



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:24 AM GMT

PDB ID : 2QJB  
Title : Crystal structure analysis of BMP-2 in complex with BMPR-IA variant IA/IB  
Authors : Kotzsch, A.; Mueller, T.D.  
Deposited on : 2007-07-06  
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](mailto: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.

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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) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
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) : trunk26865

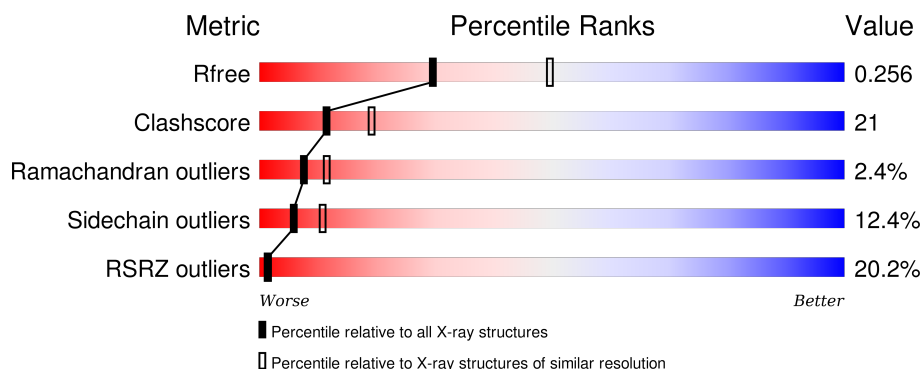
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 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))
$R_{free}$	91344	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (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  $\geq 3$ , 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	116	<div> <div>13%</div> <div>66%</div> <div>16%</div> <div>• •</div> <div>11%</div> </div>
1	B	116	<div> <div>14%</div> <div>65%</div> <div>19%</div> <div>5% •</div> <div>10%</div> </div>
2	C	135	<div> <div>24%</div> <div>37%</div> <div>21%</div> <div>5%</div> <div>37%</div> </div>
2	D	135	<div> <div>10%</div> <div>42%</div> <div>18%</div> <div>6% •</div> <div>33%</div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3052 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 Bone morphogenetic protein 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	103	Total	C	N	O	S	0	0	0
			805	510	135	151	9			
1	B	104	Total	C	N	O	S	0	0	0
			810	513	136	152	9			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MET	-	EXPRESSION TAG	UNP P12643
A	0	ALA	-	EXPRESSION TAG	UNP P12643
B	-1	MET	-	EXPRESSION TAG	UNP P12643
B	0	ALA	-	EXPRESSION TAG	UNP P12643

- Molecule 2 is a protein called Bone morphogenetic protein receptor type IA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	85	Total	C	N	O	S	0	0	0
			662	402	116	133	11			
2	D	90	Total	C	N	O	S	0	0	0
			695	425	121	138	11			

There are 42 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-5	GLY	-	EXPRESSION TAG	UNP P36894
C	-4	SER	-	EXPRESSION TAG	UNP P36894
C	-3	GLY	-	EXPRESSION TAG	UNP P36894
C	-2	ALA	-	EXPRESSION TAG	UNP P36894
C	-1	MET	-	EXPRESSION TAG	UNP P36894
C	0	ALA	-	EXPRESSION TAG	UNP P36894
C	42	HIS	GLY	ENGINEERED	UNP P36894
C	46	GLU	ASP	ENGINEERED	UNP P36894

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Chain	Residue	Modelled	Actual	Comment	Reference
C	61	THR	ALA	ENGINEERED	UNP P36894
C	62	MET	ILE	ENGINEERED	UNP P36894
C	74	THR	ALA	ENGINEERED	UNP P36894
C	78	LEU	MET	ENGINEERED	UNP P36894
C	79	GLY	LYS	ENGINEERED	UNP P36894
C	80	LEU	TYR	ENGINEERED	UNP P36894
C	88	ARG	LYS	ENGINEERED	UNP P36894
C	90	THR	SER	ENGINEERED	UNP P36894
C	92	ILE	LYS	ENGINEERED	UNP P36894
C	93	PRO	ALA	ENGINEERED	UNP P36894
C	94	HIS	GLN	ENGINEERED	UNP P36894
C	95	GLN	LEU	ENGINEERED	UNP P36894
C	98	SER	THR	ENGINEERED	UNP P36894
D	-5	GLY	-	EXPRESSION TAG	UNP P36894
D	-4	SER	-	EXPRESSION TAG	UNP P36894
D	-3	GLY	-	EXPRESSION TAG	UNP P36894
D	-2	ALA	-	EXPRESSION TAG	UNP P36894
D	-1	MET	-	EXPRESSION TAG	UNP P36894
D	0	ALA	-	EXPRESSION TAG	UNP P36894
D	42	HIS	GLY	ENGINEERED	UNP P36894
D	46	GLU	ASP	ENGINEERED	UNP P36894
D	61	THR	ALA	ENGINEERED	UNP P36894
D	62	MET	ILE	ENGINEERED	UNP P36894
D	74	THR	ALA	ENGINEERED	UNP P36894
D	78	LEU	MET	ENGINEERED	UNP P36894
D	79	GLY	LYS	ENGINEERED	UNP P36894
D	80	LEU	TYR	ENGINEERED	UNP P36894
D	88	ARG	LYS	ENGINEERED	UNP P36894
D	90	THR	SER	ENGINEERED	UNP P36894
D	92	ILE	LYS	ENGINEERED	UNP P36894
D	93	PRO	ALA	ENGINEERED	UNP P36894
D	94	HIS	GLN	ENGINEERED	UNP P36894
D	95	GLN	LEU	ENGINEERED	UNP P36894
D	98	SER	THR	ENGINEERED	UNP P36894

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total Cl 1 1	0	0

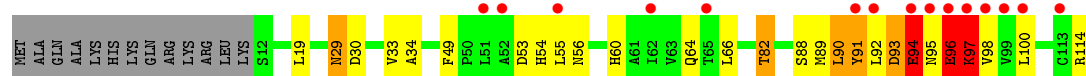
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	18	Total 18	O 18	0	0
4	B	24	Total 24	O 24	0	0
4	C	11	Total 11	O 11	0	0
4	D	26	Total 26	O 26	0	0

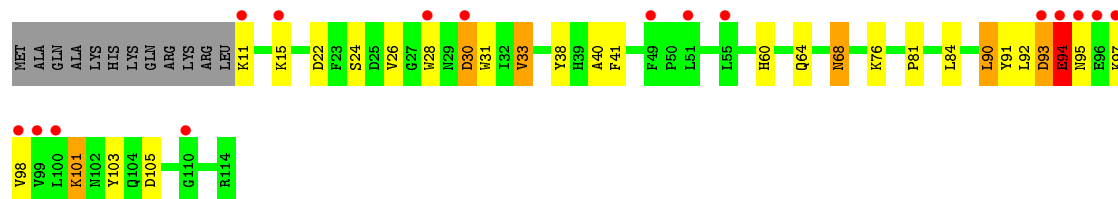
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

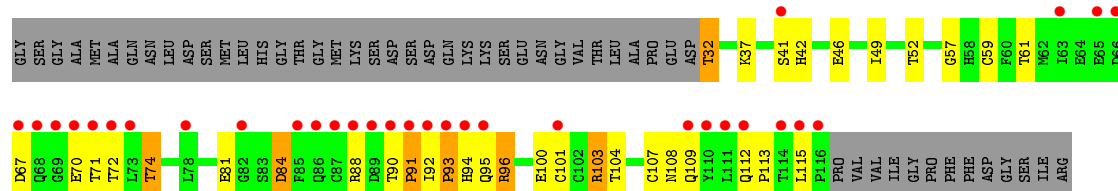
#### • Molecule 1: Bone morphogenetic protein 2



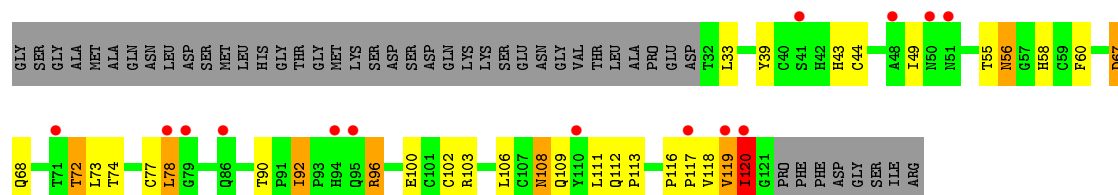
#### • Molecule 1: Bone morphogenetic protein 2



#### • Molecule 2: Bone morphogenetic protein receptor type IA



#### • Molecule 2: Bone morphogenetic protein receptor type IA



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	104.17Å 104.17Å 100.01Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	31.27 – 2.50 31.27 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.8 (31.27-2.50) 99.8 (31.27-2.50)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	0.08	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.94 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.204 , 0.263 0.201 , 0.256	Depositor DCC
$R_{free}$ test set	1097 reflections (5.41%)	DCC
Wilson B-factor (Å <sup>2</sup> )	54.5	Xtriage
Anisotropy	0.256	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 75.3	EDS
Estimated twinning fraction	0.043 for h,-h-k,-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 21357 reflections (0.005%)	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	3052	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	81.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.71% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

## 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: CL

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	0.70	0/828	0.74	0/1129
1	B	0.68	0/833	0.74	0/1136
2	C	0.59	0/676	0.74	0/919
2	D	0.65	0/710	0.79	0/967
All	All	0.66	0/3047	0.75	0/4151

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	805	0	761	28	0
1	B	810	0	763	34	0
2	C	662	0	607	28	0
2	D	695	0	646	31	0
3	B	1	0	0	0	0
4	A	18	0	0	1	0
4	B	24	0	0	5	0
4	C	11	0	0	2	0
4	D	26	0	0	5	0
All	All	3052	0	2777	118	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 21.

All (118) 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:C:32:THR:HA	4:C:135:HOH:O	1.44	1.17
1:B:11:LYS:HA	4:B:132:HOH:O	1.53	1.06
2:C:90:THR:HG22	2:C:91:PRO:HD2	1.45	0.99
2:C:37:LYS:HE2	2:C:52:THR:HG21	1.46	0.97
1:B:31:TRP:CD1	2:C:88:ARG:HG2	2.05	0.89
2:C:90:THR:HG22	2:C:91:PRO:CD	2.04	0.87
1:B:91:TYR:HE2	1:B:93:ASP:HB2	1.43	0.84
2:C:49:ILE:O	2:C:52:THR:HB	1.80	0.81
2:D:119:VAL:O	2:D:120:ILE:HG12	1.82	0.80
1:A:90:LEU:CD1	1:A:100:LEU:HD13	2.12	0.79
1:A:56:ASN:HD22	1:A:82:THR:HG21	1.47	0.78
2:C:92:ILE:HG23	2:C:96:ARG:NH1	2.00	0.76
2:C:42:HIS:HB2	2:C:74:THR:HG22	1.67	0.76
1:B:81:PRO:CB	1:B:84:LEU:HD11	2.19	0.73
1:A:56:ASN:ND2	1:A:82:THR:HG21	2.04	0.71
1:A:93:ASP:O	1:A:96:GLU:N	2.21	0.71
1:B:81:PRO:HB2	1:B:84:LEU:HD11	1.74	0.69
2:D:56:ASN:HD22	2:D:56:ASN:H	1.41	0.68
2:D:119:VAL:O	2:D:120:ILE:CG1	2.42	0.68
2:D:112:GLN:O	2:D:112:GLN:HG3	1.94	0.67
1:B:98:VAL:O	1:B:98:VAL:HG12	1.94	0.66
2:D:90:THR:HG22	2:D:92:ILE:H	1.59	0.66
1:B:30:ASP:OD2	1:B:30:ASP:N	2.25	0.65
2:D:108:ASN:HD22	2:D:108:ASN:N	1.95	0.64
2:C:90:THR:CG2	2:C:91:PRO:CD	2.76	0.64
2:C:57:GLY:O	2:C:103:ARG:HD3	1.98	0.63
2:C:109:GLN:O	2:C:109:GLN:HG2	1.97	0.62
1:B:15:LYS:HD2	4:B:127:HOH:O	2.00	0.62
1:A:54:HIS:HD2	4:A:126:HOH:O	1.84	0.59
2:D:60:PHE:CD2	2:D:78:LEU:HD13	2.37	0.59
2:D:118:VAL:O	2:D:120:ILE:HB	2.03	0.58
1:B:95:ASN:HB2	1:B:97:LYS:NZ	2.19	0.57
2:D:96:ARG:HG2	2:D:116:PRO:HD2	1.87	0.57
1:B:84:LEU:HD23	1:B:105:ASP:HB3	1.87	0.57
1:A:93:ASP:C	1:A:95:ASN:N	2.58	0.57
2:C:91:PRO:C	2:C:92:ILE:HG13	2.25	0.57
1:A:91:TYR:C	1:A:91:TYR:CD2	2.79	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:93:ASP:O	1:A:95:ASN:N	2.39	0.56
1:B:91:TYR:CE2	1:B:93:ASP:HB2	2.33	0.55
1:B:68:ASN:OD1	1:B:68:ASN:C	2.45	0.55
1:B:15:LYS:CD	4:B:127:HOH:O	2.53	0.55
1:B:92:LEU:HA	1:B:97:LYS:O	2.07	0.55
1:A:91:TYR:HD2	1:A:92:LEU:N	2.04	0.54
2:C:61:THR:OG1	2:C:108:ASN:HB2	2.07	0.53
2:D:67:ASP:N	2:D:67:ASP:OD1	2.42	0.53
1:B:81:PRO:HB2	1:B:84:LEU:CD1	2.36	0.53
1:B:81:PRO:HB3	1:B:84:LEU:HD11	1.91	0.53
2:D:72:THR:HG23	4:D:152:HOH:O	2.09	0.53
1:B:95:ASN:CB	1:B:97:LYS:NZ	2.72	0.52
2:D:102:CYS:SG	2:D:111:LEU:HD11	2.49	0.52
2:C:92:ILE:HG23	2:C:96:ARG:HH12	1.74	0.52
2:D:117:PRO:C	2:D:119:VAL:H	2.12	0.52
1:B:68:ASN:ND2	1:B:76:LYS:HG2	2.25	0.52
1:A:88:SER:O	1:A:89:MET:HG2	2.10	0.51
2:D:108:ASN:HD22	2:D:108:ASN:H	1.58	0.51
2:D:56:ASN:ND2	2:D:56:ASN:H	2.08	0.51
2:D:55:THR:HB	2:D:77:CYS:HB3	1.92	0.51
1:B:31:TRP:NE1	2:C:88:ARG:HG2	2.25	0.51
1:B:33:VAL:HG22	1:B:90:LEU:HB3	1.91	0.51
1:A:96:GLU:O	1:A:97:LYS:HG2	2.11	0.51
2:D:96:ARG:HD2	4:D:136:HOH:O	2.11	0.50
2:C:91:PRO:HD2	2:C:92:ILE:H	1.77	0.50
2:D:118:VAL:O	2:D:119:VAL:C	2.49	0.50
1:B:95:ASN:CB	1:B:97:LYS:HZ3	2.25	0.50
1:A:90:LEU:HD13	1:A:100:LEU:HD13	1.93	0.49
2:C:90:THR:CG2	2:C:91:PRO:HD2	2.31	0.49
1:B:95:ASN:HB2	1:B:97:LYS:HZ3	1.77	0.49
2:C:108:ASN:OD1	2:C:108:ASN:N	2.46	0.49
1:B:60:HIS:CE1	1:B:64:GLN:HE21	2.30	0.49
2:C:115:LEU:H	2:C:115:LEU:HD12	1.78	0.49
2:D:68:GLN:HA	2:D:68:GLN:OE1	2.13	0.48
1:A:60:HIS:CE1	1:A:64:GLN:HE21	2.32	0.48
1:B:81:PRO:CB	1:B:84:LEU:CD1	2.91	0.48
2:C:42:HIS:HB2	2:C:74:THR:CG2	2.42	0.47
2:D:56:ASN:HD22	2:D:56:ASN:N	2.03	0.47
2:D:58:HIS:N	2:D:58:HIS:CD2	2.82	0.47
1:A:49:PHE:CD2	1:A:49:PHE:C	2.88	0.47
2:C:84:ASP:OD1	2:C:84:ASP:N	2.40	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:43:HIS:CD2	2:D:74:THR:HB	2.50	0.47
1:B:40:ALA:O	1:B:41:PHE:HB2	2.15	0.47
1:A:55:LEU:O	1:A:56:ASN:C	2.52	0.46
2:C:93:PRO:HD2	2:C:96:ARG:HD3	1.97	0.46
2:C:59:CYS:O	2:C:101:CYS:HA	2.15	0.46
1:B:38:TYR:CZ	1:B:40:ALA:HB2	2.51	0.46
2:D:112:GLN:HB3	4:D:145:HOH:O	2.16	0.46
2:C:96:ARG:HD3	2:C:96:ARG:HA	1.65	0.45
1:B:22:ASP:OD1	1:B:24:SER:HB2	2.16	0.45
1:A:29:ASN:HD22	1:A:30:ASP:N	2.15	0.45
1:B:11:LYS:CB	4:B:128:HOH:O	2.64	0.45
1:B:84:LEU:HB3	1:B:105:ASP:HA	1.98	0.45
1:B:101:LYS:HD2	1:B:103:TYR:CZ	2.53	0.44
1:A:97:LYS:HE2	1:A:97:LYS:HA	1.99	0.44
2:D:112:GLN:CB	4:D:145:HOH:O	2.65	0.44
1:A:91:TYR:C	1:A:91:TYR:HD2	2.21	0.43
1:A:93:ASP:C	1:A:95:ASN:H	2.20	0.43
2:C:104:THR:O	2:C:107:CYS:HB3	2.18	0.43
2:D:100:GLU:HG3	2:D:113:PRO:HB3	1.99	0.43
2:C:100:GLU:HG3	2:C:113:PRO:HG3	2.00	0.43
1:A:93:ASP:O	1:A:94:GLU:C	2.57	0.43
2:D:39:TYR:H	2:D:108:ASN:HD21	1.66	0.43
1:A:91:TYR:CD2	1:A:92:LEU:N	2.86	0.43
1:A:33:VAL:O	1:A:34:ALA:HB2	2.19	0.43
2:D:106:LEU:O	2:D:109:GLN:HG2	2.19	0.42
1:B:76:LYS:HB3	4:B:125:HOH:O	2.18	0.42
1:B:93:ASP:O	1:B:94:GLU:C	2.57	0.42
1:A:92:LEU:HB3	1:A:96:GLU:HA	2.02	0.42
2:D:117:PRO:C	2:D:119:VAL:N	2.73	0.42
1:A:96:GLU:O	1:A:97:LYS:CG	2.67	0.42
1:B:31:TRP:O	1:B:91:TYR:HB2	2.20	0.42
1:A:60:HIS:HE1	1:A:64:GLN:HE21	1.68	0.42
1:A:66:LEU:HD13	1:B:28:TRP:HZ2	1.85	0.41
1:A:53:ASP:C	1:A:53:ASP:OD1	2.58	0.41
2:C:46:GLU:HB2	4:C:137:HOH:O	2.21	0.41
2:D:44:CYS:HB2	4:D:141:HOH:O	2.21	0.41
2:C:112:GLN:HA	2:C:113:PRO:HD3	1.86	0.41
1:A:33:VAL:O	1:A:33:VAL:HG12	2.21	0.40
2:D:56:ASN:N	2:D:56:ASN:ND2	2.69	0.40
2:D:112:GLN:HA	2:D:113:PRO:HD2	1.77	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	101/116 (87%)	92 (91%)	5 (5%)	4 (4%)	4	4
1	B	102/116 (88%)	96 (94%)	5 (5%)	1 (1%)	19	34
2	C	83/135 (62%)	76 (92%)	5 (6%)	2 (2%)	7	11
2	D	88/135 (65%)	80 (91%)	6 (7%)	2 (2%)	8	12
All	All	374/502 (74%)	344 (92%)	21 (6%)	9 (2%)	7	11

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	94	GLU
1	A	98	VAL
2	C	91	PRO
2	D	120	ILE
1	A	96	GLU
1	A	97	LYS
1	B	94	GLU
2	D	119	VAL
2	C	93	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	92/103 (89%)	82 (89%)	10 (11%)	8	15
1	B	92/103 (89%)	84 (91%)	8 (9%)	13	24

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	C	79/119 (66%)	66 (84%)	13 (16%)	3	5
2	D	83/119 (70%)	71 (86%)	12 (14%)	4	7
All	All	346/444 (78%)	303 (88%)	43 (12%)	6	11

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	LEU
1	A	29	ASN
1	A	82	THR
1	A	90	LEU
1	A	91	TYR
1	A	93	ASP
1	A	94	GLU
1	A	96	GLU
1	A	97	LYS
1	A	114	ARG
1	B	26	VAL
1	B	30	ASP
1	B	33	VAL
1	B	68	ASN
1	B	90	LEU
1	B	93	ASP
1	B	94	GLU
1	B	101	LYS
2	C	32	THR
2	C	41	SER
2	C	67	ASP
2	C	70	GLU
2	C	71	THR
2	C	72	THR
2	C	74	THR
2	C	81	GLU
2	C	84	ASP
2	C	94	HIS
2	C	95	GLN
2	C	96	ARG
2	C	103	ARG
2	D	33	LEU
2	D	49	ILE
2	D	56	ASN

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Mol	Chain	Res	Type
2	D	67	ASP
2	D	72	THR
2	D	73	LEU
2	D	78	LEU
2	D	92	ILE
2	D	96	ARG
2	D	103	ARG
2	D	108	ASN
2	D	120	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	29	ASN
1	A	56	ASN
1	A	64	GLN
1	A	95	ASN
1	B	64	GLN
2	C	58	HIS
2	C	68	GLN
2	D	43	HIS
2	D	56	ASN
2	D	58	HIS
2	D	108	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

### 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

### 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry

Of 1 ligands modelled in this entry, 1 is 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

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 ⓘ

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	103/116 (88%)	1.03	15 (14%)	3 3	72, 80, 99, 113	0
1	B	104/116 (89%)	1.21	16 (15%)	3 3	70, 80, 98, 114	0
2	C	85/135 (62%)	1.98	32 (37%)	0 0	74, 81, 110, 118	0
2	D	90/135 (66%)	1.17	14 (15%)	3 2	70, 78, 87, 91	0
All	All	382/502 (76%)	1.32	77 (20%)	1 1	70, 80, 99, 118	0

All (77) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	94	HIS	10.5
2	D	119	VAL	9.0
2	C	95	GLN	7.2
1	B	95	ASN	7.1
1	A	94	GLU	6.8
2	C	68	GLN	6.8
2	C	90	THR	6.5
2	C	115	LEU	6.4
2	C	69	GLY	5.8
2	C	110	TYR	5.7
2	D	120	ILE	5.6
1	A	95	ASN	5.3
2	C	67	ASP	5.3
2	C	92	ILE	5.1
1	B	94	GLU	4.8
1	B	30	ASP	4.8
2	C	116	PRO	4.8
2	C	93	PRO	4.8
1	B	96	GLU	4.8
2	C	114	THR	4.5
2	D	78	LEU	4.4

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Mol	Chain	Res	Type	RSRZ
1	B	93	ASP	4.3
2	C	71	THR	4.3
2	C	73	LEU	4.2
1	B	11	LYS	4.1
2	C	88	ARG	4.1
2	C	70	GLU	4.0
1	B	97	LYS	4.0
2	D	95	GLN	3.6
2	C	66	ASP	3.6
2	C	63	ILE	3.4
2	C	65	GLU	3.4
1	B	98	VAL	3.3
1	A	51	LEU	3.3
2	C	109	GLN	3.3
1	B	99	VAL	3.2
2	C	85	PHE	3.2
2	D	79	GLY	3.2
1	A	98	VAL	3.2
1	A	96	GLU	3.2
1	A	91	TYR	3.2
2	D	110	TYR	3.1
1	A	92	LEU	3.0
2	D	94	HIS	3.0
1	A	99	VAL	3.0
2	C	86	GLN	3.0
1	A	65	THR	2.9
2	C	72	THR	2.9
2	C	91	PRO	2.9
2	C	41	SER	2.9
1	B	55	LEU	2.8
1	A	97	LYS	2.7
1	B	15	LYS	2.5
1	A	62	ILE	2.5
1	A	55	LEU	2.5
2	C	111	LEU	2.5
1	A	52	ALA	2.5
2	D	50	ASN	2.5
1	B	110	GLY	2.3
1	A	100	LEU	2.3
2	D	86	GLN	2.3
2	C	87	CYS	2.3
2	D	117	PRO	2.3

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Mol	Chain	Res	Type	RSRZ
2	C	112	GLN	2.3
2	C	89	ASP	2.3
2	D	48	ALA	2.2
1	B	100	LEU	2.2
2	D	41	SER	2.2
2	C	82	GLY	2.2
2	C	101	CYS	2.2
1	B	51	LEU	2.2
1	A	113	CYS	2.1
1	B	49	PHE	2.1
1	B	28	TRP	2.0
2	C	78	LEU	2.0
2	D	71	THR	2.0
2	D	51	ASN	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 carbohydrates in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	CL	B	1	1/1	0.93	0.12	-2.24	84,84,84,84	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.