



Full wwPDB X-ray Structure Validation Report ⓘ

May 4, 2017 – 05:11 PM EDT

PDB ID : 5UNF
Title : XFEL structure of human angiotensin II type 2 receptor (Monoclinic form) in complex with compound 1 (N-benzyl-N-(2-ethyl-4-oxo-3-{[2'-(2H-tetrazol-5-yl)[1,1'-biphenyl]-4-yl])
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Deposited on : 2017-01-30
Resolution : 2.80 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	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20029077
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)	:	rb-20029077

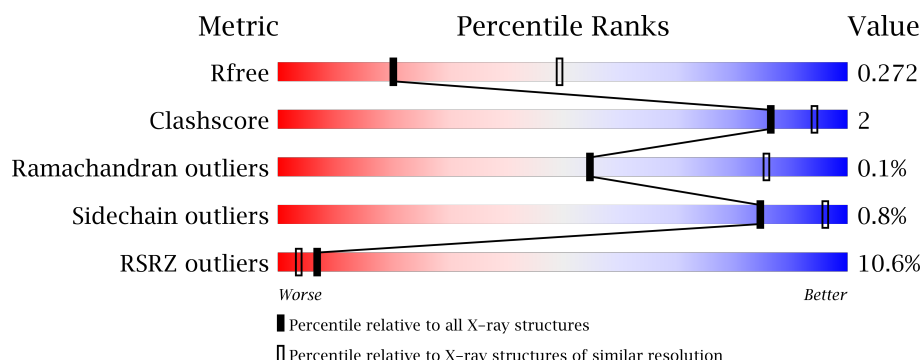
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


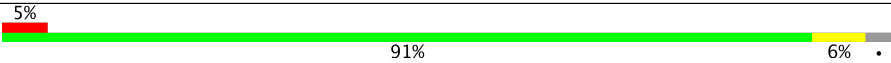
The reported resolution of this entry is 2.80 Å.

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	2583 (2.80-2.80)
Clashscore	112137	3033 (2.80-2.80)
Ramachandran outliers	110173	2983 (2.80-2.80)
Sidechain outliers	110143	2985 (2.80-2.80)
RSRZ outliers	101464	2610 (2.80-2.80)

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	411	
1	B	411	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6243 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.

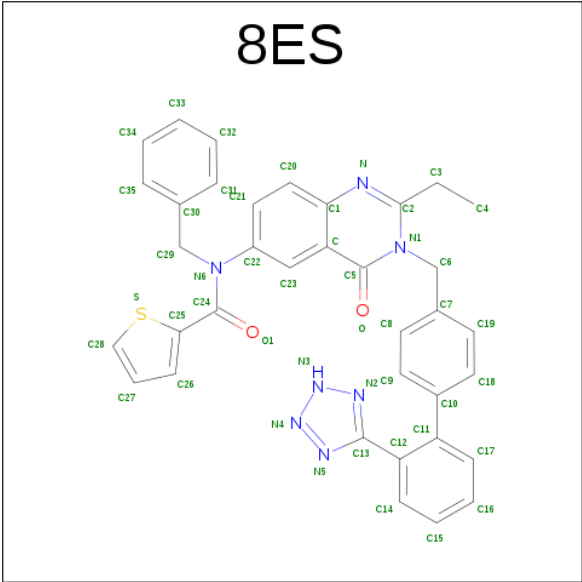
- Molecule 1 is a protein called Chimera protein of Type-2 angiotensin II receptor and Soluble cytochrome b562.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	386	Total	C	N	O	S	0	0	0
			3034	2008	482	522	22			
1	B	399	Total	C	N	O	S	0	0	0
			3117	2056	496	542	23			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1007	TRP	MET	engineered mutation	UNP P0ABE7
A	1102	ILE	HIS	engineered mutation	UNP P0ABE7
A	1106	LEU	ARG	engineered mutation	UNP P0ABE7
A	1107	GLY	-	linker	UNP P0ABE7
A	1108	SER	-	linker	UNP P0ABE7
A	1109	GLY	-	linker	UNP P0ABE7
A	1110	SER	-	linker	UNP P0ABE7
B	1007	TRP	MET	engineered mutation	UNP P0ABE7
B	1102	ILE	HIS	engineered mutation	UNP P0ABE7
B	1106	LEU	ARG	engineered mutation	UNP P0ABE7
B	1107	GLY	-	linker	UNP P0ABE7
B	1108	SER	-	linker	UNP P0ABE7
B	1109	GLY	-	linker	UNP P0ABE7
B	1110	SER	-	linker	UNP P0ABE7

- Molecule 2 is N-benzyl-N-(2-ethyl-4-oxo-3-{{[2'-(2H-tetrazol-5-yl)][1,1'-biphenyl]-4-yl]methyl}-3,4-dihydroquinazolin-6-yl)thiophene-2-carboxamide (three-letter code: 8ES) (formula: C₃₆H₂₉N₇O₂S).

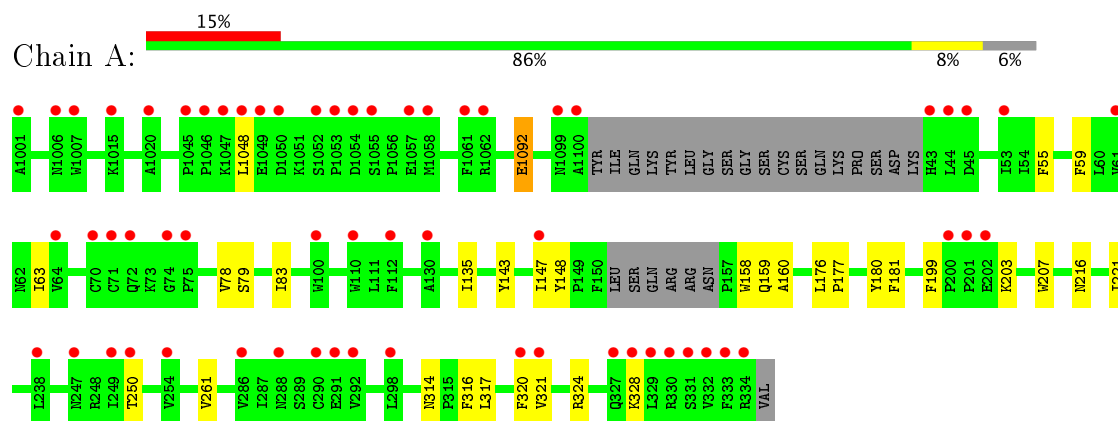


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			46	36	7	2	1		
2	B	1	Total	C	N	O	S	0	0
			46	36	7	2	1		

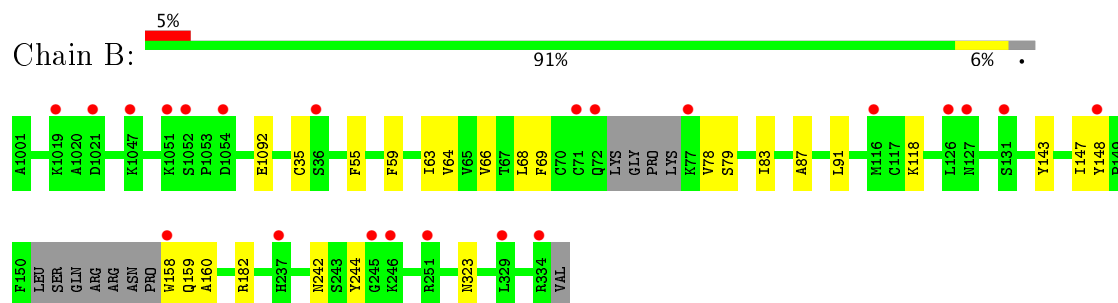
3 Residue-property plots [i](#)

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: Chimera protein of Type-2 angiotensin II receptor and Soluble cytochrome b562



- Molecule 1: Chimera protein of Type-2 angiotensin II receptor and Soluble cytochrome b562



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	77.39Å 69.11Å 90.07Å 90.00° 104.33° 90.00°	Depositor
Resolution (Å)	29.57 – 2.80 29.09 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.9 (29.57-2.80) 99.9 (29.09-2.80)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.22 (at 2.80Å)	Xtriage
Refinement program	BUSTER 2.10.2	Depositor
R, R_{free}	0.227 , 0.256 0.246 , 0.272	Depositor DCC
R_{free} test set	1117 reflections (4.88%)	DCC
Wilson B-factor (Å ²)	84.0	Xtriage
Anisotropy	0.215	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 84.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6243	wwPDB-VP
Average B, all atoms (Å ²)	111.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.32% 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

Bond lengths and bond angles in the following residue types are not validated in this section: 8ES

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	0.50	0/3112	0.55	0/4237
1	B	0.51	0/3195	0.55	0/4350
All	All	0.50	0/6307	0.55	0/8587

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

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	3034	0	3052	20	0
1	B	3117	0	3112	14	0
2	A	46	0	0	0	0
2	B	46	0	0	0	0
All	All	6243	0	6164	31	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 2.

All (31) 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:242:ASN:HD21	1:B:244:TYR:HD2	1.44	0.66
1:A:59:PHE:O	1:A:63:ILE:HG12	2.04	0.58
1:B:59:PHE:O	1:B:63:ILE:HG12	2.07	0.55
1:A:78:VAL:HG11	1:A:159:GLN:HE21	1.71	0.54
1:B:78:VAL:HG13	1:B:160:ALA:HB2	1.90	0.54
1:B:78:VAL:HG11	1:B:159:GLN:HE21	1.73	0.53
1:A:261:VAL:HA	1:A:317:LEU:HD21	1.90	0.52
1:A:176:LEU:HD13	1:B:66:VAL:HG13	1.92	0.51
1:A:78:VAL:HG13	1:A:160:ALA:HB2	1.92	0.50
1:A:135:ILE:HD11	1:A:314:ASN:HD21	1.77	0.50
1:A:1092:GLU:HG2	1:B:158:TRP:CD1	2.48	0.49
1:A:180:TYR:CE1	1:B:69:PHE:HB2	2.48	0.48
1:A:321:VAL:HG13	1:A:324:ARG:HH21	1.79	0.47
1:A:316:PHE:HA	1:A:320:PHE:HD2	1.80	0.46
1:A:250:THR:HG23	1:A:328:LYS:HD2	1.97	0.46
1:A:177:PRO:O	1:A:181:PHE:HD2	1.99	0.46
1:A:147:ILE:HG23	1:A:148:TYR:CD2	2.54	0.43
1:A:216:ASN:O	1:A:221:ILE:HG12	2.18	0.43
1:B:143:TYR:O	1:B:147:ILE:HG22	2.18	0.43
1:A:143:TYR:O	1:A:147:ILE:HG22	2.18	0.43
1:B:147:ILE:HG23	1:B:148:TYR:CD2	2.54	0.43
1:B:79:SER:O	1:B:83:ILE:HG12	2.19	0.43
1:A:261:VAL:HG12	1:A:317:LEU:HG	2.01	0.42
1:B:64:VAL:O	1:B:68:LEU:HB2	2.20	0.42
1:A:199:PHE:CD1	1:A:207:TRP:HB3	2.55	0.42
1:B:118:LYS:HG2	1:B:182:ARG:O	2.20	0.42
1:A:79:SER:O	1:A:83:ILE:HG12	2.19	0.41
1:B:87:ALA:O	1:B:91:LEU:HB2	2.21	0.41
1:A:55:PHE:O	1:A:59:PHE:HB2	2.21	0.41
1:A:203:LYS:O	1:A:207:TRP:HD1	2.05	0.40
1:B:55:PHE:O	1:B:59:PHE:HB2	2.20	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	380/411 (92%)	365 (96%)	15 (4%)	0	100	100
1	B	393/411 (96%)	381 (97%)	11 (3%)	1 (0%)	44	77
All	All	773/822 (94%)	746 (96%)	26 (3%)	1 (0%)	55	86

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	35	CYS

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	320/354 (90%)	317 (99%)	3 (1%)	82	95
1	B	327/354 (92%)	325 (99%)	2 (1%)	89	97
All	All	647/708 (91%)	642 (99%)	5 (1%)	85	96

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1048	LEU
1	A	1092	GLU
1	A	158	TRP
1	B	1092	GLU
1	B	323	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1013	ASN
1	A	72	GLN

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Mol	Chain	Res	Type
1	A	159	GLN
1	A	314	ASN
1	B	72	GLN
1	B	144	GLN
1	B	159	GLN
1	B	216	ASN
1	B	242	ASN
1	B	273	HIS
1	B	323	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	8ES	A	1501	-	51,52,52	3.67	25 (49%)	62,73,73	1.84	9 (14%)
2	8ES	B	1201	-	51,52,52	3.72	25 (49%)	62,73,73	1.81	10 (16%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral

centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	8ES	A	1501	-	-	0/26/30/30	0/7/7/7
2	8ES	B	1201	-	-	0/26/30/30	0/7/7/7

All (50) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1201	8ES	C34-C35	-3.35	1.32	1.38
2	A	1501	8ES	C34-C35	-3.35	1.32	1.38
2	B	1201	8ES	C8-C7	-3.15	1.32	1.38
2	A	1501	8ES	C9-C10	-3.15	1.32	1.39
2	A	1501	8ES	C8-C7	-3.11	1.32	1.38
2	A	1501	8ES	C31-C30	-3.08	1.32	1.38
2	B	1201	8ES	C9-C10	-3.07	1.32	1.39
2	A	1501	8ES	C18-C19	-2.95	1.33	1.38
2	B	1201	8ES	C18-C19	-2.94	1.33	1.38
2	B	1201	8ES	C31-C30	-2.74	1.33	1.38
2	A	1501	8ES	C33-C32	-2.35	1.32	1.38
2	B	1201	8ES	C33-C32	-2.09	1.33	1.38
2	A	1501	8ES	C1-N	2.04	1.41	1.37
2	A	1501	8ES	C12-C13	2.06	1.52	1.48
2	A	1501	8ES	C20-C21	2.22	1.41	1.36
2	B	1201	8ES	C12-C13	2.24	1.53	1.48
2	B	1201	8ES	C20-C21	2.25	1.41	1.36
2	B	1201	8ES	C1-N	2.29	1.41	1.37
2	A	1501	8ES	C11-C10	2.95	1.54	1.49
2	B	1201	8ES	C11-C10	3.15	1.54	1.49
2	A	1501	8ES	N3-N4	3.40	1.37	1.32
2	A	1501	8ES	C2-N1	3.45	1.44	1.36
2	B	1201	8ES	N3-N4	3.52	1.37	1.32
2	B	1201	8ES	C2-N1	3.53	1.44	1.36
2	B	1201	8ES	C27-C26	3.83	1.52	1.39
2	A	1501	8ES	C27-C26	3.86	1.52	1.39
2	A	1501	8ES	N2-N3	3.91	1.40	1.34
2	B	1201	8ES	N2-N3	4.05	1.41	1.34
2	A	1501	8ES	C5-C	4.80	1.49	1.41
2	B	1201	8ES	C5-C	5.43	1.50	1.41
2	A	1501	8ES	C24-N6	6.12	1.46	1.36
2	B	1201	8ES	C24-N6	6.24	1.47	1.36
2	A	1501	8ES	C34-C33	6.51	1.53	1.38
2	B	1201	8ES	C34-C33	6.61	1.54	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1501	8ES	O-C5	6.66	1.41	1.24
2	B	1201	8ES	O-C5	6.69	1.41	1.24
2	A	1501	8ES	C13-N5	6.79	1.42	1.33
2	B	1201	8ES	C13-N5	6.88	1.42	1.33
2	A	1501	8ES	C35-C30	7.04	1.53	1.38
2	B	1201	8ES	C18-C10	7.07	1.54	1.39
2	A	1501	8ES	C18-C10	7.17	1.54	1.39
2	B	1201	8ES	C35-C30	7.22	1.53	1.38
2	A	1501	8ES	C3-C2	7.33	1.54	1.49
2	B	1201	8ES	C32-C31	7.50	1.53	1.38
2	B	1201	8ES	C3-C2	7.68	1.54	1.49
2	A	1501	8ES	C19-C7	7.71	1.54	1.38
2	B	1201	8ES	C19-C7	7.82	1.55	1.38
2	A	1501	8ES	C32-C31	7.84	1.53	1.38
2	B	1201	8ES	C9-C8	8.06	1.53	1.38
2	A	1501	8ES	C9-C8	8.10	1.53	1.38

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1501	8ES	N5-N4-N3	-6.52	105.10	109.54
2	B	1201	8ES	N5-N4-N3	-6.31	105.25	109.54
2	A	1501	8ES	C-C1-N	-4.60	121.05	123.67
2	B	1201	8ES	C-C1-N	-4.37	121.17	123.67
2	B	1201	8ES	N2-C13-N5	-3.38	106.90	111.31
2	A	1501	8ES	N2-C13-N5	-3.13	107.23	111.31
2	A	1501	8ES	C27-C28-S	-2.59	110.40	113.18
2	A	1501	8ES	C30-C29-N6	-2.37	109.02	113.35
2	B	1201	8ES	C30-C29-N6	-2.10	109.52	113.35
2	B	1201	8ES	C12-C13-N5	2.04	126.93	124.17
2	B	1201	8ES	C23-C-C1	2.12	120.64	118.16
2	B	1201	8ES	C13-N2-N3	2.21	107.21	104.93
2	A	1501	8ES	C23-C-C1	2.40	120.97	118.16
2	A	1501	8ES	C-C5-N1	4.47	119.20	116.15
2	B	1201	8ES	C-C5-N1	4.52	119.23	116.15
2	B	1201	8ES	C25-C24-N6	4.74	123.07	118.22
2	A	1501	8ES	C13-N5-N4	5.04	110.13	104.93
2	B	1201	8ES	C13-N5-N4	5.08	110.17	104.93
2	A	1501	8ES	C25-C24-N6	5.13	123.47	118.22

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

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, 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	386/411 (93%)	0.84	61 (15%) 2 1	82, 130, 188, 226	0
1	B	399/411 (97%)	0.12	22 (5%) 26 17	59, 83, 149, 182	0
All	All	785/822 (95%)	0.47	83 (10%) 7 4	59, 108, 176, 226	0

All (83) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	75	PRO	10.3
1	A	333	PHE	10.3
1	A	331	SER	8.6
1	A	1053	PRO	8.1
1	A	1054	ASP	8.0
1	A	43	HIS	7.4
1	A	74	GLY	6.6
1	A	330	ARG	6.2
1	A	44	LEU	5.5
1	A	1048	LEU	4.7
1	B	36	SER	4.7
1	B	246	LYS	4.6
1	A	1046	PRO	4.6
1	A	71	CYS	4.6
1	A	1052	SER	4.3
1	A	249	ILE	4.2
1	A	1006	ASN	4.1
1	B	245	GLY	4.0
1	A	329	LEU	4.0
1	A	1061	PHE	4.0
1	A	290	CYS	4.0
1	A	1047	LYS	3.9
1	B	148	TYR	3.9
1	A	1058	MET	3.8

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Mol	Chain	Res	Type	RSRZ
1	A	321	VAL	3.8
1	A	1099	ASN	3.8
1	A	1045	PRO	3.8
1	A	334	ARG	3.5
1	A	201	PRO	3.4
1	A	1057	GLU	3.4
1	A	1055	SER	3.4
1	B	71	CYS	3.3
1	A	1050	ASP	3.3
1	A	288	ASN	3.2
1	A	1049	GLU	3.2
1	A	327	GLN	3.1
1	A	147	ILE	3.1
1	A	328	LYS	2.9
1	A	1007	TRP	2.9
1	A	332	VAL	2.9
1	A	1100	ALA	2.9
1	B	158	TRP	2.8
1	A	1001	ALA	2.8
1	A	202	GLU	2.8
1	A	320	PHE	2.7
1	A	130	ALA	2.7
1	A	254	VAL	2.7
1	A	1062	ARG	2.7
1	B	1021	ASP	2.7
1	A	298	LEU	2.6
1	A	53	ILE	2.6
1	A	247	ASN	2.6
1	A	1020	ALA	2.5
1	A	72	GLN	2.5
1	A	100	TRP	2.5
1	B	1054	ASP	2.4
1	A	200	PRO	2.4
1	B	1052	SER	2.4
1	B	131	SER	2.4
1	A	45	ASP	2.4
1	A	238	LEU	2.3
1	A	291	GLU	2.3
1	A	1015	LYS	2.3
1	B	72	GLN	2.3
1	B	1051	LYS	2.3
1	A	250	THR	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	237	HIS	2.2
1	A	61	VAL	2.2
1	A	112	PHE	2.2
1	A	292	VAL	2.2
1	B	1019	LYS	2.2
1	B	127	ASN	2.1
1	B	77	LYS	2.1
1	B	329	LEU	2.1
1	B	334	ARG	2.1
1	A	64	VAL	2.1
1	B	126	LEU	2.1
1	B	251	ARG	2.1
1	A	70	CYS	2.0
1	A	110	TRP	2.0
1	B	1047	LYS	2.0
1	B	116	MET	2.0
1	A	286	VAL	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	8ES	A	1501	46/46	0.88	0.27	0.73	118,123,135,136	0
2	8ES	B	1201	46/46	0.96	0.20	0.29	59,64,69,71	0

6.5 Other polymers [i](#)

There are no such residues in this entry.