



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 24, 2024 – 09:54 AM EDT

PDB ID : 3C2G  
Title : Crystal complex of SYS-1/POP-1 at 2.5A resolution  
Authors : Liu, J.; Phillips, B.T.; Amaya, M.F.; Kimble, J.; Xu, W.  
Deposited on : 2008-01-24  
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

<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>.

---

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

MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

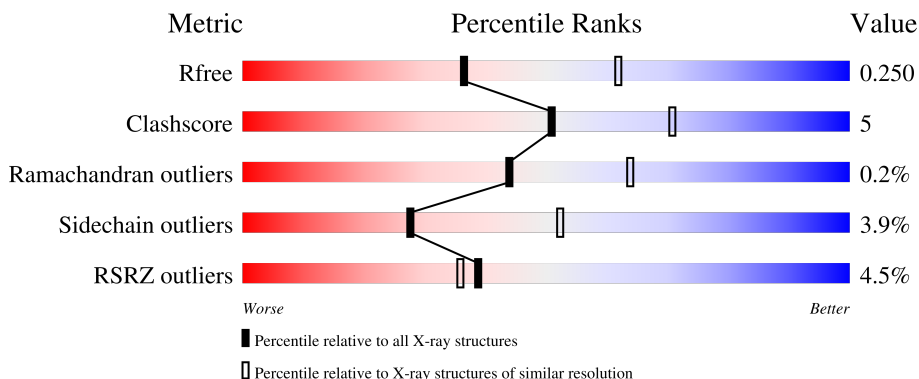
# 1 Overall quality at a glance

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))
$R_{free}$	164625	5504 (2.50-2.50)
Clashscore	180529	6282 (2.50-2.50)
Ramachandran outliers	177936	6191 (2.50-2.50)
Sidechain outliers	177891	6193 (2.50-2.50)
RSRZ outliers	164620	5504 (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 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	619	<div> <div>3%</div> <div>86%</div> <div>12%</div> <div>.</div> </div>
1	B	619	<div> <div>6%</div> <div>84%</div> <div>13%</div> <div>..</div> </div>
2	C	8	<div> <div>88%</div> <div>12%</div> </div>
2	D	8	<div> <div>12%</div> <div>88%</div> <div>12%</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9957 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 Sys-1 protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	618	Total	C	N	O	S	Se	0	0	0
			4893	3150	853	858	17	15			
1	B	615	Total	C	N	O	S	Se	0	0	0
			4864	3132	848	852	17	15			

- Molecule 2 is a protein called Pop-1 8-residue peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	8	Total	C	N	O	0	0	0
			62	40	12	10			
2	D	8	Total	C	N	O	0	0	0
			66	42	12	12			

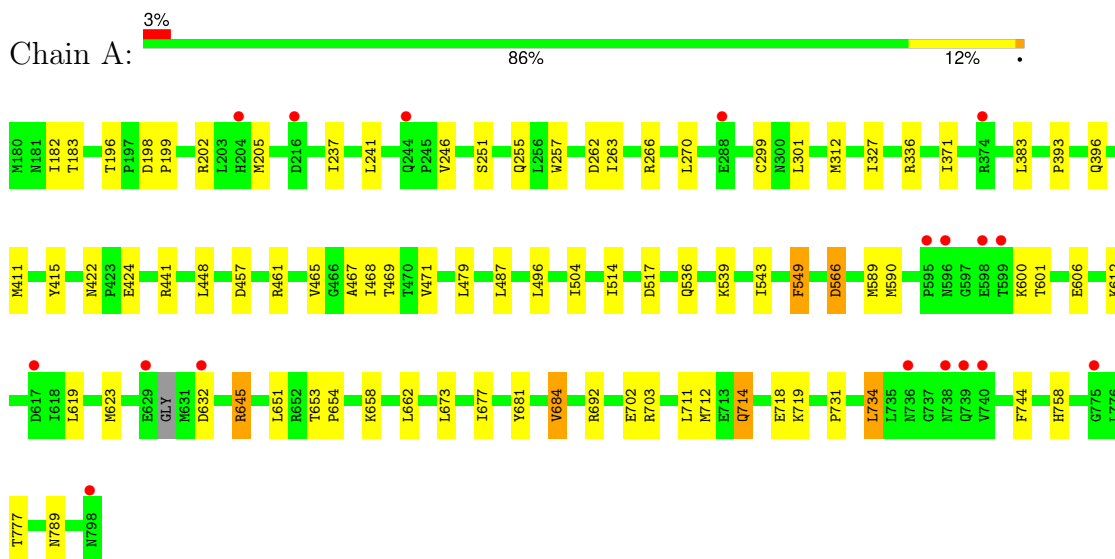
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	37	Total	O	0	0
			37	37		
3	B	35	Total	O	0	0
			35	35		

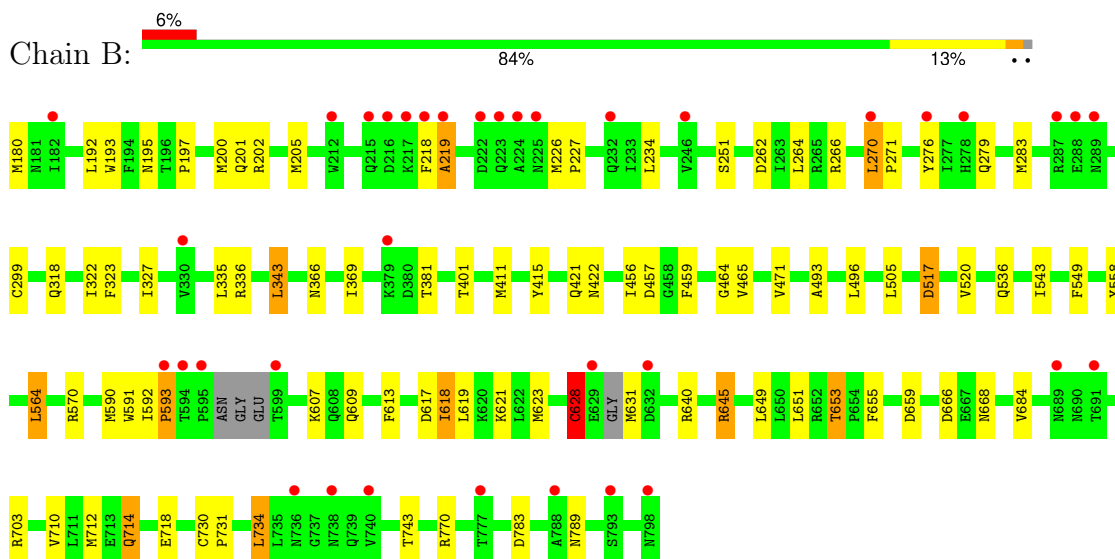
### 3 Residue-property plots [i](#)

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: Sys-1 protein



- Molecule 1: Sys-1 protein




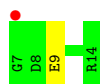
- Molecule 2: Pop-1 8-residue peptide

Chain C:  88% 12%



- Molecule 2: Pop-1 8-residue peptide

Chain D:  12% 88% 12%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	84.97Å 85.34Å 93.66Å 65.22° 78.08° 83.02°	Depositor
Resolution (Å)	48.00 – 2.50 48.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	93.7 (48.00-2.50) 90.1 (48.00-2.50)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.95 (at 2.32Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.216 , 0.261 0.209 , 0.250	Depositor DCC
$R_{free}$ test set	3629 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	42.2	Xtriage
Anisotropy	0.356	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 31.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	9957	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.34% 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	0.45	0/4992	0.55	0/6775
1	B	0.46	1/4961 (0.0%)	0.56	0/6733
2	C	0.64	0/62	0.50	0/81
2	D	0.60	0/66	0.42	0/86
All	All	0.46	1/10081 (0.0%)	0.56	0/13675

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	628	CYS	CB-SG	-5.74	1.72	1.81

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	200	MSE	Peptide

### 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	4893	0	4949	48	0
1	B	4864	0	4906	54	0
2	C	62	0	61	1	0
2	D	66	0	65	0	0
3	A	37	0	0	0	0
3	B	35	0	0	1	0
All	All	9957	0	9981	101	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 5.

All (101) 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:777:THR:HG23	1:A:789:ASN:HD22	1.38	0.89
1:A:411:MSE:HE3	1:A:471:VAL:HG21	1.59	0.82
1:A:301:LEU:HD23	1:A:312:MSE:HE1	1.61	0.82
1:B:649:LEU:O	1:B:653:THR:HG23	1.92	0.70
1:A:651:LEU:O	1:A:714:GLN:HG2	1.91	0.69
1:B:623:MSE:CE	1:B:666:ASP:HB3	2.24	0.68
1:A:411:MSE:HE3	1:A:471:VAL:CG2	2.25	0.65
1:B:234:LEU:HD13	1:B:283:MSE:HE3	1.79	0.65
1:A:327:ILE:O	1:A:336:ARG:NH2	2.30	0.64
1:A:712:MSE:HG2	1:A:718:GLU:HA	1.80	0.63
1:A:202:ARG:HA	1:A:205:MSE:HE3	1.81	0.62
1:B:651:LEU:O	1:B:714:GLN:HG2	2.02	0.60
1:B:712:MSE:HG2	1:B:718:GLU:HA	1.85	0.58
1:A:645:ARG:HG3	1:A:703:ARG:CZ	2.35	0.57
1:A:468:ILE:HG21	1:A:487:LEU:HG	1.87	0.56
1:B:543:ILE:HD11	1:B:590:MSE:HE1	1.86	0.56
1:B:564:LEU:HD23	1:B:570:ARG:HA	1.87	0.56
1:A:241:LEU:HD23	1:A:257:TRP:CD1	2.41	0.56
1:B:628:CYS:HB2	1:B:640:ARG:HH21	1.71	0.55
1:B:558:TYR:O	1:B:621:LYS:NZ	2.35	0.55
1:A:469:THR:HG21	1:A:504:ILE:HG23	1.88	0.55
1:B:202:ARG:HA	1:B:205:MSE:HE3	1.88	0.55
1:B:623:MSE:HE3	1:B:666:ASP:HB3	1.89	0.54
1:A:465:VAL:HG21	1:A:496:LEU:HD21	1.90	0.53
1:A:411:MSE:HE1	1:A:467:ALA:C	2.29	0.52
1:A:536:GLN:HE22	1:A:590:MSE:SE	2.42	0.52

*Continued on next page...*



*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:628:CYS:HB2	1:B:640:ARG:NH2	2.24	0.52
1:A:719:LYS:HE2	1:A:758:HIS:HD2	1.74	0.51
1:A:673:LEU:O	1:A:677:ILE:HG12	2.10	0.51
1:A:196:THR:HB	1:A:205:MSE:HE1	1.94	0.50
1:A:651:LEU:O	1:A:658:LYS:HE2	2.11	0.50
1:B:366:ASN:HB3	1:B:369:ILE:HD12	1.93	0.50
1:A:393:PRO:HD2	1:A:396:GLN:HG3	1.94	0.50
1:A:255:GLN:HG3	1:B:422:ASN:ND2	2.27	0.50
1:B:262:ASP:HB2	1:B:318:GLN:HE22	1.77	0.49
1:B:195:ASN:O	1:B:197:PRO:HD3	2.13	0.49
1:A:731:PRO:O	1:A:734:LEU:HB2	2.13	0.49
1:B:465:VAL:HG21	1:B:496:LEU:HD21	1.95	0.49
1:A:262:ASP:OD2	1:A:266:ARG:NH1	2.46	0.48
1:A:301:LEU:CD2	1:A:312:MSE:HE1	2.39	0.48
1:A:422:ASN:HD21	1:A:424:GLU:HB2	1.77	0.48
1:A:415:TYR:HB2	1:A:471:VAL:HG22	1.94	0.48
1:B:456:ILE:HG22	1:B:493:ALA:HB2	1.96	0.47
1:A:457:ASP:O	1:A:461:ARG:HG3	2.14	0.47
1:B:609:GLN:NE2	3:B:68:HOH:O	2.45	0.47
1:A:539:LYS:NZ	2:C:8:ASP:OD1	2.47	0.47
1:A:411:MSE:HE1	1:A:467:ALA:HB1	1.97	0.47
1:B:327:ILE:O	1:B:336:ARG:NH2	2.48	0.46
1:B:543:ILE:CD1	1:B:590:MSE:HE1	2.45	0.46
1:B:415:TYR:HB2	1:B:471:VAL:HG22	1.96	0.46
1:B:649:LEU:O	1:B:653:THR:CG2	2.62	0.46
1:B:234:LEU:HD13	1:B:283:MSE:CE	2.44	0.45
1:A:198:ASP:HA	1:A:199:PRO:HD3	1.83	0.45
1:B:730:CYS:HA	1:B:731:PRO:HD3	1.87	0.45
1:B:323:PHE:HE1	1:B:343:LEU:HD11	1.82	0.45
1:B:645:ARG:HG3	1:B:703:ARG:CZ	2.47	0.45
1:A:182:ILE:HD12	1:A:183:THR:H	1.81	0.45
1:B:549:PHE:H	1:B:609:GLN:HE21	1.64	0.45
1:B:335:LEU:HD12	1:B:335:LEU:HA	1.81	0.45
1:A:600:LYS:HE3	1:A:600:LYS:HB2	1.86	0.44
1:B:234:LEU:HB3	1:B:283:MSE:HE3	1.99	0.44
1:B:592:ILE:HA	1:B:593:PRO:HD2	1.78	0.44
1:A:479:LEU:HD23	1:A:479:LEU:C	2.37	0.44
1:B:590:MSE:HE3	1:B:591:TRP:CZ2	2.53	0.44
1:B:264:LEU:HD23	1:B:322:ILE:HD13	1.99	0.44
1:A:371:ILE:HD11	1:A:441:ARG:HH12	1.83	0.44
1:B:613:PHE:HA	1:B:618:ILE:HG21	2.00	0.44

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:662:LEU:HD21	1:A:711:LEU:HD22	1.99	0.44
1:A:383:LEU:HD12	1:A:383:LEU:HA	1.92	0.43
1:A:566:ASP:OD1	1:A:566:ASP:N	2.40	0.43
1:B:226:MSE:HG2	1:B:276:TYR:CZ	2.53	0.43
1:A:543:ILE:HG21	1:A:606:GLU:HG3	1.99	0.43
1:B:226:MSE:HB3	1:B:227:PRO:HD3	2.00	0.43
1:A:702:GLU:HG2	1:A:744:PHE:CD1	2.53	0.43
1:B:270:LEU:HA	1:B:271:PRO:HD2	1.91	0.43
1:A:549:PHE:CZ	1:A:612:LYS:HD2	2.54	0.43
1:B:459:PHE:CZ	1:B:464:GLY:HA3	2.53	0.43
1:A:411:MSE:HE2	1:A:411:MSE:HB3	1.70	0.43
1:B:279:GLN:O	1:B:283:MSE:HE2	2.19	0.43
1:A:198:ASP:O	1:A:202:ARG:HG3	2.19	0.43
1:B:218:PHE:O	1:B:219:ALA:C	2.58	0.42
1:A:182:ILE:H	1:A:182:ILE:HG13	1.58	0.42
1:B:202:ARG:HA	1:B:205:MSE:CE	2.50	0.42
1:B:457:ASP:OD1	1:B:493:ALA:HB1	2.20	0.42
1:A:681:TYR:O	1:A:684:VAL:HG12	2.20	0.42
1:A:237:ILE:HD11	1:A:263:ILE:HG22	2.02	0.42
1:A:619:LEU:O	1:A:623:MSE:HG2	2.20	0.42
1:A:653:THR:HA	1:A:654:PRO:HD3	1.88	0.41
1:B:226:MSE:HG2	1:B:276:TYR:CE2	2.55	0.41
1:B:505:LEU:HD12	1:B:505:LEU:HA	1.87	0.41
1:B:607:LYS:HG2	1:B:655:PHE:CD1	2.56	0.41
1:B:731:PRO:O	1:B:734:LEU:HB2	2.20	0.41
1:B:266:ARG:O	1:B:270:LEU:HB2	2.20	0.41
1:B:192:LEU:O	1:B:193:TRP:C	2.58	0.41
1:A:266:ARG:O	1:A:270:LEU:HB2	2.21	0.41
1:B:270:LEU:HA	1:B:270:LEU:HD12	1.98	0.41
1:B:517:ASP:OD1	1:B:520:VAL:HG13	2.21	0.40
1:B:590:MSE:CE	1:B:591:TRP:CZ2	3.04	0.40
1:B:536:GLN:HE22	1:B:590:MSE:SE	2.55	0.40
1:B:411:MSE:HG2	1:B:471:VAL:HG21	2.04	0.40
1:B:710:VAL:O	1:B:714:GLN:HB2	2.21	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	614/619 (99%)	601 (98%)	12 (2%)	1 (0%)	44	64
1	B	609/619 (98%)	581 (95%)	26 (4%)	2 (0%)	37	56
2	C	6/8 (75%)	5 (83%)	1 (17%)	0	100	100
2	D	6/8 (75%)	5 (83%)	1 (17%)	0	100	100
All	All	1235/1254 (98%)	1192 (96%)	40 (3%)	3 (0%)	44	64

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	219	ALA
1	A	632	ASP
1	B	593	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	541/548 (99%)	526 (97%)	15 (3%)	38	65
1	B	536/548 (98%)	509 (95%)	27 (5%)	20	41
2	C	6/7 (86%)	6 (100%)	0	100	100
2	D	7/7 (100%)	6 (86%)	1 (14%)	2	5
All	All	1090/1110 (98%)	1047 (96%)	43 (4%)	27	52

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

Mol	Chain	Res	Type
1	A	246	VAL
1	A	251	SER
1	A	299	CYS
1	A	448	LEU
1	A	514	ILE
1	A	517	ASP
1	A	549	PHE
1	A	566	ASP
1	A	589	MSE
1	A	601	THR
1	A	645	ARG
1	A	684	VAL
1	A	692	ARG
1	A	714	GLN
1	A	734	LEU
1	B	180	MSE
1	B	201	GLN
1	B	251	SER
1	B	270	LEU
1	B	299	CYS
1	B	343	LEU
1	B	381	THR
1	B	401	THR
1	B	421	GLN
1	B	517	ASP
1	B	564	LEU
1	B	617	ASP
1	B	618	ILE
1	B	619	LEU
1	B	628	CYS
1	B	631	MSE
1	B	645	ARG
1	B	653	THR
1	B	659	ASP
1	B	668	ASN
1	B	684	VAL
1	B	714	GLN
1	B	734	LEU
1	B	743	THR
1	B	770	ARG
1	B	783	ASP
1	B	789	ASN
2	D	9	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	247	GLN
1	A	278	HIS
1	A	515	HIS
1	A	536	GLN
1	A	714	GLN
1	A	758	HIS
1	A	789	ASN
1	B	220	GLN
1	B	259	ASN
1	B	305	ASN
1	B	307	GLN
1	B	318	GLN
1	B	342	HIS
1	B	536	GLN
1	B	546	ASN
1	B	609	GLN
1	B	699	GLN

### 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 oligosaccharides 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 ⓘ

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(Å²)	Q<0.9
1	A	603/619 (97%)	0.12	18 (2%)	52	49	23, 42, 70, 86	14 (2%)
1	B	600/619 (96%)	0.29	36 (6%)	29	27	22, 42, 73, 91	19 (3%)
2	C	8/8 (100%)	0.17	0	100	100	41, 43, 45, 55	0
2	D	8/8 (100%)	1.02	1 (12%)	9	9	57, 59, 62, 65	0
All	All	1219/1254 (97%)	0.21	55 (4%)	39	36	22, 42, 71, 91	33 (2%)

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	599	THR	6.6
1	B	215	GLN	6.3
1	B	219	ALA	4.8
1	B	595	PRO	4.6
1	B	740	VAL	4.0
1	B	225	ASN	3.9
1	A	204	HIS	3.9
1	B	217	LYS	3.7
1	B	222	ASP	3.6
1	B	594	THR	3.5
1	B	736	ASN	3.4
1	B	287	ARG	3.3
1	A	599	THR	3.3
1	B	218	PHE	3.3
1	B	276	TYR	3.1
1	A	632	ASP	3.1
1	A	740	VAL	3.1
1	A	374	ARG	2.9
1	B	798	ASN	2.9
1	A	596	ASN	2.9
1	B	689	ASN	2.9

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	B	330	VAL	2.8
1	B	289	ASN	2.7
1	A	738	ASN	2.7
1	B	629	GLU	2.6
1	B	224	ALA	2.6
1	B	777	THR	2.6
1	B	216	ASP	2.5
1	B	379	LYS	2.5
1	A	216	ASP	2.5
1	A	598	GLU	2.4
1	B	182	ILE	2.4
1	B	788	ALA	2.4
1	A	775	GLY	2.4
1	B	212	TRP	2.4
1	B	632	ASP	2.3
1	A	739	GLN	2.3
1	A	617	ASP	2.3
1	A	244	GLN	2.3
1	A	798	ASN	2.2
1	B	270	LEU	2.2
1	B	593	PRO	2.2
1	A	288	GLU	2.2
1	B	223	GLN	2.2
1	B	691	THR	2.2
1	B	288	GLU	2.1
1	B	793	SER	2.1
1	B	232	GLN	2.1
1	B	278	HIS	2.1
1	A	736	ASN	2.1
1	B	246	VAL	2.0
1	B	738	ASN	2.0
2	D	7	GLY	2.0
1	A	595	PRO	2.0
1	A	629	GLU	2.0

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

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.