	G	334	LEU	3.4
3	F	50	LEU	3.4
1	A	118	LEU	3.3
2	G	157	ILE	3.3
2	G	351	LEU	3.2
3	F	238	MET	3.2
3	F	45	PRO	3.1
1	B	153	ALA	3.1
3	F	332	MET	3.1
1	A	125	PHE	3.1
2	G	123	TRP	3.1
1	A	62	ILE	3.0
3	F	338	TRP	3.0
3	F	42	LEU	3.0
1	A	230	VAL	3.0
1	B	59	VAL	3.0
2	G	332	PRO	3.0
1	A	64	GLU	2.9
1	A	127	ILE	2.9
1	A	57	GLN	2.9
3	F	49	PHE	2.9
3	F	153	ARG	2.9
2	G	217	VAL	2.9
1	B	218	GLU	2.8

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Mol	Chain	Res	Type	RSRZ
3	F	54	PHE	2.8
1	B	217	ALA	2.8
3	F	261	GLY	2.8
3	F	195	ALA	2.8
3	F	271	LEU	2.8
3	F	149	PHE	2.8
1	B	69	HIS	2.8
2	G	349	TRP	2.7
1	A	14	LYS	2.7
3	F	358	LEU	2.7
2	G	202	GLU	2.7
2	G	90	ALA	2.7
2	G	174	TYR	2.6
2	G	192	ARG	2.6
2	G	120	VAL	2.6
2	G	199	GLN	2.6
2	G	216	GLU	2.6
3	F	196	GLU	2.6
3	F	296	PRO	2.6
1	B	210	VAL	2.5
2	G	335	ALA	2.5
1	A	162	ASP	2.5
3	F	76	TYR	2.5
1	A	85	GLN	2.5
2	G	82	ASN	2.5
3	F	272	GLN	2.5
3	F	297	ARG	2.5
1	A	217	ALA	2.5
3	F	336	LEU	2.5
1	B	44	THR	2.5
2	G	336	VAL	2.4
2	G	80	ALA	2.4
3	F	47	SER	2.4
1	A	29	GLY	2.4
2	G	154	ARG	2.4
2	G	285	LEU	2.3
1	B	22	VAL	2.3
2	G	205	THR	2.3
2	G	214	ARG	2.3
2	G	233	LEU	2.3
2	G	203	VAL	2.3
2	G	195	THR	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	7	GLN	2.3
3	F	93	MET	2.3
1	A	8	HIS	2.3
3	F	148	ARG	2.2
2	G	53	PHE	2.2
2	G	150	HIS	2.2
2	G	218	VAL	2.2
1	A	21	ASP	2.2
2	G	147	GLY	2.2
3	F	68	LEU	2.2
3	F	242	PRO	2.2
1	A	235	LEU	2.2
2	G	330	PHE	2.1
2	G	145	LYS	2.1
1	B	34	LEU	2.1
3	F	51	ILE	2.1
1	A	82	TYR	2.1
2	G	45	ILE	2.1
1	A	227	ASN	2.1
1	A	152	LEU	2.1
1	B	45	CYS	2.0
1	B	121	LEU	2.0
2	G	268	LEU	2.0
3	F	171	LEU	2.0
3	F	64	LEU	2.0
2	G	144	SER	2.0
2	G	99	VAL	2.0
1	A	58	GLY	2.0
3	F	275	LEU	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 [i](#)

There are no such residues in this entry.