



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 04:08 PM UTC

PDB ID : 1RLV / pdb_00001rlv
Title : Crystal structure of a dimeric Archaeal Splicing Endonuclease
Authors : Li, H.; Abelson, J.
Deposited on : 2003-11-26
Resolution : 3.00 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

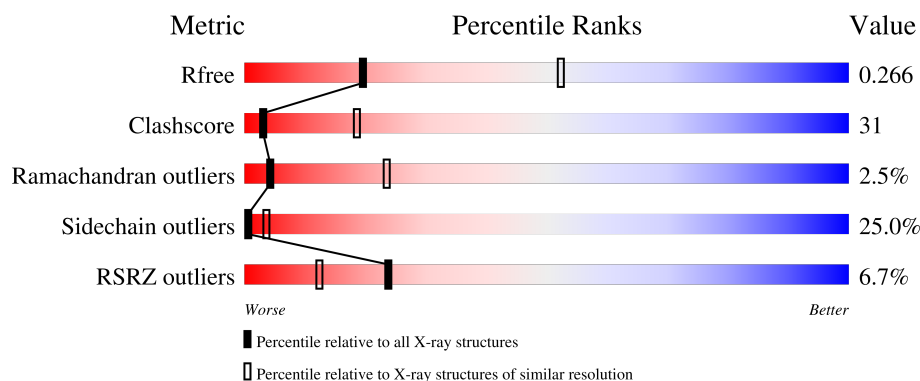
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	305	7% 41% 43% 14% .
1	B	305	7% 40% 43% 15% .

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 5058 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 Putative tRNA-intron endonuclease.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	304	Total	C	N	O	S	0	0	0
			2529	1618	433	472	6			
1	B	304	Total	C	N	O	S	0	0	0
			2529	1618	433	472	6			

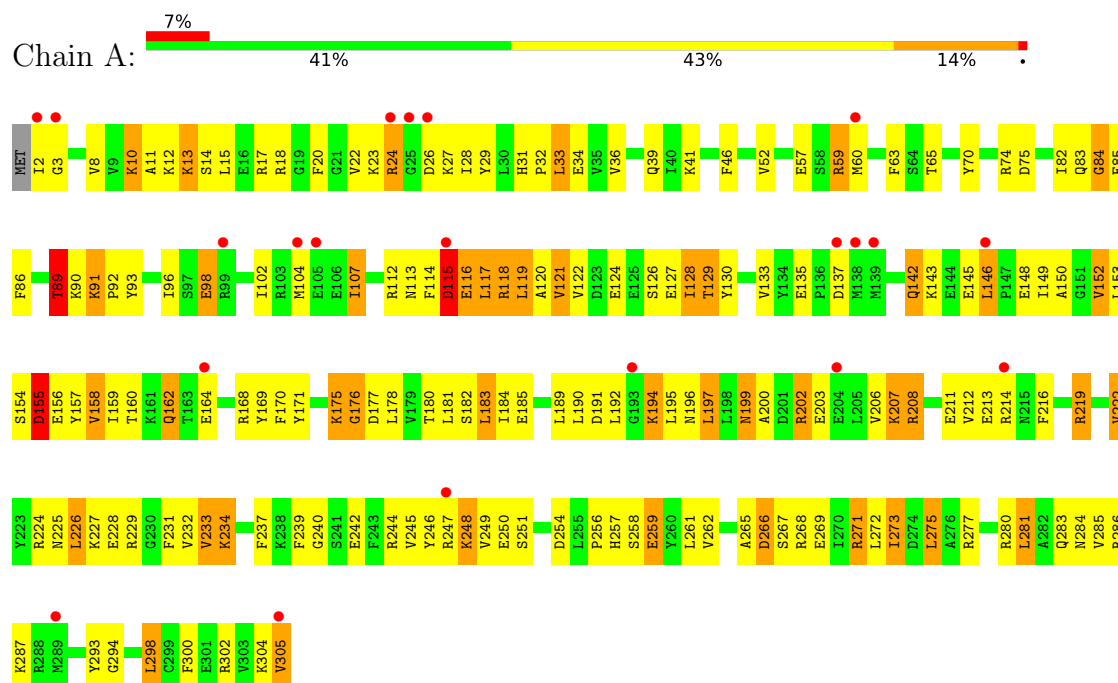
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	152	VAL	ILE	conflict	UNP O29362
B	152	VAL	ILE	conflict	UNP O29362

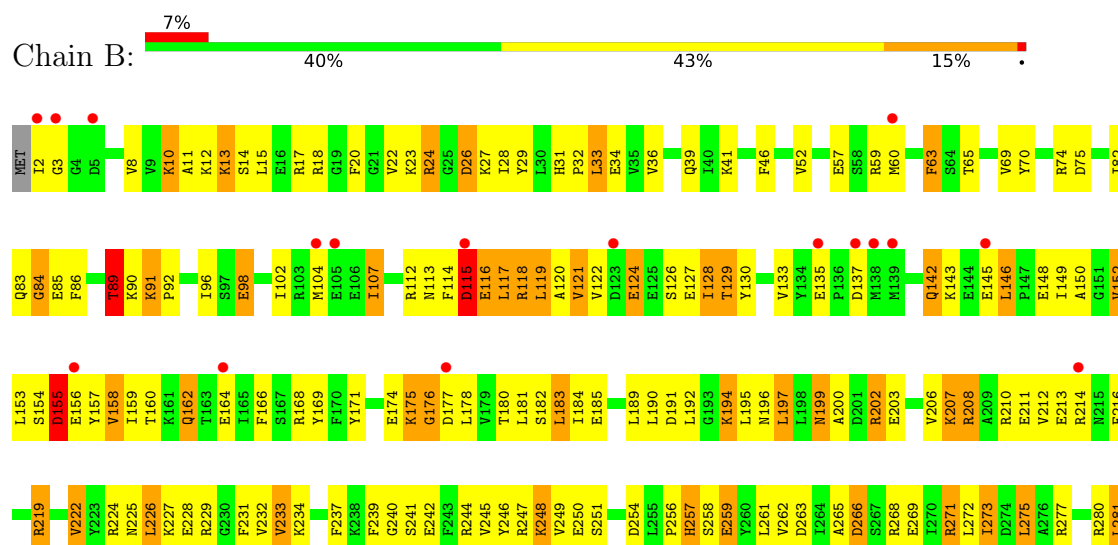
3 Residue-property plots

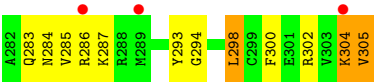
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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



• Molecule 1: Putative tRNA-intron endonuclease





4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	126.82Å 126.82Å 130.23Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.92 – 3.00 19.92 – 3.00	Depositor EDS
% Data completeness (in resolution range)	97.0 (19.92-3.00) 96.7 (19.92-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.07 (at 2.98Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.229 , 0.276 (Not available) , 0.266	Depositor DCC
R_{free} test set	1027 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å ²)	52.9	Xtriage
Anisotropy	0.514	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 82.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.016 for -h,l,k 0.014 for -l,-k,-h	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	5058	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 3.63% 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	A	0.37	0/2573	0.83	5/3451 (0.1%)
1	B	0.38	0/2573	0.83	6/3451 (0.2%)
All	All	0.37	0/5146	0.83	11/6902 (0.2%)

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	124	GLU	N-CA-C	-6.81	104.72	113.16
1	A	245	VAL	N-CA-C	6.66	118.63	108.71
1	A	124	GLU	N-CA-C	-6.64	104.93	113.16
1	B	245	VAL	N-CA-C	6.43	118.30	108.71
1	A	257	HIS	N-CA-C	6.35	118.06	110.44
1	B	257	HIS	N-CA-C	6.09	117.75	110.44
1	A	233	VAL	N-CA-C	5.75	116.14	107.80
1	A	262	VAL	N-CA-C	5.72	116.18	108.17
1	B	262	VAL	N-CA-C	5.70	116.14	108.17
1	B	233	VAL	N-CA-C	5.69	116.05	107.80
1	B	69	VAL	N-CA-C	-5.23	105.61	110.53

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2529	0	2545	177	0
1	B	2529	0	2545	168	0
All	All	5058	0	5090	319	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 31.

All (319) 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:280:ARG:HA	1:A:302:ARG:HH22	1.17	1.08
1:B:280:ARG:HA	1:B:302:ARG:HH22	1.16	1.06
1:A:284:ASN:HB3	1:B:284:ASN:HD22	1.31	0.94
1:B:273:ILE:HD12	1:B:273:ILE:H	1.32	0.94
1:A:284:ASN:HD22	1:B:284:ASN:HB3	1.32	0.93
1:A:273:ILE:H	1:A:273:ILE:HD12	1.32	0.93
1:B:39:GLN:HE22	1:B:46:PHE:H	1.20	0.90
1:B:266:ASP:HB3	1:B:293:TYR:HA	1.54	0.89
1:B:104:MET:HE1	1:B:300:PHE:HE1	1.39	0.87
1:A:39:GLN:HE22	1:A:46:PHE:H	1.23	0.86
1:A:266:ASP:HB3	1:A:293:TYR:HA	1.54	0.85
1:B:150:ALA:H	1:B:162:GLN:HE21	1.24	0.85
1:A:104:MET:HE1	1:A:300:PHE:HE1	1.42	0.84
1:A:150:ALA:H	1:A:162:GLN:HE21	1.23	0.83
1:B:104:MET:HE1	1:B:300:PHE:CE1	2.14	0.81
1:A:214:ARG:HH22	1:B:271:ARG:HG2	1.46	0.80
1:A:149:ILE:HB	1:A:162:GLN:HG3	1.64	0.80
1:A:104:MET:HE1	1:A:300:PHE:CE1	2.16	0.80
1:A:153:LEU:H	1:A:199:ASN:HD21	1.30	0.79
1:A:24:ARG:HB2	1:A:24:ARG:HH11	1.48	0.79
1:A:92:PRO:HG2	1:A:117:LEU:HA	1.66	0.78
1:A:271:ARG:HG2	1:B:214:ARG:HH22	1.47	0.78
1:B:24:ARG:HH11	1:B:24:ARG:HB2	1.49	0.78
1:B:153:LEU:H	1:B:199:ASN:HD21	1.30	0.77
1:A:39:GLN:HE21	1:A:52:VAL:HG21	1.48	0.77
1:B:149:ILE:HB	1:B:162:GLN:HG3	1.64	0.77
1:A:214:ARG:NH2	1:B:271:ARG:HG2	2.01	0.76
1:B:92:PRO:HG2	1:B:117:LEU:HA	1.67	0.76
1:B:39:GLN:HE21	1:B:52:VAL:HG21	1.52	0.75
1:A:232:VAL:HB	1:A:246:TYR:HB2	1.68	0.75
1:B:232:VAL:HB	1:B:246:TYR:HB2	1.68	0.74
1:B:39:GLN:NE2	1:B:46:PHE:H	1.84	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:271:ARG:HG2	1:B:214:ARG:NH2	2.03	0.73
1:A:212:VAL:HG23	1:A:213:GLU:HG2	1.71	0.72
1:B:280:ARG:HA	1:B:302:ARG:NH2	2.00	0.72
1:A:280:ARG:HA	1:A:302:ARG:NH2	2.00	0.72
1:A:39:GLN:NE2	1:A:46:PHE:H	1.88	0.71
1:B:212:VAL:HG23	1:B:213:GLU:HG2	1.72	0.71
1:A:284:ASN:ND2	1:B:284:ASN:HB3	2.05	0.71
1:A:107:ILE:HD11	1:A:133:VAL:HG11	1.72	0.71
1:A:150:ALA:H	1:A:162:GLN:NE2	1.89	0.71
1:B:150:ALA:H	1:B:162:GLN:NE2	1.89	0.70
1:B:249:VAL:HG13	1:B:254:ASP:HB2	1.74	0.70
1:A:249:VAL:HG13	1:A:254:ASP:HB2	1.73	0.70
1:A:180:THR:HG21	1:B:126:SER:HB3	1.72	0.70
1:A:284:ASN:HB3	1:B:284:ASN:ND2	2.04	0.70
1:B:84:GLY:O	1:B:85:GLU:HB3	1.91	0.69
1:A:146:LEU:O	1:A:194:LYS:HE3	1.92	0.69
1:B:146:LEU:O	1:B:194:LYS:HE3	1.92	0.69
1:A:39:GLN:NE2	1:A:52:VAL:HG21	2.07	0.69
1:B:107:ILE:HD11	1:B:133:VAL:HG11	1.73	0.68
1:B:158:VAL:HG22	1:B:181:LEU:HB2	1.75	0.68
1:A:142:GLN:HE22	1:A:249:VAL:H	1.40	0.68
1:B:24:ARG:O	1:B:24:ARG:HG3	1.93	0.68
1:A:24:ARG:HG3	1:A:24:ARG:O	1.95	0.67
1:A:158:VAL:HG22	1:A:181:LEU:HB2	1.76	0.67
1:B:39:GLN:NE2	1:B:52:VAL:HG21	2.10	0.67
1:A:15:LEU:HD23	1:A:18:ARG:NH2	2.10	0.66
1:A:197:LEU:HD13	1:A:200:ALA:HB3	1.77	0.66
1:B:197:LEU:HD13	1:B:200:ALA:HB3	1.77	0.66
1:B:15:LEU:HD23	1:B:18:ARG:NH2	2.10	0.66
1:B:142:GLN:HE22	1:B:249:VAL:H	1.42	0.65
1:A:126:SER:HB3	1:B:180:THR:HG21	1.77	0.64
1:B:41:LYS:HE2	1:B:75:ASP:OD2	1.98	0.64
1:A:84:GLY:O	1:A:85:GLU:HB3	1.98	0.64
1:A:2:ILE:HG12	1:A:3:GLY:H	1.63	0.63
1:B:2:ILE:HG12	1:B:3:GLY:H	1.63	0.63
1:A:237:PHE:CE1	1:B:98:GLU:HB2	2.34	0.63
1:A:41:LYS:HE2	1:A:75:ASP:OD2	1.99	0.63
1:B:89:THR:C	1:B:91:LYS:H	2.06	0.63
1:A:273:ILE:H	1:A:273:ILE:CD1	2.09	0.62
1:A:89:THR:C	1:A:91:LYS:H	2.07	0.62
1:A:98:GLU:HB2	1:B:237:PHE:CE1	2.35	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:153:LEU:N	1:B:199:ASN:HD21	1.97	0.62
1:A:153:LEU:HD12	1:A:158:VAL:HG12	1.81	0.61
1:A:107:ILE:HD11	1:A:133:VAL:HG21	1.82	0.61
1:A:226:LEU:HB3	1:A:233:VAL:HG21	1.83	0.61
1:A:281:LEU:HD21	1:B:280:ARG:HG2	1.83	0.61
1:A:280:ARG:HG2	1:B:281:LEU:HD21	1.82	0.61
1:B:226:LEU:HB3	1:B:233:VAL:HG21	1.82	0.61
1:A:160:THR:OG1	1:A:162:GLN:HG2	2.01	0.60
1:B:153:LEU:HD12	1:B:158:VAL:HG12	1.83	0.59
1:A:222:VAL:HG23	1:A:226:LEU:HD22	1.83	0.59
1:A:2:ILE:HG23	1:A:3:GLY:N	2.17	0.59
1:A:265:ALA:HA	1:A:268:ARG:NH1	2.17	0.59
1:A:197:LEU:HB3	1:A:200:ALA:O	2.03	0.59
1:A:208:ARG:O	1:A:211:GLU:HB2	2.03	0.59
1:B:107:ILE:HD11	1:B:133:VAL:HG21	1.83	0.59
1:B:142:GLN:NE2	1:B:249:VAL:H	2.01	0.59
1:A:126:SER:OG	1:B:175:LYS:HE2	2.03	0.59
1:A:153:LEU:N	1:A:199:ASN:HD21	1.98	0.58
1:B:20:PHE:HD2	1:B:34:GLU:HG2	1.67	0.58
1:B:266:ASP:CB	1:B:293:TYR:HA	2.31	0.58
1:B:2:ILE:HG23	1:B:3:GLY:N	2.18	0.58
1:B:160:THR:OG1	1:B:162:GLN:HG2	2.02	0.58
1:A:142:GLN:NE2	1:A:249:VAL:H	2.00	0.58
1:B:152:VAL:HA	1:B:199:ASN:ND2	2.19	0.58
1:B:208:ARG:O	1:B:211:GLU:HB2	2.03	0.58
1:A:102:ILE:HD11	1:A:275:LEU:HD13	1.86	0.58
1:B:102:ILE:HD11	1:B:275:LEU:HD13	1.85	0.58
1:A:199:ASN:H	1:A:199:ASN:HD22	1.52	0.57
1:B:197:LEU:HB3	1:B:200:ALA:O	2.04	0.57
1:A:20:PHE:HD2	1:A:34:GLU:HG2	1.68	0.57
1:B:102:ILE:HD11	1:B:275:LEU:CD1	2.34	0.57
1:A:102:ILE:HD11	1:A:275:LEU:CD1	2.35	0.57
1:B:222:VAL:HG23	1:B:226:LEU:HD22	1.85	0.57
1:B:117:LEU:HD22	1:B:118:ARG:N	2.20	0.56
1:A:152:VAL:HA	1:A:199:ASN:ND2	2.19	0.56
1:A:117:LEU:HD22	1:A:118:ARG:N	2.20	0.56
1:B:265:ALA:HA	1:B:268:ARG:NH1	2.19	0.56
1:A:212:VAL:HG23	1:A:213:GLU:N	2.21	0.56
1:B:92:PRO:HG3	1:B:114:PHE:HB3	1.88	0.56
1:A:207:LYS:HE3	1:A:207:LYS:O	2.06	0.55
1:B:199:ASN:HD22	1:B:199:ASN:H	1.53	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:244:ARG:HD2	1:A:246:TYR:OH	2.07	0.55
1:A:92:PRO:HG3	1:A:114:PHE:HB3	1.87	0.55
1:B:85:GLU:HG2	1:B:86:PHE:CE1	2.42	0.55
1:A:266:ASP:CB	1:A:293:TYR:HA	2.32	0.55
1:B:212:VAL:HG23	1:B:213:GLU:N	2.22	0.54
1:A:121:VAL:HG13	1:A:272:LEU:HD13	1.90	0.54
1:A:120:ALA:HB1	1:A:128:ILE:HD13	1.90	0.54
1:B:176:GLY:O	1:B:177:ASP:HB2	2.08	0.54
1:B:121:VAL:HG13	1:B:272:LEU:HD13	1.90	0.54
1:B:150:ALA:CB	1:B:196:ASN:HB3	2.38	0.54
1:A:254:ASP:C	1:A:256:PRO:HD2	2.32	0.53
1:B:280:ARG:HD2	1:B:302:ARG:NH2	2.22	0.53
1:A:280:ARG:HD2	1:A:302:ARG:NH2	2.23	0.53
1:B:207:LYS:HE3	1:B:207:LYS:O	2.07	0.53
1:A:184:ILE:HG23	1:A:185:GLU:N	2.24	0.53
1:A:244:ARG:HG3	1:A:258:SER:CB	2.39	0.53
1:A:176:GLY:O	1:A:177:ASP:HB2	2.07	0.53
1:A:228:GLU:O	1:A:228:GLU:HG2	2.09	0.53
1:B:222:VAL:O	1:B:226:LEU:HD22	2.09	0.53
1:B:254:ASP:C	1:B:256:PRO:HD2	2.33	0.53
1:A:226:LEU:O	1:A:231:PHE:HB2	2.08	0.53
1:A:85:GLU:HG2	1:A:86:PHE:CE1	2.44	0.52
1:B:244:ARG:HG3	1:B:258:SER:CB	2.39	0.52
1:A:150:ALA:CB	1:A:196:ASN:HB3	2.40	0.52
1:B:244:ARG:HD2	1:B:246:TYR:OH	2.10	0.52
1:B:226:LEU:O	1:B:231:PHE:HB2	2.10	0.51
1:A:224:ARG:HH11	1:A:224:ARG:HB3	1.75	0.51
1:B:120:ALA:HB1	1:B:128:ILE:HD13	1.92	0.51
1:B:224:ARG:HH11	1:B:224:ARG:HB3	1.76	0.51
1:B:228:GLU:O	1:B:228:GLU:HG2	2.09	0.51
1:A:24:ARG:HG2	1:A:29:TYR:CE1	2.46	0.51
1:A:240:GLY:HA3	1:B:277:ARG:HB2	1.92	0.51
1:A:277:ARG:HB2	1:B:240:GLY:HA3	1.93	0.50
1:B:273:ILE:H	1:B:273:ILE:CD1	2.08	0.50
1:B:14:SER:HB2	1:B:18:ARG:HH12	1.76	0.50
1:B:184:ILE:HG23	1:B:185:GLU:N	2.25	0.50
1:A:14:SER:HB2	1:A:18:ARG:HH12	1.77	0.50
1:A:225:ASN:OD1	1:A:229:ARG:HD2	2.11	0.50
1:B:20:PHE:CD2	1:B:34:GLU:HG2	2.45	0.50
1:B:70:TYR:CE2	1:B:74:ARG:HG3	2.47	0.50
1:A:20:PHE:CD2	1:A:34:GLU:HG2	2.46	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:168:ARG:NH2	1:B:169:TYR:OH	2.44	0.50
1:B:156:GLU:HB3	1:B:157:TYR:CE1	2.47	0.49
1:A:116:GLU:HB3	1:A:118:ARG:HH21	1.77	0.49
1:B:266:ASP:HB3	1:B:294:GLY:H	1.76	0.49
1:A:266:ASP:HB3	1:A:294:GLY:H	1.77	0.49
1:A:24:ARG:HG2	1:A:29:TYR:HE1	1.77	0.49
1:B:225:ASN:OD1	1:B:229:ARG:HD2	2.12	0.49
1:A:222:VAL:O	1:A:226:LEU:HD22	2.13	0.49
1:A:277:ARG:HG2	1:B:277:ARG:HD3	1.95	0.48
1:A:70:TYR:CE2	1:A:74:ARG:HG3	2.47	0.48
1:A:156:GLU:HB3	1:A:157:TYR:CE1	2.48	0.48
1:B:116:GLU:HB3	1:B:118:ARG:HH21	1.78	0.48
1:A:24:ARG:CG	1:A:29:TYR:HE1	2.27	0.48
1:A:96:ILE:HG21	1:A:119:LEU:HD22	1.96	0.48
1:A:168:ARG:NH2	1:A:169:TYR:OH	2.44	0.48
1:B:24:ARG:HG2	1:B:29:TYR:CE1	2.48	0.48
1:A:266:ASP:O	1:A:266:ASP:OD2	2.32	0.47
1:A:265:ALA:O	1:A:266:ASP:C	2.57	0.47
1:B:104:MET:HE3	1:B:298:LEU:HD12	1.97	0.47
1:A:104:MET:HE3	1:A:298:LEU:HD12	1.96	0.47
1:A:128:ILE:HD11	1:A:130:TYR:CD2	2.50	0.47
1:A:219:ARG:HE	1:A:242:GLU:HG3	1.79	0.47
1:B:219:ARG:HE	1:B:242:GLU:HG3	1.78	0.47
1:B:265:ALA:O	1:B:266:ASP:C	2.57	0.47
1:B:8:VAL:HG12	1:B:29:TYR:CD2	2.50	0.47
1:B:24:ARG:HG2	1:B:29:TYR:HE1	1.79	0.47
1:B:280:ARG:CA	1:B:302:ARG:HH22	2.06	0.47
1:A:171:TYR:CD2	1:A:189:LEU:HD11	2.50	0.46
1:B:190:LEU:HD13	1:B:190:LEU:C	2.41	0.46
1:A:118:ARG:HD3	1:A:305:VAL:O	2.15	0.46
1:A:13:LYS:O	1:A:17:ARG:HG3	2.16	0.46
1:A:192:LEU:HD23	1:A:194:LYS:HD2	1.97	0.46
1:A:36:VAL:HG21	1:A:63:PHE:CZ	2.51	0.46
1:A:154:SER:HB2	1:A:159:ILE:HD12	1.96	0.46
1:A:191:ASP:OD1	1:A:202:ARG:NH1	2.49	0.46
1:B:24:ARG:CG	1:B:29:TYR:HE1	2.29	0.46
1:A:33:LEU:HD13	1:A:60:MET:SD	2.56	0.46
1:A:89:THR:C	1:A:91:LYS:N	2.74	0.46
1:A:142:GLN:NE2	1:A:248:LYS:HA	2.31	0.46
1:B:171:TYR:CD2	1:B:189:LEU:HD11	2.51	0.46
1:A:152:VAL:HA	1:A:199:ASN:HD21	1.81	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:175:LYS:HE2	1:B:126:SER:OG	2.16	0.45
1:A:190:LEU:C	1:A:190:LEU:HD13	2.42	0.45
1:B:14:SER:HB2	1:B:18:ARG:NH1	2.31	0.45
1:B:266:ASP:O	1:B:266:ASP:OD2	2.33	0.45
1:A:14:SER:HB2	1:A:18:ARG:NH1	2.31	0.45
1:A:280:ARG:CA	1:A:302:ARG:HH22	2.06	0.45
1:B:150:ALA:HB2	1:B:196:ASN:HB3	1.97	0.45
1:A:150:ALA:HB2	1:A:196:ASN:HB3	1.99	0.45
1:A:277:ARG:HD3	1:B:277:ARG:HG2	1.97	0.45
1:B:154:SER:O	1:B:155:ASP:O	2.35	0.45
1:B:85:GLU:HG2	1:B:86:PHE:CD1	2.51	0.45
1:A:155:ASP:HB3	1:B:57:GLU:OE1	2.17	0.45
1:A:213:GLU:HB2	1:A:216:PHE:HB2	1.98	0.45
1:A:104:MET:CE	1:A:298:LEU:HB3	2.47	0.45
1:A:116:GLU:O	1:A:117:LEU:C	2.60	0.45
1:A:182:SER:OG	1:A:184:ILE:HG22	2.17	0.45
1:B:96:ILE:HG21	1:B:119:LEU:HD22	1.98	0.45
1:B:152:VAL:HA	1:B:199:ASN:HD21	1.80	0.45
1:A:206:VAL:HG23	1:A:207:LYS:N	2.32	0.45
1:B:83:GLN:O	1:B:84:GLY:C	2.59	0.45
1:B:199:ASN:ND2	1:B:199:ASN:H	2.15	0.45
1:A:156:GLU:O	1:A:183:LEU:HD13	2.17	0.45
1:B:213:GLU:HB2	1:B:216:PHE:HB2	1.99	0.45
1:B:258:SER:OG	1:B:287:LYS:HE3	2.17	0.45
1:A:199:ASN:HD22	1:A:199:ASN:N	2.15	0.44
1:B:247:ARG:HG3	1:B:259:GLU:CD	2.42	0.44
1:A:83:GLN:O	1:A:84:GLY:C	2.61	0.44
1:B:31:HIS:ND1	1:B:33:LEU:HB2	2.33	0.44
1:B:154:SER:HB2	1:B:159:ILE:HD12	1.98	0.44
1:A:85:GLU:HG2	1:A:86:PHE:CD1	2.52	0.44
1:B:142:GLN:NE2	1:B:248:LYS:HA	2.32	0.44
1:B:266:ASP:HA	1:B:293:TYR:HB2	2.00	0.44
1:B:24:ARG:HH11	1:B:24:ARG:CB	2.26	0.44
1:B:118:ARG:HD3	1:B:305:VAL:O	2.17	0.44
1:B:244:ARG:HG3	1:B:258:SER:HB3	2.00	0.44
1:A:244:ARG:HG3	1:A:258:SER:HB3	2.00	0.44
1:A:285:VAL:HG22	1:A:285:VAL:O	2.18	0.44
1:B:145:GLU:CD	1:B:146:LEU:N	2.76	0.44
1:B:191:ASP:OD1	1:B:202:ARG:NH1	2.51	0.44
1:B:206:VAL:HG23	1:B:207:LYS:N	2.32	0.44
1:B:206:VAL:O	1:B:210:ARG:HG3	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:226:LEU:HB3	1:A:233:VAL:CG2	2.48	0.44
1:A:184:ILE:CG2	1:A:185:GLU:N	2.80	0.44
1:A:214:ARG:HH22	1:B:271:ARG:CG	2.25	0.44
1:A:266:ASP:HA	1:A:293:TYR:HB2	1.99	0.44
1:B:239:PHE:HE2	1:B:285:VAL:HG21	1.83	0.44
1:A:145:GLU:CD	1:A:146:LEU:N	2.76	0.43
1:B:104:MET:CE	1:B:298:LEU:HB3	2.48	0.43
1:B:116:GLU:O	1:B:117:LEU:C	2.60	0.43
1:A:59:ARG:HD3	1:A:59:ARG:HA	1.84	0.43
1:A:239:PHE:HE2	1:A:285:VAL:HG21	1.83	0.43
1:A:247:ARG:HG3	1:A:259:GLU:CD	2.43	0.43
1:B:13:LYS:O	1:B:17:ARG:HG3	2.17	0.43
1:B:145:GLU:CD	1:B:146:LEU:H	2.26	0.43
1:A:145:GLU:CD	1:A:146:LEU:H	2.26	0.43
1:B:26:ASP:OD1	1:B:26:ASP:N	2.51	0.43
1:B:192:LEU:HD23	1:B:194:LYS:HD2	1.99	0.43
1:A:57:GLU:OE1	1:B:155:ASP:HB3	2.18	0.43
1:A:298:LEU:HD23	1:A:298:LEU:HA	1.82	0.43
1:A:258:SER:OG	1:A:287:LYS:HE3	2.18	0.43
1:B:182:SER:OG	1:B:184:ILE:HG22	2.19	0.43
1:A:237:PHE:CD1	1:B:98:GLU:HB2	2.53	0.43
1:A:182:SER:HA	1:B:124:GLU:O	2.18	0.43
1:B:128:ILE:HD12	1:B:129:THR:N	2.33	0.43
1:A:213:GLU:OE2	1:A:219:ARG:HD3	2.19	0.43
1:B:33:LEU:HD13	1:B:60:MET:SD	2.58	0.43
1:A:31:HIS:CG	1:A:32:PRO:HD2	2.54	0.42
1:A:89:THR:HG23	1:A:93:TYR:HE1	1.84	0.42
1:B:117:LEU:HD22	1:B:118:ARG:H	1.82	0.42
1:A:36:VAL:HG21	1:A:63:PHE:HZ	1.84	0.42
1:A:199:ASN:H	1:A:199:ASN:ND2	2.15	0.42
1:B:156:GLU:O	1:B:183:LEU:HD13	2.19	0.42
1:A:8:VAL:HG12	1:A:29:TYR:CD2	2.55	0.42
1:A:128:ILE:HD11	1:A:130:TYR:CE2	2.55	0.42
1:A:149:ILE:CB	1:A:162:GLN:HG3	2.42	0.42
1:B:285:VAL:O	1:B:285:VAL:HG22	2.19	0.42
1:A:24:ARG:HH11	1:A:24:ARG:CB	2.25	0.42
1:A:117:LEU:HD22	1:A:118:ARG:H	1.84	0.42
1:B:31:HIS:CG	1:B:32:PRO:HD2	2.54	0.42
1:B:36:VAL:HG21	1:B:63:PHE:CZ	2.55	0.42
1:A:146:LEU:HB2	1:A:194:LYS:HD3	2.02	0.42
1:A:154:SER:O	1:A:155:ASP:O	2.37	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:15:LEU:HD23	1:B:18:ARG:HH22	1.83	0.42
1:B:175:LYS:O	1:B:176:GLY:O	2.38	0.42
1:A:225:ASN:O	1:A:229:ARG:HG3	2.19	0.42
1:B:237:PHE:C	1:B:239:PHE:H	2.28	0.42
1:B:256:PRO:HG2	1:B:257:HIS:CE1	2.55	0.42
1:A:31:HIS:ND1	1:A:33:LEU:HB2	2.35	0.42
1:A:142:GLN:HE21	1:A:248:LYS:HA	1.85	0.42
1:B:128:ILE:HD11	1:B:130:TYR:CD2	2.54	0.42
1:B:213:GLU:OE2	1:B:219:ARG:HD3	2.19	0.42
1:B:184:ILE:CG2	1:B:185:GLU:N	2.82	0.42
1:A:11:ALA:HA	1:A:28:ILE:HG13	2.02	0.41
1:B:31:HIS:CG	1:B:82:ILE:HD13	2.56	0.41
1:B:142:GLN:HE22	1:B:249:VAL:N	2.14	0.41
1:A:114:PHE:O	1:A:115:ASP:C	2.63	0.41
1:A:231:PHE:HE2	1:A:247:ARG:HG2	1.86	0.41
1:B:146:LEU:HB2	1:B:194:LYS:HD3	2.02	0.41
1:A:13:LYS:HA	1:A:13:LYS:HD3	1.68	0.41
1:B:166:PHE:CE2	1:B:174:GLU:HB2	2.55	0.41
1:B:121:VAL:CG1	1:B:272:LEU:HD13	2.51	0.41
1:A:10:LYS:O	1:A:12:LYS:HG3	2.20	0.41
1:A:183:LEU:HD12	1:A:183:LEU:HA	1.86	0.41
1:A:266:ASP:O	1:A:267:SER:HB3	2.20	0.41
1:B:11:ALA:HA	1:B:28:ILE:HG13	2.03	0.41
1:B:298:LEU:HD23	1:B:298:LEU:HA	1.80	0.41
1:A:31:HIS:CG	1:A:82:ILE:HD13	2.56	0.41
1:B:156:GLU:HB3	1:B:157:TYR:CD1	2.56	0.41
1:B:226:LEU:HB3	1:B:233:VAL:CG2	2.49	0.41
1:A:107:ILE:CD1	1:A:133:VAL:HG11	2.47	0.40
1:A:155:ASP:HB3	1:B:57:GLU:CD	2.45	0.40
1:A:156:GLU:HB3	1:A:157:TYR:CD1	2.56	0.40
1:A:237:PHE:C	1:A:239:PHE:H	2.29	0.40
1:B:114:PHE:O	1:B:115:ASP:C	2.64	0.40
1:B:304:LYS:HE2	1:B:305:VAL:CG2	2.51	0.40
1:A:83:GLN:O	1:A:84:GLY:O	2.38	0.40
1:A:121:VAL:CG1	1:A:272:LEU:HD13	2.51	0.40
1:A:128:ILE:HD12	1:A:129:THR:N	2.36	0.40
1:A:153:LEU:CD1	1:A:158:VAL:HG12	2.50	0.40
1:A:170:PHE:O	1:A:234:LYS:HE2	2.22	0.40
1:B:241:SER:HA	1:B:263:ASP:OD1	2.21	0.40
1:A:46:PHE:CD1	1:A:46:PHE:N	2.89	0.40
1:B:10:LYS:O	1:B:12:LYS:HG3	2.22	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	302/305 (99%)	265 (88%)	30 (10%)	7 (2%)	5	25
1	B	302/305 (99%)	266 (88%)	28 (9%)	8 (3%)	4	23
All	All	604/610 (99%)	531 (88%)	58 (10%)	15 (2%)	4	23

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	115	ASP
1	A	266	ASP
1	B	115	ASP
1	B	155	ASP
1	B	266	ASP
1	A	84	GLY
1	A	155	ASP
1	A	176	GLY
1	B	84	GLY
1	B	176	GLY
1	A	89	THR
1	B	89	THR
1	A	146	LEU
1	B	146	LEU
1	B	63	PHE

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	272/273 (100%)	204 (75%)	68 (25%)	0	3
1	B	272/273 (100%)	204 (75%)	68 (25%)	0	3
All	All	544/546 (100%)	408 (75%)	136 (25%)	0	3

All (136) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	LYS
1	A	13	LYS
1	A	22	VAL
1	A	23	LYS
1	A	24	ARG
1	A	26	ASP
1	A	27	LYS
1	A	33	LEU
1	A	59	ARG
1	A	65	THR
1	A	89	THR
1	A	90	LYS
1	A	91	LYS
1	A	98	GLU
1	A	107	ILE
1	A	112	ARG
1	A	113	ASN
1	A	115	ASP
1	A	116	GLU
1	A	117	LEU
1	A	118	ARG
1	A	119	LEU
1	A	121	VAL
1	A	122	VAL
1	A	127	GLU
1	A	128	ILE
1	A	129	THR
1	A	135	GLU
1	A	137	ASP
1	A	142	GLN
1	A	143	LYS
1	A	148	GLU
1	A	152	VAL
1	A	155	ASP

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Mol	Chain	Res	Type
1	A	158	VAL
1	A	162	GLN
1	A	164	GLU
1	A	175	LYS
1	A	178	LEU
1	A	183	LEU
1	A	194	LYS
1	A	195	LEU
1	A	197	LEU
1	A	199	ASN
1	A	202	ARG
1	A	203	GLU
1	A	207	LYS
1	A	208	ARG
1	A	219	ARG
1	A	222	VAL
1	A	226	LEU
1	A	227	LYS
1	A	234	LYS
1	A	248	LYS
1	A	250	GLU
1	A	251	SER
1	A	259	GLU
1	A	261	LEU
1	A	269	GLU
1	A	271	ARG
1	A	273	ILE
1	A	275	LEU
1	A	281	LEU
1	A	283	GLN
1	A	286	ARG
1	A	298	LEU
1	A	304	LYS
1	A	305	VAL
1	B	10	LYS
1	B	13	LYS
1	B	22	VAL
1	B	23	LYS
1	B	24	ARG
1	B	26	ASP
1	B	27	LYS
1	B	33	LEU

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Mol	Chain	Res	Type
1	B	59	ARG
1	B	65	THR
1	B	89	THR
1	B	90	LYS
1	B	91	LYS
1	B	98	GLU
1	B	107	ILE
1	B	112	ARG
1	B	113	ASN
1	B	115	ASP
1	B	116	GLU
1	B	117	LEU
1	B	118	ARG
1	B	119	LEU
1	B	121	VAL
1	B	122	VAL
1	B	127	GLU
1	B	128	ILE
1	B	129	THR
1	B	135	GLU
1	B	137	ASP
1	B	142	GLN
1	B	143	LYS
1	B	148	GLU
1	B	152	VAL
1	B	155	ASP
1	B	158	VAL
1	B	162	GLN
1	B	164	GLU
1	B	175	LYS
1	B	178	LEU
1	B	183	LEU
1	B	194	LYS
1	B	195	LEU
1	B	197	LEU
1	B	199	ASN
1	B	202	ARG
1	B	203	GLU
1	B	207	LYS
1	B	208	ARG
1	B	219	ARG
1	B	222	VAL

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Mol	Chain	Res	Type
1	B	226	LEU
1	B	227	LYS
1	B	234	LYS
1	B	248	LYS
1	B	250	GLU
1	B	251	SER
1	B	259	GLU
1	B	261	LEU
1	B	269	GLU
1	B	271	ARG
1	B	273	ILE
1	B	275	LEU
1	B	281	LEU
1	B	283	GLN
1	B	286	ARG
1	B	298	LEU
1	B	304	LYS
1	B	305	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	39	GLN
1	A	142	GLN
1	A	162	GLN
1	A	196	ASN
1	A	199	ASN
1	A	257	HIS
1	A	284	ASN
1	B	39	GLN
1	B	142	GLN
1	B	162	GLN
1	B	196	ASN
1	B	199	ASN
1	B	257	HIS
1	B	284	ASN

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	304/305 (99%)	0.52	21 (6%) 23 12	21, 44, 70, 84	0
1	B	304/305 (99%)	0.49	20 (6%) 24 12	18, 43, 69, 84	0
All	All	608/610 (99%)	0.51	41 (6%) 24 12	18, 44, 69, 84	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	60	MET	6.5
1	B	138	MET	6.4
1	A	138	MET	5.7
1	A	60	MET	4.8
1	A	137	ASP	4.5
1	A	193	GLY	3.8
1	B	104	MET	3.7
1	A	104	MET	3.6
1	A	115	ASP	3.5
1	B	156	GLU	3.4
1	B	139	MET	3.3
1	A	289	MET	3.2
1	A	214	ARG	3.1
1	A	2	ILE	3.1
1	A	24	ARG	3.0
1	A	3	GLY	3.0
1	B	145	GLU	2.8
1	B	304	LYS	2.8
1	B	289	MET	2.7
1	A	305	VAL	2.7
1	B	135	GLU	2.6
1	A	26	ASP	2.6
1	B	2	ILE	2.6
1	B	105	GLU	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	137	ASP	2.5
1	A	247	ARG	2.4
1	B	5	ASP	2.4
1	A	99	ARG	2.4
1	A	139	MET	2.4
1	A	105	GLU	2.4
1	B	286	ARG	2.3
1	A	164	GLU	2.3
1	B	123	ASP	2.3
1	A	25	GLY	2.2
1	A	204	GLU	2.2
1	B	177	ASP	2.1
1	B	3	GLY	2.1
1	B	164	GLU	2.1
1	B	214	ARG	2.1
1	B	115	ASP	2.1
1	A	146	LEU	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.