



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:58 AM GMT

PDB ID : 3GRC
Title : Crystal structure of a sensor protein from Polaromonas sp. JS666
Authors : Palani, K.; Kumaran, D.; Burley, S.K.; Swaminathan, S.; New York SGX
Research Center for Structural Genomics (NYSGXRC)
Deposited on : 2009-03-25
Resolution : 2.21 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) : 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

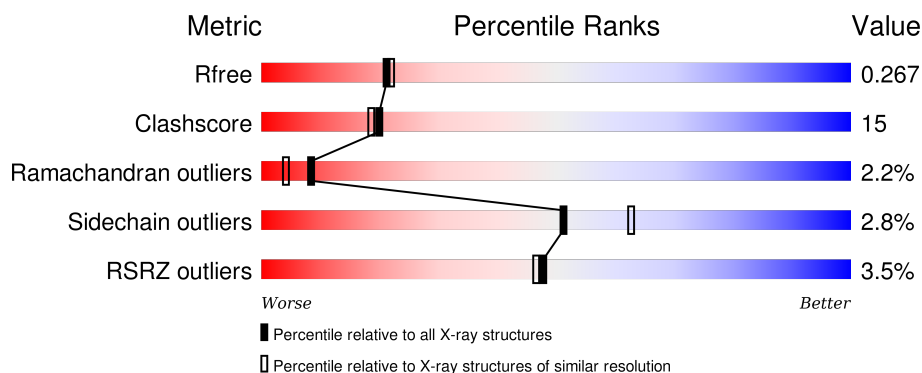
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.21 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	4405 (2.24-2.20)
Clashscore	102246	5146 (2.24-2.20)
Ramachandran outliers	100387	5065 (2.24-2.20)
Sidechain outliers	100360	5066 (2.24-2.20)
RSRZ outliers	91569	4414 (2.24-2.20)

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.

Mol	Chain	Length	Quality of chain
1	A	140	<div> <div></div> <div>67% 19% • 11%</div> </div>
1	B	140	<div> <div></div> <div>66% 20% • 12%</div> </div>
1	C	140	<div> <div>4%</div> <div>63% 21% 5% 11%</div> </div>
1	D	140	<div> <div>5%</div> <div>60% 26% • • 11%</div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4083 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 Sensor protein, Kinase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	125	Total	C	N	O	S	Se	0	0	0
			976	606	178	187	1	4			
1	B	123	Total	C	N	O	S	Se	0	0	0
			963	598	176	185	1	3			
1	C	125	Total	C	N	O	S	Se	0	0	0
			976	606	178	187	1	4			
1	D	125	Total	C	N	O	S	Se	0	0	0
			976	606	178	187	1	4			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	542	MSE	-	expression tag	UNP Q121U0
A	543	SER	-	expression tag	UNP Q121U0
A	544	LEU	-	expression tag	UNP Q121U0
A	674	GLU	-	expression tag	UNP Q121U0
A	675	GLY	-	expression tag	UNP Q121U0
A	676	HIS	-	expression tag	UNP Q121U0
A	677	HIS	-	expression tag	UNP Q121U0
A	678	HIS	-	expression tag	UNP Q121U0
A	679	HIS	-	expression tag	UNP Q121U0
A	680	HIS	-	expression tag	UNP Q121U0
A	681	HIS	-	expression tag	UNP Q121U0
B	542	MSE	-	expression tag	UNP Q121U0
B	543	SER	-	expression tag	UNP Q121U0
B	544	LEU	-	expression tag	UNP Q121U0
B	674	GLU	-	expression tag	UNP Q121U0
B	675	GLY	-	expression tag	UNP Q121U0
B	676	HIS	-	expression tag	UNP Q121U0
B	677	HIS	-	expression tag	UNP Q121U0
B	678	HIS	-	expression tag	UNP Q121U0
B	679	HIS	-	expression tag	UNP Q121U0
B	680	HIS	-	expression tag	UNP Q121U0

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Chain	Residue	Modelled	Actual	Comment	Reference
B	681	HIS	-	expression tag	UNP Q121U0
C	542	MSE	-	expression tag	UNP Q121U0
C	543	SER	-	expression tag	UNP Q121U0
C	544	LEU	-	expression tag	UNP Q121U0
C	674	GLU	-	expression tag	UNP Q121U0
C	675	GLY	-	expression tag	UNP Q121U0
C	676	HIS	-	expression tag	UNP Q121U0
C	677	HIS	-	expression tag	UNP Q121U0
C	678	HIS	-	expression tag	UNP Q121U0
C	679	HIS	-	expression tag	UNP Q121U0
C	680	HIS	-	expression tag	UNP Q121U0
C	681	HIS	-	expression tag	UNP Q121U0
D	542	MSE	-	expression tag	UNP Q121U0
D	543	SER	-	expression tag	UNP Q121U0
D	544	LEU	-	expression tag	UNP Q121U0
D	674	GLU	-	expression tag	UNP Q121U0
D	675	GLY	-	expression tag	UNP Q121U0
D	676	HIS	-	expression tag	UNP Q121U0
D	677	HIS	-	expression tag	UNP Q121U0
D	678	HIS	-	expression tag	UNP Q121U0
D	679	HIS	-	expression tag	UNP Q121U0
D	680	HIS	-	expression tag	UNP Q121U0
D	681	HIS	-	expression tag	UNP Q121U0

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	50	Total O 50 50	0	0
2	B	51	Total O 51 51	0	0
2	C	42	Total O 42 42	0	0
2	D	49	Total O 49 49	0	0

L660	S661	L662	H663	H664	H669	A670	GLU	GLY	LYS	GLU	GLY	HIS	HIS	HIS	HIS	HIS	HIS
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4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	64.36Å 87.81Å 103.81Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.44 – 2.21 46.43 – 2.21	Depositor EDS
% Data completeness (in resolution range)	95.9 (40.44-2.21) 96.0 (46.43-2.21)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.50 (at 2.20Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.221 , 0.265 0.221 , 0.267	Depositor DCC
R_{free} test set	1424 reflections (4.91%)	DCC
Wilson B-factor (Å ²)	29.7	Xtriage
Anisotropy	0.464	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 46.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 29717 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4083	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 22.97 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 5.1150e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical 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](#)

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	0/985	0.70	0/1329
1	B	0.39	0/973	0.67	0/1315
1	C	0.39	0/985	0.71	0/1329
1	D	0.35	0/985	0.63	0/1329
All	All	0.38	0/3928	0.68	0/5302

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	976	0	988	26	0
1	B	963	0	974	26	0
1	C	976	0	988	29	0
1	D	976	0	988	35	0
2	A	50	0	0	1	0
2	B	51	0	0	0	0
2	C	42	0	0	0	0
2	D	49	0	0	1	0
All	All	4083	0	3938	116	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 15.

All (116) 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:551:LEU:HD23	1:B:595:MSE:HE2	1.49	0.93
1:A:595:MSE:HE1	1:A:609:LEU:HD21	1.48	0.93
1:B:585:GLU:HG2	1:B:589:ARG:HH12	1.35	0.91
1:D:547:ARG:HB3	1:D:548:PRO:HD3	1.53	0.89
1:C:551:LEU:HD23	1:C:595:MSE:HE2	1.60	0.84
1:D:623:ALA:HB2	1:D:669:MSE:HE1	1.60	0.84
1:D:595:MSE:HE3	1:D:597:VAL:CG2	2.08	0.82
1:A:645:VAL:HG11	1:A:648:TRP:CZ2	2.14	0.81
1:A:553:CYS:HB2	1:A:595:MSE:HE3	1.63	0.80
1:D:580:ALA:HB3	1:D:604:GLN:HB3	1.64	0.78
1:D:621:ASP:O	1:D:669:MSE:HE2	1.84	0.76
1:B:566:MSE:HE2	1:B:655:GLU:HB2	1.67	0.76
1:D:623:ALA:N	1:D:669:MSE:HE1	2.04	0.72
1:A:616:ASP:OD1	1:A:618:ARG:HG2	1.90	0.71
1:B:561:ARG:HH11	1:B:564:ASN:HD22	1.38	0.71
1:C:668:ASN:C	1:C:670:ALA:N	2.42	0.71
1:A:547:ARG:HB3	1:A:548:PRO:HD3	1.72	0.71
1:C:669:MSE:HE2	1:C:669:MSE:N	2.06	0.71
1:D:613:LEU:O	1:D:619:THR:HG23	1.93	0.69
1:D:623:ALA:CB	1:D:669:MSE:HE1	2.22	0.68
1:C:668:ASN:C	1:C:670:ALA:H	1.96	0.68
1:B:585:GLU:HG2	1:B:589:ARG:NH1	2.07	0.67
1:D:562:LEU:HD11	1:D:566:MSE:HE3	1.78	0.66
1:D:595:MSE:HE3	1:D:597:VAL:HG21	1.78	0.66
1:B:547:ARG:HG3	1:B:547:ARG:HH11	1.60	0.65
1:D:623:ALA:CA	1:D:669:MSE:HE1	2.27	0.65
1:A:562:LEU:HD11	1:A:566:MSE:HE3	1.78	0.65
1:C:547:ARG:HB3	1:C:548:PRO:HD3	1.82	0.62
1:A:621:ASP:O	1:A:669:MSE:HG2	1.99	0.62
1:A:595:MSE:HE2	1:A:597:VAL:HG22	1.80	0.62
1:B:665:ALA:O	1:B:666:ILE:HB	2.00	0.60
1:C:669:MSE:H	1:C:669:MSE:HE2	1.65	0.59
1:A:553:CYS:CB	1:A:595:MSE:HE3	2.33	0.59
1:C:630:ASN:HB3	1:C:633:GLU:HG2	1.84	0.58
1:D:616:ASP:HB3	1:D:619:THR:HG22	1.84	0.58
1:B:546:PRO:N	1:B:571:GLY:O	2.36	0.58
1:D:623:ALA:HB2	1:D:669:MSE:CE	2.32	0.57
1:C:607:VAL:HG21	1:C:638:PHE:CE1	2.40	0.57
1:A:641:GLN:HA	1:A:641:GLN:NE2	2.20	0.56
1:C:635:GLU:HB2	1:C:648:TRP:CD2	2.41	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:595:MSE:HE1	1:D:613:LEU:CD1	2.37	0.54
1:C:639:ASN:O	1:C:641:GLN:N	2.38	0.53
1:B:549:ARG:HE	1:B:590:ARG:NH1	2.06	0.53
1:C:651:LYS:HA	1:C:652:PRO:C	2.29	0.53
1:D:579:SER:HB2	1:D:603:ASP:OD1	2.09	0.52
1:A:546:PRO:HD2	1:A:571:GLY:C	2.30	0.52
1:C:638:PHE:O	1:C:640:SER:N	2.42	0.51
1:B:566:MSE:HE2	1:B:655:GLU:CB	2.39	0.51
1:D:550:ILE:HD13	1:D:662:LEU:HD13	1.93	0.51
1:D:617:SER:HA	1:D:620:ARG:HB2	1.92	0.51
1:B:547:ARG:HG3	1:B:547:ARG:NH1	2.26	0.50
1:D:563:LEU:O	1:D:567:LEU:HG	2.11	0.50
1:A:611:ARG:HH11	1:A:611:ARG:HG2	1.77	0.50
1:C:562:LEU:HD11	1:C:566:MSE:HE3	1.93	0.50
1:C:668:ASN:O	1:C:670:ALA:N	2.45	0.50
1:B:666:ILE:C	1:B:668:ASN:H	2.15	0.50
1:C:664:ARG:HH11	1:C:664:ARG:HG3	1.77	0.50
1:B:561:ARG:HH11	1:B:564:ASN:ND2	2.06	0.50
1:B:549:ARG:HD2	1:B:590:ARG:HD2	1.94	0.49
1:B:563:LEU:O	1:B:567:LEU:HG	2.11	0.49
1:D:660:LEU:HD21	1:D:664:ARG:NH2	2.28	0.48
1:B:549:ARG:HE	1:B:590:ARG:CZ	2.27	0.47
1:A:605:ASP:HB2	2:A:27:HOH:O	2.14	0.47
1:C:607:VAL:HG21	1:C:638:PHE:CZ	2.48	0.47
1:D:598:ASP:O	1:D:601:LEU:HG	2.15	0.47
1:D:601:LEU:HB3	1:D:602:PRO:HD2	1.97	0.47
1:C:645:VAL:HG22	1:C:648:TRP:NE1	2.30	0.47
1:B:547:ARG:H	1:B:548:PRO:HD2	1.79	0.47
1:B:601:LEU:HB2	1:B:604:GLN:O	2.14	0.47
1:D:603:ASP:CG	1:D:604:GLN:H	2.17	0.47
1:C:641:GLN:N	1:C:642:PRO:HD2	2.30	0.47
1:C:645:VAL:CG2	1:C:648:TRP:NE1	2.79	0.46
1:D:550:ILE:CD1	1:D:662:LEU:HD13	2.46	0.46
1:A:553:CYS:CB	1:A:595:MSE:CE	2.94	0.45
1:D:636:LEU:C	1:D:636:LEU:HD23	2.36	0.45
1:A:547:ARG:HB3	1:A:548:PRO:CD	2.45	0.44
1:C:549:ARG:NH1	1:C:575:ASP:OD1	2.51	0.44
1:A:547:ARG:HA	1:A:547:ARG:HD2	1.75	0.44
1:A:567:LEU:HA	1:A:567:LEU:HD23	1.74	0.44
1:A:599:LEU:H	1:A:628:SER:HB2	1.82	0.44
1:D:611:ARG:HG2	1:D:611:ARG:HH11	1.83	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:645:VAL:HG11	1:A:648:TRP:HZ2	1.73	0.43
1:C:563:LEU:HD21	1:C:653:ILE:HD13	2.00	0.43
1:B:599:LEU:HA	1:B:606:GLY:HA3	1.99	0.43
1:B:631:ALA:HB3	1:B:650:GLU:HG2	1.99	0.43
1:A:599:LEU:HD13	1:A:626:VAL:CG1	2.49	0.43
1:A:651:LYS:HA	1:A:652:PRO:C	2.39	0.43
1:D:547:ARG:CB	1:D:548:PRO:HD3	2.36	0.43
1:B:550:ILE:HD12	1:B:594:ALA:HB3	2.01	0.43
1:C:645:VAL:HG21	1:C:648:TRP:CE2	2.54	0.43
1:D:645:VAL:HG22	1:D:648:TRP:NE1	2.33	0.43
1:C:633:GLU:O	1:C:637:GLU:HB2	2.18	0.42
1:D:546:PRO:HD3	2:D:44:HOH:O	2.18	0.42
1:D:547:ARG:HB3	1:D:548:PRO:CD	2.36	0.42
1:A:641:GLN:HA	1:A:641:GLN:HE21	1.85	0.42
1:D:628:SER:HA	1:D:651:LYS:HE3	2.01	0.42
1:A:602:PRO:HG2	1:A:603:ASP:H	1.84	0.42
1:D:669:MSE:HE3	1:D:669:MSE:HB2	1.68	0.42
1:B:547:ARG:HB3	1:B:548:PRO:CD	2.49	0.42
1:A:662:LEU:O	1:A:666:ILE:HG13	2.20	0.42
1:B:633:GLU:OE1	1:B:633:GLU:N	2.53	0.42
1:B:623:ALA:HB1	1:B:665:ALA:HB1	2.02	0.41
1:C:632:ARG:HG3	1:C:632:ARG:HH11	1.86	0.41
1:D:623:ALA:H	1:D:669:MSE:HE1	1.84	0.41
1:C:639:ASN:C	1:C:641:GLN:H	2.22	0.41
1:D:596:THR:HA	1:D:625:VAL:O	2.21	0.41
1:C:586:GLN:HA	1:C:589:ARG:HD3	2.03	0.41
1:C:643:LEU:HD12	1:C:643:LEU:N	2.35	0.41
1:B:665:ALA:O	1:B:667:ASP:N	2.54	0.40
1:B:590:ARG:HB2	1:B:591:PRO:HD2	2.02	0.40
1:D:549:ARG:HG2	1:D:592:TYR:CE1	2.55	0.40
1:A:641:GLN:HB2	1:A:642:PRO:HD3	2.03	0.40
1:A:599:LEU:HD13	1:A:626:VAL:HG13	2.03	0.40
1:C:599:LEU:HD13	1:C:626:VAL:CG1	2.50	0.40
1:C:601:LEU:HA	1:C:602:PRO:HD3	1.88	0.40
1:D:607:VAL:HG13	1:D:638:PHE:CE2	2.57	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	123/140 (88%)	119 (97%)	1 (1%)	3 (2%)	7	4
1	B	121/140 (86%)	118 (98%)	1 (1%)	2 (2%)	11	7
1	C	123/140 (88%)	112 (91%)	8 (6%)	3 (2%)	7	4
1	D	123/140 (88%)	118 (96%)	2 (2%)	3 (2%)	7	4
All	All	490/560 (88%)	467 (95%)	12 (2%)	11 (2%)	8	4

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	603	ASP
1	D	603	ASP
1	A	547	ARG
1	C	640	SER
1	D	602	PRO
1	C	547	ARG
1	D	547	ARG
1	A	602	PRO
1	C	602	PRO
1	B	547	ARG
1	B	666	ILE

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	107/114 (94%)	104 (97%)	3 (3%)	51	63
1	B	106/114 (93%)	105 (99%)	1 (1%)	84	92
1	C	107/114 (94%)	102 (95%)	5 (5%)	32	38
1	D	107/114 (94%)	104 (97%)	3 (3%)	51	63
All	All	427/456 (94%)	415 (97%)	12 (3%)	51	63

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	599	LEU
1	A	645	VAL
1	A	664	ARG
1	B	549	ARG
1	C	599	LEU
1	C	638	PHE
1	C	650	GLU
1	C	664	ARG
1	C	669	MSE
1	D	547	ARG
1	D	599	LEU
1	D	645	VAL

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

Mol	Chain	Res	Type
1	A	641	GLN
1	B	564	ASN
1	B	600	ASN
1	D	604	GLN
1	D	630	ASN
1	D	656	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 [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 ⓘ

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, 95th 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 ²)	Q<0.9
1	A	121/140 (86%)	-0.25	2 (1%) 73 72	12, 30, 54, 60	0
1	B	120/140 (85%)	-0.30	2 (1%) 73 72	15, 27, 53, 72	0
1	C	121/140 (86%)	-0.05	6 (4%) 32 32	13, 34, 58, 66	0
1	D	121/140 (86%)	0.07	7 (5%) 26 26	14, 38, 66, 77	0
All	All	483/560 (86%)	-0.13	17 (3%) 48 46	12, 32, 59, 77	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	604	GLN	4.6
1	C	638	PHE	3.8
1	B	546	PRO	3.7
1	D	546	PRO	3.6
1	D	603	ASP	3.3
1	C	546	PRO	3.2
1	D	636	LEU	2.8
1	C	615	ARG	2.6
1	C	632	ARG	2.6
1	A	603	ASP	2.5
1	A	546	PRO	2.5
1	B	666	ILE	2.5
1	C	639	ASN	2.3
1	D	600	ASN	2.3
1	D	618	ARG	2.3
1	C	603	ASP	2.2
1	D	617	SER	2.2

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 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.