



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 18, 2026 – 05:58 PM UTC

PDB ID : 4RNS / pdb_00004rns
Title : PcpR inducer binding domain (apo-form)
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Deposited on : 2014-10-25
Resolution : 2.70 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

<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	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

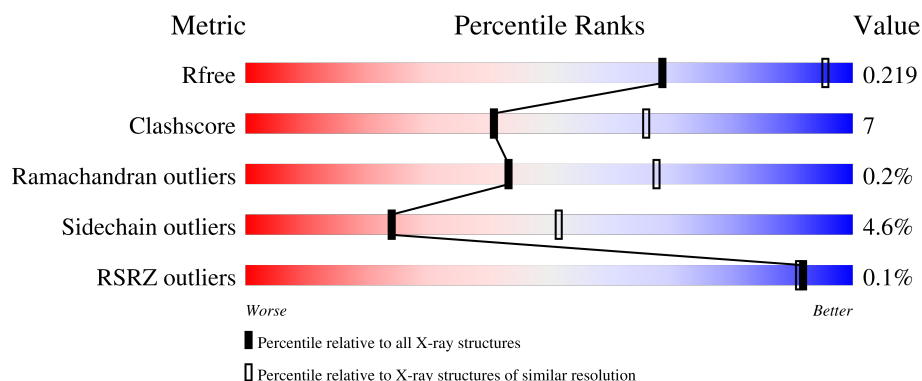
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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}	180053	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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 ≥ 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	223	 77% 18% . .
1	B	223	 71% 23% . .
1	C	223	 81% 15% . .
1	D	223	 80% 16% . .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7013 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 PCP degradation transcriptional activation protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	214	Total	C	N	O	S	0	0	0
			1720	1101	306	309	4			
1	B	213	Total	C	N	O	S	0	0	0
			1709	1095	302	308	4			
1	C	217	Total	C	N	O	S	0	0	0
			1736	1112	307	313	4			
1	D	217	Total	C	N	O	S	0	0	0
			1736	1112	307	313	4			

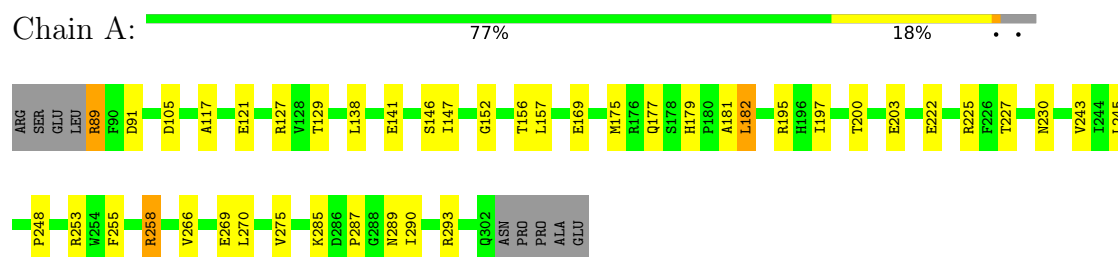
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	32	Total	O	0	0
			32	32		
2	B	18	Total	O	0	0
			18	18		
2	C	33	Total	O	0	0
			33	33		
2	D	29	Total	O	0	0
			29	29		

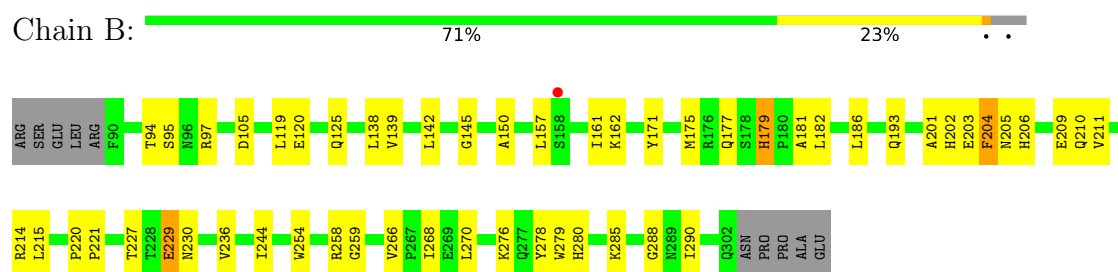
3 Residue-property plots

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.

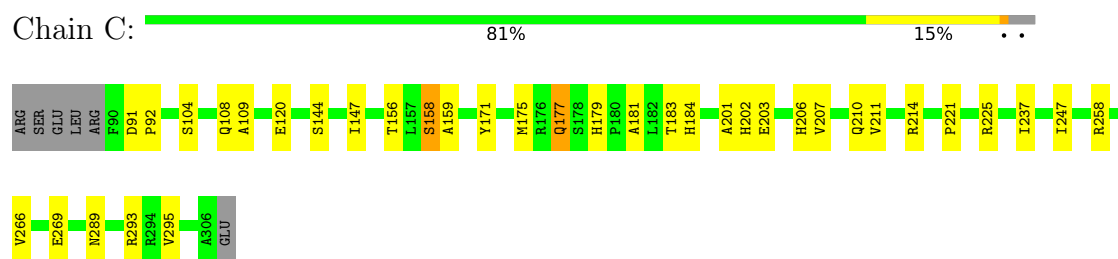
- Molecule 1: PCP degradation transcriptional activation protein



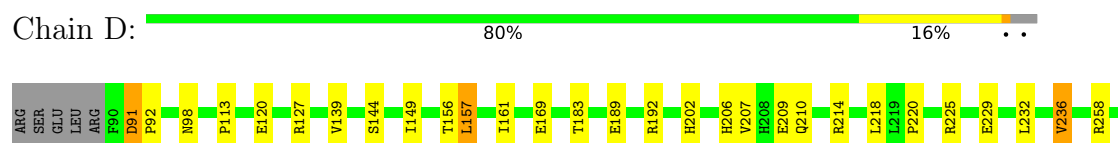
- Molecule 1: PCP degradation transcriptional activation protein



- Molecule 1: PCP degradation transcriptional activation protein



- Molecule 1: PCP degradation transcriptional activation protein



I268	E269	I273	D264	I290	R291	L292	R293	R294	Q302	A306	GLU
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4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	169.45Å 169.45Å 109.84Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	42.36 – 2.70 42.36 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.9 (42.36-2.70) 96.3 (42.36-2.70)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	-1.06 (at 2.22Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, R_{free}	0.161 , 0.216 0.172 , 0.219	Depositor DCC
R_{free} test set	2005 reflections (3.49%)	wwPDB-VP
Wilson B-factor (Å ²)	31.9	Xtriage
Anisotropy	0.169	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 27.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.025 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7013	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.01% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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

5.1 Standard geometry

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	0.66	0/1763	0.84	2/2397 (0.1%)
1	B	0.65	0/1752	0.90	4/2383 (0.2%)
1	C	0.65	0/1781	0.83	2/2425 (0.1%)
1	D	0.64	0/1781	0.86	2/2425 (0.1%)
All	All	0.65	0/7077	0.86	10/9630 (0.1%)

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	259	GLY	N-CA-C	8.80	121.89	110.45
1	D	91	ASP	CA-C-N	6.99	126.51	119.24
1	D	91	ASP	C-N-CA	6.99	126.51	119.24
1	B	145	GLY	N-CA-C	-5.74	107.75	115.21
1	C	179	HIS	CA-C-N	5.52	124.98	119.24
1	C	179	HIS	C-N-CA	5.52	124.98	119.24
1	A	91	ASP	CA-C-N	5.32	126.49	119.84
1	A	91	ASP	C-N-CA	5.32	126.49	119.84
1	B	179	HIS	CA-C-N	5.22	125.27	119.32
1	B	179	HIS	C-N-CA	5.22	125.27	119.32

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	1720	0	1711	30	0
1	B	1709	0	1698	30	0
1	C	1736	0	1723	19	0
1	D	1736	0	1723	23	0
2	A	32	0	0	3	0
2	B	18	0	0	0	0
2	C	33	0	0	0	0
2	D	29	0	0	1	0
All	All	7013	0	6855	90	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 7.

All (90) 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:225:ARG:NH2	1:B:120:GLU:OE2	2.11	0.83
1:C:120:GLU:OE2	1:D:225:ARG:NH2	2.19	0.76
1:A:179:HIS:HD2	1:A:181:ALA:H	1.33	0.74
1:D:202:HIS:HA	1:D:209:GLU:HG2	1.70	0.72
1:A:89:ARG:HB2	1:A:287:PRO:HG2	1.77	0.66
1:D:206:HIS:O	1:D:209:GLU:HG3	1.98	0.64
1:B:175:MET:HE1	1:B:181:ALA:HB3	1.80	0.62
1:A:117:ALA:O	1:A:121:GLU:HG2	2.00	0.62
1:C:214:ARG:NH2	1:C:269:GLU:O	2.32	0.62
1:A:270:LEU:O	2:A:413:HOH:O	2.16	0.60
1:D:149:ILE:HD12	1:D:292:LEU:HD23	1.85	0.58
1:B:266:VAL:HG11	1:B:270:LEU:HD12	1.86	0.58
1:D:139:VAL:HG13	1:D:161:ILE:HD11	1.87	0.56
1:A:175:MET:HE2	1:A:179:HIS:CD2	2.40	0.56
1:A:225:ARG:HD3	2:A:422:HOH:O	2.06	0.54
1:C:158:SER:OG	1:C:159:ALA:N	2.39	0.54
1:B:138:LEU:HD23	1:B:157:LEU:HD11	1.89	0.54
1:A:253:ARG:O	1:A:253:ARG:HD3	2.08	0.54
1:A:179:HIS:CD2	1:A:181:ALA:H	2.20	0.53
1:B:139:VAL:HG13	1:B:161:ILE:HD11	1.90	0.53
1:D:192:ARG:NH1	1:D:220:PRO:HD3	2.24	0.52
1:A:197:ILE:HG23	1:A:245:LEU:HD13	1.92	0.51
1:A:225:ARG:NH1	2:A:422:HOH:O	2.31	0.51
1:C:104:SER:O	1:C:108:GLN:HG3	2.11	0.50
1:B:210:GLN:O	1:B:214:ARG:HG2	2.12	0.50
1:C:221:PRO:O	1:D:127:ARG:NH2	2.43	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:138:LEU:HB3	1:B:157:LEU:HD21	1.93	0.49
1:A:138:LEU:HD23	1:A:157:LEU:HD11	1.94	0.49
1:B:162:LYS:HB2	1:B:279:TRP:CE2	2.48	0.49
1:A:225:ARG:HD2	1:B:119:LEU:HD13	1.95	0.49
1:B:201:ALA:HB1	1:B:206:HIS:HB3	1.95	0.48
1:B:142:LEU:HB3	1:B:280:HIS:HB2	1.94	0.48
1:C:109:ALA:HB1	1:D:236:VAL:HG21	1.94	0.48
1:B:202:HIS:HA	1:B:209:GLU:HG2	1.95	0.48
1:C:225:ARG:HH21	1:D:120:GLU:HG2	1.79	0.48
1:C:177:GLN:N	1:C:177:GLN:OE1	2.47	0.48
1:C:289:ASN:OD1	1:C:293:ARG:NH1	2.46	0.48
1:A:258:ARG:NH2	1:B:254:TRP:CH2	2.82	0.48
1:A:89:ARG:HH11	1:A:89:ARG:HG2	1.78	0.47
1:B:276:LYS:HB2	1:B:278:TYR:CE1	2.48	0.47
1:B:285:LYS:HA	1:B:290:ILE:HD11	1.96	0.47
1:A:177:GLN:HA	1:A:182:LEU:HD13	1.97	0.47
1:A:289:ASN:O	1:A:293:ARG:HG3	2.14	0.47
1:B:105:ASP:OD2	1:B:230:ASN:HB3	2.15	0.46
1:B:171:TYR:CZ	1:B:211:VAL:HG21	2.49	0.46
1:A:105:ASP:OD2	1:A:230:ASN:HB3	2.15	0.46
1:D:290:ILE:O	1:D:294:ARG:HG2	2.14	0.46
1:C:201:ALA:HB1	1:C:206:HIS:HB3	1.96	0.46
1:C:207:VAL:O	1:C:210:GLN:HG2	2.16	0.46
1:C:237:ILE:HG12	1:D:113:PRO:HG3	1.98	0.46
1:B:95:SER:O	1:B:125:GLN:HB3	2.16	0.45
1:A:152:GLY:HA2	1:A:275:VAL:HG13	1.99	0.45
1:D:207:VAL:HG13	1:D:273:ILE:HD13	1.99	0.45
1:C:175:MET:HE1	1:C:181:ALA:HB3	2.00	0.44
1:C:225:ARG:NH2	1:D:120:GLU:OE2	2.50	0.44
1:A:141:GLU:OE1	1:A:146:SER:OG	2.33	0.44
1:D:202:HIS:CA	1:D:209:GLU:HG2	2.42	0.44
1:D:214:ARG:NH1	1:D:269:GLU:O	2.51	0.44
1:C:183:THR:OG1	1:C:184:HIS:N	2.50	0.44
1:A:195:ARG:HB3	1:A:243:VAL:HG12	2.00	0.43
1:C:171:TYR:CE1	1:C:211:VAL:HG21	2.52	0.43
1:B:175:MET:HE3	1:B:179:HIS:CG	2.53	0.43
1:A:285:LYS:HA	1:A:290:ILE:HD11	2.00	0.43
1:C:289:ASN:O	1:C:293:ARG:HG3	2.18	0.43
1:B:220:PRO:HA	1:B:221:PRO:HD3	1.85	0.43
1:D:210:GLN:O	1:D:214:ARG:HG2	2.19	0.43
1:D:91:ASP:HA	1:D:92:PRO:HD3	1.89	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:284:ASP:HA	1:D:293:ARG:HH12	1.83	0.42
1:A:255:PHE:O	1:A:258:ARG:HG3	2.19	0.42
1:B:177:GLN:O	1:B:182:LEU:HD22	2.19	0.42
1:A:147:ILE:HD13	1:A:147:ILE:HA	1.93	0.42
1:B:276:LYS:HD2	1:B:278:TYR:OH	2.20	0.42
1:A:89:ARG:HH11	1:A:89:ARG:CG	2.33	0.42
1:D:98:ASN:ND2	2:D:409:HOH:O	2.49	0.42
1:B:215:LEU:HD23	1:B:268:ILE:HD12	2.01	0.41
1:D:218:LEU:HD22	1:D:268:ILE:HD13	2.02	0.41
1:A:127:ARG:NH2	1:B:221:PRO:O	2.52	0.41
1:B:171:TYR:CE1	1:B:211:VAL:HG21	2.54	0.41
1:B:97:ARG:NH1	1:B:288:GLY:HA3	2.35	0.41
1:A:269:GLU:OE1	1:D:183:THR:OG1	2.37	0.41
1:B:204:PHE:HZ	1:B:229:GLU:OE1	2.03	0.41
1:A:177:GLN:O	1:A:182:LEU:HD22	2.21	0.41
1:D:232:LEU:O	1:D:236:VAL:HG13	2.20	0.41
1:A:169:GLU:HG2	1:A:248:PRO:HB3	2.02	0.40
1:A:129:THR:OG1	1:B:227:THR:HB	2.22	0.40
1:B:186:LEU:HA	1:B:186:LEU:HD12	1.88	0.40
1:D:157:LEU:HG	1:D:161:ILE:HG13	2.03	0.40
1:B:150:ALA:HB3	1:B:278:TYR:HB2	2.02	0.40
1:C:202:HIS:CD2	1:C:203:GLU:HG3	2.56	0.40
1:C:91:ASP:HA	1:C:92:PRO:HD3	1.96	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	212/223 (95%)	202 (95%)	10 (5%)	0	100	100
1	B	211/223 (95%)	197 (93%)	13 (6%)	1 (0%)	24	48

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	215/223 (96%)	206 (96%)	8 (4%)	1 (0%)	24	48
1	D	215/223 (96%)	207 (96%)	8 (4%)	0	100	100
All	All	853/892 (96%)	812 (95%)	39 (5%)	2 (0%)	43	68

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	158	SER
1	B	205	ASN

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	185/193 (96%)	176 (95%)	9 (5%)	22	49
1	B	184/193 (95%)	176 (96%)	8 (4%)	26	54
1	C	187/193 (97%)	179 (96%)	8 (4%)	26	54
1	D	187/193 (97%)	178 (95%)	9 (5%)	23	50
All	All	743/772 (96%)	709 (95%)	34 (5%)	24	51

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	89	ARG
1	A	156	THR
1	A	182	LEU
1	A	200	THR
1	A	203	GLU
1	A	222	GLU
1	A	227	THR
1	A	258	ARG
1	A	266	VAL
1	B	94	THR

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Mol	Chain	Res	Type
1	B	193	GLN
1	B	203	GLU
1	B	204	PHE
1	B	229	GLU
1	B	236	VAL
1	B	244	ILE
1	B	258	ARG
1	C	144	SER
1	C	147	ILE
1	C	156	THR
1	C	177	GLN
1	C	247	ILE
1	C	258	ARG
1	C	266	VAL
1	C	295	VAL
1	D	144	SER
1	D	156	THR
1	D	157	LEU
1	D	169	GLU
1	D	189	GLU
1	D	229	GLU
1	D	236	VAL
1	D	258	ARG
1	D	302	GLN

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

Mol	Chain	Res	Type
1	A	135	HIS
1	A	179	HIS
1	A	277	GLN
1	C	98	ASN
1	C	108	GLN
1	C	135	HIS
1	C	164	GLN
1	C	202	HIS
1	C	277	GLN
1	D	202	HIS
1	D	210	GLN
1	D	303	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 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 [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, 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	214/223 (95%)	-0.58	0 100 100	27, 40, 67, 101	0
1	B	213/223 (95%)	-0.34	1 (0%) 87 86	29, 47, 81, 91	0
1	C	217/223 (97%)	-0.52	0 100 100	27, 43, 67, 82	0
1	D	217/223 (97%)	-0.55	0 100 100	28, 41, 65, 81	0
All	All	861/892 (96%)	-0.50	1 (0%) 92 91	27, 43, 70, 101	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	158	SER	2.1

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 oligosaccharides 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.