



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 07:37 am GMT

PDB ID : 3CXB  
Title : Crystal Structure of sifa and skip  
Authors : Huang, Z.; Chai, J.  
Deposited on : 2008-04-24  
Resolution : 2.60 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  
Xtriage (Phenix) : 1.9-1692  
EDS : trunk28620  
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) : recalc28949

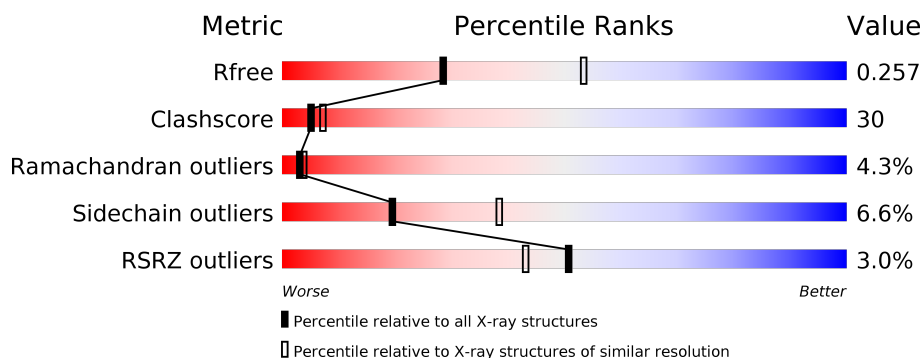
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.60 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	2542 (2.60-2.60)
Clashscore	112137	2895 (2.60-2.60)
Ramachandran outliers	110173	2848 (2.60-2.60)
Sidechain outliers	110143	2848 (2.60-2.60)
RSRZ outliers	101464	2550 (2.60-2.60)

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	336	<div> <div>2%</div> <div> <div></div> <div>46%</div> <div>39%</div> <div>• • 10%</div> </div> </div>
2	B	112	<div> <div>4%</div> <div> <div></div> <div>46%</div> <div>40%</div> <div>5% 8%</div> </div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3335 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 Protein sifA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	302	Total	C	N	O	S	0	0	0
			2442	1549	421	456	16			

- Molecule 2 is a protein called Pleckstrin homology domain-containing family M member 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	103	Total	C	N	O	S	0	0	0
			773	478	140	146	9			

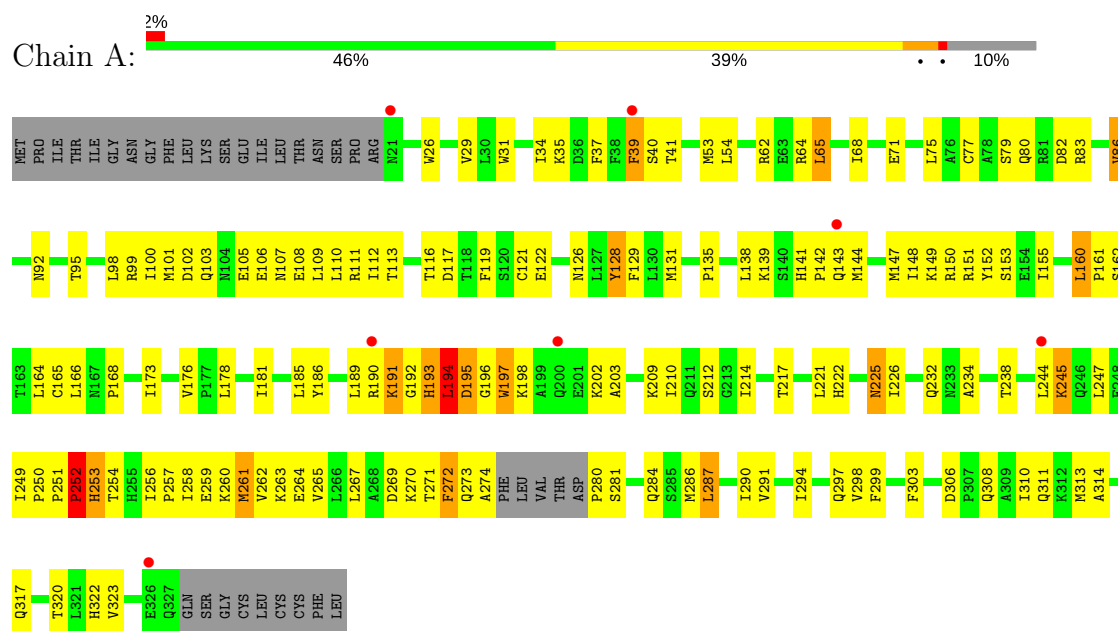
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	93	Total	O	0	0
			93	93		
3	B	27	Total	O	0	0
			27	27		

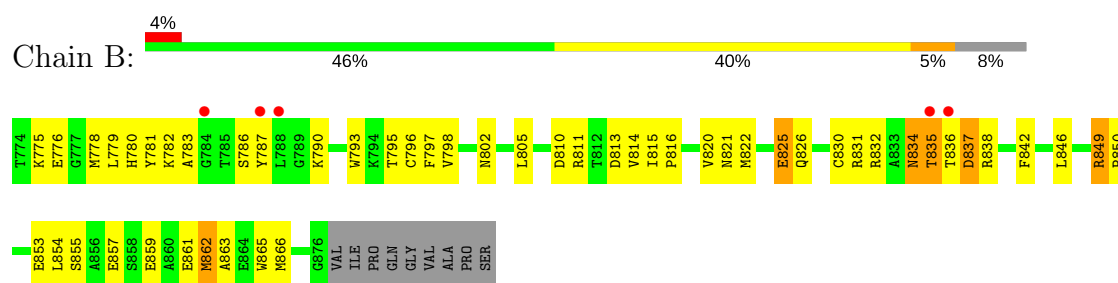
### 3 Residue-property plots

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



#### • Molecule 2: Pleckstrin homology domain-containing family M member 2



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	91.69Å 110.68Å 45.53Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.60 35.30 – 2.50	Depositor EDS
% Data completeness (in resolution range)	95.3 (20.00-2.60) 86.6 (35.30-2.50)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	0.03	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.77 (at 2.51Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.258 , 0.288 0.261 , 0.257	Depositor DCC
$R_{free}$ test set	628 reflections (5.05%)	DCC
Wilson B-factor (Å <sup>2</sup> )	49.0	Xtriage
Anisotropy	0.747	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.29 , 58.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	3335	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.99% 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, 0.333 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	0.43	0/2493	0.67	1/3369 (0.0%)
2	B	0.49	1/788 (0.1%)	0.65	0/1060
All	All	0.45	1/3281 (0.0%)	0.67	1/4429 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	825	GLU	CD-OE2	6.91	1.33	1.25

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	281	SER	N-CA-C	8.73	134.58	111.00

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	2442	0	2441	144	0
2	B	773	0	731	48	0
3	A	93	0	0	9	0
3	B	27	0	0	5	0
All	All	3335	0	3172	189	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 30.

All (189) 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:A:191:LYS:HG3	1:A:192:GLY:H	1.32	0.94
1:A:202:LYS:HG2	1:A:244:LEU:HD21	1.49	0.93
1:A:203:ALA:HA	1:A:244:LEU:HD23	1.53	0.90
1:A:153:SER:HB2	1:A:160:LEU:HD12	1.58	0.83
2:B:830:CYS:HB3	2:B:866:MET:HE1	1.63	0.81
1:A:238:THR:HG22	1:A:317:GLN:HG3	1.62	0.80
1:A:267:LEU:HD22	1:A:272:PHE:HE2	1.45	0.80
2:B:825:GLU:CA	2:B:826:GLN:N	2.46	0.79
1:A:297:GLN:HB3	3:A:414:HOH:O	1.84	0.76
1:A:65:LEU:HD22	1:A:112:ILE:HD13	1.66	0.76
1:A:191:LYS:C	1:A:193:HIS:H	1.87	0.75
1:A:191:LYS:HG3	1:A:192:GLY:N	2.01	0.74
1:A:194:LEU:HD23	1:A:195:ASP:N	2.03	0.74
1:A:164:LEU:HD11	1:A:166:LEU:HG	1.71	0.73
1:A:267:LEU:HD22	1:A:272:PHE:CE2	2.24	0.72
1:A:234:ALA:O	1:A:238:THR:HG23	1.89	0.71
1:A:267:LEU:HA	1:A:272:PHE:CZ	2.28	0.68
1:A:164:LEU:HB2	1:A:178:LEU:HD11	1.76	0.67
1:A:31:TRP:O	1:A:35:LYS:HB2	1.94	0.67
1:A:164:LEU:CD1	1:A:166:LEU:HG	2.24	0.66
1:A:113:THR:HG21	3:A:350:HOH:O	1.94	0.66
1:A:191:LYS:CG	1:A:192:GLY:H	2.01	0.66
2:B:830:CYS:HB3	2:B:866:MET:CE	2.24	0.66
1:A:287:LEU:O	1:A:291:VAL:HG23	1.96	0.65
1:A:164:LEU:HD13	1:A:164:LEU:C	2.16	0.65
1:A:210:ILE:O	1:A:214:ILE:HG12	1.96	0.65
1:A:250:PRO:O	1:A:252:PRO:HD3	1.96	0.65
1:A:259:GLU:O	1:A:263:LYS:HG3	1.95	0.65
2:B:853:GLU:HG3	3:B:36:HOH:O	1.96	0.64
2:B:862:MET:HG3	2:B:863:ALA:N	2.13	0.63
1:A:286:MET:O	1:A:290:ILE:HG12	1.98	0.63
1:A:101:MET:HE1	1:A:108:GLU:HG2	1.82	0.62
1:A:26:TRP:HA	1:A:29:VAL:CG1	2.30	0.61
2:B:778:MET:O	2:B:779:LEU:HD12	2.00	0.61
1:A:26:TRP:HA	1:A:29:VAL:HG12	1.81	0.61
1:A:251:PRO:O	1:A:252:PRO:C	2.39	0.61
2:B:782:LYS:HD3	3:B:36:HOH:O	2.00	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:202:LYS:HB2	1:A:303:PHE:CE2	2.35	0.61
1:A:317:GLN:O	1:A:320:THR:HG22	2.01	0.61
1:A:82:ASP:HA	3:A:362:HOH:O	1.99	0.61
2:B:835:THR:O	2:B:838:ARG:HG2	2.01	0.61
1:A:151:ARG:HG2	1:A:162:SER:HB3	1.83	0.61
1:A:99:ARG:HD2	1:A:101:MET:HE1	1.84	0.60
1:A:232:GLN:OE1	1:A:232:GLN:HA	2.02	0.59
1:A:86:VAL:HG22	1:A:152:TYR:HB2	1.83	0.59
2:B:778:MET:C	2:B:779:LEU:HD12	2.23	0.59
2:B:793:TRP:CZ3	2:B:855:SER:HB2	2.37	0.59
2:B:805:LEU:HB3	2:B:820:VAL:HG23	1.85	0.59
1:A:226:ILE:HG12	1:A:322:HIS:ND1	2.17	0.58
1:A:191:LYS:C	1:A:193:HIS:N	2.56	0.58
1:A:238:THR:HG21	1:A:314:ALA:HA	1.85	0.58
1:A:153:SER:HB2	1:A:160:LEU:CD1	2.32	0.58
1:A:251:PRO:O	1:A:253:HIS:N	2.37	0.57
2:B:813:ASP:O	2:B:815:ILE:N	2.36	0.57
1:A:53:MET:HG2	1:A:54:LEU:HG	1.86	0.57
1:A:65:LEU:CD2	1:A:112:ILE:HD13	2.33	0.57
1:A:116:THR:O	1:A:135:PRO:HG3	2.04	0.57
2:B:838:ARG:HG3	2:B:838:ARG:O	2.04	0.56
1:A:267:LEU:HB3	1:A:272:PHE:HZ	1.69	0.56
1:A:267:LEU:CA	1:A:272:PHE:CZ	2.89	0.56
1:A:257:PRO:O	1:A:261:MET:HB2	2.05	0.56
1:A:264:GLU:HG3	3:A:375:HOH:O	2.05	0.55
1:A:317:GLN:HA	1:A:320:THR:HG22	1.88	0.55
1:A:287:LEU:HD22	1:A:287:LEU:O	2.07	0.55
2:B:779:LEU:HD11	2:B:865:TRP:NE1	2.23	0.54
1:A:54:LEU:HD21	1:A:121:CYS:SG	2.49	0.53
1:A:274:ALA:C	1:A:280:PRO:O	2.47	0.53
1:A:256:ILE:HG23	1:A:260:LYS:HD2	1.91	0.53
1:A:155:ILE:HD11	1:A:186:TYR:CE1	2.44	0.52
1:A:39:PHE:H	1:A:39:PHE:HD1	1.57	0.52
1:A:77:CYS:HB3	3:A:397:HOH:O	2.09	0.52
1:A:53:MET:HE1	1:A:54:LEU:HD11	1.92	0.52
2:B:781:TYR:HA	2:B:853:GLU:O	2.09	0.51
1:A:53:MET:CE	1:A:110:LEU:HD11	2.40	0.51
1:A:147:MET:HG3	1:A:165:CYS:HB2	1.91	0.51
1:A:122:GLU:HA	1:A:126:ASN:O	2.10	0.51
1:A:209:LYS:HE3	1:A:308:GLN:NE2	2.25	0.51
1:A:245:LYS:NZ	1:A:247:LEU:HD12	2.25	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:138:LEU:HD21	1:A:290:ILE:CD1	2.41	0.50
1:A:166:LEU:O	1:A:173:ILE:HG13	2.12	0.50
1:A:287:LEU:C	1:A:287:LEU:HD13	2.32	0.50
1:A:103:GLN:C	1:A:105:GLU:H	2.15	0.50
1:A:149:LYS:HE3	1:A:165:CYS:SG	2.52	0.49
2:B:834:ASN:H	2:B:834:ASN:ND2	2.09	0.49
2:B:815:ILE:O	2:B:815:ILE:HG23	2.13	0.49
1:A:160:LEU:HD13	1:A:160:LEU:O	2.12	0.49
2:B:835:THR:HG22	2:B:836:THR:N	2.26	0.49
1:A:274:ALA:O	1:A:280:PRO:O	2.30	0.49
2:B:805:LEU:HB3	2:B:820:VAL:CG2	2.43	0.49
1:A:225:ASN:HB2	3:A:409:HOH:O	2.13	0.48
1:A:284:GLN:N	1:A:284:GLN:OE1	2.46	0.48
1:A:64:ARG:O	1:A:68:ILE:HG13	2.13	0.48
1:A:267:LEU:O	1:A:272:PHE:CE1	2.66	0.48
1:A:197:TRP:CD1	1:A:198:LYS:N	2.82	0.48
1:A:245:LYS:HZ2	1:A:247:LEU:HD12	1.79	0.48
2:B:849:ARG:HD2	2:B:850:PRO:O	2.14	0.48
1:A:86:VAL:CG2	1:A:152:TYR:HB2	2.44	0.48
1:A:105:GLU:O	1:A:105:GLU:HG2	2.13	0.47
2:B:835:THR:HG22	2:B:837:ASP:H	1.79	0.47
1:A:139:LYS:HA	3:A:405:HOH:O	2.14	0.47
1:A:71:GLU:O	1:A:75:LEU:HG	2.14	0.47
1:A:99:ARG:HD2	1:A:101:MET:CE	2.44	0.47
2:B:781:TYR:CE2	2:B:796:CYS:HB3	2.50	0.47
1:A:191:LYS:O	1:A:193:HIS:N	2.45	0.47
2:B:779:LEU:HD11	2:B:865:TRP:HE1	1.80	0.46
1:A:142:PRO:HG2	1:A:143:GLN:H	1.80	0.46
1:A:86:VAL:HB	1:A:98:LEU:CD2	2.45	0.46
1:A:269:ASP:O	1:A:272:PHE:CE1	2.68	0.46
2:B:813:ASP:C	2:B:815:ILE:H	2.18	0.46
1:A:267:LEU:HB3	1:A:272:PHE:CZ	2.50	0.46
1:A:186:TYR:CD1	1:A:189:LEU:HD12	2.51	0.46
1:A:101:MET:CE	1:A:108:GLU:HG2	2.46	0.46
1:A:217:THR:HG22	1:A:221:LEU:HD12	1.97	0.46
2:B:813:ASP:O	2:B:815:ILE:HG22	2.16	0.46
1:A:108:GLU:OE2	1:A:111:ARG:HD2	2.17	0.46
1:A:191:LYS:CG	1:A:192:GLY:N	2.69	0.45
1:A:34:ILE:O	1:A:37:PHE:HB2	2.15	0.45
1:A:129:PHE:HB3	2:B:831:ARG:CB	2.47	0.45
1:A:80:GLN:O	1:A:83:ARG:HB2	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:129:PHE:HB3	2:B:831:ARG:HB2	1.99	0.45
2:B:832:ARG:NH1	2:B:859:GLU:OE2	2.49	0.45
2:B:793:TRP:CZ2	2:B:838:ARG:HD2	2.52	0.45
1:A:128:TYR:O	1:A:129:PHE:HB3	2.16	0.45
1:A:147:MET:CG	1:A:165:CYS:HB2	2.46	0.45
2:B:853:GLU:HG3	3:B:96:HOH:O	2.16	0.45
1:A:259:GLU:HB3	1:A:263:LYS:HE3	1.99	0.45
1:A:317:GLN:CA	1:A:320:THR:HG22	2.46	0.45
2:B:810:ASP:OD1	2:B:813:ASP:HB2	2.17	0.44
1:A:79:SER:HB2	1:A:80:GLN:HE21	1.82	0.44
1:A:294:ILE:O	1:A:298:VAL:HG23	2.16	0.44
1:A:138:LEU:HD21	1:A:290:ILE:HD13	2.00	0.44
1:A:164:LEU:HD13	1:A:165:CYS:N	2.32	0.44
1:A:226:ILE:HG12	1:A:322:HIS:CE1	2.52	0.44
1:A:267:LEU:O	1:A:272:PHE:CZ	2.70	0.44
2:B:849:ARG:HH11	2:B:850:PRO:HG2	1.83	0.44
1:A:262:VAL:HA	1:A:265:VAL:HG22	1.99	0.44
1:A:310:ILE:HD12	1:A:310:ILE:N	2.33	0.44
1:A:212:SER:HB2	1:A:311:GLN:HE21	1.83	0.44
1:A:131:MET:HE2	3:A:356:HOH:O	2.18	0.43
1:A:139:LYS:HD3	3:A:413:HOH:O	2.17	0.43
1:A:160:LEU:C	1:A:160:LEU:HD22	2.39	0.43
2:B:842:PHE:CE1	2:B:854:LEU:HD12	2.54	0.43
1:A:141:HIS:HA	1:A:142:PRO:HD2	1.86	0.43
1:A:62:ARG:HB3	1:A:119:PHE:CE2	2.54	0.43
1:A:164:LEU:C	1:A:164:LEU:CD1	2.85	0.43
1:A:258:ILE:O	1:A:262:VAL:HG12	2.19	0.43
2:B:778:MET:HE2	2:B:811:ARG:HE	1.82	0.43
2:B:798:VAL:CG1	2:B:805:LEU:HD11	2.49	0.43
2:B:781:TYR:CD2	2:B:796:CYS:HB3	2.54	0.43
1:A:153:SER:CB	1:A:160:LEU:HD12	2.39	0.43
1:A:269:ASP:C	1:A:271:THR:H	2.22	0.43
1:A:209:LYS:HE3	1:A:308:GLN:HE21	1.83	0.43
2:B:780:HIS:CD2	2:B:795:THR:HG22	2.54	0.43
2:B:861:GLU:HG3	3:B:4:HOH:O	2.18	0.43
1:A:194:LEU:O	1:A:196:GLY:N	2.52	0.42
2:B:857:GLU:HG2	2:B:861:GLU:OE2	2.18	0.42
2:B:778:MET:CE	2:B:811:ARG:HE	2.32	0.42
2:B:775:LYS:HE3	2:B:775:LYS:HB2	1.89	0.42
1:A:306:ASP:O	1:A:310:ILE:HD13	2.20	0.42
1:A:129:PHE:HB2	2:B:831:ARG:HB3	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:129:PHE:CD2	1:A:129:PHE:C	2.92	0.42
1:A:194:LEU:HD23	1:A:195:ASP:H	1.83	0.42
1:A:299:PHE:C	1:A:299:PHE:CD1	2.93	0.42
1:A:86:VAL:HB	1:A:98:LEU:HD22	2.00	0.42
2:B:782:LYS:HB2	2:B:793:TRP:CZ2	2.54	0.42
2:B:776:GLU:OE1	2:B:797:PHE:HZ	2.02	0.42
1:A:62:ARG:HB3	1:A:119:PHE:CD2	2.55	0.42
1:A:176:VAL:O	1:A:176:VAL:HG22	2.18	0.42
1:A:323:VAL:O	1:A:323:VAL:HG12	2.19	0.42
1:A:41:THR:O	1:A:41:THR:HG22	2.19	0.42
1:A:258:ILE:HA	1:A:261:MET:HB2	2.02	0.41
1:A:106:GLU:O	1:A:107:ASN:HB2	2.20	0.41
1:A:181:ILE:CG1	1:A:265:VAL:HG21	2.50	0.41
1:A:148:ILE:HD12	1:A:297:GLN:NE2	2.34	0.41
1:A:185:LEU:HD12	1:A:261:MET:HE3	2.02	0.41
1:A:252:PRO:O	1:A:254:THR:N	2.53	0.41
1:A:79:SER:HB2	1:A:80:GLN:NE2	2.36	0.41
1:A:39:PHE:N	1:A:39:PHE:CD1	2.89	0.41
1:A:270:LYS:HA	1:A:272:PHE:HD1	1.85	0.41
2:B:820:VAL:HG11	2:B:846:LEU:HD21	2.03	0.41
2:B:782:LYS:HB2	2:B:793:TRP:CE2	2.56	0.41
2:B:853:GLU:CG	3:B:96:HOH:O	2.69	0.41
1:A:150:ARG:HA	1:A:162:SER:H	1.85	0.40
1:A:310:ILE:O	1:A:313:MET:HG2	2.21	0.40
1:A:100:ILE:O	1:A:109:LEU:HB2	2.21	0.40
2:B:815:ILE:HA	2:B:816:PRO:HD2	1.80	0.40
1:A:144:MET:SD	1:A:168:PRO:HA	2.61	0.40
1:A:249:ILE:HA	1:A:250:PRO:HA	1.86	0.40
1:A:102:ASP:OD1	1:A:103:GLN:N	2.51	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	298/336 (89%)	257 (86%)	31 (10%)	10 (3%)	4	6
2	B	99/112 (88%)	84 (85%)	8 (8%)	7 (7%)	1	1
All	All	397/448 (89%)	341 (86%)	39 (10%)	17 (4%)	3	4

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	40	SER
1	A	193	HIS
1	A	252	PRO
2	B	814	VAL
1	A	39	PHE
1	A	191	LYS
1	A	194	LEU
1	A	253	HIS
1	A	272	PHE
2	B	783	ALA
2	B	790	LYS
2	B	786	SER
2	B	787	TYR
2	B	835	THR
1	A	161	PRO
2	B	837	ASP
1	A	195	ASP

### 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	269/300 (90%)	252 (94%)	17 (6%)	21	42
2	B	81/94 (86%)	75 (93%)	6 (7%)	16	32
All	All	350/394 (89%)	327 (93%)	23 (7%)	19	38

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

Mol	Chain	Res	Type
1	A	65	LEU
1	A	86	VAL
1	A	92	ASN
1	A	95	THR
1	A	117	ASP
1	A	128	TYR
1	A	160	LEU
1	A	190	ARG
1	A	194	LEU
1	A	197	TRP
1	A	222	HIS
1	A	225	ASN
1	A	245	LYS
1	A	252	PRO
1	A	261	MET
1	A	273	GLN
1	A	287	LEU
2	B	802	ASN
2	B	821	ASN
2	B	822	MET
2	B	834	ASN
2	B	849	ARG
2	B	862	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	21	ASN
1	A	80	GLN
1	A	92	ASN
1	A	114	GLN
1	A	225	ASN
1	A	246	GLN
1	A	308	GLN
1	A	311	GLN
2	B	780	HIS
2	B	802	ASN
2	B	821	ASN
2	B	868	HIS

### 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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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	302/336 (89%)	-0.00	7 (2%) 61 54	24, 60, 130, 153	0
2	B	103/112 (91%)	0.08	5 (4%) 30 24	31, 59, 139, 156	0
All	All	405/448 (90%)	0.02	12 (2%) 51 43	24, 60, 135, 156	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	21	ASN	4.2
1	A	39	PHE	3.8
2	B	788	LEU	3.4
1	A	326	GLU	2.9
1	A	244	LEU	2.8
1	A	200	GLN	2.5
2	B	784	GLY	2.2
1	A	190	ARG	2.2
2	B	836	THR	2.2
1	A	143	GLN	2.1
2	B	835	THR	2.1
2	B	787	TYR	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](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.