



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 04:19 AM UTC

PDB ID : 2ESN / pdb_00002esn
Title : The crystal structure of probable transcriptional regulator PA0477 from *Pseudomonas aeruginosa*
Authors : Lunin, V.V.; Chang, C.; Skarina, T.; Gorodischenskaya, E.; Edwards, A.M.; Joachimiak, A.; Savchenko, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on : 2005-10-26
Resolution : 2.10 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

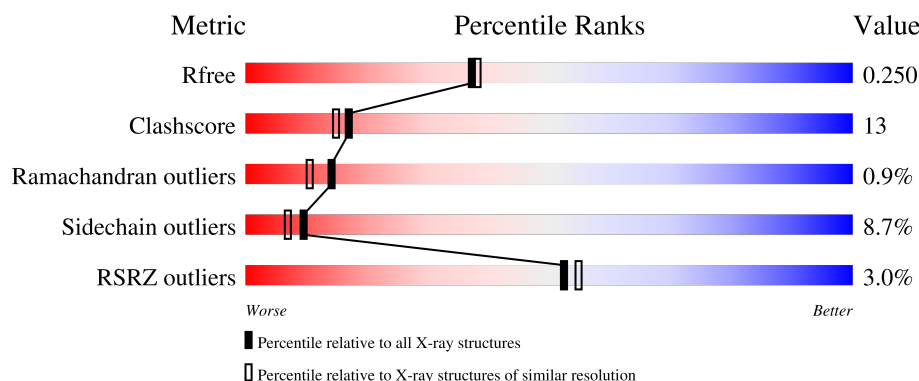
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	310	2% 72% 21% . .
1	B	310	% 71% 21% . . .
1	C	310	2% 71% 21% . .
1	D	310	6% 68% 23% 5% .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10202 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 probable transcriptional regulator.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	299	Total	C	N	O	S	0	2	0
			2360	1498	445	414	3			
1	C	298	Total	C	N	O	S	0	3	0
			2363	1502	449	409	3			
1	B	300	Total	C	N	O	S	0	0	0
			2347	1493	437	414	3			
1	D	298	Total	C	N	O	S	0	3	0
			2362	1500	444	415	3			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	cloning artifact	GB 9946339
A	0	HIS	-	cloning artifact	GB 9946339
B	-1	GLY	-	cloning artifact	GB 9946339
B	0	HIS	-	cloning artifact	GB 9946339
C	-1	GLY	-	cloning artifact	GB 9946339
C	0	HIS	-	cloning artifact	GB 9946339
D	-1	GLY	-	cloning artifact	GB 9946339
D	0	HIS	-	cloning artifact	GB 9946339

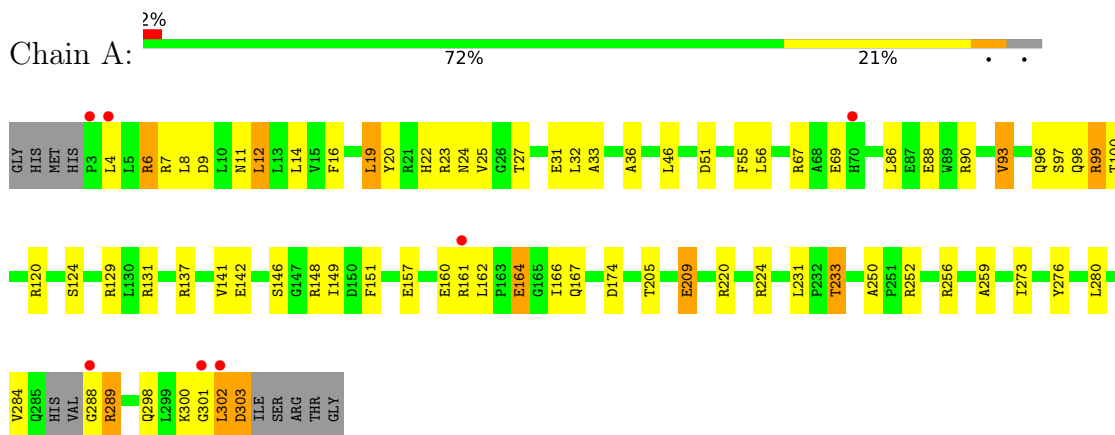
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	224	Total	O	0	0
			224	224		
2	C	204	Total	O	0	0
			204	204		
2	B	201	Total	O	0	0
			201	201		
2	D	141	Total	O	0	0
			141	141		

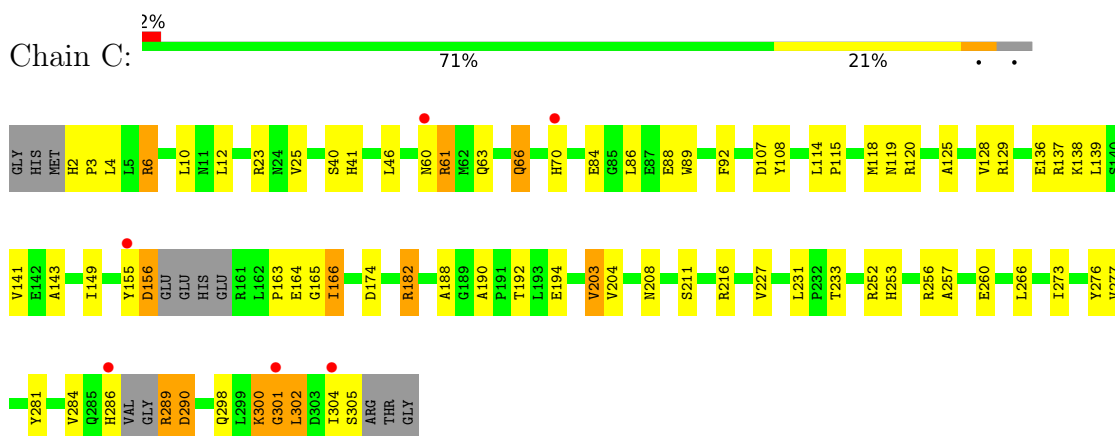
3 Residue-property plots [i](#)

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.

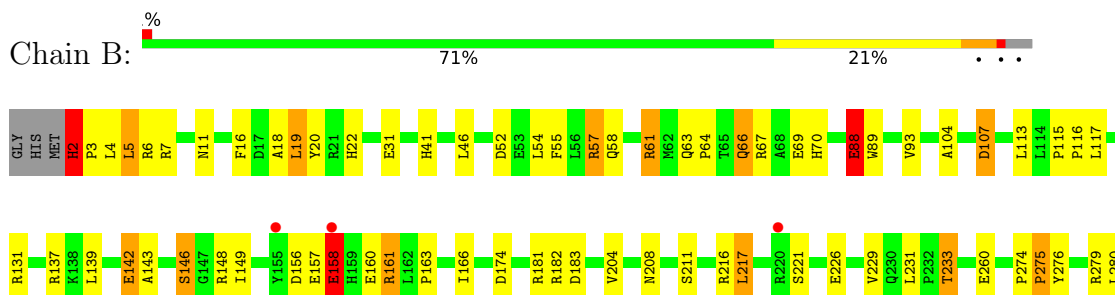
- Molecule 1: probable transcriptional regulator

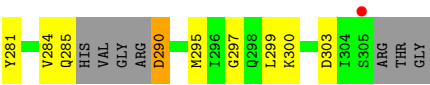


- Molecule 1: probable transcriptional regulator

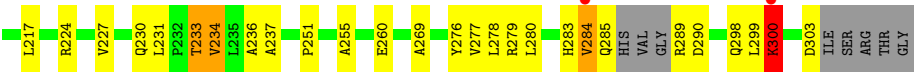
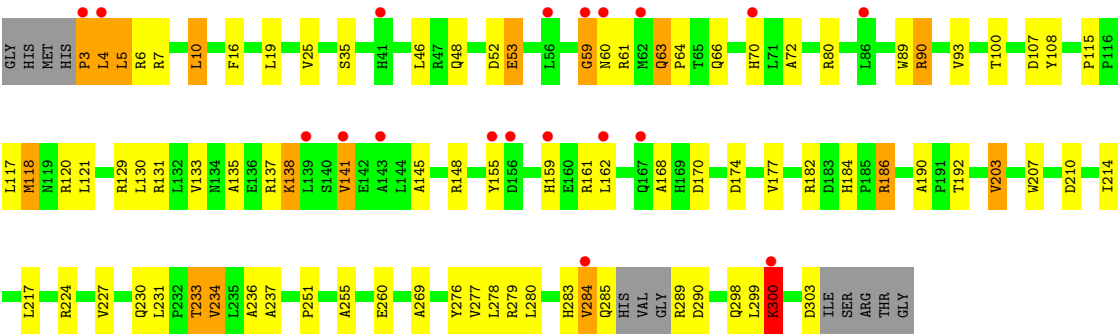


- Molecule 1: probable transcriptional regulator





● Molecule 1: probable transcriptional regulator



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	97.03Å 171.21Å 86.78Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.10 50.00 – 2.10	Depositor EDS
% Data completeness (in resolution range)	97.1 (50.00-2.10) 97.1 (50.00-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.81 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.194 , 0.251 0.193 , 0.250	Depositor DCC
R_{free} test set	2466 reflections (2.91%)	wwPDB-VP
Wilson B-factor (Å ²)	38.6	Xtriage
Anisotropy	0.352	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 58.0	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	0.96	EDS
Total number of atoms	10202	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 12.19% 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	1.28	5/2414 (0.2%)	1.23	10/3280 (0.3%)
1	B	1.25	2/2402 (0.1%)	1.25	15/3268 (0.5%)
1	C	1.35	6/2418 (0.2%)	1.21	9/3286 (0.3%)
1	D	1.17	6/2417 (0.2%)	1.21	9/3286 (0.3%)
All	All	1.26	19/9651 (0.2%)	1.23	43/13120 (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	B	0	3
1	C	0	1
1	D	0	2
All	All	0	6

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	118	MET	SD-CE	-7.78	1.60	1.79
1	C	125	ALA	C-O	-7.27	1.15	1.24
1	A	250	ALA	CA-CB	6.43	1.61	1.54
1	C	273	ILE	CA-CB	6.43	1.59	1.54
1	C	273	ILE	CA-C	-6.25	1.47	1.52
1	D	115	PRO	CA-C	5.86	1.57	1.52
1	A	273	ILE	C-O	-5.79	1.18	1.23
1	A	233	THR	CA-CB	5.47	1.61	1.53
1	A	93	VAL	CA-CB	5.39	1.59	1.54
1	D	93	VAL	CA-CB	5.24	1.59	1.54
1	D	269	ALA	CA-CB	-5.20	1.44	1.52
1	A	259	ALA	N-CA	5.19	1.52	1.46
1	B	18	ALA	CA-CB	-5.15	1.45	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	118	MET	SD-CE	-5.13	1.66	1.79
1	B	295	MET	SD-CE	-5.13	1.66	1.79
1	C	128	VAL	C-O	5.12	1.30	1.24
1	D	184	HIS	CA-C	5.06	1.58	1.53
1	C	257	ALA	CA-CB	5.04	1.61	1.53
1	D	255	ALA	CA-CB	5.01	1.61	1.53

All (43) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	53	GLU	N-CA-C	-9.82	101.65	114.31
1	B	233	THR	CB-CA-C	-8.83	95.71	110.19
1	B	2	HIS	CA-C-N	7.46	128.45	120.47
1	B	2	HIS	C-N-CA	7.46	128.45	120.47
1	B	58	GLN	CB-CA-C	-6.93	101.23	111.77
1	C	115	PRO	CA-C-N	-6.72	111.44	119.84
1	C	115	PRO	C-N-CA	-6.72	111.44	119.84
1	C	192	THR	N-CA-C	-6.43	101.87	110.55
1	C	114	LEU	CA-C-N	-6.33	113.86	120.38
1	C	114	LEU	C-N-CA	-6.33	113.86	120.38
1	D	234	VAL	CB-CA-C	-6.29	103.64	112.14
1	A	162	LEU	CA-C-N	-5.96	114.13	120.03
1	A	162	LEU	C-N-CA	-5.96	114.13	120.03
1	D	269	ALA	CA-C-N	-5.79	114.92	120.83
1	D	269	ALA	C-N-CA	-5.79	114.92	120.83
1	D	161	ARG	N-CA-C	5.79	118.36	111.71
1	C	107	ASP	N-CA-C	5.66	117.45	111.28
1	B	181	ARG	NE-CZ-NH1	-5.64	115.86	121.50
1	A	36	ALA	CA-C-N	5.57	128.00	120.65
1	A	36	ALA	C-N-CA	5.57	128.00	120.65
1	A	209	GLU	CA-CB-CG	5.56	125.21	114.10
1	D	192	THR	N-CA-C	-5.51	102.75	110.35
1	A	205	THR	CA-C-N	-5.48	114.03	119.56
1	A	205	THR	C-N-CA	-5.48	114.03	119.56
1	D	186	ARG	N-CA-C	5.45	119.71	112.30
1	A	99	ARG	N-CA-C	5.41	117.67	110.53
1	C	190	ALA	CA-C-N	-5.39	114.15	119.92
1	C	190	ALA	C-N-CA	-5.39	114.15	119.92
1	B	52	ASP	CA-C-N	5.36	127.46	120.28
1	B	52	ASP	C-N-CA	5.36	127.46	120.28
1	A	25	VAL	N-CA-C	5.30	115.51	110.53
1	C	188	ALA	N-CA-C	-5.21	105.21	112.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	181	ARG	NE-CZ-NH2	5.21	123.89	119.20
1	B	5	LEU	N-CA-C	-5.17	105.71	112.23
1	B	63	GLN	CA-C-N	5.17	125.33	119.90
1	B	63	GLN	C-N-CA	5.17	125.33	119.90
1	B	221	SER	N-CA-C	-5.13	105.38	113.02
1	A	124	SER	N-CA-C	5.13	117.77	111.82
1	B	104	ALA	N-CA-C	-5.11	100.91	109.24
1	D	25	VAL	N-CA-C	5.10	115.31	110.42
1	D	60	ASN	N-CA-C	-5.07	105.81	113.51
1	B	275	PRO	N-CA-C	-5.06	103.17	111.21
1	B	158	GLU	N-CA-C	5.02	121.49	110.80

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	158	GLU	Peptide
1	B	2	HIS	Peptide
1	B	284	VAL	Peptide
1	C	60	ASN	Peptide
1	D	3	PRO	Peptide
1	D	59	GLY	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	2360	0	2364	50	1
1	B	2347	0	2346	69	1
1	C	2363	0	2370	68	1
1	D	2362	0	2361	71	0
2	A	224	0	0	13	0
2	B	201	0	0	17	0
2	C	204	0	0	12	0
2	D	141	0	0	13	0
All	All	10202	0	9441	239	2

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 13.

All (239) 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:D:3:PRO:N	1:D:4:LEU:HB2	1.63	1.12
1:C:23:ARG:HH21	1:C:61:ARG:HH22	1.04	1.01
1:B:174:ASP:HB3	1:B:276:TYR:CZ	2.08	0.89
1:C:182:ARG:HH22	1:C:260:GLU:CD	1.81	0.87
1:B:285:GLN:HA	2:B:407:HOH:O	1.75	0.86
1:C:120[A]:ARG:NH2	1:C:298:GLN:O	2.08	0.85
1:D:117:LEU:HG	1:D:118:MET:HE2	1.58	0.84
1:D:70:HIS:NE2	2:D:440:HOH:O	2.10	0.83
1:C:25:VAL:HB	2:C:507:HOH:O	1.78	0.83
1:B:158:GLU:O	1:B:158:GLU:HG3	1.76	0.83
1:C:300:LYS:C	1:C:302:LEU:H	1.87	0.82
1:D:138:LYS:H	1:D:138:LYS:HD2	1.43	0.82
1:C:182:ARG:NH2	1:C:260:GLU:OE2	2.13	0.81
1:B:66:GLN:CD	1:B:66:GLN:H	1.86	0.81
1:B:107:ASP:HB2	1:D:233[B]:THR:HG21	1.64	0.80
1:A:67:ARG:HH12	1:D:90:ARG:HG3	1.46	0.79
1:C:86:LEU:HD12	1:B:5:LEU:HD23	1.64	0.79
1:C:141:VAL:HG22	1:C:166:ILE:HD11	1.63	0.79
1:A:51:ASP:OD2	1:D:90:ARG:NH2	2.17	0.78
1:B:67:ARG:HG2	2:B:395:HOH:O	1.83	0.77
1:C:23:ARG:NH2	1:C:61:ARG:HH22	1.82	0.77
1:B:2:HIS:N	1:B:3:PRO:HA	2.00	0.77
1:D:138:LYS:HD2	1:D:138:LYS:N	1.97	0.76
1:D:138:LYS:H	1:D:138:LYS:CD	1.99	0.76
1:A:9:ASP:HB3	1:A:12:LEU:HD22	1.67	0.75
1:A:302:LEU:HB3	2:A:521:HOH:O	1.85	0.75
1:B:290:ASP:N	2:B:504:HOH:O	2.20	0.75
1:C:300:LYS:O	1:C:302:LEU:N	2.21	0.74
1:C:10:LEU:H	1:B:11:ASN:HD21	1.35	0.74
1:C:86:LEU:CD1	1:B:5:LEU:HD23	2.18	0.74
1:C:41:HIS:HB2	2:C:431:HOH:O	1.86	0.73
1:A:20:TYR:O	1:A:23:ARG:HD3	1.88	0.73
1:C:66:GLN:OE1	1:C:70[B]:HIS:CE1	2.42	0.72
1:D:233[B]:THR:CG2	2:D:447:HOH:O	2.38	0.72
1:D:117:LEU:HD23	1:D:118:MET:HE3	1.71	0.72
1:C:165:GLY:HA2	1:C:284:VAL:HG22	1.72	0.72
1:B:157:GLU:O	1:B:160:GLU:HG3	1.89	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:66:GLN:OE1	1:C:70[B]:HIS:HE1	1.72	0.71
1:A:141:VAL:HG13	1:A:166:ILE:HD11	1.71	0.71
1:D:70:HIS:CE1	2:D:440:HOH:O	2.44	0.70
1:C:174:ASP:HB3	1:C:276:TYR:CZ	2.26	0.70
1:C:141:VAL:HG22	1:C:166:ILE:CD1	2.21	0.70
1:D:170:ASP:OD1	1:D:279:ARG:HD3	1.92	0.70
1:A:256:ARG:NH1	2:A:520:HOH:O	2.25	0.69
1:D:117:LEU:HD23	1:D:118:MET:CE	2.22	0.69
1:A:24:ASN:OD1	1:A:27:THR:HG22	1.93	0.69
2:A:443:HOH:O	1:D:66:GLN:HB3	1.91	0.69
1:A:67:ARG:NH1	1:D:90:ARG:HG3	2.07	0.69
1:C:86:LEU:HD12	1:B:5:LEU:CD2	2.24	0.68
1:B:66:GLN:HG2	2:B:395:HOH:O	1.95	0.66
1:A:174:ASP:HB3	1:A:276:TYR:CZ	2.30	0.66
1:D:233[B]:THR:HG21	2:D:447:HOH:O	1.96	0.66
1:C:119:ASN:ND2	1:C:120[B]:ARG:HH21	1.95	0.65
1:D:120:ARG:HD3	1:D:298:GLN:HE22	1.60	0.65
1:B:260:GLU:HG2	2:B:505:HOH:O	1.96	0.64
1:D:80:ARG:HD3	2:D:444:HOH:O	1.98	0.64
1:D:117:LEU:CD2	1:D:118:MET:CE	2.75	0.64
1:A:4:LEU:HB2	2:A:428:HOH:O	1.98	0.63
1:B:139:LEU:HD21	1:B:166:ILE:HG13	1.78	0.63
1:B:161:ARG:HH11	1:B:161:ARG:HG2	1.62	0.63
1:C:286:HIS:HA	1:C:289:ARG:HD3	1.80	0.62
1:A:11:ASN:HD21	1:D:10:LEU:H	1.47	0.62
1:D:117:LEU:CG	1:D:118:MET:HE2	2.30	0.62
1:D:59:GLY:HA3	1:D:61:ARG:H	1.65	0.61
1:D:155:TYR:HB3	1:D:277:VAL:O	2.00	0.61
1:B:57:ARG:NH1	1:B:57:ARG:HG2	2.16	0.61
1:C:23:ARG:HH21	1:C:61:ARG:NH2	1.87	0.61
1:C:3:PRO:HB2	2:C:482:HOH:O	2.00	0.61
1:A:302:LEU:CD2	1:A:303:ASP:H	2.15	0.60
1:D:117:LEU:HG	1:D:118:MET:CE	2.31	0.60
1:D:35:SER:HB2	2:D:389:HOH:O	2.01	0.60
1:B:161:ARG:HH11	1:B:161:ARG:CG	2.14	0.60
1:D:108:TYR:HB2	1:D:234:VAL:HG21	1.84	0.60
1:C:300:LYS:C	1:C:302:LEU:N	2.55	0.59
1:D:3:PRO:N	1:D:4:LEU:CB	2.53	0.59
1:D:284:VAL:O	1:D:285:GLN:HB2	2.01	0.59
1:B:66:GLN:CD	1:B:66:GLN:N	2.60	0.58
1:B:70:HIS:HB2	2:B:414:HOH:O	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:2:HIS:CD2	1:C:3:PRO:HD2	2.39	0.57
1:D:190:ALA:HB1	2:D:443:HOH:O	2.03	0.57
1:A:148:ARG:HD2	2:A:500:HOH:O	2.02	0.57
1:C:252:ARG:NH2	1:C:256:ARG:HD2	2.19	0.57
1:B:57:ARG:CG	1:B:57:ARG:HH11	2.17	0.57
1:B:217:LEU:HD23	1:B:217:LEU:C	2.29	0.57
1:B:2:HIS:N	1:B:3:PRO:CA	2.68	0.57
1:C:12:LEU:HB3	1:C:46:LEU:HD13	1.87	0.57
1:B:88:GLU:HB3	1:B:89:TRP:CE3	2.40	0.57
1:A:19:LEU:HD13	1:A:55:PHE:CD1	2.40	0.57
1:D:16:PHE:HB2	1:D:46:LEU:HD21	1.87	0.57
1:A:22:HIS:NE2	1:A:31:GLU:OE2	2.24	0.56
1:C:84:GLU:HG2	2:C:492:HOH:O	2.05	0.56
1:D:174:ASP:HB3	1:D:276:TYR:CZ	2.41	0.56
1:A:157:GLU:O	1:A:160:GLU:HG3	2.06	0.56
1:D:120:ARG:HD3	1:D:298:GLN:NE2	2.20	0.56
1:D:155:TYR:CB	1:D:277:VAL:O	2.54	0.56
1:A:93:VAL:HG21	2:A:529:HOH:O	2.06	0.55
1:C:10:LEU:H	1:B:11:ASN:ND2	2.01	0.55
1:B:57:ARG:HG2	1:B:57:ARG:HH11	1.72	0.55
1:A:146:SER:HB2	1:A:148:ARG:H	1.72	0.55
1:B:19:LEU:HD13	1:B:55:PHE:CD1	2.42	0.54
1:A:302:LEU:HD23	1:A:303:ASP:H	1.72	0.54
1:D:217:LEU:HD12	1:D:217:LEU:O	2.08	0.54
1:B:217:LEU:HD23	1:B:217:LEU:O	2.07	0.54
1:D:107:ASP:OD2	1:D:233[B]:THR:HG23	2.07	0.54
1:A:164:GLU:OE2	1:A:164:GLU:HA	2.07	0.54
1:D:203:VAL:HG22	1:D:230:GLN:HG2	1.90	0.54
1:A:303:ASP:OD2	1:A:303:ASP:C	2.51	0.53
1:B:66:GLN:NE2	2:B:452:HOH:O	2.42	0.53
1:D:118:MET:HE1	1:D:121:LEU:HD12	1.90	0.53
1:C:163:PRO:HB2	1:C:166:ILE:HD13	1.90	0.53
1:D:203:VAL:HG13	1:D:227:VAL:HG13	1.90	0.53
1:C:108:TYR:HH	1:C:276:TYR:HH	1.57	0.53
1:A:97:SER:OG	1:A:99:ARG:HG2	2.08	0.52
1:D:168:ALA:HA	1:D:280:LEU:O	2.09	0.52
1:A:16:PHE:HB2	1:A:46:LEU:HD21	1.91	0.52
1:B:156:ASP:OD2	1:B:279:ARG:NH1	2.42	0.52
1:D:182:ARG:HH12	1:D:260:GLU:CD	2.18	0.52
1:C:120[A]:ARG:HH12	1:C:301:GLY:HA3	1.75	0.51
1:B:107:ASP:HB2	2:D:447:HOH:O	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:165:GLY:HA2	1:C:284:VAL:CG2	2.38	0.51
1:B:208:ASN:HB3	2:D:442:HOH:O	2.09	0.51
1:A:100:THR:HA	1:A:129[B]:ARG:O	2.10	0.51
1:B:142:GLU:O	1:B:146:SER:OG	2.28	0.51
1:A:22:HIS:HB3	1:A:27:THR:HG23	1.93	0.50
1:C:155:TYR:HE1	2:C:392:HOH:O	1.93	0.50
1:B:22:HIS:NE2	1:B:31:GLU:OE2	2.28	0.50
1:D:145:ALA:HA	1:D:283:HIS:CD2	2.46	0.50
1:C:203:VAL:HG13	1:C:227:VAL:HG13	1.93	0.50
1:C:165:GLY:CA	1:C:284:VAL:HG22	2.40	0.50
1:B:61:ARG:HD2	2:B:503:HOH:O	2.10	0.50
1:A:67:ARG:HB3	1:D:89:TRP:CZ3	2.47	0.50
1:D:289:ARG:HA	2:D:410:HOH:O	2.11	0.50
1:B:174:ASP:HB3	1:B:276:TYR:CE1	2.45	0.49
1:A:120:ARG:HH22	1:A:301:GLY:HA3	1.77	0.49
1:B:226:GLU:HB3	2:B:350:HOH:O	2.12	0.49
1:A:14:LEU:HD12	2:A:421:HOH:O	2.12	0.49
1:A:141:VAL:HG13	1:A:166:ILE:CD1	2.42	0.48
1:D:300:LYS:HG3	2:D:366:HOH:O	2.13	0.48
1:B:16:PHE:HB2	1:B:46:LEU:CD2	2.43	0.48
1:A:6:ARG:HD2	2:A:380:HOH:O	2.12	0.48
1:C:156:ASP:HB3	2:C:460:HOH:O	2.14	0.48
1:A:220:ARG:NH2	2:A:460:HOH:O	2.45	0.48
1:C:252:ARG:O	1:C:256:ARG:HG3	2.13	0.48
1:B:161:ARG:HG2	1:B:161:ARG:NH1	2.28	0.48
1:C:174:ASP:HB3	1:C:276:TYR:OH	2.13	0.48
1:B:208:ASN:O	1:D:131:ARG:NH2	2.47	0.47
1:C:63:GLN:HG2	2:C:496:HOH:O	2.14	0.47
1:C:6:ARG:HH21	1:C:6:ARG:HG3	1.79	0.47
1:B:20:TYR:OH	1:B:69:GLU:OE2	2.24	0.47
1:D:16:PHE:HB2	1:D:46:LEU:CD2	2.44	0.47
1:D:135:ALA:HB1	2:D:395:HOH:O	2.14	0.47
1:D:251:PRO:HB3	1:D:276:TYR:OH	2.14	0.47
1:B:279:ARG:HG3	1:B:281:TYR:CZ	2.50	0.46
1:D:280:LEU:HD22	1:D:299:LEU:HD22	1.97	0.46
1:D:138:LYS:H	1:D:138:LYS:CE	2.28	0.46
1:C:163:PRO:HB2	1:C:166:ILE:CD1	2.44	0.46
1:C:304:ILE:HG23	2:C:319:HOH:O	2.14	0.46
1:A:131:ARG:NH2	1:C:208:ASN:O	2.48	0.46
1:B:16:PHE:HB2	1:B:46:LEU:HD21	1.98	0.46
1:B:285:GLN:HB2	2:B:487:HOH:O	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:302:LEU:HA	1:C:302:LEU:HD12	1.82	0.46
1:C:6:ARG:HG3	1:C:6:ARG:NH2	2.30	0.46
1:A:137:ARG:HD2	2:A:525:HOH:O	2.16	0.45
1:A:9:ASP:O	1:A:12:LEU:HB2	2.16	0.45
1:D:133:VAL:HG12	2:D:442:HOH:O	2.16	0.45
1:A:98:GLN:OE1	1:A:129[B]:ARG:NH1	2.49	0.45
1:C:290:ASP:HB2	2:C:382:HOH:O	2.17	0.45
1:A:88:GLU:OE1	1:D:70:HIS:ND1	2.50	0.45
1:C:119:ASN:ND2	1:C:120[B]:ARG:NH2	2.62	0.45
1:B:285:GLN:CG	2:B:407:HOH:O	2.65	0.45
1:C:89:TRP:O	1:B:67:ARG:NH1	2.50	0.45
1:A:12:LEU:HB3	1:A:46:LEU:HD13	1.99	0.45
1:B:157:GLU:H	1:B:160:GLU:CD	2.25	0.45
1:A:146:SER:CB	1:A:148:ARG:H	2.30	0.44
1:C:204:VAL:HA	1:C:231:LEU:O	2.17	0.44
1:C:211:SER:O	1:C:216:ARG:NH1	2.50	0.44
1:A:164:GLU:OE2	1:A:164:GLU:CA	2.65	0.44
1:B:54:LEU:O	1:B:64:PRO:HA	2.18	0.44
1:D:120:ARG:HH11	1:D:298:GLN:NE2	2.15	0.44
1:D:233[B]:THR:HG22	1:D:236:ALA:H	1.82	0.44
1:B:280:LEU:HD22	1:B:299:LEU:HD22	1.99	0.44
1:B:57:ARG:HH11	1:B:57:ARG:HB3	1.82	0.44
1:A:86:LEU:HD12	1:D:5:LEU:HD13	2.00	0.44
1:A:252:ARG:NH1	1:A:256:ARG:HD3	2.33	0.44
1:C:61:ARG:HE	1:C:61:ARG:HB3	1.20	0.44
1:C:253:HIS:HE1	1:C:305:SER:HB2	1.83	0.44
1:C:304:ILE:O	1:C:305:SER:C	2.61	0.44
1:B:211:SER:O	1:B:216:ARG:NH2	2.51	0.44
1:C:119:ASN:HD21	1:C:120[B]:ARG:HH21	1.62	0.43
1:C:155:TYR:HB2	1:C:277:VAL:O	2.18	0.43
1:A:288:GLY:HA3	2:A:434:HOH:O	2.16	0.43
1:B:137:ARG:HD3	2:B:368:HOH:O	2.19	0.43
1:B:303:ASP:OD2	2:B:498:HOH:O	2.21	0.43
1:D:284:VAL:O	1:D:285:GLN:CB	2.66	0.43
1:C:86:LEU:CD1	1:B:5:LEU:CD2	2.89	0.43
1:A:32:LEU:O	1:A:33:ALA:HB3	2.18	0.43
1:A:174:ASP:HB3	1:A:276:TYR:OH	2.17	0.43
1:C:300:LYS:HD3	1:C:300:LYS:N	2.34	0.43
1:B:115:PRO:HB2	1:B:116:PRO:HD3	1.99	0.43
1:B:229:VAL:HA	1:D:130:LEU:O	2.18	0.43
1:D:118:MET:HE3	1:D:130:LEU:HD22	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:299:LEU:O	1:D:300:LYS:C	2.62	0.43
1:C:2:HIS:HD2	1:C:4:LEU:H	1.67	0.42
1:B:182:ARG:NH2	2:B:370:HOH:O	2.52	0.42
1:D:117:LEU:CG	1:D:118:MET:CE	2.93	0.42
1:A:151:PHE:CG	1:A:280:LEU:HD11	2.54	0.42
1:B:7:ARG:NH1	2:B:439:HOH:O	2.32	0.42
1:B:297:GLY:HA2	1:B:300:LYS:HD2	2.02	0.42
1:D:203:VAL:HG22	1:D:230:GLN:CG	2.50	0.42
1:B:217:LEU:C	1:B:217:LEU:CD2	2.93	0.42
1:C:143:ALA:HB1	1:C:149:ILE:HD12	2.02	0.42
1:A:131:ARG:NH2	2:A:418:HOH:O	2.53	0.41
1:D:231:LEU:HD21	1:D:237:ALA:HB2	2.01	0.41
1:C:70[A]:HIS:HE2	1:B:88:GLU:HB2	1.84	0.41
1:C:40:SER:HB2	2:C:463:HOH:O	2.20	0.41
1:D:59:GLY:CA	1:D:61:ARG:H	2.33	0.41
1:C:86:LEU:O	1:C:89:TRP:HB3	2.20	0.41
1:B:274:PRO:HA	1:B:275:PRO:HD3	1.93	0.41
1:A:4:LEU:HD12	2:A:428:HOH:O	2.20	0.41
1:A:120:ARG:HH11	1:A:298:GLN:NE2	2.19	0.41
1:A:284:VAL:O	1:A:284:VAL:HG22	2.21	0.41
1:B:143:ALA:HB1	1:B:149:ILE:HD12	2.02	0.41
1:B:163:PRO:HA	2:B:477:HOH:O	2.21	0.41
1:D:100:THR:HA	1:D:129[A]:ARG:O	2.20	0.41
1:D:278:LEU:HD23	1:D:278:LEU:HA	1.89	0.41
1:C:138:LYS:HE3	1:C:281:TYR:CZ	2.56	0.41
1:B:157:GLU:OE2	1:B:216:ARG:HD2	2.21	0.41
1:B:161:ARG:HH11	1:B:161:ARG:CB	2.34	0.41
1:D:63:GLN:HA	1:D:64:PRO:HD3	1.95	0.40
1:C:129:ARG:NH1	2:C:409:HOH:O	2.54	0.40
1:C:141:VAL:HG13	1:C:166:ILE:HD11	2.04	0.40
1:B:137:ARG:HB2	2:B:447:HOH:O	2.21	0.40
1:D:16:PHE:HE1	1:D:72:ALA:HA	1.86	0.40
1:D:138:LYS:HB3	1:D:207:TRP:CH2	2.56	0.40
1:D:138:LYS:HB3	1:D:207:TRP:CZ2	2.57	0.40
1:C:260:GLU:HG2	2:C:478:HOH:O	2.20	0.40
1:C:92:PHE:HB2	1:C:290:ASP:HB3	2.03	0.40
1:B:204:VAL:HA	1:B:231:LEU:O	2.21	0.40
1:D:52:ASP:CG	1:D:53:GLU:H	2.28	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:41:HIS:ND1	1:B:183:ASP:OD1[1_556]	2.12	0.08
1:A:161[A]:ARG:NH1	1:C:289:ARG:N[4_456]	2.13	0.07

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	297/310 (96%)	286 (96%)	8 (3%)	3 (1%)	12	9
1	B	296/310 (96%)	286 (97%)	8 (3%)	2 (1%)	18	15
1	C	295/310 (95%)	283 (96%)	10 (3%)	2 (1%)	18	15
1	D	297/310 (96%)	281 (95%)	12 (4%)	4 (1%)	9	6
All	All	1185/1240 (96%)	1136 (96%)	38 (3%)	11 (1%)	14	10

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	300	LYS
1	D	141	VAL
1	D	159	HIS
1	A	164	GLU
1	C	301	GLY
1	B	88	GLU
1	A	289	ARG
1	B	158	GLU
1	D	162	LEU
1	D	300	LYS
1	C	88	GLU

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	234/241 (97%)	215 (92%)	19 (8%)	11	8
1	B	234/241 (97%)	214 (92%)	20 (8%)	10	7
1	C	235/241 (98%)	217 (92%)	18 (8%)	12	9
1	D	235/241 (98%)	210 (89%)	25 (11%)	6	4
All	All	938/964 (97%)	856 (91%)	82 (9%)	9	7

All (82) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	ARG
1	A	7	ARG
1	A	8	LEU
1	A	12	LEU
1	A	19	LEU
1	A	56	LEU
1	A	69	GLU
1	A	90	ARG
1	A	96	GLN
1	A	142	GLU
1	A	149	ILE
1	A	167	GLN
1	A	209	GLU
1	A	224	ARG
1	A	231	LEU
1	A	233	THR
1	A	289	ARG
1	A	302	LEU
1	A	303	ASP
1	C	6	ARG
1	C	61	ARG
1	C	66	GLN
1	C	136	GLU
1	C	137	ARG
1	C	139	LEU
1	C	156	ASP
1	C	164	GLU
1	C	166	ILE

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Mol	Chain	Res	Type
1	C	182	ARG
1	C	194	GLU
1	C	203	VAL
1	C	233	THR
1	C	266	LEU
1	C	289	ARG
1	C	290	ASP
1	C	300	LYS
1	C	302	LEU
1	B	2	HIS
1	B	4	LEU
1	B	6	ARG
1	B	19	LEU
1	B	57	ARG
1	B	61	ARG
1	B	66	GLN
1	B	88	GLU
1	B	93	VAL
1	B	107	ASP
1	B	113	LEU
1	B	117	LEU
1	B	131	ARG
1	B	142	GLU
1	B	146	SER
1	B	148	ARG
1	B	161	ARG
1	B	217	LEU
1	B	233	THR
1	B	290	ASP
1	D	4	LEU
1	D	5	LEU
1	D	6	ARG
1	D	7	ARG
1	D	10	LEU
1	D	19	LEU
1	D	48	GLN
1	D	63	GLN
1	D	90	ARG
1	D	137	ARG
1	D	138	LYS
1	D	141	VAL
1	D	148	ARG

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Mol	Chain	Res	Type
1	D	177	VAL
1	D	186	ARG
1	D	203	VAL
1	D	210	ASP
1	D	214	ILE
1	D	224	ARG
1	D	233[A]	THR
1	D	233[B]	THR
1	D	284	VAL
1	D	290	ASP
1	D	300	LYS
1	D	303	ASP

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	11	ASN
1	A	167	GLN
1	A	230	GLN
1	A	298	GLN
1	C	2	HIS
1	C	41	HIS
1	C	48	GLN
1	C	119	ASN
1	B	11	ASN
1	B	41	HIS
1	B	169	HIS
1	B	230	GLN
1	D	63	GLN
1	D	123	HIS
1	D	159	HIS
1	D	169	HIS
1	D	298	GLN

5.3.3 RNA ⓘ

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

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	299/310 (96%)	-0.02	7 (2%) 61 64	18, 39, 64, 86	2 (0%)
1	B	300/310 (96%)	0.07	4 (1%) 75 77	25, 41, 64, 79	0
1	C	298/310 (96%)	-0.02	6 (2%) 65 67	17, 37, 63, 81	3 (1%)
1	D	298/310 (96%)	0.39	19 (6%) 25 27	16, 50, 78, 96	3 (1%)
All	All	1195/1240 (96%)	0.11	36 (3%) 52 55	16, 41, 70, 96	8 (0%)

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	60	ASN	4.2
1	A	302	LEU	4.2
1	D	41[A]	HIS	3.8
1	A	3	PRO	3.4
1	A	4	LEU	3.3
1	C	155	TYR	3.3
1	D	56	LEU	3.2
1	D	141	VAL	3.1
1	C	60	ASN	2.9
1	B	220	ARG	2.8
1	D	156	ASP	2.8
1	D	284	VAL	2.8
1	D	159	HIS	2.7
1	D	3	PRO	2.6
1	A	161[A]	ARG	2.6
1	C	301	GLY	2.5
1	D	4	LEU	2.5
1	D	62	MET	2.5
1	B	305	SER	2.5
1	B	155	TYR	2.4
1	D	139	LEU	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	286	HIS	2.4
1	D	155	TYR	2.3
1	A	70	HIS	2.3
1	D	167	GLN	2.3
1	D	59	GLY	2.3
1	D	162	LEU	2.3
1	D	86	LEU	2.2
1	D	143	ALA	2.2
1	C	70[A]	HIS	2.2
1	D	300	LYS	2.2
1	C	304	ILE	2.1
1	B	158	GLU	2.1
1	D	70	HIS	2.1
1	A	288	GLY	2.0
1	A	301	GLY	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 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.