



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 02:10 AM UTC

PDB ID : 1M8Y / pdb_00001m8y
Title : CRYSTAL STRUCTURE OF THE PUMILIO-HOMOLOGY DOMAIN
FROM HUMAN PUMILIO1 IN COMPLEX WITH NRE2-10 RNA
Authors : Wang, X.; McLachlan, J.; Zamore, P.D.; Hall, T.M.T.
Deposited on : 2002-07-26
Resolution : 2.60 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

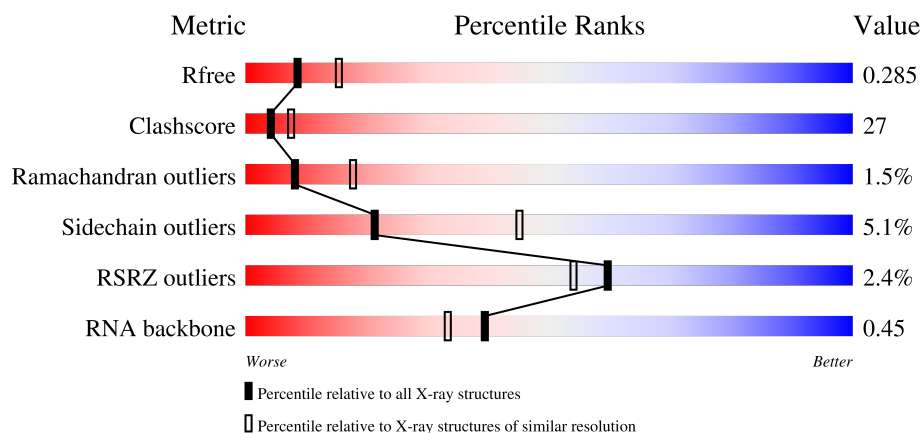
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)
RNA backbone	3983	1014 (2.84-2.36)

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	C	10	10% 30% 50% 10% 10%
1	D	10	10% 20% 60% 20%
2	A	349	% 58% 38% ..
2	B	349	3% 44% 46% 8% .

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6303 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 RNA chain called 5'-R(P*AP*UP*UP*GP*UP*AP*CP*AP*UP*A)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	10	Total	C	N	O	P	0	0	0
			211	95	36	70	10			
1	D	10	Total	C	N	O	P	0	0	0
			211	95	36	70	10			

- Molecule 2 is a protein called Pumilio 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	341	Total	C	N	O	S	0	0	0
			2768	1752	501	498	17			
2	B	342	Total	C	N	O	S	0	0	0
			2773	1755	502	499	17			

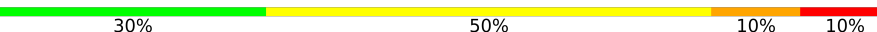
- Molecule 3 is water.

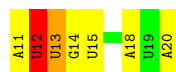
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	30	Total	O	0	0
			30	30		
3	D	16	Total	O	0	0
			16	16		
3	A	190	Total	O	0	0
			190	190		
3	B	104	Total	O	0	0
			104	104		

3 Residue-property plots

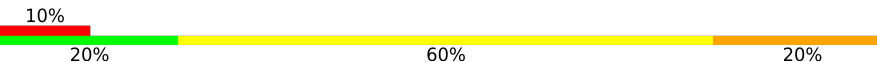
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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.

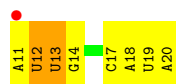
- Molecule 1: 5'-R(P*AP*UP*UP*GP*UP*AP*CP*AP*UP*A)-3'

Chain C: 



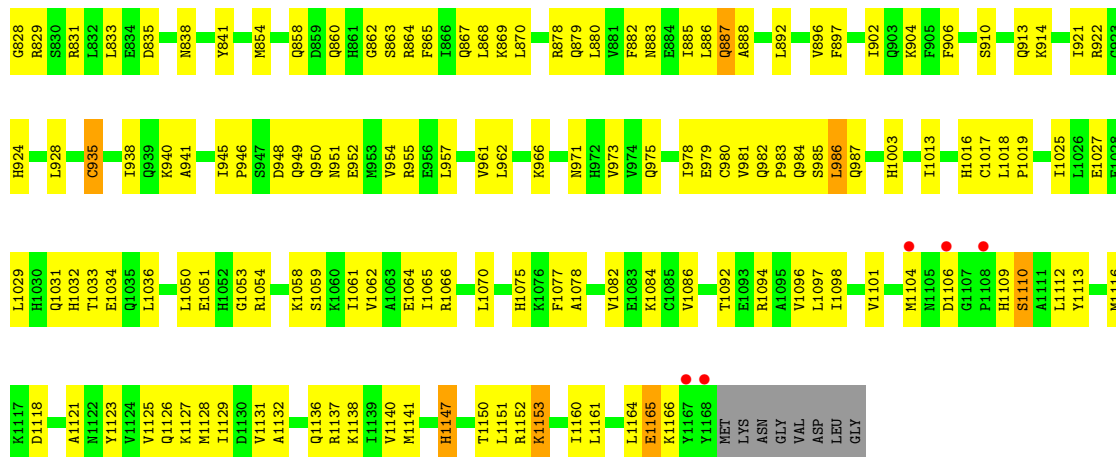
- Molecule 1: 5'-R(P*AP*UP*UP*GP*UP*AP*CP*AP*UP*A)-3'

Chain D: 



- Molecule 2: Pumilio 1

Chain A: 



- Molecule 2: Pumilio 1

Chain B: 





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	35.84Å 59.73Å 340.85Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.74 – 2.60 37.74 – 2.60	Depositor EDS
% Data completeness (in resolution range)	83.4 (37.74-2.60) 83.4 (37.74-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.08 (at 2.61Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.214 , 0.286 0.214 , 0.285	Depositor DCC
R_{free} test set	1193 reflections (5.84%)	wwPDB-VP
Wilson B-factor (Å ²)	29.0	Xtriage
Anisotropy	0.486	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 60.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6303	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.92% 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 [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	C	0.46	0/235	0.86	1/363 (0.3%)
1	D	0.40	0/235	0.78	0/363
2	A	0.48	0/2821	0.97	14/3804 (0.4%)
2	B	0.46	0/2826	0.98	11/3811 (0.3%)
All	All	0.47	0/6117	0.96	26/8341 (0.3%)

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	D	0	1

There are no bond length outliers.

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1091	ARG	N-CA-C	-12.99	97.90	113.88
2	B	1111	ALA	N-CA-C	-7.57	103.04	111.82
2	A	869	LYS	N-CA-C	7.29	120.27	111.82
2	B	1160	ILE	N-CA-C	-6.85	103.63	110.62
2	B	944	PHE	N-CA-C	6.84	121.91	113.50
2	A	1025	ILE	N-CA-C	-6.55	104.38	110.53
2	B	1071	VAL	N-CA-C	6.32	117.10	110.72
2	A	1018	LEU	CA-C-N	6.07	127.42	119.84
2	A	1018	LEU	C-N-CA	6.07	127.42	119.84
2	B	997	VAL	N-CA-C	5.95	116.60	110.36
1	C	12	U	C2'-C3'-O3'	5.91	118.36	109.50
2	A	1016	HIS	N-CA-C	5.84	121.28	114.04
2	A	1017	CYS	N-CA-C	5.73	118.09	110.53
2	B	1005	TYR	N-CA-C	-5.67	106.03	113.12

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1029	LEU	N-CA-C	-5.63	105.44	112.93
2	B	1099	ASP	N-CA-C	-5.49	106.56	113.20
2	A	1054	ARG	CA-C-N	5.42	126.61	119.84
2	A	1054	ARG	C-N-CA	5.42	126.61	119.84
2	A	973	VAL	N-CA-C	-5.35	105.50	110.53
2	B	1148	ILE	N-CA-C	5.34	116.07	110.62
2	A	1036	LEU	N-CA-C	5.31	117.07	111.28
2	A	980	CYS	N-CA-C	5.26	119.97	113.50
2	A	1053	GLY	N-CA-C	5.10	121.06	112.85
2	A	887	GLN	N-CA-C	5.07	118.61	112.23
2	B	851	GLY	N-CA-C	-5.06	108.30	115.43
2	A	1147	HIS	N-CA-C	5.05	118.47	112.72

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	19	U	Sidechain

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	C	211	0	107	14	0
1	D	211	0	107	7	0
2	A	2768	0	2783	117	0
2	B	2773	0	2785	197	0
3	A	190	0	0	4	0
3	B	104	0	0	6	0
3	C	30	0	0	0	0
3	D	16	0	0	0	0
All	All	6303	0	5782	322	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 27.

All (322) 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)
2:A:1141:MET:HE1	2:A:1164:LEU:HA	1.39	1.04
2:B:831:ARG:H	2:B:831:ARG:HD3	1.21	1.04
2:B:1124:VAL:HG12	2:B:1128:MET:HE2	1.43	1.01
2:B:975:GLN:HG3	2:B:1012:ARG:HG3	1.45	0.96
1:D:11:A:H5'	1:D:14:G:OP2	1.67	0.95
1:C:11:A:H5'	1:C:12:U:H3'	1.50	0.92
2:B:1113:TYR:HA	2:B:1116:MET:HE3	1.50	0.90
2:B:951:ASN:HB3	2:B:955:ARG:HH22	1.37	0.87
2:B:1089:ALA:HB3	2:B:1094:ARG:HB2	1.58	0.85
2:B:831:ARG:HD3	2:B:831:ARG:N	1.96	0.80
2:A:1113:TYR:HA	2:A:1116:MET:HE2	1.63	0.79
2:B:1138:LYS:HB3	2:B:1138:LYS:NZ	1.98	0.78
2:B:919:GLU:HG2	2:B:922:ARG:NH2	1.99	0.77
2:A:864:ARG:O	2:A:868:LEU:HG	1.84	0.77
2:B:921:ILE:HG21	2:B:953:MET:HE2	1.68	0.74
2:A:1141:MET:HE3	2:A:1164:LEU:HD12	1.70	0.73
2:B:1144:ILE:HG23	2:B:1164:LEU:HD11	1.69	0.73
2:A:829:ARG:HD3	2:A:833:LEU:HD23	1.69	0.73
2:A:885:ILE:HD12	2:A:886:LEU:N	2.04	0.73
2:B:990:ILE:HG22	2:B:994:LYS:HE3	1.71	0.73
1:C:11:A:H5''	1:C:12:U:C6	2.24	0.73
2:A:966:LYS:HE3	3:A:92:HOH:O	1.89	0.72
2:B:1094:ARG:HD3	2:B:1131:VAL:HG23	1.70	0.72
2:B:883:ASN:HD22	2:B:883:ASN:C	1.97	0.72
2:B:1166:LYS:HA	2:B:1169:MET:O	1.89	0.71
2:B:866:ILE:O	2:B:870:LEU:HB2	1.91	0.71
2:B:1018:LEU:H	2:B:1021:GLN:NE2	1.89	0.71
2:B:946:PRO:HB2	2:B:949:GLN:NE2	2.06	0.70
2:A:952:GLU:HA	2:A:955:ARG:HH11	1.56	0.70
2:B:949:GLN:NE2	2:B:949:GLN:H	1.89	0.70
2:B:910:SER:OG	2:B:913:GLN:HG3	1.93	0.69
2:B:1084:LYS:HA	2:B:1087:THR:HG22	1.74	0.69
2:B:874:THR:O	2:B:878:ARG:HG3	1.93	0.69
2:A:1050:LEU:HD13	2:A:1084:LYS:HB3	1.73	0.69
2:A:882:PHE:CD2	2:A:885:ILE:HD11	2.29	0.68
2:B:1125:VAL:O	2:B:1129:ILE:HG13	1.93	0.68
2:B:1033:THR:O	2:B:1037:VAL:HG13	1.94	0.68
1:C:20:A:H1'	2:A:897:PHE:HE2	1.60	0.67
2:A:1109:HIS:HD2	2:A:1110:SER:H	1.42	0.67
1:D:20:A:H1'	2:B:860:GLN:HE21	1.60	0.67
1:C:12:U:O2'	1:C:13:U:OP1	2.12	0.67
2:A:1118:ASP:HB3	2:A:1121:ALA:HB3	1.75	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:831:ARG:H	2:B:831:ARG:CD	2.03	0.66
2:B:949:GLN:H	2:B:949:GLN:HE21	1.41	0.66
2:B:874:THR:OG1	2:B:877:GLU:HG3	1.95	0.66
2:B:847:ARG:H	2:B:847:ARG:HD3	1.60	0.66
2:B:1116:MET:HB2	2:B:1151:LEU:HD21	1.76	0.66
2:A:1152:ARG:NH1	2:A:1161:LEU:HD21	2.11	0.66
2:B:979:GLU:HB3	2:B:1016:HIS:CD2	2.29	0.66
2:A:828:GLY:HA2	3:A:230:HOH:O	1.95	0.66
2:A:1113:TYR:CE1	2:A:1150:THR:HG21	2.30	0.66
2:B:949:GLN:NE2	2:B:949:GLN:N	2.43	0.66
2:B:883:ASN:HD22	2:B:884:GLU:N	1.94	0.65
2:A:1066:ARG:HG3	2:A:1066:ARG:HH11	1.61	0.65
2:A:978:ILE:HD13	2:A:1013:ILE:HG13	1.79	0.65
2:B:860:GLN:HE22	2:B:864:ARG:NH1	1.94	0.65
2:B:854:MET:HE2	2:B:888:ALA:CB	2.27	0.65
2:B:1145:ARG:N	2:B:1146:PRO:HD2	2.12	0.65
2:B:1018:LEU:H	2:B:1021:GLN:HE21	1.45	0.65
2:A:952:GLU:HA	2:A:955:ARG:NH1	2.12	0.64
2:B:950:GLN:HB3	2:B:981:VAL:HG22	1.80	0.64
2:A:910:SER:H	2:A:913:GLN:NE2	1.96	0.63
2:B:919:GLU:HG2	2:B:922:ARG:HH22	1.62	0.63
2:B:896:VAL:HB	2:B:933:TYR:CZ	2.34	0.63
2:A:1086:VAL:HG12	2:A:1131:VAL:HG21	1.80	0.62
2:B:1118:ASP:HB3	2:B:1121:ALA:HB3	1.80	0.62
2:A:921:ILE:HG23	2:A:928:LEU:CD1	2.30	0.61
2:B:907:GLU:HG2	2:B:944:PHE:CE1	2.35	0.61
2:B:978:ILE:HD11	2:B:1009:VAL:HG13	1.81	0.61
2:B:1062:VAL:HG12	2:B:1085:CYS:SG	2.40	0.61
2:B:889:ALA:O	2:B:893:MET:HG3	1.99	0.61
2:A:1033:THR:HG21	2:A:1064:GLU:HG3	1.80	0.61
2:B:1141:MET:O	2:B:1145:ARG:HG3	2.01	0.61
1:D:11:A:H5''	1:D:12:U:O5'	2.00	0.61
2:A:1051:GLU:OE1	2:A:1084:LYS:HE2	1.99	0.61
2:B:1018:LEU:O	2:B:1021:GLN:HB2	2.00	0.61
1:C:11:A:H5''	1:C:12:U:H6	1.66	0.60
1:C:20:A:H2'	2:A:860:GLN:OE1	2.01	0.60
2:B:888:ALA:HB1	2:B:891:GLN:NE2	2.16	0.60
2:A:1086:VAL:CG1	2:A:1131:VAL:HG21	2.31	0.60
2:B:885:ILE:C	2:B:885:ILE:HD12	2.27	0.60
2:B:941:ALA:O	2:B:945:ILE:HG12	2.02	0.60
2:A:1109:HIS:CD2	2:A:1110:SER:H	2.19	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:838:ASN:HB3	2:B:1019:PRO:HG2	1.85	0.59
2:A:1127:LYS:HE3	2:A:1127:LYS:HA	1.85	0.59
2:A:831:ARG:HD2	2:A:831:ARG:C	2.28	0.59
2:B:946:PRO:O	2:B:950:GLN:HG3	2.02	0.59
2:A:1136:GLN:O	2:A:1140:VAL:HG23	2.03	0.58
2:B:838:ASN:O	2:B:840:ARG:N	2.35	0.58
2:B:1074:GLN:HG2	2:B:1118:ASP:CG	2.28	0.58
2:A:1126:GLN:HG3	2:A:1160:ILE:HD13	1.85	0.58
2:B:951:ASN:HA	2:B:954:VAL:HG22	1.84	0.58
2:B:1004:PRO:HB3	2:B:1041:TYR:CZ	2.39	0.58
2:B:999:ALA:O	2:B:1002:THR:HG22	2.03	0.57
2:A:1125:VAL:O	2:A:1129:ILE:HG13	2.04	0.57
2:B:1094:ARG:HD3	2:B:1131:VAL:CG2	2.35	0.57
2:B:829:ARG:HG3	2:B:829:ARG:HH11	1.69	0.57
2:A:854:MET:O	2:A:858:GLN:HG3	2.05	0.57
2:A:882:PHE:HA	2:A:885:ILE:HG13	1.87	0.57
2:B:951:ASN:C	2:B:953:MET:H	2.12	0.57
2:B:967:ASP:O	2:B:971:ASN:HB3	2.04	0.57
2:B:916:ALA:O	2:B:919:GLU:HB3	2.04	0.56
2:A:951:ASN:O	2:A:955:ARG:HG3	2.04	0.56
2:B:1047:GLN:HG2	2:B:1084:LYS:HG3	1.87	0.56
2:B:849:ILE:O	2:B:852:HIS:HB2	2.06	0.56
2:B:1040:GLN:HG2	3:B:103:HOH:O	2.05	0.56
2:B:912:GLU:HG3	3:B:112:HOH:O	2.06	0.56
1:D:17:C:C6	2:B:1008:ARG:NH2	2.74	0.56
2:A:1033:THR:HG21	2:A:1064:GLU:CG	2.35	0.56
2:A:885:ILE:HD12	2:A:885:ILE:C	2.31	0.55
2:B:874:THR:HB	2:B:875:PRO:CD	2.36	0.55
1:C:12:U:O2'	1:C:13:U:P	2.65	0.55
2:A:1113:TYR:HE1	2:A:1150:THR:HG21	1.71	0.55
2:B:971:ASN:HD21	2:B:972:HIS:CE1	2.25	0.55
2:B:1138:LYS:HB3	2:B:1138:LYS:HZ3	1.72	0.55
2:B:984:GLN:H	2:B:984:GLN:CD	2.13	0.55
2:B:1019:PRO:C	2:B:1021:GLN:H	2.13	0.55
2:B:1124:VAL:O	2:B:1128:MET:HG3	2.07	0.55
2:A:882:PHE:HA	2:A:885:ILE:CG1	2.37	0.55
2:A:1033:THR:HG23	2:A:1034:GLU:H	1.72	0.55
1:D:12:U:O2'	1:D:13:U:OP1	2.17	0.55
2:B:1091:ARG:HA	2:B:1091:ARG:HH11	1.72	0.55
2:A:975:GLN:O	2:A:979:GLU:HG3	2.06	0.55
2:B:1018:LEU:N	2:B:1021:GLN:HE21	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:966:LYS:HG2	2:A:1003:HIS:NE2	2.21	0.54
2:A:1138:LYS:HE3	3:A:217:HOH:O	2.06	0.54
2:B:987:GLN:HG3	2:B:991:ASP:OD2	2.07	0.54
2:B:1155:THR:HA	2:B:1158:LYS:HG3	1.88	0.54
2:B:997:VAL:HB	2:B:1028:GLU:HG2	1.89	0.54
2:A:950:GLN:HB3	2:A:981:VAL:HG22	1.88	0.54
2:A:1152:ARG:HH11	2:A:1161:LEU:HD21	1.72	0.54
2:B:1108:PRO:O	2:B:1109:HIS:HB2	2.08	0.54
2:B:1098:ILE:HD13	2:B:1132:ALA:HB2	1.89	0.54
2:A:838:ASN:CG	2:B:1019:PRO:HG3	2.32	0.54
2:A:1109:HIS:CE1	2:A:1147:HIS:NE2	2.76	0.54
2:A:1033:THR:HG23	2:A:1034:GLU:N	2.22	0.54
2:B:984:GLN:H	2:B:984:GLN:NE2	2.05	0.54
2:B:1052:HIS:O	2:B:1053:GLY:O	2.26	0.53
2:B:1162:ALA:O	2:B:1166:LYS:HG2	2.08	0.53
2:B:1075:HIS:O	2:B:1079:SER:HB2	2.07	0.53
2:B:831:ARG:HH11	2:B:832:LEU:H	1.55	0.53
2:B:981:VAL:CG1	2:B:982:GLN:N	2.71	0.53
2:A:1050:LEU:O	2:A:1058:LYS:NZ	2.38	0.53
2:A:854:MET:SD	2:A:888:ALA:HB3	2.49	0.53
2:B:907:GLU:CG	2:B:940:LYS:HD2	2.38	0.53
2:B:1013:ILE:HG23	2:B:1017:CYS:SG	2.49	0.53
2:A:906:PHE:O	2:A:914:LYS:HD3	2.09	0.53
2:B:975:GLN:HG3	2:B:1012:ARG:CG	2.30	0.52
2:A:854:MET:CE	2:A:888:ALA:HB3	2.39	0.52
2:B:935:CYS:O	2:B:939:GLN:HG3	2.09	0.52
2:A:941:ALA:O	2:A:945:ILE:HG12	2.10	0.52
2:B:1139:ILE:O	2:B:1142:HIS:HB3	2.09	0.52
2:B:1122:ASN:HA	2:B:1160:ILE:HD11	1.92	0.52
2:A:910:SER:H	2:A:913:GLN:HE21	1.58	0.52
2:A:880:LEU:C	2:A:880:LEU:HD23	2.36	0.51
2:A:966:LYS:HG2	2:A:1003:HIS:CD2	2.45	0.51
2:B:874:THR:HB	2:B:875:PRO:HD2	1.91	0.51
2:B:882:PHE:CE1	2:B:886:LEU:HD22	2.46	0.51
2:B:972:HIS:HE1	3:B:41:HOH:O	1.93	0.51
2:A:879:GLN:HE21	2:A:883:ASN:ND2	2.08	0.51
2:B:1070:LEU:HD13	2:B:1111:ALA:HB1	1.93	0.51
2:A:1113:TYR:HA	2:A:1116:MET:CE	2.38	0.51
2:A:1116:MET:HB3	2:A:1151:LEU:HD21	1.93	0.51
1:C:11:A:H5'	1:C:14:G:OP2	2.11	0.50
2:B:1076:LYS:HD3	2:B:1077:PHE:CE2	2.47	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:1076:LYS:HA	2:B:1120:TYR:CD1	2.47	0.50
2:B:950:GLN:HB3	2:B:981:VAL:CG2	2.40	0.50
2:B:1141:MET:O	2:B:1145:ARG:NH1	2.44	0.50
2:A:854:MET:HE3	2:A:892:LEU:HD21	1.93	0.50
2:B:961:VAL:O	2:B:965:VAL:HG23	2.11	0.50
2:B:886:LEU:HA	2:B:889:ALA:HB2	1.94	0.50
2:B:1144:ILE:HD11	2:B:1151:LEU:HD11	1.94	0.50
1:C:15:U:H1'	2:A:1077:PHE:HE2	1.77	0.49
2:A:948:ASP:O	2:A:952:GLU:HG3	2.12	0.49
2:B:829:ARG:HG3	2:B:829:ARG:NH1	2.27	0.49
2:B:954:VAL:HG12	2:B:977:CYS:SG	2.52	0.49
2:B:1048:HIS:CE1	2:B:1052:HIS:HD2	2.30	0.49
2:B:1084:LYS:HA	2:B:1087:THR:CG2	2.41	0.49
2:A:862:GLY:O	2:A:865:PHE:HB3	2.12	0.49
2:A:951:ASN:HA	2:A:954:VAL:HG12	1.93	0.49
2:B:883:ASN:C	2:B:883:ASN:ND2	2.70	0.49
2:B:1005:TYR:O	2:B:1008:ARG:HB2	2.13	0.49
2:B:978:ILE:HG21	2:B:1012:ARG:C	2.38	0.49
1:C:11:A:P	1:C:14:G:N7	2.86	0.49
2:B:1117:LYS:HD2	2:B:1154:TYR:CZ	2.48	0.49
2:B:873:ALA:HA	2:B:877:GLU:OE1	2.13	0.49
2:B:882:PHE:HE1	2:B:886:LEU:HD22	1.77	0.49
2:B:846:LEU:HD12	2:B:846:LEU:HA	1.70	0.48
2:A:1137:ARG:CB	2:A:1137:ARG:HH11	2.25	0.48
2:B:907:GLU:HG2	2:B:944:PHE:CZ	2.47	0.48
2:A:854:MET:CE	2:A:892:LEU:HG	2.43	0.48
2:B:1121:ALA:O	2:B:1125:VAL:HG23	2.13	0.48
2:B:906:PHE:O	2:B:914:LYS:HD3	2.13	0.48
2:B:982:GLN:HB3	2:B:984:GLN:HE21	1.78	0.48
2:B:1138:LYS:HB3	2:B:1138:LYS:HZ2	1.77	0.48
2:B:1148:ILE:HG23	2:B:1149:ALA:N	2.28	0.48
1:C:11:A:C5'	1:C:14:G:OP2	2.62	0.48
2:B:1122:ASN:CA	2:B:1160:ILE:HD11	2.44	0.48
2:B:860:GLN:HE22	2:B:864:ARG:HH12	1.62	0.47
2:B:879:GLN:O	2:B:880:LEU:C	2.56	0.47
2:B:1119:GLN:HG3	2:B:1120:TYR:CD2	2.49	0.47
2:B:1028:GLU:O	2:B:1032:HIS:HD2	1.97	0.47
2:B:978:ILE:HD13	2:B:1013:ILE:HG13	1.95	0.47
2:A:1132:ALA:O	2:A:1137:ARG:NH1	2.47	0.47
2:B:1050:LEU:O	2:B:1058:LYS:HE2	2.14	0.47
2:A:838:ASN:HB3	2:B:1019:PRO:CG	2.44	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:1059:SER:HA	2:B:1062:VAL:HG22	1.97	0.47
2:B:1067:GLY:N	2:B:1100:GLU:OE2	2.39	0.47
2:A:928:LEU:HD13	2:A:938:ILE:HD11	1.97	0.46
2:A:949:GLN:H	2:A:949:GLN:NE2	2.14	0.46
2:A:1066:ARG:HG3	2:A:1066:ARG:NH1	2.29	0.46
2:B:1086:VAL:HG11	2:B:1128:MET:HG2	1.97	0.46
2:A:1153:LYS:NZ	2:A:1153:LYS:HB3	2.31	0.46
2:B:972:HIS:CD2	3:B:56:HOH:O	2.68	0.46
2:B:1087:THR:HB	2:B:1127:LYS:HE2	1.97	0.46
2:A:983:PRO:HG2	2:A:984:GLN:NE2	2.31	0.46
2:A:946:PRO:HD2	2:A:949:GLN:NE2	2.31	0.46
2:B:1003:HIS:CG	2:B:1004:PRO:HD2	2.51	0.46
2:B:1145:ARG:N	2:B:1146:PRO:CD	2.79	0.46
2:A:879:GLN:HE21	2:A:883:ASN:HD21	1.63	0.46
2:B:1023:LEU:HB3	2:B:1024:PRO:CD	2.46	0.46
1:C:20:A:H1'	2:A:897:PHE:CE2	2.45	0.45
2:B:1013:ILE:C	2:B:1015:GLU:H	2.24	0.45
2:B:971:ASN:ND2	2:B:972:HIS:CE1	2.84	0.45
2:A:863:SER:O	2:A:867:GLN:HG3	2.16	0.45
2:A:1029:LEU:HD23	2:A:1029:LEU:HA	1.79	0.45
2:B:907:GLU:HG3	2:B:940:LYS:HD2	1.98	0.45
2:A:854:MET:HE1	2:A:892:LEU:HG	1.99	0.45
2:A:887:GLN:HA	2:A:887:GLN:NE2	2.32	0.45
2:A:981:VAL:HG12	2:A:982:GLN:N	2.32	0.45
2:B:838:ASN:O	2:B:840:ARG:HG3	2.17	0.45
2:B:1026:LEU:HD13	2:B:1057:ASP:OD1	2.17	0.45
2:A:1075:HIS:HB3	2:A:1078:ALA:HB3	1.99	0.44
2:A:1061:ILE:O	2:A:1065:ILE:HD13	2.17	0.44
2:B:1116:MET:SD	2:B:1151:LEU:HG	2.58	0.44
2:A:1027:GLU:HA	2:A:1027:GLU:OE1	2.17	0.44
2:A:1098:ILE:HD13	2:A:1132:ALA:HB2	2.00	0.44
2:A:1098:ILE:HG12	2:A:1128:MET:HG2	1.99	0.44
2:B:1003:HIS:HB3	2:B:1006:GLY:HA3	1.98	0.44
2:B:1089:ALA:HB3	2:B:1094:ARG:CB	2.39	0.44
2:B:1089:ALA:CB	2:B:1094:ARG:HB2	2.40	0.44
2:B:880:LEU:C	2:B:880:LEU:HD12	2.42	0.44
2:B:1144:ILE:C	2:B:1146:PRO:HD2	2.43	0.44
2:B:978:ILE:CD1	2:B:1012:ARG:HB2	2.48	0.44
2:B:1094:ARG:O	2:B:1094:ARG:HG2	2.18	0.44
2:B:1155:THR:HA	2:B:1158:LYS:HE3	2.00	0.44
2:A:978:ILE:HD13	2:A:1013:ILE:CG1	2.45	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:832:LEU:HD21	2:B:849:ILE:HG22	2.00	0.43
2:B:981:VAL:HG13	2:B:982:GLN:N	2.34	0.43
2:B:1164:LEU:O	2:B:1167:TYR:HB3	2.18	0.43
2:B:906:PHE:CE1	2:B:941:ALA:HB2	2.52	0.43
2:B:935:CYS:SG	2:B:969:ASN:HB3	2.58	0.43
2:B:1086:VAL:HG12	2:B:1131:VAL:HG21	2.00	0.43
2:B:1104:MET:O	2:B:1104:MET:HG2	2.18	0.43
2:A:1137:ARG:NH1	2:A:1137:ARG:HB2	2.32	0.43
2:B:957:LEU:HB3	2:B:988:PHE:HE1	1.83	0.43
2:B:1109:HIS:ND1	2:B:1110:SER:HB2	2.32	0.43
2:B:972:HIS:HD2	3:B:56:HOH:O	2.00	0.43
1:C:18:A:H2	2:A:935:CYS:SG	2.42	0.43
2:A:1126:GLN:HG3	2:A:1160:ILE:CD1	2.49	0.43
2:A:1070:LEU:HD21	2:A:1106:ASP:HB2	2.01	0.43
2:A:870:LEU:HD23	2:A:878:ARG:HG2	2.00	0.43
1:C:14:G:C6	2:A:1123:TYR:CD2	3.06	0.43
2:A:860:GLN:HB2	2:A:897:PHE:CZ	2.54	0.43
2:B:951:ASN:C	2:B:953:MET:N	2.73	0.43
2:B:951:ASN:HB3	2:B:955:ARG:NH2	2.20	0.43
2:B:882:PHE:O	2:B:885:ILE:HG13	2.19	0.43
2:B:1043:ASN:OD1	2:B:1047:GLN:NE2	2.52	0.43
2:A:971:ASN:O	2:A:975:GLN:HG3	2.19	0.42
2:A:1098:ILE:O	2:A:1101:VAL:HG22	2.18	0.42
2:B:1021:GLN:C	2:B:1023:LEU:H	2.27	0.42
2:A:1092:THR:O	2:A:1096:VAL:HG23	2.18	0.42
2:B:1021:GLN:O	2:B:1024:PRO:HD2	2.18	0.42
2:B:1111:ALA:O	2:B:1115:MET:HG3	2.19	0.42
1:D:20:A:O2'	2:B:860:GLN:NE2	2.52	0.42
2:A:1109:HIS:CD2	2:A:1110:SER:N	2.86	0.42
2:A:1161:LEU:O	2:A:1165:GLU:HB2	2.20	0.42
2:B:978:ILE:HD12	2:B:1012:ARG:HB2	2.01	0.42
2:B:1055:PRO:HA	2:B:1058:LYS:HB2	2.00	0.42
2:B:1087:THR:HG23	2:B:1088:HIS:CD2	2.54	0.42
2:B:1110:SER:O	2:B:1113:TYR:HB3	2.19	0.42
1:D:18:A:C5	2:B:972:HIS:CD2	3.08	0.42
2:A:924:HIS:O	2:A:928:LEU:HD12	2.19	0.42
2:A:978:ILE:HA	2:A:986:LEU:HD21	2.02	0.42
2:B:1061:ILE:O	2:B:1065:ILE:HG12	2.20	0.42
2:A:924:HIS:HA	3:A:5:HOH:O	2.20	0.42
2:B:1010:ILE:HG22	2:B:1011:GLN:N	2.35	0.42
2:B:1119:GLN:HA	2:B:1156:TYR:CG	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:981:VAL:CG1	2:A:982:GLN:N	2.83	0.41
2:B:1050:LEU:HD21	2:B:1061:ILE:HB	2.03	0.41
2:A:1027:GLU:O	2:A:1031:GLN:HG3	2.20	0.41
2:B:1050:LEU:HD21	2:B:1061:ILE:HG21	2.01	0.41
2:B:945:ILE:HD12	2:B:949:GLN:HB3	2.02	0.41
2:B:1145:ARG:HH22	2:B:1167:TYR:HD2	1.69	0.41
2:A:1086:VAL:O	2:A:1094:ARG:HD3	2.20	0.41
2:B:886:LEU:O	2:B:920:ARG:NH1	2.54	0.41
2:A:1062:VAL:HG23	2:A:1097:LEU:HD11	2.03	0.41
2:B:971:ASN:O	2:B:975:GLN:HB2	2.20	0.41
2:B:1117:LYS:HD2	2:B:1154:TYR:CE2	2.55	0.41
2:A:921:ILE:O	2:A:922:ARG:C	2.63	0.41
2:A:1129:ILE:O	2:A:1129:ILE:HG22	2.20	0.41
2:A:1032:HIS:O	2:A:1033:THR:C	2.62	0.41
2:A:1112:LEU:O	2:A:1116:MET:HG3	2.21	0.41
2:B:951:ASN:N	2:B:951:ASN:HD22	2.17	0.41
2:B:958:ASP:HA	2:B:988:PHE:HD1	1.86	0.41
2:B:982:GLN:HG2	3:B:195:HOH:O	2.21	0.41
2:B:986:LEU:HA	2:B:988:PHE:CE2	2.56	0.41
2:B:1007:CYS:O	2:B:1011:GLN:HG3	2.21	0.41
2:B:1141:MET:SD	2:B:1164:LEU:HD23	2.60	0.41
2:A:882:PHE:HA	2:A:885:ILE:HD11	2.02	0.41
2:B:1019:PRO:C	2:B:1021:GLN:N	2.77	0.40
2:A:829:ARG:HD3	2:A:833:LEU:CD2	2.45	0.40
2:A:1126:GLN:CG	2:A:1160:ILE:HD13	2.50	0.40
2:B:988:PHE:H	2:B:988:PHE:HD2	1.69	0.40
2:B:1050:LEU:HD21	2:B:1061:ILE:CG2	2.51	0.40
2:A:835:ASP:HB2	2:A:841:TYR:HE2	1.86	0.40
2:B:1089:ALA:O	2:B:1090:SER:O	2.39	0.40
2:A:896:VAL:HG13	2:A:897:PHE:CD1	2.56	0.40
2:A:1059:SER:HA	2:A:1062:VAL:HG12	2.03	0.40
2:B:845:GLN:HE22	2:B:869:LYS:NZ	2.19	0.40
2:B:1076:LYS:HA	2:B:1120:TYR:CE1	2.57	0.40
2:B:876:ALA:O	2:B:877:GLU:C	2.64	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	
2	A	339/349 (97%)	325 (96%)	10 (3%)	4 (1%)	10	23
2	B	340/349 (97%)	306 (90%)	28 (8%)	6 (2%)	6	14
All	All	679/698 (97%)	631 (93%)	38 (6%)	10 (2%)	8	18

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	1110	SER
2	A	1166	LYS
2	B	839	ASN
2	B	1090	SER
2	B	1105	ASN
2	B	1109	HIS
2	B	909	GLY
2	B	1053	GLY
2	A	1104	MET
2	A	961	VAL

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	
2	A	303/310 (98%)	290 (96%)	13 (4%)	26	51
2	B	303/310 (98%)	285 (94%)	18 (6%)	18	38
All	All	606/620 (98%)	575 (95%)	31 (5%)	21	45

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	902	ILE
2	A	904	LYS
2	A	935	CYS
2	A	940	LYS
2	A	957	LEU
2	A	962	LEU
2	A	985	SER
2	A	986	LEU
2	A	987	GLN
2	A	1019	PRO
2	A	1082	VAL
2	A	1153	LYS
2	A	1165	GLU
2	B	831	ARG
2	B	854	MET
2	B	870	LEU
2	B	880	LEU
2	B	883	ASN
2	B	896	VAL
2	B	949	GLN
2	B	955	ARG
2	B	971	ASN
2	B	981	VAL
2	B	984	GLN
2	B	988	PHE
2	B	996	GLN
2	B	1002	THR
2	B	1010	ILE
2	B	1018	LEU
2	B	1105	ASN
2	B	1138	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (34) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	858	GLN
2	A	879	GLN
2	A	883	ASN
2	A	887	GLN
2	A	891	GLN
2	A	913	GLN

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Mol	Chain	Res	Type
2	A	949	GLN
2	A	968	GLN
2	A	969	ASN
2	A	984	GLN
2	A	1031	GLN
2	A	1032	HIS
2	A	1074	GLN
2	A	1109	HIS
2	A	1126	GLN
2	A	1159	HIS
2	B	843	ASN
2	B	858	GLN
2	B	860	GLN
2	B	861	HIS
2	B	879	GLN
2	B	883	ASN
2	B	891	GLN
2	B	949	GLN
2	B	951	ASN
2	B	972	HIS
2	B	984	GLN
2	B	1021	GLN
2	B	1032	HIS
2	B	1038	GLN
2	B	1040	GLN
2	B	1052	HIS
2	B	1088	HIS
2	B	1136	GLN

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	C	9/10 (90%)	2 (22%)	1 (11%)
1	D	9/10 (90%)	2 (22%)	1 (11%)
All	All	18/20 (90%)	4 (22%)	2 (11%)

All (4) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	C	12	U
1	C	13	U

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Mol	Chain	Res	Type
1	D	12	U
1	D	13	U

All (2) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	C	12	U
1	D	12	U

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 ⓘ

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	C	10/10 (100%)	-0.18	0 100 100	15, 34, 53, 57	0
1	D	10/10 (100%)	0.16	1 (10%) 12 9	29, 45, 87, 98	0
2	A	341/349 (97%)	-0.41	5 (1%) 72 68	6, 20, 56, 94	0
2	B	342/349 (97%)	0.26	11 (3%) 50 44	14, 38, 73, 90	0
All	All	703/718 (97%)	-0.07	17 (2%) 59 54	6, 31, 70, 98	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	1168	TYR	3.4
2	A	1106	ASP	3.4
2	A	1168	TYR	3.1
2	B	1108	PRO	2.9
2	A	1167	TYR	2.9
2	B	1167	TYR	2.9
2	B	1091	ARG	2.8
2	B	1104	MET	2.8
2	A	1104	MET	2.6
2	B	1107	GLY	2.6
2	B	1106	ASP	2.4
2	B	1031	GLN	2.4
2	B	1169	MET	2.3
1	D	11	A	2.3
2	B	1109	HIS	2.2
2	A	1108	PRO	2.0
2	B	888	ALA	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 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.