



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:06 PM UTC

PDB ID : 4DFC / pdb_00004dfc
Title : Core UvrA/TRCF complex
Authors : Deaconescu, A.M.; Grigorieff, N.
Deposited on : 2012-01-23
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

<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
Xtrriage (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

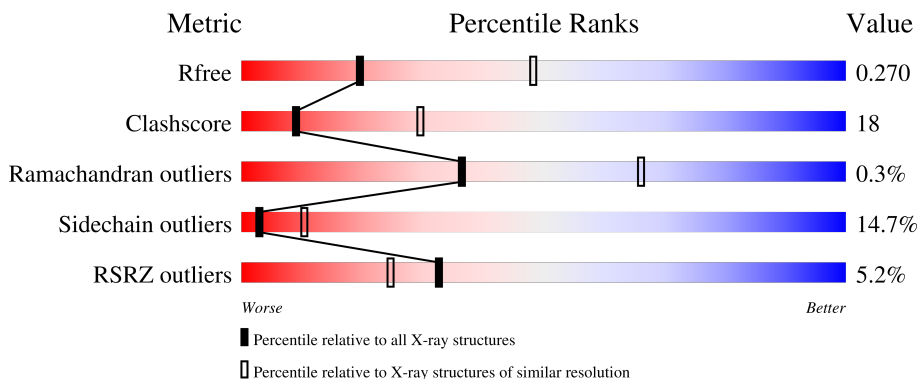
1 Overall quality at a glance i

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}	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (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 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	93	
1	C	93	
2	B	126	
2	D	126	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3143 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 Transcription-repair-coupling factor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	85	680	425	119	133	3	0	0	0
1	C	79	608	378	108	119	3	0	0	0

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	121	GLY	-	expression tag	UNP P30958
A	122	PRO	-	expression tag	UNP P30958
A	123	HIS	-	expression tag	UNP P30958
A	124	MET	-	expression tag	UNP P30958
A	125	ALA	-	expression tag	UNP P30958
A	126	SER	-	expression tag	UNP P30958
C	121	GLY	-	expression tag	UNP P30958
C	122	PRO	-	expression tag	UNP P30958
C	123	HIS	-	expression tag	UNP P30958
C	124	MET	-	expression tag	UNP P30958
C	125	ALA	-	expression tag	UNP P30958
C	126	SER	-	expression tag	UNP P30958

- Molecule 2 is a protein called UvrABC system protein A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	121	925	578	158	185	4	0	0	0
2	D	122	930	579	158	188	5	0	0	0

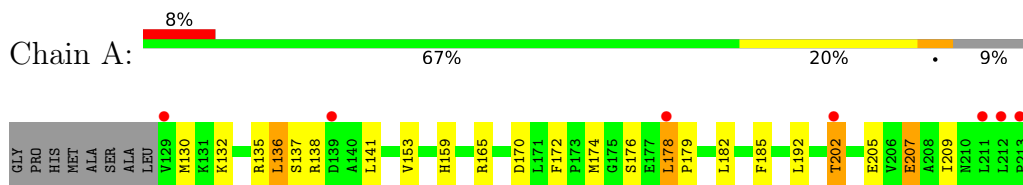
There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	125	GLY	-	expression tag	UNP P0A698
B	126	PRO	-	expression tag	UNP P0A698
B	127	HIS	-	expression tag	UNP P0A698
B	128	MET	-	expression tag	UNP P0A698
B	129	ALA	-	expression tag	UNP P0A698
B	130	SER	-	expression tag	UNP P0A698
D	125	GLY	-	expression tag	UNP P0A698
D	126	PRO	-	expression tag	UNP P0A698
D	127	HIS	-	expression tag	UNP P0A698
D	128	MET	-	expression tag	UNP P0A698
D	129	ALA	-	expression tag	UNP P0A698
D	130	SER	-	expression tag	UNP P0A698

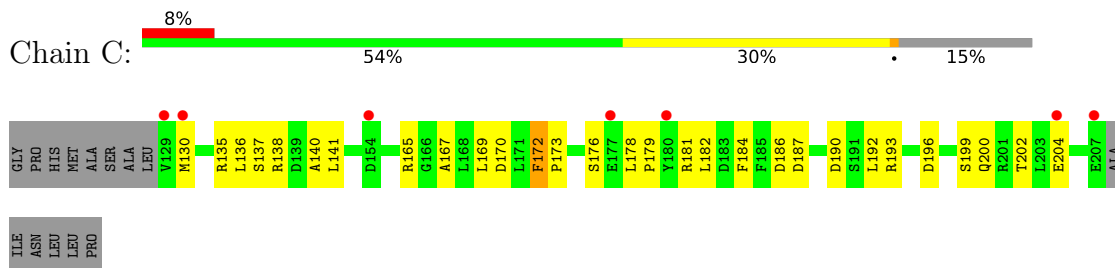
3 Residue-property plots [i](#)

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.

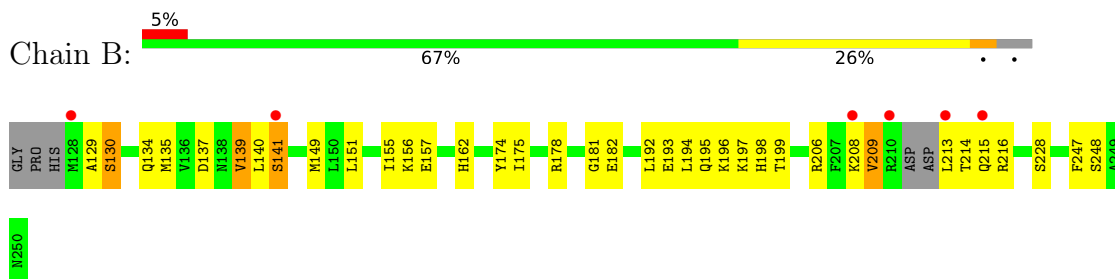
- Molecule 1: Transcription-repair-coupling factor



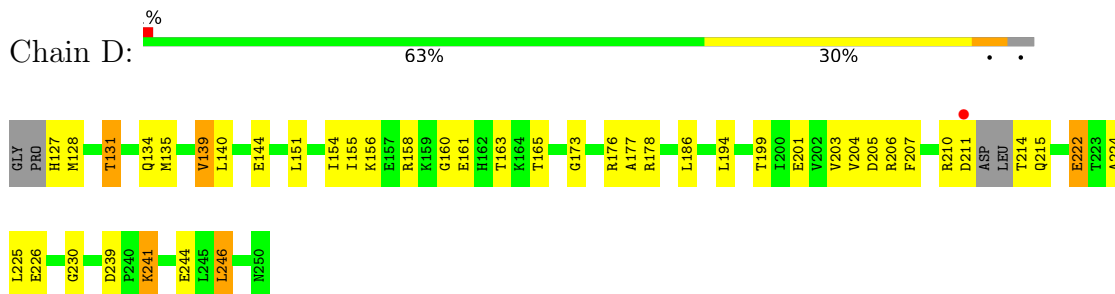
- Molecule 1: Transcription-repair-coupling factor



- Molecule 2: UvrABC system protein A



- Molecule 2: UvrABC system protein A



4 Data and refinement statistics

Property	Value	Source
Space group	P 62 2 2	Depositor
Cell constants a, b, c, α , β , γ	119.17Å 119.17Å 234.33Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.98 – 2.80 29.98 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.2 (29.98-2.80) 91.0 (29.98-2.80)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.59 (at 2.80Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: 1.7.1_743)	Depositor
R, R_{free}	0.238 , 0.282 0.233 , 0.270	Depositor DCC
R_{free} test set	1244 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	69.9	Xtrriage
Anisotropy	0.647	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 58.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3143	wwPDB-VP
Average B, all atoms (Å ²)	94.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.06% 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	0.51	0/691	0.91	1/933 (0.1%)
1	C	0.46	0/618	0.90	4/834 (0.5%)
2	B	0.65	0/934	1.05	6/1260 (0.5%)
2	D	0.60	0/940	0.97	3/1269 (0.2%)
All	All	0.57	0/3183	0.97	14/4296 (0.3%)

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	128	MET	N-CA-C	9.72	122.28	110.91
2	B	208	LYS	N-CA-C	7.11	122.12	113.38
2	B	130	SER	N-CA-C	6.58	119.01	110.53
2	D	139	VAL	N-CA-C	-6.46	100.40	110.09
2	B	141	SER	N-CA-C	-6.19	105.99	112.93
1	C	178	LEU	CA-C-N	6.02	126.22	119.90
1	C	178	LEU	C-N-CA	6.02	126.22	119.90
2	B	139	VAL	N-CA-C	-6.01	101.07	110.09
1	A	136	LEU	N-CA-C	5.65	116.82	107.73
2	B	209	VAL	N-CA-C	5.55	120.89	109.34
2	D	230	GLY	N-CA-C	5.48	122.32	114.64
2	B	228	SER	N-CA-C	-5.46	101.40	109.86
1	C	172	PHE	CA-C-N	5.19	125.44	119.83
1	C	172	PHE	C-N-CA	5.19	125.44	119.83

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	680	0	653	24	0
1	C	608	0	552	22	0
2	B	925	0	929	43	0
2	D	930	0	921	35	0
All	All	3143	0	3055	114	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 18.

All (114) 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:C:137:SER:HB2	1:C:140:ALA:HB3	1.33	1.10
2:B:214:THR:HG22	2:B:215:GLN:H	0.96	1.08
2:B:206:ARG:HG3	2:B:206:ARG:HH11	1.22	0.99
1:C:173:PRO:HG2	1:C:176:SER:HB2	1.45	0.95
2:D:155:ILE:HD11	2:D:165:THR:HG21	1.48	0.94
2:B:214:THR:HG22	2:B:215:GLN:N	1.74	0.92
2:B:214:THR:CG2	2:B:215:GLN:H	1.85	0.89
2:D:160:GLY:HA2	2:D:194:LEU:HD12	1.54	0.88
2:B:213:LEU:HG	2:B:214:THR:H	1.38	0.86
1:C:186:ASP:HB2	2:D:207:PHE:HA	1.57	0.86
1:A:130:MET:HE1	1:A:141:LEU:HD21	1.60	0.83
2:B:193:GLU:OE2	2:B:196:LYS:NZ	2.16	0.78
2:B:130:SER:HB2	2:B:134:GLN:HB2	1.67	0.76
1:A:178:LEU:H	1:A:178:LEU:HD23	1.49	0.75
1:C:137:SER:CB	1:C:140:ALA:HB3	2.16	0.73
2:D:178:ARG:NH2	2:D:244:GLU:OE1	2.23	0.72
2:D:135:MET:CE	2:D:225:LEU:HD11	2.20	0.71
2:B:206:ARG:HH11	2:B:206:ARG:CG	2.02	0.70
2:B:206:ARG:HG3	2:B:206:ARG:NH1	2.02	0.69
1:C:182:LEU:HG	1:C:192:LEU:HD13	1.75	0.69
2:D:139:VAL:O	2:D:140:LEU:HB2	1.91	0.68
2:D:176:ARG:HG2	2:D:203:VAL:HB	1.76	0.68
2:D:239:ASP:OD1	2:D:241:LYS:HG3	1.96	0.66
1:A:178:LEU:HD23	1:A:178:LEU:N	2.11	0.65
2:B:214:THR:O	2:B:216:ARG:N	2.30	0.65
2:D:206:ARG:HH11	2:D:206:ARG:HG3	1.61	0.65
2:D:214:THR:HG23	2:D:215:GLN:H	1.61	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:182:LEU:HG	1:A:192:LEU:HD12	1.77	0.64
1:A:202:THR:HG21	2:B:175:ILE:HG22	1.80	0.64
2:B:214:THR:O	2:B:215:GLN:C	2.40	0.63
2:D:222:GLU:OE2	2:D:222:GLU:HA	2.01	0.60
2:B:194:LEU:HD13	2:B:195:GLN:HG2	1.84	0.59
1:C:172:PHE:CZ	1:C:179:PRO:HD3	2.38	0.58
2:D:131:THR:HG23	2:D:134:GLN:HG3	1.83	0.58
2:B:139:VAL:O	2:B:140:LEU:HB2	2.04	0.58
1:C:167:ALA:HB2	2:D:205:ASP:OD1	2.04	0.58
2:B:130:SER:OG	2:B:135:MET:HG3	2.04	0.57
1:A:182:LEU:HG	1:A:192:LEU:CD1	2.34	0.57
1:A:172:PHE:CD1	1:A:179:PRO:HB3	2.40	0.56
2:D:151:LEU:HD13	2:D:201:GLU:OE1	2.06	0.56
2:D:206:ARG:HG3	2:D:206:ARG:NH1	2.21	0.56
2:B:192:LEU:HB3	2:B:198:HIS:NE2	2.22	0.55
1:C:135:ARG:HG3	1:C:187:ASP:OD1	2.06	0.55
2:D:214:THR:HG23	2:D:215:GLN:N	2.20	0.55
2:B:135:MET:HG2	2:B:247:PHE:HB3	1.89	0.55
1:C:187:ASP:OD1	1:C:187:ASP:C	2.49	0.55
1:A:136:LEU:HD12	1:A:137:SER:H	1.72	0.54
2:D:177:ALA:HB2	2:D:186:LEU:HD11	1.89	0.54
2:B:213:LEU:HG	2:B:214:THR:N	2.17	0.54
1:C:169:LEU:C	1:C:169:LEU:HD12	2.33	0.53
2:D:154:ILE:HG22	2:D:155:ILE:HD12	1.91	0.52
1:C:138:ARG:O	1:C:141:LEU:N	2.34	0.52
1:A:138:ARG:HH11	2:B:216:ARG:NH2	2.08	0.52
1:A:136:LEU:HD12	1:A:137:SER:N	2.25	0.52
1:A:202:THR:CG2	2:B:175:ILE:HG22	2.39	0.52
2:D:135:MET:HE2	2:D:225:LEU:HD11	1.90	0.52
2:B:214:THR:CG2	2:B:215:GLN:N	2.49	0.51
1:C:199:SER:O	1:C:200:GLN:HB2	2.10	0.51
2:D:161:GLU:HG3	2:D:194:LEU:HD13	1.92	0.51
2:B:178:ARG:CD	2:B:181:GLY:HA2	2.40	0.51
2:B:192:LEU:HB3	2:B:198:HIS:CD2	2.47	0.50
1:C:130:MET:HE3	1:C:136:LEU:HD13	1.94	0.50
2:B:156:LYS:HA	2:B:199:THR:HB	1.94	0.49
2:D:144:GLU:OE1	2:D:210:ARG:HA	2.11	0.49
1:C:193:ARG:HD3	1:C:204:GLU:O	2.13	0.48
2:D:151:LEU:HD22	2:D:203:VAL:HG22	1.95	0.48
2:B:135:MET:O	2:B:139:VAL:HG23	2.13	0.48
1:C:169:LEU:HD12	1:C:170:ASP:N	2.29	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:149:MET:HE2	2:B:151:LEU:HD21	1.96	0.47
1:A:165:ARG:NH2	1:A:170:ASP:OD2	2.47	0.47
1:C:181:ARG:NH2	1:C:200:GLN:O	2.46	0.46
2:D:204:VAL:HG21	2:D:224:ALA:HB2	1.97	0.46
2:B:155:ILE:HD13	2:B:162:HIS:CE1	2.52	0.45
2:B:194:LEU:CD1	2:B:195:GLN:HG2	2.45	0.45
1:A:185:PHE:HD2	1:A:185:PHE:C	2.24	0.45
1:A:135:ARG:O	1:A:136:LEU:HB3	2.16	0.45
1:C:165:ARG:NH1	2:D:173:GLY:HA3	2.32	0.45
2:D:156:LYS:O	2:D:158:ARG:NH1	2.49	0.45
2:D:161:GLU:HB3	2:D:163:THR:HG23	1.98	0.45
2:D:178:ARG:NH1	2:D:201:GLU:OE1	2.50	0.45
2:B:137:ASP:O	2:B:141:SER:HB3	2.17	0.45
2:B:178:ARG:HD2	2:B:181:GLY:HA2	1.97	0.45
2:D:131:THR:HG23	2:D:134:GLN:CG	2.46	0.45
2:D:160:GLY:HA2	2:D:194:LEU:CD1	2.37	0.45
1:A:185:PHE:C	1:A:185:PHE:CD2	2.95	0.45
1:A:165:ARG:HG2	2:B:174:TYR:CE1	2.52	0.44
1:A:165:ARG:HG2	2:B:174:TYR:CZ	2.52	0.44
2:D:214:THR:CG2	2:D:215:GLN:H	2.30	0.44
1:A:138:ARG:NH1	2:B:216:ARG:NH2	2.66	0.44
1:A:207:GLU:N	1:A:207:GLU:OE1	2.50	0.44
2:D:210:ARG:CG	2:D:211:ASP:N	2.81	0.43
1:A:178:LEU:N	1:A:178:LEU:CD2	2.79	0.43
2:B:135:MET:HG2	2:B:247:PHE:CB	2.48	0.43
2:B:157:GLU:OE2	2:B:157:GLU:HA	2.19	0.43
2:B:135:MET:HG2	2:B:247:PHE:CG	2.53	0.43
1:A:132:LYS:HD2	1:A:205:GLU:OE2	2.17	0.43
1:A:153:VAL:C	2:D:246:LEU:HD22	2.43	0.42
2:D:131:THR:HG23	2:D:134:GLN:CD	2.44	0.42
2:D:151:LEU:CD2	2:D:203:VAL:HG22	2.49	0.42
1:C:136:LEU:HD12	1:C:136:LEU:HA	1.68	0.42
2:B:178:ARG:HD3	2:B:181:GLY:HA2	2.01	0.42
1:A:172:PHE:HE2	1:A:176:SER:O	2.02	0.42
2:B:129:ALA:HA	2:B:248:SER:O	2.21	0.41
1:C:182:LEU:HD12	1:C:182:LEU:N	2.34	0.41
2:D:158:ARG:HH11	2:D:158:ARG:HG2	1.85	0.41
1:C:167:ALA:HB1	1:C:184:PHE:O	2.20	0.41
2:B:193:GLU:HB2	2:B:196:LYS:HD2	2.01	0.41
1:A:159:HIS:CE1	1:A:172:PHE:CE2	3.08	0.41
2:B:178:ARG:HA	2:B:182:GLU:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:192:LEU:HB3	2:B:198:HIS:CE1	2.56	0.41
2:B:206:ARG:CG	2:B:206:ARG:NH1	2.72	0.40
2:B:157:GLU:O	2:B:197:LYS:HG2	2.22	0.40
1:C:186:ASP:O	1:C:186:ASP:OD1	2.39	0.40
1:C:196:ASP:O	1:C:200:GLN:N	2.50	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	83/93 (89%)	80 (96%)	3 (4%)	0	100	100
1	C	77/93 (83%)	68 (88%)	9 (12%)	0	100	100
2	B	117/126 (93%)	110 (94%)	6 (5%)	1 (1%)	14	41
2	D	118/126 (94%)	116 (98%)	2 (2%)	0	100	100
All	All	395/438 (90%)	374 (95%)	20 (5%)	1 (0%)	36	66

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	209	VAL

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	36/80 (45%)	31 (86%)	5 (14%)	3	12
1	C	9/80 (11%)	7 (78%)	2 (22%)	1	3
2	B	6/109 (6%)	6 (100%)	0	100	100
2	D	44/109 (40%)	37 (84%)	7 (16%)	2	9
All	All	95/378 (25%)	81 (85%)	14 (15%)	3	10

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	174	MET
1	A	178	LEU
1	A	202	THR
1	A	207	GLU
1	A	209	ILE
1	C	190	ASP
1	C	202	THR
2	D	127	HIS
2	D	131	THR
2	D	199	THR
2	D	222	GLU
2	D	226	GLU
2	D	241	LYS
2	D	246	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

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	85/93 (91%)	0.22	7 (8%) 17 13	60, 91, 133, 150	0
1	C	79/93 (84%)	0.61	7 (8%) 15 11	76, 123, 165, 179	0
2	B	121/126 (96%)	-0.03	6 (4%) 34 26	59, 81, 115, 160	0
2	D	122/126 (96%)	-0.04	1 (0%) 82 75	57, 81, 115, 140	0
All	All	407/438 (92%)	0.14	21 (5%) 33 25	57, 89, 141, 179	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	204	GLU	5.2
1	C	129	VAL	5.1
2	B	128	MET	4.2
2	D	211	ASP	3.7
1	A	129	VAL	3.4
2	B	210	ARG	3.3
1	C	180	TYR	3.3
1	C	207	GLU	3.1
1	A	212	LEU	3.1
1	C	177	GLU	2.8
2	B	215	GLN	2.8
2	B	208	LYS	2.7
1	A	202	THR	2.7
1	C	130	MET	2.6
1	C	154	ASP	2.6
2	B	141	SER	2.4
2	B	213	LEU	2.3
1	A	213	PRO	2.2
1	A	139	ASP	2.2
1	A	178	LEU	2.2
1	A	211	LEU	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.