Full wwPDB X-ray Structure Validation Report ⓘ

Jan 30, 2024 – 10:05 AM EST

PDB ID	:	1FO1
Title	:	CRYSTAL STRUCTURE OF THE RNA-BINDING DOMAIN OF THE MRNA EXPORT FACTOR TAP
Authors	:	Liker, E.; Fernandez, E.; Izaurralde, E.; Conti, E.
Deposited on	:	2000-08-24
Resolution	:	2.90 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:

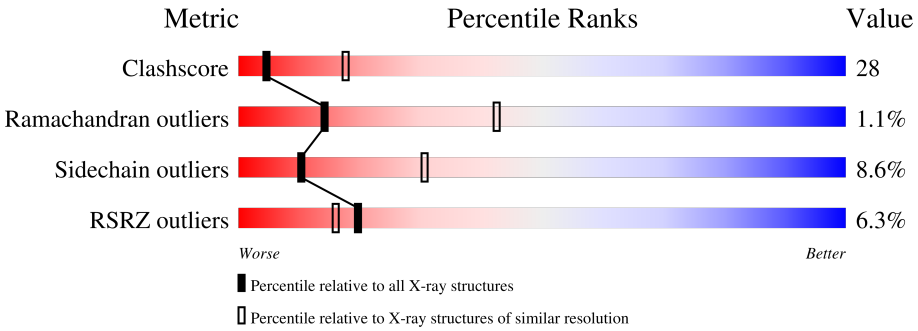
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 2.90 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))
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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 <=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	271	8% 41% 32% 6% 21%
1	B	271	. 37% 21% . 40%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 3079 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 NUCLEAR RNA EXPORT FACTOR 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	215	Total	C	N	O	S	0	0	0
			1751	1110	309	325	7			
1	B	163	Total	C	N	O	S	0	0	0
			1311	823	230	253	5			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	226	VAL	ALA	conflict	UNP Q9UBU9
B	226	VAL	ALA	conflict	UNP Q9UBU9

- Molecule 2 is water.

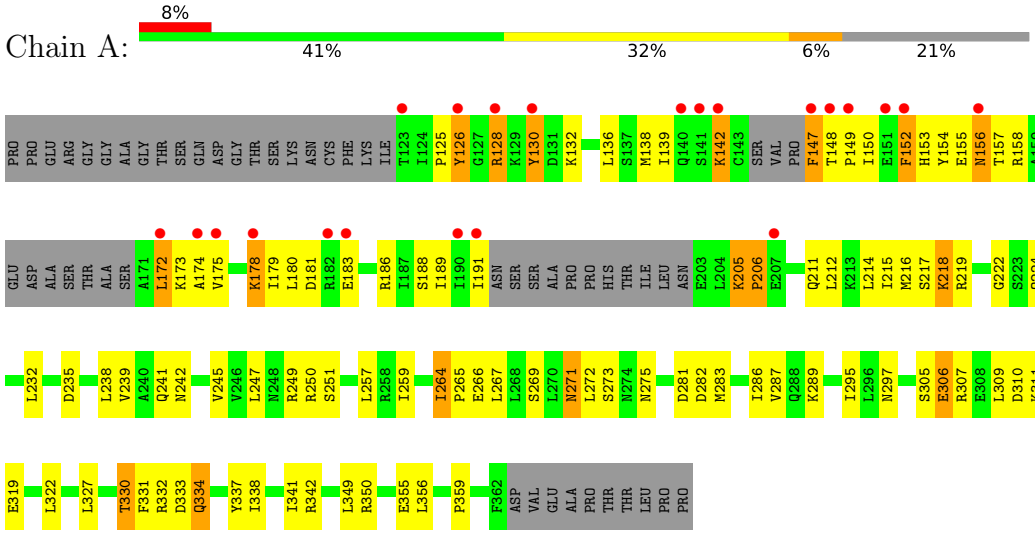
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	10	Total	O	0	0
			10	10		
2	B	7	Total	O	0	0
			7	7		

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.

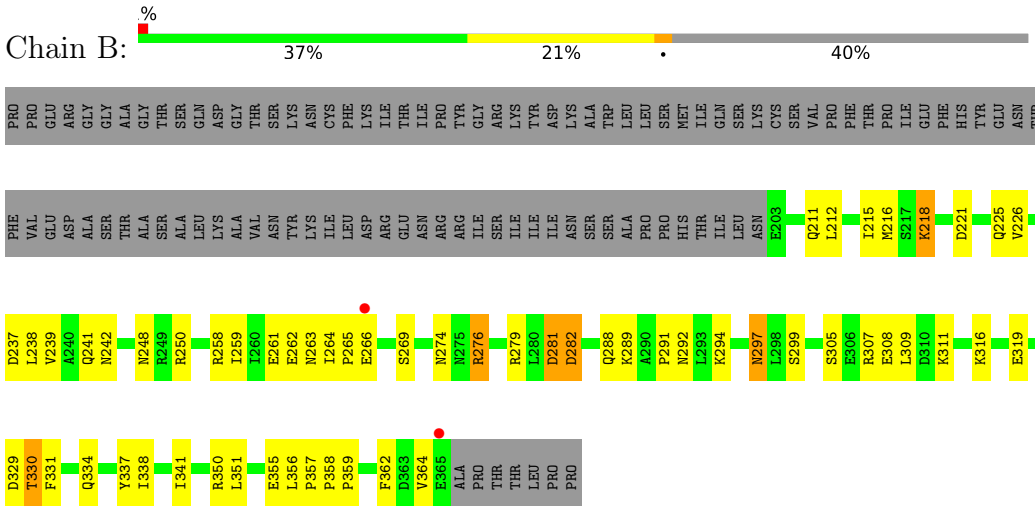
• Molecule 1: NUCLEAR RNA EXPORT FACTOR 1

Chain A:



• Molecule 1: NUCLEAR RNA EXPORT FACTOR 1

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α, β, γ	96.35Å 96.35Å 152.23Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.90 29.88 – 2.90	Depositor EDS
% Data completeness (in resolution range)	(Not available) (30.00-2.90) 98.7 (29.88-2.90)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.79 (at 2.90Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.250 , 0.278 0.242 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å^2)	65.0	Xtriage
Anisotropy	0.283	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 45.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.91	EDS
Total number of atoms	3079	wwPDB-VP
Average B, all atoms (Å^2)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.55% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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	0.38	0/1775	0.62	0/2388
1	B	0.39	0/1327	0.61	0/1789
All	All	0.38	0/3102	0.62	0/4177

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1751	0	1810	118	0
1	B	1311	0	1351	58	0
2	A	10	0	0	0	0
2	B	7	0	0	4	0
All	All	3079	0	3161	176	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 28.

All (176) 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:172:LEU:O	1:A:175:VAL:HG12	1.36	1.24

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:142:LYS:HG2	1:A:178:LYS:HD3	1.25	1.14
1:A:126:TYR:HA	1:A:157:THR:HG22	1.34	1.09
1:A:211:GLN:HG3	1:A:241:GLN:HE22	1.11	1.07
1:A:172:LEU:HD12	1:A:191:ILE:HD11	1.41	1.00
1:A:211:GLN:HG3	1:A:241:GLN:NE2	1.78	0.95
1:A:212:LEU:HG	1:A:216:MET:HE3	1.50	0.93
1:A:142:LYS:CG	1:A:178:LYS:HD3	2.01	0.90
1:A:211:GLN:CG	1:A:241:GLN:HE22	1.85	0.89
1:A:175:VAL:HG13	1:A:189:ILE:HD12	1.53	0.88
1:A:175:VAL:HG23	1:A:178:LYS:CE	2.06	0.86
1:A:342:ARG:HH21	1:A:356:LEU:HB3	1.43	0.83
1:A:273:SER:HB3	1:A:297:ASN:ND2	1.94	0.83
1:A:218:LYS:HE3	1:A:218:LYS:HA	1.63	0.80
1:A:128:ARG:HA	1:A:154:TYR:CE2	2.17	0.79
1:A:180:LEU:HD21	1:A:186:ARG:CZ	2.12	0.79
1:B:226:VAL:HG22	1:B:269:SER:HB3	1.64	0.79
1:A:334:GLN:NE2	1:A:334:GLN:H	1.81	0.79
1:A:172:LEU:O	1:A:175:VAL:CG1	2.27	0.77
1:B:288:GLN:HG2	2:B:26:HOH:O	1.83	0.77
1:A:128:ARG:HA	1:A:154:TYR:HE2	1.50	0.76
1:A:175:VAL:HA	1:A:178:LYS:HZ3	1.51	0.75
1:A:235:ASP:O	1:A:239:VAL:HG23	1.89	0.71
1:A:180:LEU:CD2	1:A:186:ARG:HG2	2.21	0.71
1:A:219:ARG:HD3	1:A:228:ASP:O	1.91	0.70
1:B:322:LEU:HD12	1:B:341:ILE:HD11	1.72	0.70
1:B:212:LEU:HG	1:B:216:MET:HE2	1.74	0.69
1:A:250:ARG:HG3	1:A:282:ASP:OD1	1.93	0.69
1:A:175:VAL:CG1	1:A:189:ILE:HD12	2.23	0.68
1:B:212:LEU:HG	1:B:216:MET:CE	2.24	0.68
1:A:327:LEU:O	1:A:330:THR:HG22	1.95	0.66
1:A:271:ASN:ND2	1:A:273:SER:H	1.94	0.66
1:A:311:LYS:HD3	2:B:25:HOH:O	1.94	0.66
1:A:310:ASP:HA	1:A:313:LYS:HG3	1.77	0.66
1:A:175:VAL:HG23	1:A:178:LYS:NZ	2.11	0.66
1:A:179:ILE:O	1:A:180:LEU:HD23	1.96	0.66
1:A:150:ILE:HG22	1:A:150:ILE:O	1.96	0.64
1:A:212:LEU:HG	1:A:216:MET:CE	2.26	0.63
1:A:136:LEU:HD11	1:A:147:PHE:CE1	2.33	0.63
1:B:276:ARG:HH11	1:B:276:ARG:HB3	1.63	0.63
1:A:175:VAL:HG23	1:A:178:LYS:HE2	1.81	0.62
1:A:180:LEU:HD22	1:A:186:ARG:HG2	1.81	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:289:LYS:C	1:B:291:PRO:HD3	2.19	0.62
1:A:132:LYS:HG3	1:A:154:TYR:OH	2.00	0.62
1:B:215:ILE:HD13	1:B:238:LEU:HG	1.80	0.61
1:B:230:LYS:HG3	1:B:274:ASN:HD22	1.66	0.61
1:A:205:LYS:HG3	1:A:206:PRO:HD2	1.84	0.59
1:B:316:LYS:NZ	1:B:316:LYS:HB2	2.17	0.59
1:B:319:GLU:HG3	1:B:350:ARG:HB2	1.85	0.59
1:A:214:LEU:O	1:A:217:SER:HB3	2.02	0.58
1:B:309:LEU:HD11	1:B:322:LEU:HD11	1.84	0.58
1:A:132:LYS:HB2	1:A:132:LYS:NZ	2.18	0.58
1:B:250:ARG:HA	1:B:282:ASP:OD2	2.03	0.58
1:A:142:LYS:HB3	1:A:178:LYS:HZ2	1.68	0.58
1:B:289:LYS:O	1:B:291:PRO:HD3	2.04	0.58
1:B:216:MET:HE1	1:B:259:ILE:HD12	1.85	0.57
1:B:305:SER:OG	1:B:307:ARG:HG2	2.04	0.57
1:A:271:ASN:ND2	1:A:297:ASN:HD22	2.03	0.57
1:A:264:ILE:N	1:A:265:PRO:HD3	2.19	0.57
1:A:178:LYS:HD2	1:A:178:LYS:C	2.26	0.56
1:A:271:ASN:HD22	1:A:272:LEU:N	2.03	0.56
1:A:175:VAL:CA	1:A:178:LYS:HZ3	2.18	0.56
1:A:309:LEU:HD11	1:A:322:LEU:HD11	1.85	0.56
1:A:216:MET:HE1	1:A:259:ILE:HD12	1.88	0.56
1:A:313:LYS:NZ	1:A:313:LYS:HB3	2.22	0.55
1:A:273:SER:HB3	1:A:297:ASN:HD22	1.66	0.55
1:B:322:LEU:HD12	1:B:341:ILE:CD1	2.36	0.55
1:A:273:SER:HB3	1:A:297:ASN:HD21	1.72	0.55
1:A:211:GLN:O	1:A:215:ILE:HD13	2.08	0.54
1:B:225:GLN:NE2	1:B:266:GLU:HG3	2.22	0.54
1:A:147:PHE:CG	1:A:147:PHE:O	2.61	0.53
1:A:216:MET:CE	1:A:259:ILE:HD12	2.38	0.53
1:A:271:ASN:HD21	1:A:297:ASN:HD22	1.55	0.53
1:B:288:GLN:NE2	2:B:26:HOH:O	2.38	0.53
1:A:205:LYS:HE3	1:A:206:PRO:HD2	1.91	0.53
1:A:235:ASP:HB3	1:A:238:LEU:HD12	1.92	0.52
1:A:342:ARG:NH2	1:A:356:LEU:HB3	2.19	0.52
1:A:342:ARG:NH1	1:A:359:PRO:HG3	2.24	0.52
1:B:316:LYS:HB2	1:B:316:LYS:HZ2	1.75	0.52
1:A:205:LYS:HA	1:A:205:LYS:NZ	2.25	0.52
1:A:128:ARG:C	1:A:130:TYR:H	2.11	0.51
1:B:292:ASN:O	1:B:294:LYS:HG3	2.10	0.51
1:A:152:PHE:C	1:A:152:PHE:CD1	2.84	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:211:GLN:HG3	1:B:241:GLN:NE2	2.24	0.51
1:A:142:LYS:CB	1:A:178:LYS:NZ	2.75	0.50
1:B:216:MET:HE1	1:B:259:ILE:CD1	2.41	0.50
1:B:218:LYS:HA	1:B:218:LYS:NZ	2.25	0.50
1:A:139:ILE:HA	1:A:179:ILE:HD11	1.93	0.50
1:A:211:GLN:OE1	1:A:241:GLN:OE1	2.30	0.49
1:A:178:LYS:HE3	1:A:179:ILE:HG13	1.93	0.49
1:A:334:GLN:O	1:A:338:ILE:HG13	2.12	0.49
1:A:128:ARG:HG2	1:A:154:TYR:CD2	2.47	0.49
1:A:267:LEU:HD11	1:A:269:SER:O	2.12	0.49
1:B:211:GLN:CD	1:B:241:GLN:HE22	2.14	0.49
1:B:279:ARG:O	1:B:279:ARG:HG3	2.11	0.49
1:A:142:LYS:HB3	1:A:178:LYS:NZ	2.28	0.48
1:A:126:TYR:HA	1:A:157:THR:CG2	2.25	0.48
1:A:295:ILE:HG12	1:A:319:GLU:HB3	1.94	0.48
1:A:319:GLU:OE1	1:A:350:ARG:NH1	2.46	0.48
1:A:281:ASP:HB3	1:A:311:LYS:HG3	1.95	0.48
1:A:247:LEU:C	1:A:249:ARG:H	2.16	0.48
1:A:306:GLU:HG2	1:A:331:PHE:CZ	2.49	0.48
1:A:286:ILE:HG23	1:A:287:VAL:N	2.29	0.47
1:A:126:TYR:CG	1:A:126:TYR:O	2.67	0.47
1:A:319:GLU:HG3	1:A:350:ARG:HB2	1.97	0.46
1:A:333:ASP:OD2	1:A:333:ASP:O	2.33	0.46
1:A:178:LYS:CE	1:A:179:ILE:HG13	2.46	0.46
1:B:216:MET:CE	1:B:259:ILE:CD1	2.93	0.46
1:B:334:GLN:O	1:B:338:ILE:HG13	2.16	0.46
1:B:211:GLN:CG	1:B:241:GLN:HE22	2.28	0.46
1:A:136:LEU:HD21	1:A:147:PHE:CE1	2.51	0.46
1:B:221:ASP:O	1:B:225:GLN:N	2.48	0.46
1:B:226:VAL:HG22	1:B:269:SER:CB	2.40	0.46
1:B:261:GLU:O	1:B:265:PRO:HG3	2.15	0.46
1:A:130:TYR:HE1	1:A:181:ASP:OD1	1.99	0.46
1:B:288:GLN:CG	2:B:26:HOH:O	2.53	0.46
1:B:212:LEU:CG	1:B:216:MET:HE2	2.45	0.45
1:A:147:PHE:O	1:A:148:THR:OG1	2.32	0.45
1:B:258:ARG:O	1:B:262:GLU:HG3	2.17	0.45
1:A:132:LYS:HB2	1:A:132:LYS:HZ2	1.82	0.45
1:A:157:THR:HG22	1:A:157:THR:O	2.17	0.45
1:A:175:VAL:O	1:A:178:LYS:HB3	2.16	0.45
1:A:205:LYS:HA	1:A:205:LYS:HZ1	1.80	0.45
1:A:318:GLU:HG2	1:A:349:LEU:HD12	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:358:PRO:HA	1:B:359:PRO:HD3	1.92	0.44
1:A:306:GLU:H	1:A:306:GLU:HG3	1.36	0.44
1:B:308:GLU:OE1	1:B:308:GLU:HA	2.17	0.44
1:A:172:LEU:O	1:A:174:ALA:N	2.51	0.44
1:A:172:LEU:C	1:A:174:ALA:H	2.21	0.44
1:A:305:SER:OG	1:A:307:ARG:HG2	2.16	0.44
1:A:211:GLN:O	1:A:214:LEU:HB3	2.18	0.44
1:A:232:LEU:HD23	1:A:245:VAL:HG11	1.99	0.44
1:A:272:LEU:C	1:A:275:ASN:HD22	2.22	0.44
1:B:216:MET:CE	1:B:259:ILE:HD12	2.47	0.44
1:B:281:ASP:HA	1:B:311:LYS:HB3	2.00	0.44
1:A:156:ASN:HB3	1:A:157:THR:H	1.57	0.43
1:B:327:LEU:O	1:B:330:THR:HB	2.18	0.43
1:A:136:LEU:C	1:A:136:LEU:HD13	2.37	0.43
1:A:181:ASP:C	1:A:183:GLU:H	2.20	0.43
1:A:172:LEU:C	1:A:174:ALA:N	2.72	0.43
1:A:337:TYR:O	1:A:341:ILE:HG12	2.19	0.43
1:B:350:ARG:NH1	1:B:355:GLU:HB2	2.33	0.43
1:A:126:TYR:HH	1:A:181:ASP:CG	2.22	0.43
1:A:152:PHE:CD1	1:A:153:HIS:N	2.87	0.43
1:A:142:LYS:HB2	1:A:178:LYS:NZ	2.34	0.43
1:A:173:LYS:HB3	1:A:191:ILE:HD12	2.00	0.43
1:B:237:ASP:O	1:B:241:GLN:HG3	2.19	0.43
1:B:328:CYS:C	1:B:330:THR:H	2.22	0.42
1:A:178:LYS:HE2	1:A:178:LYS:HB3	1.75	0.42
1:A:283:MET:O	1:A:286:ILE:HB	2.19	0.42
1:A:331:PHE:CE2	1:A:337:TYR:HA	2.55	0.42
1:B:297:ASN:HD22	1:B:299:SER:H	1.68	0.42
1:A:128:ARG:C	1:A:130:TYR:N	2.72	0.42
1:B:276:ARG:HB3	1:B:276:ARG:NH1	2.34	0.42
1:A:125:PRO:O	1:A:126:TYR:CD1	2.73	0.42
1:A:153:HIS:CD2	1:A:155:GLU:HG3	2.55	0.41
1:A:224:GLN:HB2	1:A:226:VAL:HG23	2.02	0.41
1:B:235:ASP:O	1:B:239:VAL:HG23	2.19	0.41
1:B:263:ASN:C	1:B:265:PRO:HD3	2.40	0.41
1:B:337:TYR:CE2	1:B:351:LEU:HD21	2.55	0.41
1:B:211:GLN:HG3	1:B:241:GLN:HE22	1.85	0.41
1:B:216:MET:HE3	1:B:259:ILE:HD13	2.02	0.41
1:B:250:ARG:CA	1:B:282:ASP:OD2	2.67	0.41
1:B:264:ILE:N	1:B:265:PRO:HD3	2.35	0.41
1:A:257:LEU:HB3	1:A:289:LYS:HD2	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:138:MET:HG2	1:A:179:ILE:CG2	2.51	0.41
1:A:225:GLN:HE22	1:A:266:GLU:HB3	1.84	0.41
1:A:332:ARG:HG3	1:A:332:ARG:HH11	1.85	0.41
1:B:338:ILE:HD13	1:B:356:LEU:HD22	2.03	0.40
1:B:356:LEU:HA	1:B:357:PRO:HD3	1.91	0.40
1:A:175:VAL:CB	1:A:178:LYS:HZ3	2.34	0.40
1:B:266:GLU:HG3	1:B:266:GLU:O	2.21	0.40
1:B:330:THR:HG22	1:B:331:PHE:CD1	2.57	0.40
1:B:331:PHE:CD2	1:B:337:TYR:HA	2.57	0.40
1:A:156:ASN:HD22	1:A:156:ASN:HA	1.56	0.40
1:B:248:ASN:C	1:B:248:ASN:OD1	2.60	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	207/271 (76%)	181 (87%)	23 (11%)	3 (1%)	11	36
1	B	161/271 (59%)	147 (91%)	13 (8%)	1 (1%)	25	58
All	All	368/542 (68%)	328 (89%)	36 (10%)	4 (1%)	14	42

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	149	PRO
1	B	364	VAL
1	A	222	GLY
1	A	206	PRO

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	198/245 (81%)	177 (89%)	21 (11%)	6 20
1	B	151/245 (62%)	142 (94%)	9 (6%)	19 49
All	All	349/490 (71%)	319 (91%)	30 (9%)	10 30

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	126	TYR
1	A	128	ARG
1	A	130	TYR
1	A	142	LYS
1	A	147	PHE
1	A	152	PHE
1	A	156	ASN
1	A	158	ARG
1	A	172	LEU
1	A	178	LYS
1	A	188	SER
1	A	205	LYS
1	A	218	LYS
1	A	242	ASN
1	A	251	SER
1	A	264	ILE
1	A	271	ASN
1	A	306	GLU
1	A	330	THR
1	A	334	GLN
1	A	355	GLU
1	B	218	LYS
1	B	242	ASN
1	B	276	ARG
1	B	281	ASP
1	B	282	ASP
1	B	297	ASN

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Mol	Chain	Res	Type
1	B	329	ASP
1	B	330	THR
1	B	362	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	156	ASN
1	A	225	GLN
1	A	241	GLN
1	A	242	ASN
1	A	271	ASN
1	A	297	ASN
1	A	334	GLN
1	B	225	GLN
1	B	241	GLN
1	B	242	ASN
1	B	274	ASN
1	B	297	ASN

5.3.3 RNA

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains

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

5.5 Carbohydrates

There are no monosaccharides in this entry.

5.6 Ligand geometry

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, 95^{th} 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	215/271 (79%)	0.35	22 (10%) 6 5	36, 60, 101, 101	0
1	B	163/271 (60%)	-0.19	2 (1%) 79 79	31, 49, 78, 99	0
All	All	378/542 (69%)	0.11	24 (6%) 20 16	31, 54, 101, 101	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	141	SER	6.4
1	A	149	PRO	5.4
1	A	147	PHE	5.1
1	A	172	LEU	5.0
1	A	142	LYS	5.0
1	A	126	TYR	4.7
1	A	151	GLU	4.2
1	A	174	ALA	3.3
1	A	178	LYS	3.2
1	A	156	ASN	3.1
1	A	130	TYR	3.0
1	A	175	VAL	3.0
1	A	182	ARG	3.0
1	B	365	GLU	3.0
1	A	190	ILE	2.8
1	A	123	THR	2.7
1	A	140	GLN	2.7
1	A	191	ILE	2.5
1	A	152	PHE	2.3
1	A	128	ARG	2.3
1	A	148	THR	2.3
1	A	207	GLU	2.2
1	A	183	GLU	2.1
1	B	266	GLU	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 monosaccharides 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.