



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 11:29 PM GMT

PDB ID : 1XB1
Title : The Structure of the BIR domain of IAP-like protein 2
Authors : Shin, H.; Renatus, M.; Eckelman, B.P.; Nunes, V.A.; Sampaio, C.A.M.;
Salvesen, G.S.
Deposited on : 2004-08-27
Resolution : 2.70 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
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

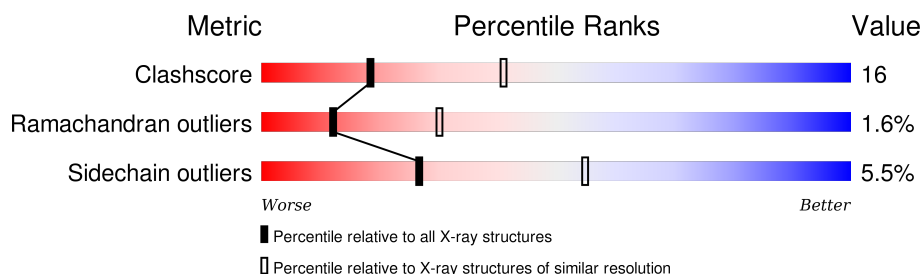
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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))
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)

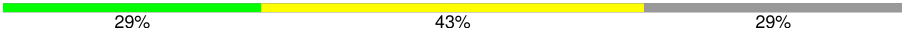


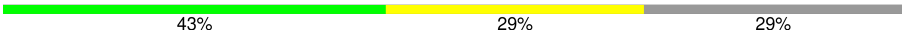

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey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

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	108	
1	B	108	
1	C	108	
1	D	108	
1	E	108	
1	F	108	
2	G	7	

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Mol	Chain	Length	Quality of chain
2	H	7	
2	I	7	
2	J	7	
2	K	7	
2	L	7	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5005 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 Baculoviral IAP repeat-containing protein 8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	90	Total	C	N	O	S	21	0	0
			739	475	128	131	5			
1	B	102	Total	C	N	O	S	13	0	0
			827	528	144	150	5			
1	C	90	Total	C	N	O	S	32	0	0
			739	475	128	131	5			
1	D	90	Total	C	N	O	S	15	0	0
			739	475	128	131	5			
1	E	102	Total	C	N	O	S	39	0	0
			827	528	144	150	5			
1	F	89	Total	C	N	O	S	38	0	0
			731	469	127	130	5			

There are 78 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	249	GLY	-	CLONING ARTIFACT	UNP Q96P09
A	250	SER	-	CLONING ARTIFACT	UNP Q96P09
A	251	HIS	-	CLONING ARTIFACT	UNP Q96P09
A	252	MET	-	CLONING ARTIFACT	UNP Q96P09
A	253	SER	-	SEE REMARK 999	UNP Q96P09
A	254	THR	-	SEE REMARK 999	UNP Q96P09
A	255	ASN	-	SEE REMARK 999	UNP Q96P09
A	256	LEU	-	SEE REMARK 999	UNP Q96P09
A	257	PRO	-	SEE REMARK 999	UNP Q96P09
A	258	ARG	-	SEE REMARK 999	UNP Q96P09
A	259	ASN	-	SEE REMARK 999	UNP Q96P09
A	260	PRO	-	SEE REMARK 999	UNP Q96P09
A	261	SER	-	SEE REMARK 999	UNP Q96P09
B	249	GLY	-	CLONING ARTIFACT	UNP Q96P09
B	250	SER	-	CLONING ARTIFACT	UNP Q96P09
B	251	HIS	-	CLONING ARTIFACT	UNP Q96P09
B	252	MET	-	CLONING ARTIFACT	UNP Q96P09

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Chain	Residue	Modelled	Actual	Comment	Reference
B	253	SER	-	SEE REMARK 999	UNP Q96P09
B	254	THR	-	SEE REMARK 999	UNP Q96P09
B	255	ASN	-	SEE REMARK 999	UNP Q96P09
B	256	LEU	-	SEE REMARK 999	UNP Q96P09
B	257	PRO	-	SEE REMARK 999	UNP Q96P09
B	258	ARG	-	SEE REMARK 999	UNP Q96P09
B	259	ASN	-	SEE REMARK 999	UNP Q96P09
B	260	PRO	-	SEE REMARK 999	UNP Q96P09
B	261	SER	-	SEE REMARK 999	UNP Q96P09
C	249	GLY	-	CLONING ARTIFACT	UNP Q96P09
C	250	SER	-	CLONING ARTIFACT	UNP Q96P09
C	251	HIS	-	CLONING ARTIFACT	UNP Q96P09
C	252	MET	-	CLONING ARTIFACT	UNP Q96P09
C	253	SER	-	SEE REMARK 999	UNP Q96P09
C	254	THR	-	SEE REMARK 999	UNP Q96P09
C	255	ASN	-	SEE REMARK 999	UNP Q96P09
C	256	LEU	-	SEE REMARK 999	UNP Q96P09
C	257	PRO	-	SEE REMARK 999	UNP Q96P09
C	258	ARG	-	SEE REMARK 999	UNP Q96P09
C	259	ASN	-	SEE REMARK 999	UNP Q96P09
C	260	PRO	-	SEE REMARK 999	UNP Q96P09
C	261	SER	-	SEE REMARK 999	UNP Q96P09
D	249	GLY	-	CLONING ARTIFACT	UNP Q96P09
D	250	SER	-	CLONING ARTIFACT	UNP Q96P09
D	251	HIS	-	CLONING ARTIFACT	UNP Q96P09
D	252	MET	-	CLONING ARTIFACT	UNP Q96P09
D	253	SER	-	SEE REMARK 999	UNP Q96P09
D	254	THR	-	SEE REMARK 999	UNP Q96P09
D	255	ASN	-	SEE REMARK 999	UNP Q96P09
D	256	LEU	-	SEE REMARK 999	UNP Q96P09
D	257	PRO	-	SEE REMARK 999	UNP Q96P09
D	258	ARG	-	SEE REMARK 999	UNP Q96P09
D	259	ASN	-	SEE REMARK 999	UNP Q96P09
D	260	PRO	-	SEE REMARK 999	UNP Q96P09
D	261	SER	-	SEE REMARK 999	UNP Q96P09
E	249	GLY	-	CLONING ARTIFACT	UNP Q96P09
E	250	SER	-	CLONING ARTIFACT	UNP Q96P09
E	251	HIS	-	CLONING ARTIFACT	UNP Q96P09
E	252	MET	-	CLONING ARTIFACT	UNP Q96P09
E	253	SER	-	SEE REMARK 999	UNP Q96P09
E	254	THR	-	SEE REMARK 999	UNP Q96P09
E	255	ASN	-	SEE REMARK 999	UNP Q96P09

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Chain	Residue	Modelled	Actual	Comment	Reference
E	256	LEU	-	SEE REMARK 999	UNP Q96P09
E	257	PRO	-	SEE REMARK 999	UNP Q96P09
E	258	ARG	-	SEE REMARK 999	UNP Q96P09
E	259	ASN	-	SEE REMARK 999	UNP Q96P09
E	260	PRO	-	SEE REMARK 999	UNP Q96P09
E	261	SER	-	SEE REMARK 999	UNP Q96P09
F	249	GLY	-	CLONING ARTIFACT	UNP Q96P09
F	250	SER	-	CLONING ARTIFACT	UNP Q96P09
F	251	HIS	-	CLONING ARTIFACT	UNP Q96P09
F	252	MET	-	CLONING ARTIFACT	UNP Q96P09
F	253	SER	-	SEE REMARK 999	UNP Q96P09
F	254	THR	-	SEE REMARK 999	UNP Q96P09
F	255	ASN	-	SEE REMARK 999	UNP Q96P09
F	256	LEU	-	SEE REMARK 999	UNP Q96P09
F	257	PRO	-	SEE REMARK 999	UNP Q96P09
F	258	ARG	-	SEE REMARK 999	UNP Q96P09
F	259	ASN	-	SEE REMARK 999	UNP Q96P09
F	260	PRO	-	SEE REMARK 999	UNP Q96P09
F	261	SER	-	SEE REMARK 999	UNP Q96P09

- Molecule 2 is a protein called Diablo homolog, mitochondrial.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	G	5	Total	C	N	O	0	0	0
			32	22	5	5			
2	H	5	Total	C	N	O	0	0	0
			32	22	5	5			
2	I	5	Total	C	N	O	0	0	0
			32	22	5	5			
2	J	5	Total	C	N	O	0	0	0
			32	22	5	5			
2	K	5	Total	C	N	O	0	0	0
			32	22	5	5			
2	L	5	Total	C	N	O	0	0	0
			32	22	5	5			

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	3	Total	Zn	0	0
			3	3		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	E	4	Total 4	Zn 4	0	0
3	B	4	Total 4	Zn 4	0	0
3	C	4	Total 4	Zn 4	0	0
3	A	4	Total 4	Zn 4	0	0
3	F	4	Total 4	Zn 4	0	0

- Molecule 4 is water.

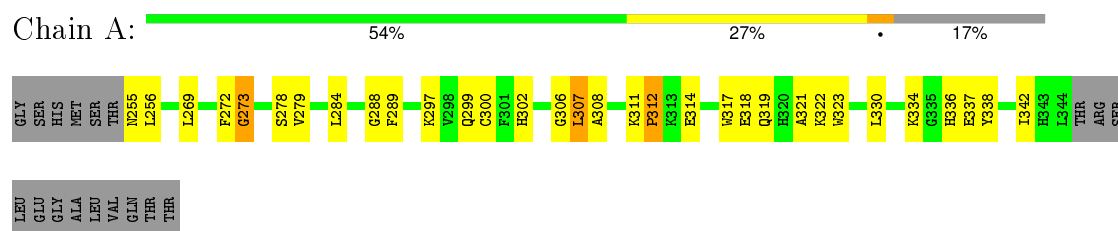
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	24	Total 24	O 24	0	0
4	B	42	Total 42	O 42	0	0
4	C	28	Total 28	O 28	1	0
4	D	39	Total 39	O 39	0	0
4	E	25	Total 25	O 25	0	0
4	F	22	Total 22	O 22	0	0
4	H	3	Total 3	O 3	0	0
4	I	4	Total 4	O 4	0	0
4	J	1	Total 1	O 1	0	0

3 Residue-property plots

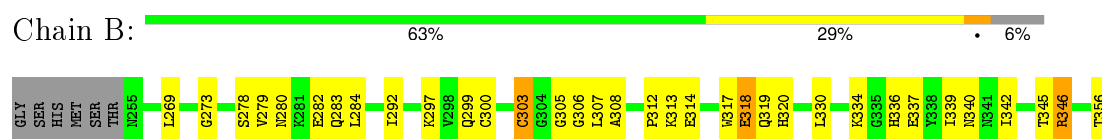
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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.

Note EDS was not executed.

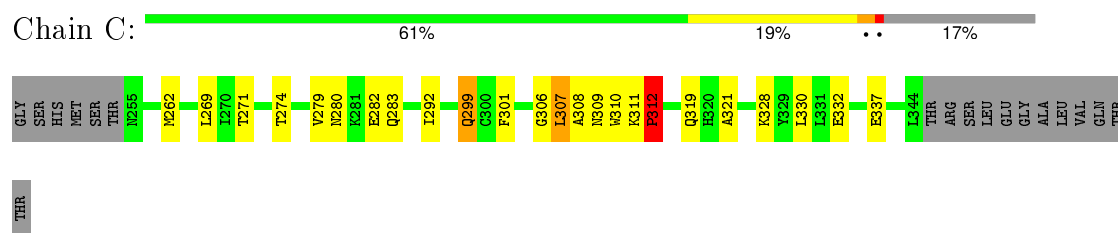
- Molecule 1: Baculoviral IAP repeat-containing protein 8



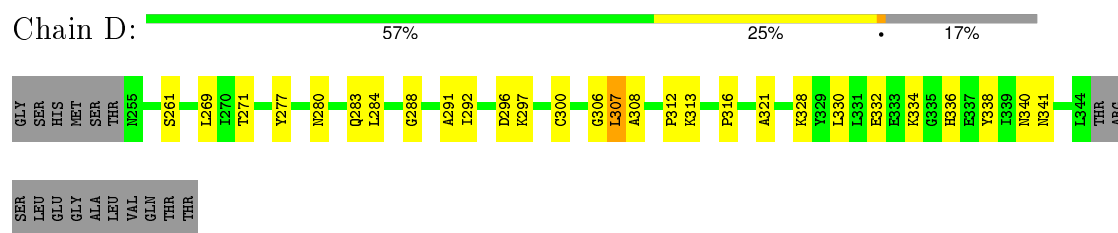
- Molecule 1: Baculoviral IAP repeat-containing protein 8



- Molecule 1: Baculoviral IAP repeat-containing protein 8

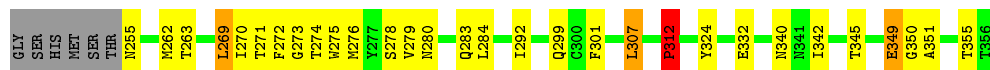


- Molecule 1: Baculoviral IAP repeat-containing protein 8



- Molecule 1: Baculoviral IAP repeat-containing protein 8

Chain E:  67% 24% • • 6%



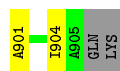
- Molecule 1: Baculoviral IAP repeat-containing protein 8

Chain F:  51% 28% • 18%



- Molecule 2: Diablo homolog, mitochondrial

Chain G:  43% 29% 29%



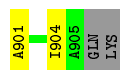
- Molecule 2: Diablo homolog, mitochondrial

Chain H:  29% 43% 29%



- Molecule 2: Diablo homolog, mitochondrial

Chain I:  43% 29% 29%



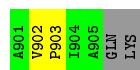
- Molecule 2: Diablo homolog, mitochondrial

Chain J:  57% 14% 29%



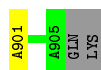
- Molecule 2: Diablo homolog, mitochondrial

Chain K:  43% 29% 29%



- Molecule 2: Diablo homolog, mitochondrial

Chain L:  57% 14% 29%



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	88.69 Å 88.69 Å 191.35 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	40.23 – 2.70	Depositor
% Data completeness (in resolution range)	93.0 (40.23-2.70)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.232 , 0.277	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5005	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section:
ZN

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	0.35	0/764	0.53	0/1034
1	B	0.38	0/852	0.55	0/1154
1	C	0.38	0/764	0.56	0/1034
1	D	0.39	0/764	0.57	0/1034
1	E	0.34	0/852	0.55	0/1154
1	F	0.33	0/756	0.53	0/1023
2	G	0.24	0/32	0.59	0/44
2	H	0.27	0/32	0.53	0/44
2	I	0.24	0/32	0.55	0/44
2	J	0.28	0/32	0.62	0/44
2	K	0.42	0/32	0.62	0/44
2	L	0.26	0/32	0.56	0/44
All	All	0.36	0/4944	0.55	0/6697

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	739	0	694	24	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	827	0	789	26	0
1	C	739	0	694	21	0
1	D	739	0	694	20	0
1	E	827	0	786	22	0
1	F	731	0	683	22	0
2	G	32	0	36	2	0
2	H	32	0	36	3	0
2	I	32	0	36	2	0
2	J	32	0	36	3	0
2	K	32	0	36	1	0
2	L	32	0	36	1	0
3	A	4	0	0	0	0
3	B	4	0	0	0	0
3	C	4	0	0	0	0
3	D	3	0	0	0	0
3	E	4	0	0	0	0
3	F	4	0	0	0	0
4	A	24	0	0	1	0
4	B	42	0	0	0	0
4	C	28	0	0	2	0
4	D	39	0	0	0	0
4	E	25	0	0	0	0
4	F	22	0	0	0	0
4	H	3	0	0	0	0
4	I	4	0	0	0	0
4	J	1	0	0	0	0
All	All	5005	0	4556	137	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 16.

All (137) 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:B:292:ILE:HD11	1:B:299:GLN:HB2	1.50	0.93
1:B:303:CYS:SG	1:B:320:HIS:NE2	2.44	0.90
1:B:345:THR:HG21	1:B:356:THR:HG21	1.56	0.88
1:B:280:ASN:HD21	1:B:282:GLU:HB2	1.43	0.83
1:E:278:SER:HB2	1:E:312:PRO:HA	1.64	0.80
1:C:280:ASN:HD22	1:C:283:GLN:H	1.31	0.78
1:C:280:ASN:ND2	1:C:283:GLN:H	1.81	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:280:ASN:ND2	1:B:283:GLN:H	1.85	0.74
1:C:311:LYS:HB3	1:C:312:PRO:HD2	1.70	0.74
1:F:334:LYS:HE2	1:F:334:LYS:HA	1.71	0.72
1:A:321:ALA:HA	1:A:330:LEU:HD21	1.73	0.70
1:C:321:ALA:HA	1:C:330:LEU:HD21	1.74	0.69
1:E:279:VAL:HG23	1:E:284:LEU:HD11	1.75	0.69
1:E:280:ASN:HD22	1:E:283:GLN:HG3	1.57	0.68
1:E:262:MET:HE3	1:E:271:THR:HG21	1.76	0.68
1:F:281:LYS:HG3	1:F:282:GLU:OE1	1.94	0.67
1:C:280:ASN:HD21	1:C:282:GLU:HB2	1.60	0.67
1:B:346:ARG:HG3	1:B:346:ARG:HH21	1.58	0.66
2:I:904:ILE:HD12	2:I:904:ILE:O	1.95	0.66
1:E:340:ASN:HB3	1:E:345:THR:O	1.96	0.66
1:D:308:ALA:HB2	2:J:904:ILE:HG21	1.78	0.64
1:D:280:ASN:HD22	1:D:283:GLN:HG3	1.63	0.63
1:F:308:ALA:O	2:L:901:ALA:HB1	1.99	0.63
1:F:265:TYR:CE2	1:F:269:LEU:HD12	2.34	0.62
1:C:274:THR:HG22	1:C:274:THR:O	2.02	0.60
1:A:317:TRP:CD2	1:A:334:LYS:HD2	2.36	0.60
1:F:272:PHE:O	1:F:274:THR:N	2.35	0.60
1:C:319:GLN:OE1	2:I:901:ALA:N	2.36	0.59
1:A:314:GLU:OE2	2:G:901:ALA:N	2.36	0.59
1:E:269:LEU:HD22	1:E:269:LEU:O	2.02	0.59
1:F:279:VAL:HG21	1:F:310:TRP:CE3	2.38	0.58
1:B:317:TRP:CG	1:B:334:LYS:HE3	2.38	0.58
1:E:274:THR:O	1:E:276:MET:HG3	2.03	0.58
1:E:263:THR:HA	1:E:301:PHE:CE1	2.38	0.58
1:A:318:GLU:HG2	1:A:342:ILE:HD12	1.84	0.58
1:A:311:LYS:HB2	1:A:314:GLU:HG3	1.84	0.57
1:B:280:ASN:ND2	1:B:283:GLN:N	2.52	0.55
1:B:330:LEU:HG	1:B:339:ILE:HD11	1.88	0.55
1:A:272:PHE:O	1:A:273:GLY:C	2.44	0.55
1:C:328:LYS:O	1:C:332:GLU:HG3	2.07	0.55
1:D:297:LYS:HD3	2:J:904:ILE:HD12	1.89	0.54
1:A:288:GLY:O	1:A:300:CYS:HA	2.07	0.54
1:C:279:VAL:HG21	1:C:310:TRP:CB	2.37	0.54
1:C:311:LYS:HB3	1:C:312:PRO:CD	2.38	0.53
1:B:280:ASN:ND2	1:B:282:GLU:HB2	2.17	0.53
1:C:279:VAL:HG21	1:C:310:TRP:HB2	1.91	0.53
1:A:279:VAL:HG23	1:A:284:LEU:HD11	1.91	0.53
1:E:262:MET:CE	1:E:271:THR:HG21	2.38	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:328:LYS:O	1:D:332:GLU:HG3	2.09	0.53
1:D:321:ALA:HA	1:D:330:LEU:HD21	1.91	0.53
1:F:273:GLY:C	1:F:275:TRP:H	2.12	0.52
1:E:307:LEU:HD13	1:E:324:TYR:HE1	1.74	0.52
1:D:277:TYR:CD2	1:D:296:ASP:HB3	2.45	0.52
1:F:280:ASN:HB3	1:F:283:GLN:HB2	1.93	0.51
1:C:337:GLU:HG3	4:C:146:HOH:O	2.09	0.51
1:C:292:ILE:HD11	1:C:299:GLN:HB2	1.93	0.51
1:E:292:ILE:HD11	1:E:299:GLN:HB2	1.92	0.51
1:F:298:VAL:HG22	1:F:310:TRP:HE1	1.75	0.51
1:E:355:THR:HG22	1:E:355:THR:O	2.10	0.50
1:B:297:LYS:HG2	1:B:308:ALA:HB2	1.94	0.50
1:B:346:ARG:HG3	1:B:346:ARG:NH2	2.26	0.50
1:D:312:PRO:O	1:D:313:LYS:HB2	2.12	0.50
1:B:319:GLN:OE1	2:H:901:ALA:N	2.46	0.49
1:A:255:ASN:HA	4:A:49:HOH:O	2.12	0.49
1:D:280:ASN:HD22	1:D:283:GLN:CG	2.23	0.49
1:D:280:ASN:ND2	1:D:283:GLN:H	2.10	0.49
1:F:284:LEU:HD23	1:F:316:PRO:HB3	1.95	0.49
1:E:263:THR:HA	1:E:301:PHE:CD1	2.48	0.49
1:F:277:TYR:CE2	1:F:296:ASP:HB3	2.47	0.48
1:E:272:PHE:HB3	1:E:275:TRP:HB2	1.95	0.48
1:D:277:TYR:CE2	1:D:296:ASP:HB3	2.48	0.48
1:A:278:SER:HB2	1:A:312:PRO:HA	1.94	0.48
1:C:280:ASN:HD22	1:C:283:GLN:N	2.06	0.48
1:F:330:LEU:HD12	1:F:334:LYS:HG3	1.96	0.47
1:D:292:ILE:HD12	1:D:297:LYS:HB3	1.96	0.47
1:D:306:GLY:O	1:D:307:LEU:HD13	2.13	0.47
1:B:306:GLY:HA3	2:H:904:ILE:HD12	1.96	0.47
1:C:321:ALA:HA	1:C:330:LEU:CD2	2.44	0.47
1:B:336:HIS:CD2	1:B:340:ASN:ND2	2.82	0.47
1:B:312:PRO:O	1:B:313:LYS:HB2	2.14	0.47
1:A:330:LEU:HD12	1:A:334:LYS:HG3	1.97	0.47
1:A:322:LYS:HD3	1:A:323:TRP:CE2	2.50	0.47
1:C:328:LYS:HA	1:C:328:LYS:HD3	1.66	0.46
1:A:336:HIS:O	1:A:337:GLU:C	2.54	0.46
1:D:334:LYS:HD3	1:D:338:TYR:CE2	2.50	0.46
1:F:312:PRO:O	1:F:313:LYS:HB2	2.16	0.46
1:F:292:ILE:O	1:F:292:ILE:HG22	2.15	0.46
1:B:292:ILE:HG22	1:B:292:ILE:O	2.15	0.46
1:C:280:ASN:ND2	1:C:283:GLN:N	2.59	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:279:VAL:CG2	1:E:284:LEU:HD11	2.45	0.45
1:F:265:TYR:HE1	1:F:285:ALA:HB3	1.81	0.45
1:B:314:GLU:OE2	2:H:901:ALA:N	2.49	0.45
1:A:306:GLY:HA3	2:G:904:ILE:HD12	1.98	0.45
1:F:265:TYR:CE1	1:F:285:ALA:HB3	2.50	0.45
1:A:255:ASN:HD22	1:A:302:HIS:HE1	1.64	0.45
1:E:349:GLU:C	1:E:351:ALA:H	2.20	0.45
1:D:288:GLY:O	1:D:300:CYS:HA	2.16	0.45
1:F:266:GLU:OE1	1:F:266:GLU:N	2.48	0.45
1:F:330:LEU:HG	1:F:339:ILE:HD11	1.99	0.45
1:B:318:GLU:HG2	1:B:342:ILE:HD12	1.98	0.45
1:A:297:LYS:HG2	1:A:308:ALA:HB2	1.99	0.44
1:C:262:MET:HB2	1:C:301:PHE:HB2	1.99	0.44
1:A:279:VAL:CG2	1:A:284:LEU:HD11	2.47	0.44
1:A:334:LYS:HE2	1:A:334:LYS:HA	1.99	0.44
1:D:308:ALA:CB	2:J:904:ILE:HG21	2.46	0.44
1:A:337:GLU:O	1:A:338:TYR:C	2.55	0.44
1:E:270:ILE:O	1:E:270:ILE:HG12	2.17	0.44
1:F:268:ARG:HD2	1:F:289:PHE:O	2.18	0.44
1:A:318:GLU:OE1	1:A:319:GLN:NE2	2.43	0.44
1:A:289:PHE:HA	1:A:299:GLN:O	2.17	0.44
1:A:255:ASN:HD22	1:A:302:HIS:CE1	2.36	0.44
1:F:317:TRP:HA	1:F:317:TRP:CE3	2.52	0.43
1:B:300:CYS:CB	1:B:303:CYS:SG	3.07	0.43
1:E:349:GLU:O	1:E:351:ALA:N	2.51	0.43
1:A:256:LEU:HD12	1:A:256:LEU:N	2.33	0.43
1:D:336:HIS:O	1:D:340:ASN:ND2	2.51	0.43
1:B:279:VAL:HG23	1:B:284:LEU:HD11	2.00	0.43
1:D:321:ALA:HA	1:D:330:LEU:CD2	2.49	0.42
1:E:292:ILE:HG22	1:E:292:ILE:O	2.18	0.42
1:C:308:ALA:O	1:C:309:ASN:HB2	2.19	0.42
1:B:280:ASN:HD22	1:B:283:GLN:H	1.62	0.42
1:E:292:ILE:HD11	1:E:299:GLN:CB	2.50	0.42
1:B:336:HIS:O	1:B:337:GLU:C	2.58	0.42
1:B:280:ASN:OD1	1:B:282:GLU:OE2	2.38	0.41
1:B:336:HIS:CD2	1:B:340:ASN:HD21	2.38	0.41
1:C:306:GLY:O	1:C:307:LEU:HD13	2.21	0.41
1:E:274:THR:O	1:E:275:TRP:C	2.59	0.41
1:D:328:LYS:HD3	1:D:328:LYS:HA	1.86	0.41
1:B:303:CYS:HB2	1:B:305:GLY:H	1.84	0.41
2:K:902:VAL:HG13	2:K:903:PRO:HD2	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:327:CYS:SG	1:F:329:TYR:HB3	2.61	0.41
1:E:342:ILE:HD13	1:E:342:ILE:HA	1.88	0.41
1:A:307:LEU:HA	1:A:307:LEU:HD12	1.89	0.41
1:D:284:LEU:HD23	1:D:316:PRO:HB3	2.04	0.41
1:C:271:THR:HA	4:C:175:HOH:O	2.20	0.40
1:D:271:THR:O	1:D:291:ALA:HB3	2.21	0.40
1:F:284:LEU:O	1:F:287:ALA:N	2.47	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	88/108 (82%)	80 (91%)	7 (8%)	1 (1%)	17	42
1	B	100/108 (93%)	90 (90%)	8 (8%)	2 (2%)	9	24
1	C	88/108 (82%)	80 (91%)	7 (8%)	1 (1%)	17	42
1	D	88/108 (82%)	85 (97%)	3 (3%)	0	100	100
1	E	100/108 (93%)	88 (88%)	8 (8%)	4 (4%)	4	8
1	F	87/108 (81%)	72 (83%)	14 (16%)	1 (1%)	17	42
2	G	3/7 (43%)	3 (100%)	0	0	100	100
2	H	3/7 (43%)	3 (100%)	0	0	100	100
2	I	3/7 (43%)	3 (100%)	0	0	100	100
2	J	3/7 (43%)	3 (100%)	0	0	100	100
2	K	3/7 (43%)	3 (100%)	0	0	100	100
2	L	3/7 (43%)	3 (100%)	0	0	100	100
All	All	569/690 (82%)	513 (90%)	47 (8%)	9 (2%)	12	30

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	349	GLU
1	F	273	GLY
1	A	273	GLY
1	B	273	GLY
1	B	278	SER
1	C	312	PRO
1	E	350	GLY
1	E	273	GLY
1	E	312	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	A	75/90 (83%)	72 (96%)	3 (4%)	38	69
1	B	85/90 (94%)	80 (94%)	5 (6%)	24	51
1	C	75/90 (83%)	71 (95%)	4 (5%)	28	57
1	D	75/90 (83%)	71 (95%)	4 (5%)	28	57
1	E	85/90 (94%)	80 (94%)	5 (6%)	24	51
1	F	74/90 (82%)	69 (93%)	5 (7%)	20	43
2	G	3/5 (60%)	3 (100%)	0	100	100
2	H	3/5 (60%)	2 (67%)	1 (33%)	0	0
2	I	3/5 (60%)	3 (100%)	0	100	100
2	J	3/5 (60%)	3 (100%)	0	100	100
2	K	3/5 (60%)	3 (100%)	0	100	100
2	L	3/5 (60%)	3 (100%)	0	100	100
All	All	487/570 (85%)	460 (94%)	27 (6%)	27	55

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	269	LEU
1	A	307	LEU

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Mol	Chain	Res	Type
1	A	312	PRO
1	B	269	LEU
1	B	303	CYS
1	B	307	LEU
1	B	318	GLU
1	B	346	ARG
2	H	902	VAL
1	C	269	LEU
1	C	299	GLN
1	C	307	LEU
1	C	312	PRO
1	D	261	SER
1	D	269	LEU
1	D	307	LEU
1	D	341	ASN
1	E	255	ASN
1	E	269	LEU
1	E	307	LEU
1	E	312	PRO
1	E	332	GLU
1	F	269	LEU
1	F	282	GLU
1	F	307	LEU
1	F	317	TRP
1	F	341	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	280	ASN
1	A	294	GLN
1	A	302	HIS
1	A	309	ASN
1	B	280	ASN
1	B	309	ASN
1	B	336	HIS
1	B	341	ASN
1	C	280	ASN
1	C	294	GLN
1	C	299	GLN
1	C	309	ASN
1	C	341	ASN

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Mol	Chain	Res	Type
1	D	280	ASN
1	D	309	ASN
1	E	280	ASN
1	E	283	GLN
1	E	302	HIS
1	E	309	ASN
1	E	341	ASN
1	F	280	ASN
1	F	309	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](#)

Of 23 ligands modelled in this entry, 23 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

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 ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.