



Full wwPDB X-ray Structure Validation Report ⓘ

Jul 19, 2017 – 05:21 PM EDT

PDB ID : 2P1L
Title : Structure of the Bcl-XL:Beclin 1 complex
Authors : Jeffrey, P.D.; Shi, Y.; Oberstein, A.L.
Deposited on : unknown
Resolution : 2.50 Å(reported)

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

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

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.9-1692
EDS : rb-20029824
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20029824

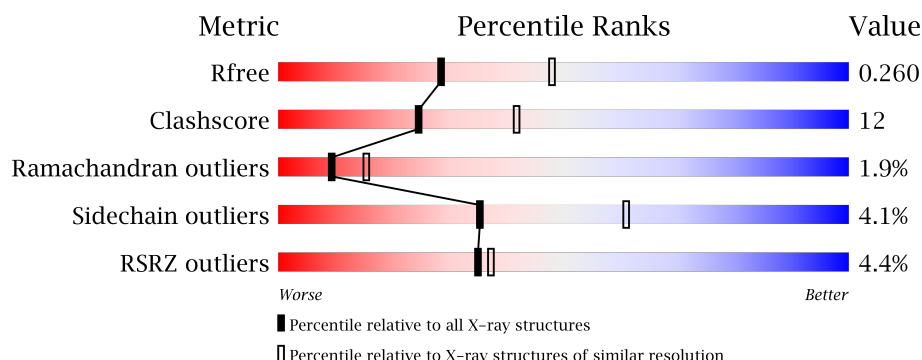
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.






Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	100719	3846 (2.50-2.50)
Clashscore	112137	4554 (2.50-2.50)
Ramachandran outliers	110173	4463 (2.50-2.50)
Sidechain outliers	110143	4465 (2.50-2.50)
RSRZ outliers	101464	3876 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	153	<div> <div>3%</div> <div>65% 22% 5% 8%</div> </div>
1	C	153	<div> <div>3%</div> <div>72% 18% • 8%</div> </div>
1	E	153	<div> <div>5%</div> <div>76% 11% • • 8%</div> </div>
1	G	153	<div> <div>5%</div> <div>63% 24% 5% • 8%</div> </div>
2	B	31	<div> <div>6%</div> <div>58% 19% 23%</div> </div>

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Mol	Chain	Length	Quality of chain
2	D	31	 68% 10% 23%
2	F	31	 6% 55% 23% 23%
2	H	31	 6% 52% 26% 23%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5300 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 Apoptosis regulator Bcl-X.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	141	Total	C	N	O	S	0	0	0
			1143	729	194	216	4			
1	C	141	Total	C	N	O	S	0	0	0
			1143	729	194	216	4			
1	E	141	Total	C	N	O	S	0	0	0
			1143	729	194	216	4			
1	G	141	Total	C	N	O	S	0	0	0
			1143	729	194	216	4			

- Molecule 2 is a protein called Beclin 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	24	Total	C	N	O	S	0	0	0
			178	108	32	36	2			
2	D	24	Total	C	N	O	S	0	0	0
			178	108	32	36	2			
2	F	24	Total	C	N	O	S	0	0	0
			178	108	32	36	2			
2	H	24	Total	C	N	O	S	0	0	0
			178	108	32	36	2			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	105	GLY	-	CLONING ARTIFACT	UNP Q14457
B	106	SER	-	CLONING ARTIFACT	UNP Q14457
D	105	GLY	-	CLONING ARTIFACT	UNP Q14457
D	106	SER	-	CLONING ARTIFACT	UNP Q14457
F	105	GLY	-	CLONING ARTIFACT	UNP Q14457
F	106	SER	-	CLONING ARTIFACT	UNP Q14457
H	105	GLY	-	CLONING ARTIFACT	UNP Q14457
H	106	SER	-	CLONING ARTIFACT	UNP Q14457

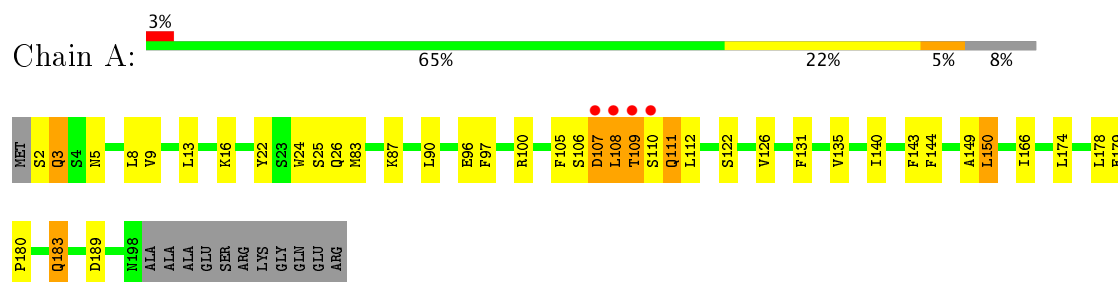
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	5	Total 5	O 5	0	0
3	C	4	Total 4	O 4	0	0
3	E	5	Total 5	O 5	0	0
3	G	2	Total 2	O 2	0	0

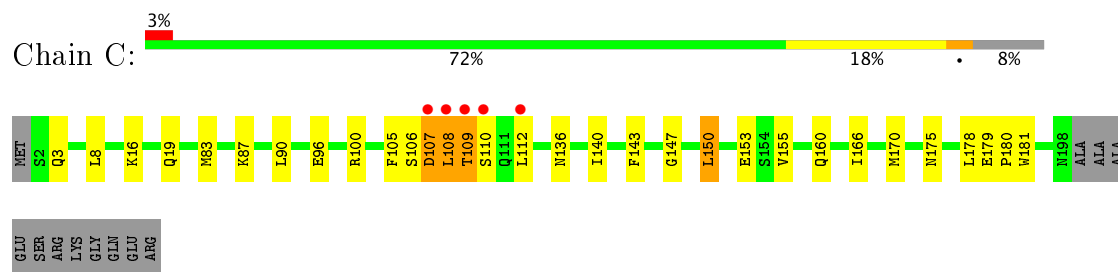
3 Residue-property plots [i](#)

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.

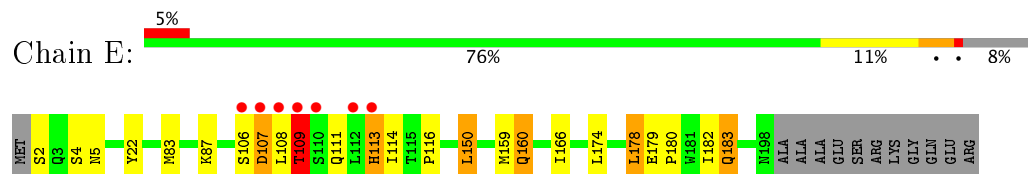
• Molecule 1: Apoptosis regulator Bcl-X



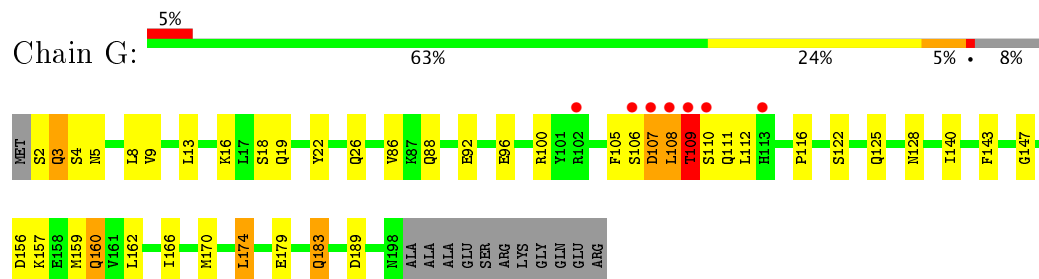
• Molecule 1: Apoptosis regulator Bcl-X



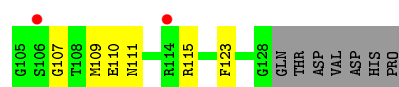
• Molecule 1: Apoptosis regulator Bcl-X



• Molecule 1: Apoptosis regulator Bcl-X



• Molecule 2: Beclin 1



• Molecule 2: Beclin 1



• Molecule 2: Beclin 1



• Molecule 2: Beclin 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	107.94Å 110.59Å 100.52Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.73 – 2.50 45.56 – 2.40	Depositor EDS
% Data completeness (in resolution range)	97.6 (29.73-2.50) 97.5 (45.56-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.56 (at 2.39Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.226 , 0.262 0.225 , 0.260	Depositor DCC
R_{free} test set	2107 reflections (5.11%)	DCC
Wilson B-factor (Å ²)	53.3	Xtriage
Anisotropy	0.504	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 40.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.055 for k,h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5300	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.42% 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	0.54	0/1171	0.66	0/1586
1	C	0.52	0/1171	0.65	1/1586 (0.1%)
1	E	0.55	0/1171	0.68	0/1586
1	G	0.53	0/1171	0.69	2/1586 (0.1%)
2	B	0.47	0/178	0.61	0/235
2	D	0.43	0/178	0.63	0/235
2	F	0.47	0/178	0.65	0/235
2	H	0.44	0/178	0.62	0/235
All	All	0.52	0/5396	0.67	3/7284 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	174	LEU	CA-CB-CG	6.34	129.89	115.30
1	G	108	LEU	CA-CB-CG	5.29	127.47	115.30
1	C	108	LEU	CA-CB-CG	5.05	126.92	115.30

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	A	1143	0	1083	36	0
1	C	1143	0	1083	22	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	1143	0	1083	29	0
1	G	1143	0	1083	41	0
2	B	178	0	179	5	0
2	D	178	0	179	3	0
2	F	178	0	179	6	0
2	H	178	0	179	7	0
3	A	5	0	0	0	0
3	C	4	0	0	0	0
3	E	5	0	0	0	0
3	G	2	0	0	0	0
All	All	5300	0	5048	127	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 12.

All (127) 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)
1:E:83:MET:HE3	1:E:87:LYS:HE3	1.53	0.89
1:A:109:THR:HG23	1:A:110:SER:H	1.39	0.88
1:C:109:THR:HG23	1:C:110:SER:H	1.39	0.87
1:E:83:MET:CE	1:E:87:LYS:HE3	2.05	0.85
1:E:2:SER:HB2	1:G:86:VAL:HG21	1.62	0.81
1:E:2:SER:CB	1:G:86:VAL:HG21	2.13	0.78
1:E:107:ASP:OD2	1:E:109:THR:HG23	1.85	0.76
2:F:111:ASN:HD21	2:F:115:ARG:CZ	2.01	0.73
1:G:107:ASP:HB2	1:G:109:THR:HG22	1.71	0.72
1:C:107:ASP:HB2	1:C:109:THR:HG22	1.70	0.72
1:A:8:LEU:HD22	1:C:90:LEU:HD23	1.73	0.71
1:G:107:ASP:HB2	1:G:109:THR:CG2	2.20	0.71
2:D:111:ASN:O	2:D:115:ARG:HG3	1.90	0.70
1:A:108:LEU:HD23	1:A:108:LEU:H	1.57	0.70
1:A:107:ASP:HB2	1:A:109:THR:HG22	1.73	0.69
1:E:174:LEU:HD12	1:E:178:LEU:HD23	1.77	0.66
1:E:183:GLN:CA	1:E:183:GLN:HE21	2.10	0.64
1:A:150:LEU:HB3	1:A:166:ILE:HD13	1.78	0.64
1:E:179:GLU:HB3	1:E:180:PRO:HD3	1.81	0.63
1:G:116:PRO:HA	1:G:162:LEU:HD21	1.81	0.63
1:E:114:ILE:HG22	1:E:159:MET:HE3	1.80	0.62
1:G:179:GLU:OE1	1:G:179:GLU:HA	2.00	0.62
1:E:174:LEU:HD23	1:G:5:ASN:HB3	1.80	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:183:GLN:HA	1:E:183:GLN:HE21	1.64	0.62
1:A:96:GLU:OE2	1:A:100:ARG:HD2	2.00	0.61
1:A:2:SER:HB3	1:C:175:ASN:OD1	2.01	0.61
1:E:160:GLN:CD	1:E:160:GLN:H	2.04	0.61
1:E:150:LEU:HB3	1:E:166:ILE:HD13	1.84	0.60
2:H:107:GLY:HA2	2:H:110:GLU:OE1	2.02	0.60
1:G:109:THR:OG1	1:G:153:GLU:HG2	2.02	0.59
1:G:16:LYS:HD3	1:G:19:GLN:OE1	2.03	0.59
1:E:113:HIS:NE2	2:F:105:GLY:HA2	2.18	0.59
1:E:178:LEU:O	1:E:182:ILE:HG13	2.02	0.59
1:E:183:GLN:NE2	1:E:183:GLN:HA	2.17	0.58
2:B:111:ASN:O	2:B:115:ARG:HG3	2.04	0.58
1:C:150:LEU:HB3	1:C:166:ILE:HD13	1.86	0.58
2:F:114:ARG:O	2:F:118:VAL:HG23	2.04	0.58
1:C:109:THR:HG23	1:C:110:SER:N	2.16	0.57
1:G:150:LEU:HB3	1:G:166:ILE:HD13	1.84	0.57
1:C:179:GLU:HB3	1:C:180:PRO:HD3	1.87	0.56
1:A:109:THR:HG23	1:A:110:SER:N	2.16	0.56
1:C:112:LEU:HD21	2:D:109:MET:HG2	1.88	0.56
1:G:112:LEU:HD23	2:H:108:THR:HB	1.88	0.54
1:A:183:GLN:HE21	1:A:183:GLN:CA	2.20	0.54
1:G:107:ASP:OD2	1:G:109:THR:HG23	2.08	0.53
1:E:107:ASP:HB2	1:E:109:THR:CG2	2.38	0.53
1:E:160:GLN:CD	1:E:160:GLN:N	2.62	0.53
1:A:8:LEU:HD22	1:C:90:LEU:CD2	2.39	0.53
2:F:111:ASN:HD21	2:F:115:ARG:NE	2.07	0.53
1:E:2:SER:HB3	1:G:86:VAL:HG21	1.91	0.52
1:A:83:MET:CE	1:A:87:LYS:HE3	2.40	0.51
1:A:140:ILE:O	1:A:143:PHE:HB3	2.11	0.51
1:G:116:PRO:HD3	1:G:159:MET:SD	2.51	0.51
1:E:108:LEU:HD23	1:E:108:LEU:H	1.75	0.51
1:A:105:PHE:CD2	2:B:115:ARG:HD3	2.46	0.51
1:G:160:GLN:H	1:G:160:GLN:CD	2.13	0.51
1:A:108:LEU:HA	1:A:111:GLN:HE21	1.77	0.50
1:G:160:GLN:N	1:G:160:GLN:CD	2.66	0.49
2:F:111:ASN:HD21	2:F:115:ARG:NH2	2.11	0.49
1:E:183:GLN:NE2	1:E:183:GLN:CA	2.76	0.48
1:A:22:TYR:CG	1:C:155:VAL:HG11	2.49	0.48
1:A:26:GLN:NE2	1:A:189:ASP:OD1	2.47	0.48
1:C:108:LEU:H	1:C:108:LEU:HD23	1.78	0.48
1:A:179:GLU:HB3	1:A:180:PRO:HD3	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:96:GLU:OE2	1:C:100:ARG:HD2	2.14	0.48
1:A:112:LEU:HD21	2:B:109:MET:HG2	1.96	0.48
1:G:105:PHE:CE1	2:H:115:ARG:HB3	2.49	0.47
1:A:83:MET:HE1	1:A:87:LYS:HE3	1.97	0.47
1:G:18:SER:HA	1:G:22:TYR:O	2.14	0.47
1:G:2:SER:HB2	1:G:3:GLN:NE2	2.29	0.47
1:G:96:GLU:OE2	1:G:100:ARG:HD2	2.14	0.47
1:G:153:GLU:OE2	1:G:157:LYS:HE2	2.15	0.46
1:G:109:THR:HG23	1:G:110:SER:H	1.80	0.46
1:G:96:GLU:HG3	2:H:123:PHE:HE1	1.78	0.46
1:A:90:LEU:HD21	1:A:144:PHE:CE1	2.50	0.46
1:E:5:ASN:HB3	1:G:174:LEU:HD23	1.96	0.46
1:A:131:PHE:CD1	1:A:135:VAL:HG22	2.50	0.46
1:E:174:LEU:CD1	1:E:178:LEU:HD23	2.45	0.46
1:G:140:ILE:O	1:G:143:PHE:HB3	2.16	0.46
1:A:96:GLU:HG3	2:B:123:PHE:HE1	1.81	0.46
1:G:183:GLN:HE21	1:G:183:GLN:CA	2.29	0.45
1:A:183:GLN:HA	1:A:183:GLN:HE21	1.82	0.45
1:G:183:GLN:NE2	1:G:183:GLN:HA	2.32	0.45
1:A:16:LYS:HA	1:A:16:LYS:HD3	1.78	0.45
1:E:106:SER:O	1:E:107:ASP:O	2.35	0.45
1:G:9:VAL:O	1:G:13:LEU:HG	2.17	0.45
1:C:106:SER:O	1:C:107:ASP:O	2.36	0.44
1:G:128:ASN:HA	1:G:128:ASN:HD22	1.68	0.44
1:A:106:SER:O	1:A:107:ASP:C	2.56	0.44
1:G:108:LEU:HD23	1:G:108:LEU:H	1.82	0.44
1:A:106:SER:O	1:A:107:ASP:O	2.35	0.44
1:E:106:SER:O	1:E:107:ASP:C	2.56	0.44
1:A:9:VAL:O	1:A:13:LEU:HG	2.18	0.44
1:E:107:ASP:HB2	1:E:109:THR:HG22	2.00	0.43
1:A:109:THR:HA	1:A:149:ALA:HB1	1.99	0.43
1:A:97:PHE:CD1	1:A:97:PHE:C	2.92	0.43
2:F:121:ASP:O	2:F:125:ILE:HG13	2.19	0.43
1:A:5:ASN:O	1:A:9:VAL:HG23	2.19	0.43
1:G:26:GLN:NE2	1:G:189:ASP:OD1	2.52	0.43
1:G:112:LEU:HD13	1:G:122:SER:HB3	2.01	0.42
1:E:116:PRO:HD3	1:E:159:MET:SD	2.60	0.42
1:C:105:PHE:CD2	2:D:115:ARG:HD3	2.54	0.42
2:B:107:GLY:HA2	2:B:110:GLU:OE2	2.20	0.42
1:C:109:THR:OG1	1:C:153:GLU:HG2	2.20	0.42
1:A:108:LEU:HA	1:A:111:GLN:NE2	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:107:ASP:OD2	1:A:109:THR:HG23	2.20	0.41
1:A:122:SER:O	1:A:126:VAL:HG23	2.20	0.41
1:G:88:GLN:HE21	1:G:92:GLU:CD	2.21	0.41
1:C:136:ASN:HA	1:C:181:TRP:CZ2	2.55	0.41
1:C:8:LEU:HD23	1:C:8:LEU:HA	1.89	0.41
1:G:150:LEU:HD12	1:G:150:LEU:HA	1.92	0.41
1:A:183:GLN:HA	1:A:183:GLN:NE2	2.35	0.41
1:A:3:GLN:OE1	1:A:3:GLN:HA	2.20	0.41
1:G:183:GLN:NE2	1:G:183:GLN:CA	2.82	0.41
1:E:22:TYR:OH	1:G:156:ASP:OD1	2.32	0.41
1:C:166:ILE:O	1:C:170:MET:HG3	2.21	0.41
1:E:2:SER:HB2	1:G:86:VAL:CG2	2.39	0.41
1:C:140:ILE:O	1:C:143:PHE:HB3	2.20	0.41
1:C:147:GLY:HA3	1:C:170:MET:SD	2.61	0.41
1:C:16:LYS:HD3	1:C:19:GLN:OE1	2.21	0.41
1:C:83:MET:O	1:C:87:LYS:HG3	2.21	0.41
1:G:125:GLN:OE1	2:H:109:MET:SD	2.79	0.40
1:G:147:GLY:HA3	1:G:170:MET:SD	2.61	0.40
1:G:112:LEU:HD21	2:H:109:MET:HG3	2.04	0.40
2:H:121:ASP:O	2:H:125:ILE:HG13	2.21	0.40
1:A:24:TRP:CD1	1:A:25:SER:HB3	2.56	0.40
1:G:106:SER:O	1:G:107:ASP:O	2.40	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	A	139/153 (91%)	135 (97%)	1 (1%)	3 (2%)	8	12
1	C	139/153 (91%)	134 (96%)	3 (2%)	2 (1%)	13	23
1	E	139/153 (91%)	131 (94%)	4 (3%)	4 (3%)	5	7

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	G	139/153 (91%)	132 (95%)	4 (3%)	3 (2%)	8	12
2	B	22/31 (71%)	22 (100%)	0	0	100	100
2	D	22/31 (71%)	22 (100%)	0	0	100	100
2	F	22/31 (71%)	22 (100%)	0	0	100	100
2	H	22/31 (71%)	22 (100%)	0	0	100	100
All	All	644/736 (88%)	620 (96%)	12 (2%)	12 (2%)	9	15

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	107	ASP
1	A	109	THR
1	C	107	ASP
1	C	109	THR
1	E	107	ASP
1	E	109	THR
1	G	107	ASP
1	G	109	THR
1	E	4	SER
1	A	111	GLN
1	G	111	GLN
1	E	111	GLN

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	120/128 (94%)	114 (95%)	6 (5%)	28	51
1	C	120/128 (94%)	116 (97%)	4 (3%)	43	70
1	E	120/128 (94%)	114 (95%)	6 (5%)	28	51
1	G	120/128 (94%)	113 (94%)	7 (6%)	23	43
2	B	20/27 (74%)	20 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	D	20/27 (74%)	20 (100%)	0	100	100
2	F	20/27 (74%)	20 (100%)	0	100	100
2	H	20/27 (74%)	20 (100%)	0	100	100
All	All	560/620 (90%)	537 (96%)	23 (4%)	35	61

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	GLN
1	A	108	LEU
1	A	150	LEU
1	A	174	LEU
1	A	178	LEU
1	A	183	GLN
1	C	3	GLN
1	C	150	LEU
1	C	160	GLN
1	C	178	LEU
1	E	109	THR
1	E	113	HIS
1	E	150	LEU
1	E	160	GLN
1	E	178	LEU
1	E	183	GLN
1	G	3	GLN
1	G	4	SER
1	G	8	LEU
1	G	109	THR
1	G	153	GLU
1	G	160	GLN
1	G	183	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	111	GLN
1	A	125	GLN
1	A	128	ASN
1	A	183	GLN
1	A	197	ASN

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Mol	Chain	Res	Type
1	C	111	GLN
1	C	128	ASN
1	C	160	GLN
1	C	183	GLN
1	E	128	ASN
1	E	183	GLN
2	F	111	ASN
1	G	3	GLN
1	G	128	ASN
1	G	183	GLN
1	G	185	ASN

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 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	141/153 (92%)	-0.10	4 (2%) 53 56	36, 51, 93, 127	0
1	C	141/153 (92%)	0.08	5 (3%) 44 47	36, 54, 99, 131	0
1	E	141/153 (92%)	-0.02	7 (4%) 30 31	34, 48, 100, 127	0
1	G	141/153 (92%)	0.05	7 (4%) 30 31	36, 54, 103, 130	0
2	B	24/31 (77%)	0.57	2 (8%) 12 12	51, 68, 107, 115	0
2	D	24/31 (77%)	0.24	0 100 100	49, 66, 106, 112	0
2	F	24/31 (77%)	0.42	2 (8%) 12 12	45, 65, 108, 117	0
2	H	24/31 (77%)	0.44	2 (8%) 12 12	55, 72, 112, 117	0
All	All	660/736 (89%)	0.06	29 (4%) 35 37	34, 54, 109, 131	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	108	LEU	8.8
1	C	109	THR	8.7
1	G	107	ASP	6.9
1	C	110	SER	6.6
1	G	109	THR	6.5
1	G	108	LEU	6.1
1	E	109	THR	6.1
1	A	108	LEU	5.2
1	C	107	ASP	4.8
1	A	109	THR	4.8
1	E	113	HIS	4.8
1	A	110	SER	4.5
1	E	108	LEU	4.5
2	H	106	SER	4.3
2	B	106	SER	4.2
1	G	110	SER	4.2

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Mol	Chain	Res	Type	RSRZ
2	F	106	SER	4.0
1	E	112	LEU	4.0
1	A	107	ASP	3.8
1	G	106	SER	3.7
1	G	113	HIS	3.5
1	E	107	ASP	3.0
1	E	110	SER	3.0
2	F	107	GLY	2.7
2	B	114	ARG	2.6
2	H	107	GLY	2.4
1	E	106	SER	2.3
1	G	102	ARG	2.1
1	C	112	LEU	2.1

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.