



# Full wwPDB X-ray Structure Validation Report ⓘ

May 28, 2024 – 07:24 pm BST

PDB ID : 8BRH  
Title : Co-crystal structure of She4 with Myo4 peptide  
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Deposited on : 2022-11-23  
Resolution : 2.40 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

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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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) : 1.13  
EDS : 2.36.2  
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) : 2.36.2

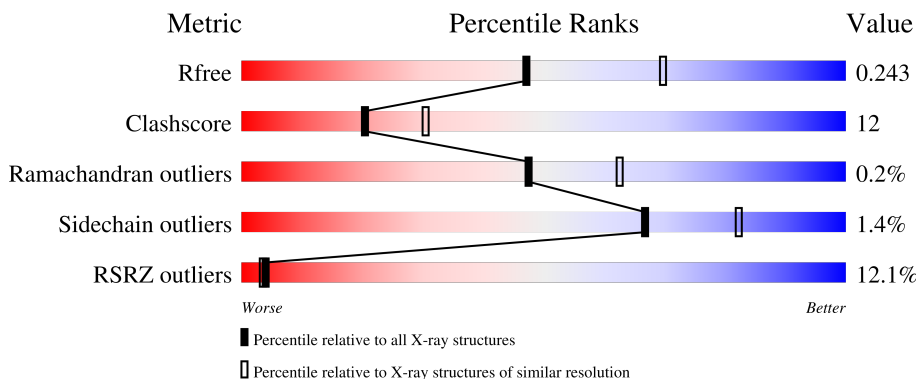
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 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	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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 of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	669	
2	B	7	

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 5215 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 KLLA0E16699p.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	632	5020	3190	849	958	23	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	662	LEU	-	expression tag	UNP Q6CMY7
A	663	GLU	-	expression tag	UNP Q6CMY7
A	664	HIS	-	expression tag	UNP Q6CMY7
A	665	HIS	-	expression tag	UNP Q6CMY7
A	666	HIS	-	expression tag	UNP Q6CMY7
A	667	HIS	-	expression tag	UNP Q6CMY7
A	668	HIS	-	expression tag	UNP Q6CMY7
A	669	HIS	-	expression tag	UNP Q6CMY7

- Molecule 2 is a protein called Myo4 peptide (LYS-PHE-ILE-VAL-SER-HIS-TYR).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	B	7	63	44	10	9	0	0	0

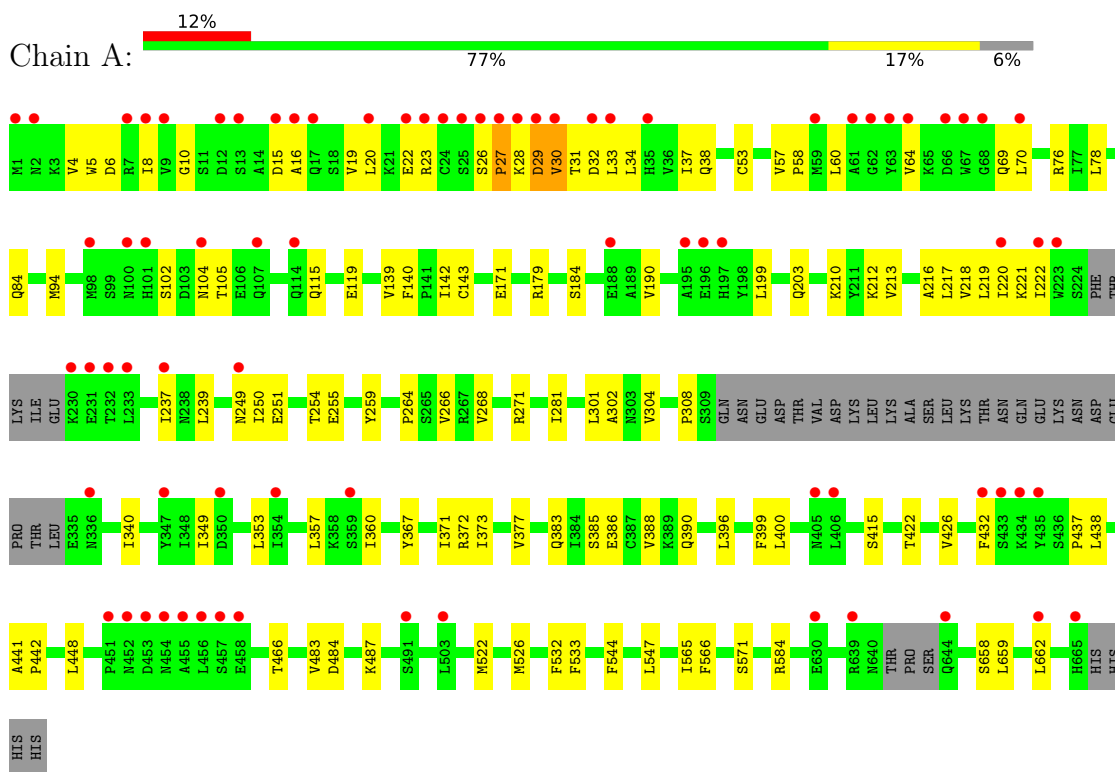
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	130	Total 130	O 130	0	0
3	B	2	Total 2	O 2	0	0

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

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: KLLA0E16699p



- Molecule 2: Myo4 peptide (LYS-PHE-ILE-VAL-SER-HIS-TYR)

Chain B: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	128.48Å 128.48Å 77.54Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.47 – 2.40 49.47 – 2.26	Depositor EDS
% Data completeness (in resolution range)	99.9 (49.47-2.40) 99.9 (49.47-2.26)	Depositor EDS
$R_{merge}$	0.14	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.02 (at 2.27Å)	Xtrriage
Refinement program	PHENIX 1.21.1_5286	Depositor
R, $R_{free}$	0.212 , 0.250 0.208 , 0.243	Depositor DCC
$R_{free}$ test set	1748 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	56.7	Xtrriage
Anisotropy	0.021	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 44.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.029 for -h,-k,l	Xtrriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	5215	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	68.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.18% 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	0.39	0/5093	0.53	0/6897
2	B	0.34	0/65	0.51	0/86
All	All	0.39	0/5158	0.53	0/6983

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	5020	0	5135	127	0
2	B	63	0	62	2	0
3	A	130	0	0	2	0
3	B	2	0	0	0	0
All	All	5215	0	5197	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:A:281:ILE:CD1	1:A:353:LEU:HD12	1.69	1.21

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:217:LEU:HD12	1:A:220:ILE:HD11	1.37	1.06
1:A:8:ILE:CD1	1:A:19:VAL:HG21	1.86	1.06
1:A:212:LYS:HB3	1:A:239:LEU:CD1	1.89	1.03
1:A:212:LYS:CB	1:A:239:LEU:HD11	1.86	1.02
1:A:281:ILE:HD12	1:A:353:LEU:HD12	1.40	1.02
1:A:217:LEU:O	1:A:220:ILE:HG12	1.62	1.00
1:A:281:ILE:HD13	1:A:353:LEU:HD12	1.48	0.95
1:A:8:ILE:HD11	1:A:19:VAL:HG21	1.49	0.95
1:A:658:SER:O	1:A:662:LEU:HD13	1.70	0.90
1:A:212:LYS:HB3	1:A:239:LEU:HD11	0.93	0.90
1:A:219:LEU:HD23	1:A:222:ILE:HD11	1.53	0.89
1:A:34:LEU:CD1	1:A:60:LEU:HD21	2.04	0.88
1:A:4:VAL:HG13	1:A:19:VAL:CG1	2.06	0.86
1:A:281:ILE:CD1	1:A:353:LEU:CD1	2.54	0.86
1:A:34:LEU:HD12	1:A:60:LEU:HD21	1.59	0.83
1:A:658:SER:O	1:A:662:LEU:CD1	2.27	0.83
1:A:8:ILE:HD13	1:A:19:VAL:HG21	1.60	0.81
1:A:8:ILE:HD11	1:A:19:VAL:CG2	2.13	0.79
1:A:349:ILE:CD1	1:A:383:GLN:HB3	2.14	0.77
1:A:438:LEU:CD1	1:A:484:ASP:HB3	2.15	0.77
1:A:217:LEU:O	1:A:220:ILE:CG1	2.33	0.77
1:A:33:LEU:HD23	1:A:60:LEU:HD22	1.66	0.76
1:A:237:ILE:HD13	1:A:266:VAL:HG22	1.68	0.76
1:A:522:MET:CE	1:A:533:PHE:HE2	1.98	0.76
1:A:4:VAL:HG13	1:A:19:VAL:HG11	1.68	0.74
1:A:213:VAL:HG21	1:A:249:ASN:HD22	1.54	0.72
1:A:349:ILE:HD13	1:A:383:GLN:HB3	1.70	0.72
1:A:386:GLU:HG2	3:A:728:HOH:O	1.90	0.70
1:A:64:VAL:HB	1:A:70:LEU:HD11	1.73	0.70
1:A:84:GLN:HE22	1:A:179:ARG:HH22	1.41	0.68
1:A:522:MET:HE2	1:A:533:PHE:HE2	1.59	0.68
1:A:385:SER:O	1:A:388:VAL:HG22	1.94	0.67
1:A:34:LEU:HD11	1:A:60:LEU:HD21	1.77	0.66
1:A:217:LEU:HD12	1:A:220:ILE:CD1	2.22	0.66
1:A:8:ILE:HD11	1:A:19:VAL:CB	2.26	0.65
1:A:448:LEU:HD22	1:A:466:THR:HG23	1.79	0.63
1:A:216:ALA:O	1:A:220:ILE:HG23	1.98	0.63
1:A:57:VAL:N	1:A:58:PRO:HD2	2.13	0.63
1:A:139:VAL:O	1:A:139:VAL:HG12	1.98	0.63
1:A:349:ILE:HD11	1:A:377:VAL:CG1	2.28	0.62
1:A:237:ILE:CD1	1:A:266:VAL:HG22	2.30	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:27:PRO:HA	1:A:30:VAL:HG12	1.82	0.61
1:A:4:VAL:CG1	1:A:19:VAL:CG1	2.78	0.60
1:A:281:ILE:HD13	1:A:353:LEU:CD1	2.26	0.59
1:A:8:ILE:CD1	1:A:19:VAL:CG2	2.69	0.59
1:A:5:TRP:CE3	1:A:37:ILE:HD11	2.38	0.58
1:A:34:LEU:HD23	1:A:69:GLN:O	2.04	0.58
1:A:4:VAL:CG1	1:A:19:VAL:HG11	2.32	0.58
1:A:271:ARG:HG3	1:A:304:VAL:CG1	2.34	0.57
1:A:8:ILE:HD12	1:A:16:ALA:HA	1.87	0.57
1:A:33:LEU:CD2	1:A:60:LEU:HD22	2.33	0.57
1:A:219:LEU:HA	1:A:222:ILE:HG12	1.87	0.57
1:A:139:VAL:HG12	1:A:143:CYS:HB2	1.85	0.56
1:A:571:SER:OG	2:B:74:LYS:HD3	2.06	0.56
1:A:522:MET:CE	1:A:533:PHE:CE2	2.85	0.56
1:A:522:MET:HE1	1:A:533:PHE:CE2	2.41	0.56
1:A:547:LEU:HG	1:A:566:PHE:CZ	2.40	0.56
1:A:271:ARG:HG3	1:A:304:VAL:HG13	1.88	0.56
1:A:385:SER:O	1:A:388:VAL:CG2	2.54	0.55
1:A:522:MET:HE1	1:A:533:PHE:HE2	1.69	0.55
1:A:33:LEU:O	1:A:37:ILE:HG12	2.07	0.55
1:A:220:ILE:HG13	1:A:221:LYS:N	2.22	0.54
1:A:658:SER:O	1:A:662:LEU:HD12	2.08	0.54
1:A:27:PRO:O	1:A:28:LYS:C	2.45	0.53
1:A:199:LEU:O	1:A:203:GLN:HG3	2.09	0.53
1:A:29:ASP:O	1:A:33:LEU:HB3	2.09	0.52
1:A:349:ILE:HD11	1:A:383:GLN:HB3	1.89	0.52
1:A:78:LEU:HD21	1:A:139:VAL:HG21	1.92	0.52
1:A:27:PRO:O	1:A:31:THR:N	2.43	0.51
1:A:522:MET:HG3	1:A:565:ILE:HG23	1.93	0.51
1:A:171:GLU:HB3	1:A:210:LYS:NZ	2.24	0.51
1:A:5:TRP:CZ2	1:A:37:ILE:HD12	2.46	0.51
1:A:251:GLU:O	1:A:254:THR:HG22	2.11	0.51
1:A:281:ILE:HD12	1:A:353:LEU:CD1	2.28	0.51
1:A:20:LEU:HA	1:A:23:ARG:HE	1.75	0.51
1:A:19:VAL:O	1:A:22:GLU:N	2.44	0.50
1:A:5:TRP:CE2	1:A:37:ILE:CD1	2.95	0.49
1:A:171:GLU:HB3	1:A:210:LYS:CE	2.42	0.49
1:A:217:LEU:C	1:A:220:ILE:HG12	2.32	0.49
1:A:5:TRP:CZ2	1:A:53:CYS:SG	3.03	0.49
1:A:441:ALA:HB3	1:A:442:PRO:HD3	1.94	0.49
1:A:254:THR:HG23	1:A:255:GLU:N	2.26	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:139:VAL:O	1:A:139:VAL:CG1	2.61	0.48
1:A:94:MET:CG	1:A:142:ILE:HD11	2.44	0.48
1:A:522:MET:HE2	1:A:533:PHE:CE2	2.44	0.48
1:A:526:MET:HB2	3:A:727:HOH:O	2.14	0.48
1:A:571:SER:OG	2:B:74:LYS:CD	2.62	0.48
1:A:102:SER:O	1:A:105:THR:HG22	2.14	0.48
1:A:422:THR:O	1:A:426:VAL:HG22	2.15	0.47
1:A:432:PHE:CD2	1:A:437:PRO:HA	2.50	0.47
1:A:371:ILE:HD12	1:A:396:LEU:HD23	1.97	0.47
1:A:94:MET:HG2	1:A:142:ILE:HD11	1.97	0.46
1:A:115:GLN:O	1:A:119:GLU:HG3	2.16	0.46
1:A:5:TRP:HZ2	1:A:53:CYS:SG	2.39	0.46
1:A:438:LEU:CD1	1:A:484:ASP:CB	2.91	0.46
1:A:301:LEU:HD23	1:A:373:ILE:HD13	1.99	0.45
1:A:31:THR:O	1:A:32:ASP:C	2.54	0.45
1:A:184:SER:HB2	1:A:190:VAL:HG11	1.99	0.45
1:A:28:LYS:HE2	1:A:28:LYS:HB3	1.77	0.44
1:A:5:TRP:CZ2	1:A:37:ILE:CD1	2.99	0.44
1:A:371:ILE:HG21	1:A:415:SER:OG	2.17	0.44
1:A:218:VAL:O	1:A:222:ILE:HG12	2.18	0.44
1:A:302:ALA:HB1	1:A:372:ARG:HG2	1.99	0.44
1:A:38:GLN:O	1:A:76:ARG:NH2	2.51	0.44
1:A:308:PRO:HD3	1:A:340:ILE:HD12	1.98	0.44
1:A:6:ASP:O	1:A:10:GLY:N	2.51	0.43
1:A:250:ILE:O	1:A:254:THR:HG22	2.19	0.43
1:A:483:VAL:HG12	1:A:487:LYS:HE3	2.01	0.43
1:A:5:TRP:CH2	1:A:37:ILE:HD12	2.53	0.43
1:A:5:TRP:CZ3	1:A:37:ILE:HD11	2.53	0.43
1:A:532:PHE:O	1:A:544:PHE:HB2	2.19	0.42
1:A:217:LEU:CD1	1:A:220:ILE:HD11	2.26	0.42
1:A:26:SER:O	1:A:30:VAL:HB	2.20	0.42
1:A:399:PHE:CD2	1:A:415:SER:OG	2.71	0.42
1:A:23:ARG:CD	1:A:33:LEU:HD22	2.49	0.42
1:A:264:PRO:O	1:A:268:VAL:HG23	2.20	0.42
1:A:5:TRP:CD2	1:A:37:ILE:HD11	2.55	0.41
1:A:27:PRO:O	1:A:30:VAL:N	2.53	0.41
1:A:659:LEU:HA	1:A:662:LEU:HD13	2.02	0.41
1:A:4:VAL:O	1:A:8:ILE:HG12	2.21	0.41
1:A:15:ASP:O	1:A:19:VAL:HG23	2.20	0.41
1:A:8:ILE:HD11	1:A:19:VAL:HB	2.02	0.40
1:A:386:GLU:O	1:A:390:GLN:HG3	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:349:ILE:HD11	1:A:377:VAL:HG11	2.01	0.40
1:A:357:LEU:HA	1:A:360:ILE:HG22	2.03	0.40
1:A:438:LEU:HD13	1:A:484:ASP:HB3	2.00	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	624/669 (93%)	604 (97%)	19 (3%)	1 (0%)	47	62
2	B	5/7 (71%)	5 (100%)	0	0	100	100
All	All	629/676 (93%)	609 (97%)	19 (3%)	1 (0%)	47	62

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	27	PRO

### 5.3.2 Protein sidechains [i](#)

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	576/612 (94%)	568 (99%)	8 (1%)	67	82
2	B	7/7 (100%)	7 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	583/619 (94%)	575 (99%)	8 (1%)	67 82

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	ASP
1	A	30	VAL
1	A	104	ASN
1	A	140	PHE
1	A	259	TYR
1	A	367	TYR
1	A	400	LEU
1	A	584	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	35	HIS
1	A	84	GLN
1	A	104	ASN
1	A	249	ASN
1	A	460	GLN
1	A	480	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 monosaccharides 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, 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	632/669 (94%)	0.61	77 (12%) <b>4</b>   <b>3</b>	36, 62, 121, 214	0
2	B	7/7 (100%)	0.40	0 <b>100</b>   <b>100</b>	48, 51, 60, 67	0
All	All	639/676 (94%)	0.61	77 (12%) <b>4</b>   <b>3</b>	36, 62, 121, 214	0

All (77) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	452	ASN	9.1
1	A	61	ALA	8.4
1	A	23	ARG	7.4
1	A	17	GLN	7.0
1	A	32	ASP	6.4
1	A	434	LYS	6.4
1	A	453	ASP	6.3
1	A	644	GLN	5.5
1	A	101	HIS	5.4
1	A	639	ARG	5.2
1	A	233	LEU	5.2
1	A	29	ASP	5.1
1	A	451	PRO	5.0
1	A	230	LYS	4.7
1	A	433	SER	4.5
1	A	30	VAL	4.4
1	A	457	SER	4.1
1	A	222	ILE	4.1
1	A	223	TRP	4.0
1	A	68	GLY	3.9
1	A	347	TYR	3.8
1	A	1	MET	3.8
1	A	405	ASN	3.8
1	A	456	LEU	3.8

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>	<b>RSRZ</b>
1	A	100	ASN	3.8
1	A	20	LEU	3.8
1	A	7	ARG	3.8
1	A	237	ILE	3.7
1	A	107	GLN	3.6
1	A	66	ASP	3.5
1	A	27	PRO	3.5
1	A	8	ILE	3.5
1	A	232	THR	3.4
1	A	98	MET	3.4
1	A	249	ASN	3.2
1	A	63	TYR	3.2
1	A	188	GLU	3.0
1	A	503	LEU	3.0
1	A	662	LEU	2.9
1	A	195	ALA	2.8
1	A	24	CYS	2.8
1	A	26	SER	2.7
1	A	454	ASN	2.7
1	A	33	LEU	2.7
1	A	12	ASP	2.6
1	A	455	ALA	2.6
1	A	13	SER	2.6
1	A	67	TRP	2.5
1	A	104	ASN	2.5
1	A	336	ASN	2.5
1	A	458	GLU	2.4
1	A	406	LEU	2.4
1	A	22	GLU	2.4
1	A	70	LEU	2.4
1	A	62	GLY	2.3
1	A	35	HIS	2.3
1	A	435	TYR	2.3
1	A	9	VAL	2.3
1	A	25	SER	2.3
1	A	220	ILE	2.3
1	A	28	LYS	2.2
1	A	491	SER	2.2
1	A	231	GLU	2.2
1	A	197	HIS	2.2
1	A	630	GLU	2.2
1	A	2	ASN	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	15	ASP	2.2
1	A	196	GLU	2.2
1	A	114	GLN	2.1
1	A	59	MET	2.1
1	A	350	ASP	2.1
1	A	665	HIS	2.1
1	A	64	VAL	2.1
1	A	354	ILE	2.1
1	A	432	PHE	2.1
1	A	359	SER	2.1
1	A	16	ALA	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 monosaccharides 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.