



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:53 PM GMT

PDB ID : 1R1R
Title : RIBONUCLEOTIDE REDUCTASE R1 PROTEIN MUTANT Y730F WITH
A REDUCED ACTIVE SITE FROM ESCHERICHIA COLI
Authors : Eriksson, M.; Eklund, H.
Deposited on : 1997-07-15
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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

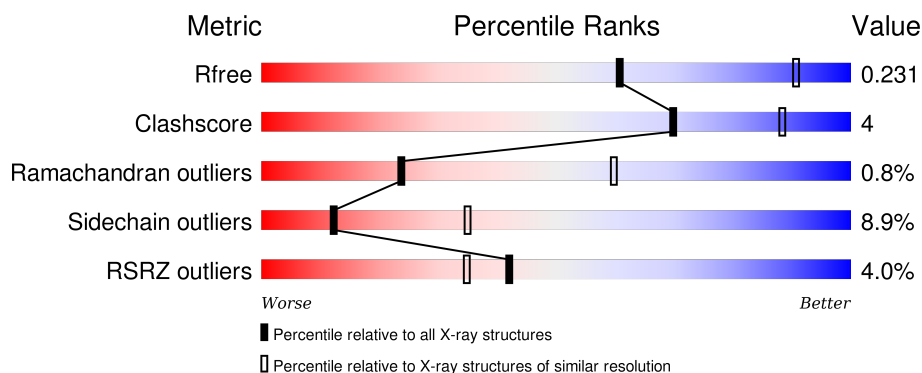
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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))
R_{free}	91344	1451 (2.90-2.90)
Clashscore	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (2.90-2.90)
RSRZ outliers	91569	1456 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	761	<div> <div>4%</div> <div>79%</div> <div>16%</div> <div>..</div> </div>
1	B	761	<div> <div>4%</div> <div>79%</div> <div>16%</div> <div>..</div> </div>
1	C	761	<div> <div>3%</div> <div>79%</div> <div>16%</div> <div>..</div> </div>
2	D	20	<div> <div>15%</div> <div>60%</div> <div>20%</div> <div>20%</div> </div>
2	E	20	<div> <div>10%</div> <div>60%</div> <div>20%</div> <div>20%</div> </div>

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Mol	Chain	Length	Quality of chain
2	F	20	<div><div>5%</div><div>60%</div><div>20%</div><div>20%</div></div>
2	P	20	<div><div>5%</div><div>5%</div><div>5%</div><div>85%</div></div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 18258 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 RIBONUCLEOTIDE REDUCTASE R1 PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	734	Total	C	N	O	S	0	0	0
			5844	3712	1004	1104	24			
1	B	734	Total	C	N	O	S	0	0	0
			5844	3712	1004	1104	24			
1	C	734	Total	C	N	O	S	0	0	0
			5844	3712	1004	1104	24			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	730	PHE	TYR	ENGINEERED	UNP P00452
B	730	PHE	TYR	ENGINEERED	UNP P00452
C	730	PHE	TYR	ENGINEERED	UNP P00452

- Molecule 2 is a protein called RIBONUCLEOTIDE REDUCTASE R2 PROTEIN.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	D	16	Total	C	N	O	0	0	0
			129	77	19	33			
2	E	16	Total	C	N	O	0	0	0
			129	77	19	33			
2	F	16	Total	C	N	O	0	0	0
			129	77	19	33			
2	P	3	Total	C	N	O	0	0	0
			27	20	3	4			

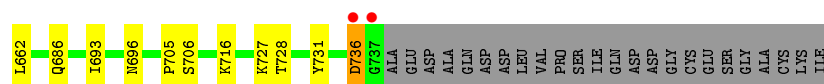
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	104	Total	O	0	0
			104	104		

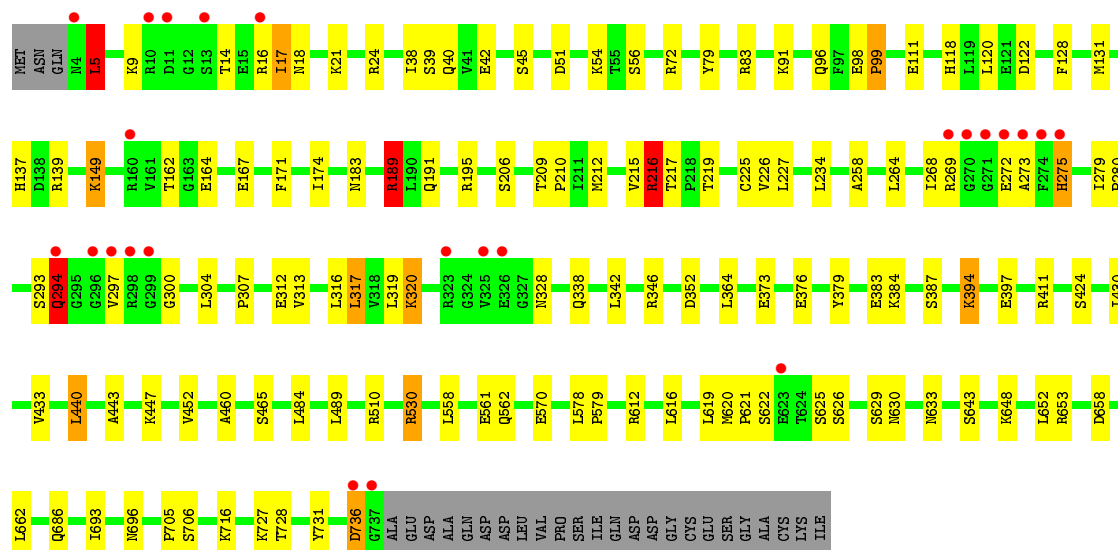
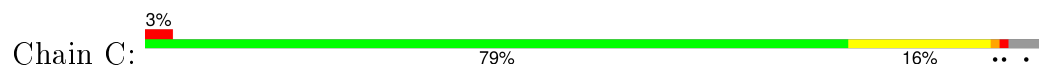
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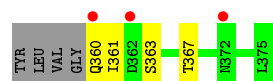
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	104	Total 104	O 104	0	0
3	C	104	Total 104	O 104	0	0



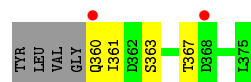
• Molecule 1: RIBONUCLEOTIDE REDUCTASE R1 PROTEIN



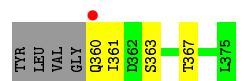
• Molecule 2: RIBONUCLEOTIDE REDUCTASE R2 PROTEIN



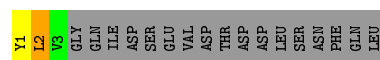
• Molecule 2: RIBONUCLEOTIDE REDUCTASE R2 PROTEIN



• Molecule 2: RIBONUCLEOTIDE REDUCTASE R2 PROTEIN



• Molecule 2: RIBONUCLEOTIDE REDUCTASE R2 PROTEIN



4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	227.82Å 227.82Å 343.47Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.90 20.00 – 2.90	Depositor EDS
% Data completeness (in resolution range)	96.1 (20.00-2.90) 96.1 (20.00-2.90)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.92 (at 2.88Å)	Xtriage
Refinement program	REFMAC, TNT	Depositor
R, R_{free}	0.210 , 0.245 0.216 , 0.231	Depositor DCC
R_{free} test set	2008 reflections (2.85%)	DCC
Wilson B-factor (Å ²)	46.3	Xtriage
Anisotropy	0.045	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 46.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 72549 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	18258	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.90% 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, 0.375 respectively for untwinned datasets, and 0.333, 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	0.36	0/5972	0.88	5/8088 (0.1%)
1	B	0.36	0/5972	0.88	5/8088 (0.1%)
1	C	0.36	0/5972	0.88	5/8088 (0.1%)
2	D	0.36	0/129	0.75	0/173
2	E	0.36	0/129	0.75	0/173
2	F	0.36	0/129	0.75	0/173
2	P	0.72	0/27	1.81	1/36 (2.8%)
All	All	0.36	0/18330	0.88	16/24819 (0.1%)

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	C	653	ARG	NE-CZ-NH1	7.81	124.20	120.30
1	B	653	ARG	NE-CZ-NH1	7.75	124.17	120.30
1	A	653	ARG	NE-CZ-NH1	7.74	124.17	120.30
1	B	653	ARG	NE-CZ-NH2	-6.20	117.20	120.30
1	C	653	ARG	NE-CZ-NH2	-6.17	117.22	120.30
1	A	653	ARG	NE-CZ-NH2	-6.07	117.27	120.30
2	P	2	LEU	CA-CB-CG	5.87	128.79	115.30
1	A	530	ARG	NE-CZ-NH1	5.48	123.04	120.30
1	C	189	ARG	CD-NE-CZ	5.42	131.19	123.60
1	B	189	ARG	CD-NE-CZ	5.42	131.18	123.60
1	B	530	ARG	NE-CZ-NH1	5.41	123.01	120.30
1	C	530	ARG	NE-CZ-NH1	5.41	123.00	120.30
1	A	189	ARG	CD-NE-CZ	5.40	131.16	123.60
1	B	736	ASP	CA-CB-CG	5.21	124.86	113.40
1	A	736	ASP	CA-CB-CG	5.20	124.84	113.40
1	C	736	ASP	CA-CB-CG	5.19	124.83	113.40

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	5844	0	5770	54	0
1	B	5844	0	5770	53	0
1	C	5844	0	5770	54	0
2	D	129	0	111	0	0
2	E	129	0	111	0	0
2	F	129	0	111	0	0
2	P	27	0	31	1	0
3	A	104	0	0	0	0
3	B	104	0	0	0	0
3	C	104	0	0	0	0
All	All	18258	0	17674	160	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 4.

All (160) 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:394:LYS:HB3	1:A:397:GLU:HG3	1.68	0.76
1:C:215:VAL:O	1:C:216:ARG:HB3	1.85	0.76
1:B:394:LYS:HB3	1:B:397:GLU:HG3	1.68	0.75
1:C:394:LYS:HB3	1:C:397:GLU:HG3	1.68	0.75
1:A:215:VAL:O	1:A:216:ARG:HB3	1.85	0.75
1:B:215:VAL:O	1:B:216:ARG:HB3	1.85	0.74
1:A:279:ILE:HB	1:A:280:PRO:HD3	1.70	0.73
1:B:279:ILE:HB	1:B:280:PRO:HD3	1.70	0.73
1:C:279:ILE:HB	1:C:280:PRO:HD3	1.70	0.73
1:A:149:LYS:HG2	1:A:652:LEU:HD21	1.74	0.69
1:B:227:LEU:HB2	1:B:460:ALA:HB3	1.75	0.69
1:C:227:LEU:HB2	1:C:460:ALA:HB3	1.75	0.68
1:B:149:LYS:HG2	1:B:652:LEU:HD21	1.74	0.68
1:C:149:LYS:HG2	1:C:652:LEU:HD21	1.74	0.68
1:B:313:VAL:HG22	1:B:317:LEU:HD22	1.77	0.67
1:A:313:VAL:HG22	1:A:317:LEU:HD22	1.77	0.67
1:A:227:LEU:HB2	1:A:460:ALA:HB3	1.75	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:313:VAL:HG22	1:C:317:LEU:HD22	1.77	0.66
1:A:5:LEU:HD22	1:A:17:ILE:HG13	1.79	0.64
1:A:619:LEU:HD12	1:A:693:ILE:HG12	1.80	0.63
1:B:561:GLU:HG2	1:B:562:GLN:HG3	1.81	0.63
1:C:5:LEU:HD22	1:C:17:ILE:HG13	1.79	0.63
1:C:561:GLU:HG2	1:C:562:GLN:HG3	1.81	0.63
1:B:619:LEU:HD12	1:B:693:ILE:HG12	1.80	0.63
1:B:5:LEU:HD22	1:B:17:ILE:HG13	1.79	0.62
1:A:658:ASP:HB3	1:A:662:LEU:HD12	1.82	0.62
1:B:658:ASP:HB3	1:B:662:LEU:HD12	1.82	0.62
1:C:619:LEU:HD12	1:C:693:ILE:HG12	1.80	0.62
1:C:658:ASP:HB3	1:C:662:LEU:HD12	1.82	0.62
1:A:561:GLU:HG2	1:A:562:GLN:HG3	1.81	0.61
1:A:279:ILE:HD12	1:A:319:LEU:HD21	1.83	0.60
1:A:558:LEU:HD23	1:A:612:ARG:HG2	1.83	0.60
1:B:279:ILE:HD12	1:B:319:LEU:HD21	1.84	0.59
1:A:5:LEU:O	1:A:17:ILE:HB	2.03	0.59
1:B:5:LEU:O	1:B:17:ILE:HB	2.03	0.59
1:C:279:ILE:HD12	1:C:319:LEU:HD21	1.84	0.59
1:C:5:LEU:O	1:C:17:ILE:HB	2.02	0.59
1:B:558:LEU:HD23	1:B:612:ARG:HG2	1.84	0.59
1:C:558:LEU:HD23	1:C:612:ARG:HG2	1.84	0.58
1:B:342:LEU:HD13	1:B:376:GLU:HG3	1.86	0.58
1:C:342:LEU:HD13	1:C:376:GLU:HG3	1.86	0.56
1:A:342:LEU:HD13	1:A:376:GLU:HG3	1.86	0.56
1:A:578:LEU:HB3	1:A:579:PRO:HD2	1.88	0.56
1:B:578:LEU:HB3	1:B:579:PRO:HD2	1.88	0.55
1:C:578:LEU:HB3	1:C:579:PRO:HD2	1.88	0.55
1:C:167:GLU:OE2	1:C:216:ARG:NH1	2.41	0.54
1:A:167:GLU:OE2	1:A:216:ARG:NH1	2.41	0.53
1:B:167:GLU:OE2	1:B:216:ARG:NH1	2.41	0.53
1:B:430:ILE:HG21	1:B:570:GLU:HG2	1.89	0.53
1:A:430:ILE:HG21	1:A:570:GLU:HG2	1.90	0.53
1:C:430:ILE:HG21	1:C:570:GLU:HG2	1.90	0.53
1:C:120:LEU:HD13	2:P:1:TYR:CG	2.44	0.52
1:C:258:ALA:HB3	1:C:304:LEU:HD21	1.91	0.52
1:B:258:ALA:HB3	1:B:304:LEU:HD21	1.91	0.52
1:C:264:LEU:HD12	1:C:275:HIS:O	2.10	0.52
1:A:258:ALA:HB3	1:A:304:LEU:HD21	1.91	0.52
1:A:264:LEU:HD12	1:A:275:HIS:O	2.10	0.51
1:B:264:LEU:HD12	1:B:275:HIS:O	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:268:ILE:HB	1:B:275:HIS:CD2	2.47	0.50
1:C:268:ILE:HB	1:C:275:HIS:CD2	2.47	0.50
1:B:465:SER:HB2	1:B:489:LEU:HD11	1.93	0.50
1:C:234:LEU:HG	1:C:272:GLU:HB3	1.94	0.50
1:C:465:SER:HB2	1:C:489:LEU:HD11	1.93	0.50
1:B:234:LEU:HG	1:B:272:GLU:HB3	1.94	0.50
1:A:234:LEU:HG	1:A:272:GLU:HB3	1.93	0.49
1:A:465:SER:HB2	1:A:489:LEU:HD11	1.93	0.49
1:A:268:ILE:HB	1:A:275:HIS:CD2	2.47	0.48
1:A:128:PHE:HA	1:A:131:MET:HE3	1.96	0.48
1:A:79:TYR:O	1:A:83:ARG:HG3	2.14	0.48
1:C:268:ILE:HB	1:C:275:HIS:HD2	1.79	0.47
1:C:79:TYR:O	1:C:83:ARG:HG3	2.14	0.47
1:B:79:TYR:O	1:B:83:ARG:HG3	2.14	0.47
1:C:128:PHE:HA	1:C:131:MET:HE3	1.97	0.47
1:C:622:SER:O	1:C:633:ASN:HB3	2.15	0.47
1:B:128:PHE:HA	1:B:131:MET:HE3	1.97	0.47
1:A:510:ARG:NH2	1:A:570:GLU:OE1	2.48	0.47
1:A:18:ASN:ND2	1:A:21:LYS:HG3	2.31	0.46
1:C:18:ASN:ND2	1:C:21:LYS:HG3	2.31	0.46
1:B:622:SER:O	1:B:633:ASN:HB3	2.15	0.46
1:B:268:ILE:HB	1:B:275:HIS:HD2	1.79	0.46
1:B:18:ASN:ND2	1:B:21:LYS:HG3	2.31	0.46
1:C:510:ARG:NH2	1:C:570:GLU:OE1	2.48	0.46
1:C:705:PRO:O	1:C:706:SER:HB2	2.16	0.46
1:A:622:SER:O	1:A:633:ASN:HB3	2.15	0.46
1:B:705:PRO:O	1:B:706:SER:HB2	2.15	0.46
1:A:705:PRO:O	1:A:706:SER:HB2	2.15	0.46
1:A:122:ASP:O	1:A:189:ARG:NH2	2.49	0.45
1:A:268:ILE:HB	1:A:275:HIS:HD2	1.80	0.45
1:A:433:VAL:HG11	1:A:443:ALA:HB1	1.97	0.45
1:B:510:ARG:NH2	1:B:570:GLU:OE1	2.48	0.45
1:C:122:ASP:O	1:C:189:ARG:NH2	2.49	0.45
1:B:122:ASP:O	1:B:189:ARG:NH2	2.49	0.45
1:C:212:MET:O	1:C:216:ARG:NH2	2.50	0.45
1:A:620:MET:HB2	1:A:621:PRO:HD2	1.99	0.45
1:B:620:MET:HB2	1:B:621:PRO:HD2	1.99	0.45
1:C:433:VAL:HG11	1:C:443:ALA:HB1	1.98	0.44
1:C:686:GLN:NE2	1:C:727:LYS:HE3	2.33	0.44
1:B:686:GLN:NE2	1:B:727:LYS:HE3	2.33	0.44
1:C:620:MET:HB2	1:C:621:PRO:HD2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:433:VAL:HG11	1:B:443:ALA:HB1	1.97	0.44
1:B:212:MET:O	1:B:216:ARG:NH2	2.50	0.44
1:B:17:ILE:HD13	1:B:18:ASN:N	2.33	0.44
1:B:293:SER:O	1:B:294:GLN:C	2.56	0.44
1:A:217:THR:OG1	1:A:219:THR:HG22	2.18	0.44
1:C:217:THR:OG1	1:C:219:THR:HG22	2.18	0.44
1:A:212:MET:O	1:A:216:ARG:NH2	2.50	0.43
1:C:293:SER:O	1:C:294:GLN:C	2.57	0.43
1:A:686:GLN:NE2	1:A:727:LYS:HE3	2.33	0.43
1:A:328:ASN:HA	1:A:328:ASN:HD22	1.64	0.43
1:C:312:GLU:O	1:C:316:LEU:HG	2.19	0.43
1:A:17:ILE:HD13	1:A:18:ASN:N	2.33	0.43
1:B:312:GLU:O	1:B:316:LEU:HG	2.19	0.43
1:A:162:THR:OG1	1:A:164:GLU:HG3	2.18	0.43
1:A:293:SER:O	1:A:294:GLN:C	2.56	0.43
1:B:217:THR:OG1	1:B:219:THR:HG22	2.18	0.43
1:B:162:THR:OG1	1:B:164:GLU:HG3	2.18	0.43
1:A:558:LEU:CD2	1:A:612:ARG:HG2	2.49	0.43
1:C:17:ILE:HD13	1:C:18:ASN:N	2.33	0.42
1:B:558:LEU:CD2	1:B:612:ARG:HG2	2.49	0.42
1:C:346:ARG:HD2	1:C:352:ASP:O	2.19	0.42
1:B:379:TYR:O	1:B:383:GLU:HG3	2.19	0.42
1:A:379:TYR:O	1:A:383:GLU:HG3	2.19	0.42
1:A:312:GLU:O	1:A:316:LEU:HG	2.19	0.42
1:C:162:THR:OG1	1:C:164:GLU:HG3	2.18	0.42
1:C:328:ASN:HD22	1:C:328:ASN:HA	1.64	0.42
1:C:191:GLN:O	1:C:195:ARG:HG3	2.20	0.42
1:A:171:PHE:HA	1:A:174:ILE:HG22	2.02	0.42
1:B:346:ARG:HD2	1:B:352:ASP:O	2.19	0.42
1:B:98:GLU:HA	1:B:99:PRO:HD3	1.91	0.42
1:A:346:ARG:HD2	1:A:352:ASP:O	2.19	0.42
1:C:558:LEU:CD2	1:C:612:ARG:HG2	2.49	0.42
1:C:379:TYR:O	1:C:383:GLU:HG3	2.19	0.42
1:B:191:GLN:O	1:B:195:ARG:HG3	2.20	0.42
1:C:209:THR:N	1:C:210:PRO:CD	2.83	0.42
1:C:171:PHE:HA	1:C:174:ILE:HG22	2.02	0.41
1:B:209:THR:HB	1:B:210:PRO:HD3	2.02	0.41
1:C:209:THR:HB	1:C:210:PRO:HD3	2.02	0.41
1:B:209:THR:N	1:B:210:PRO:CD	2.83	0.41
1:C:307:PRO:HA	1:C:338:GLN:HB2	2.02	0.41
1:A:672:MET:HA	1:A:673:PRO:HD3	1.87	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:440:LEU:HD12	1:B:728:THR:HB	2.02	0.41
1:B:171:PHE:HA	1:B:174:ILE:HG22	2.02	0.41
1:C:440:LEU:HD12	1:C:728:THR:HB	2.02	0.41
1:A:209:THR:HB	1:A:210:PRO:HD3	2.02	0.41
1:C:98:GLU:HA	1:C:99:PRO:HD3	1.92	0.41
1:C:99:PRO:HG2	1:C:137:HIS:CD2	2.56	0.41
1:A:191:GLN:O	1:A:195:ARG:HG3	2.20	0.41
1:B:99:PRO:HG2	1:B:137:HIS:CD2	2.56	0.41
1:A:209:THR:N	1:A:210:PRO:CD	2.83	0.41
1:A:284:HIS:CE1	1:B:284:HIS:CE1	3.09	0.41
1:A:307:PRO:HA	1:A:338:GLN:HB2	2.02	0.41
1:A:99:PRO:HG2	1:A:137:HIS:CD2	2.56	0.41
1:B:307:PRO:HA	1:B:338:GLN:HB2	2.02	0.41
1:A:203:PHE:HB3	1:A:629:SER:HB3	2.03	0.41
1:B:434:ARG:HD3	1:B:434:ARG:HH11	1.75	0.41
1:A:320:LYS:HE2	1:A:411:ARG:HB2	2.03	0.40
1:C:320:LYS:HE2	1:C:411:ARG:HB2	2.03	0.40
1:A:440:LEU:HD12	1:A:728:THR:HB	2.03	0.40
1:B:320:LYS:HE2	1:B:411:ARG:HB2	2.03	0.40
1:C:38:ILE:O	1:C:42:GLU: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	732/761 (96%)	697 (95%)	29 (4%)	6 (1%)	24 60
1	B	732/761 (96%)	697 (95%)	29 (4%)	6 (1%)	24 60
1	C	732/761 (96%)	697 (95%)	29 (4%)	6 (1%)	24 60
2	D	14/20 (70%)	13 (93%)	1 (7%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	E	14/20 (70%)	13 (93%)	1 (7%)	0	100	100
2	F	14/20 (70%)	13 (93%)	1 (7%)	0	100	100
2	P	1/20 (5%)	0	1 (100%)	0	100	100
All	All	2239/2363 (95%)	2130 (95%)	91 (4%)	18 (1%)	24	60

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	294	GLN
1	B	294	GLN
1	C	294	GLN
1	A	216	ARG
1	A	273	ALA
1	B	216	ARG
1	B	273	ALA
1	C	216	ARG
1	C	273	ALA
1	A	5	LEU
1	A	300	GLY
1	B	5	LEU
1	B	300	GLY
1	C	5	LEU
1	C	300	GLY
1	A	731	TYR
1	B	731	TYR
1	C	731	TYR

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	629/651 (97%)	576 (92%)	53 (8%)	14	37
1	B	629/651 (97%)	576 (92%)	53 (8%)	14	37
1	C	629/651 (97%)	576 (92%)	53 (8%)	14	37

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	D	16/19 (84%)	12 (75%)	4 (25%)	1	2
2	E	16/19 (84%)	12 (75%)	4 (25%)	1	2
2	F	16/19 (84%)	12 (75%)	4 (25%)	1	2
2	P	3/19 (16%)	2 (67%)	1 (33%)	0	1
All	All	1938/2029 (96%)	1766 (91%)	172 (9%)	12	35

All (172) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	LEU
1	A	9	LYS
1	A	14	THR
1	A	16	ARG
1	A	17	ILE
1	A	24	ARG
1	A	39	SER
1	A	40	GLN
1	A	45	SER
1	A	51	ASP
1	A	54	LYS
1	A	56	SER
1	A	72	ARG
1	A	91	LYS
1	A	96	GLN
1	A	99	PRO
1	A	111	GLU
1	A	118	HIS
1	A	139	ARG
1	A	149	LYS
1	A	183	ASN
1	A	189	ARG
1	A	206	SER
1	A	216	ARG
1	A	225	CYS
1	A	226	VAL
1	A	269	ARG
1	A	275	HIS
1	A	294	GLN
1	A	297	VAL
1	A	317	LEU
1	A	320	LYS

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Mol	Chain	Res	Type
1	A	364	LEU
1	A	373	GLU
1	A	384	LYS
1	A	387	SER
1	A	394	LYS
1	A	424	SER
1	A	440	LEU
1	A	447	LYS
1	A	452	VAL
1	A	484	LEU
1	A	530	ARG
1	A	616	LEU
1	A	625	SER
1	A	626	SER
1	A	629	SER
1	A	630	ASN
1	A	643	SER
1	A	648	LYS
1	A	696	ASN
1	A	716	LYS
1	A	736	ASP
2	D	360	GLN
2	D	361	ILE
2	D	363	SER
2	D	367	THR
1	B	5	LEU
1	B	9	LYS
1	B	14	THR
1	B	16	ARG
1	B	17	ILE
1	B	24	ARG
1	B	39	SER
1	B	40	GLN
1	B	45	SER
1	B	51	ASP
1	B	54	LYS
1	B	56	SER
1	B	72	ARG
1	B	91	LYS
1	B	96	GLN
1	B	99	PRO
1	B	111	GLU

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Mol	Chain	Res	Type
1	B	118	HIS
1	B	139	ARG
1	B	149	LYS
1	B	183	ASN
1	B	189	ARG
1	B	206	SER
1	B	216	ARG
1	B	225	CYS
1	B	226	VAL
1	B	269	ARG
1	B	275	HIS
1	B	294	GLN
1	B	297	VAL
1	B	317	LEU
1	B	320	LYS
1	B	364	LEU
1	B	373	GLU
1	B	384	LYS
1	B	387	SER
1	B	394	LYS
1	B	424	SER
1	B	440	LEU
1	B	447	LYS
1	B	452	VAL
1	B	484	LEU
1	B	530	ARG
1	B	616	LEU
1	B	625	SER
1	B	626	SER
1	B	629	SER
1	B	630	ASN
1	B	643	SER
1	B	648	LYS
1	B	696	ASN
1	B	716	LYS
1	B	736	ASP
2	E	360	GLN
2	E	361	ILE
2	E	363	SER
2	E	367	THR
1	C	5	LEU
1	C	9	LYS

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Mol	Chain	Res	Type
1	C	14	THR
1	C	16	ARG
1	C	17	ILE
1	C	24	ARG
1	C	39	SER
1	C	40	GLN
1	C	45	SER
1	C	51	ASP
1	C	54	LYS
1	C	56	SER
1	C	72	ARG
1	C	91	LYS
1	C	96	GLN
1	C	99	PRO
1	C	111	GLU
1	C	118	HIS
1	C	139	ARG
1	C	149	LYS
1	C	183	ASN
1	C	189	ARG
1	C	206	SER
1	C	216	ARG
1	C	225	CYS
1	C	226	VAL
1	C	269	ARG
1	C	275	HIS
1	C	294	GLN
1	C	297	VAL
1	C	317	LEU
1	C	320	LYS
1	C	364	LEU
1	C	373	GLU
1	C	384	LYS
1	C	387	SER
1	C	394	LYS
1	C	424	SER
1	C	440	LEU
1	C	447	LYS
1	C	452	VAL
1	C	484	LEU
1	C	530	ARG
1	C	616	LEU

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Mol	Chain	Res	Type
1	C	625	SER
1	C	626	SER
1	C	629	SER
1	C	630	ASN
1	C	643	SER
1	C	648	LYS
1	C	696	ASN
1	C	716	LYS
1	C	736	ASP
2	F	360	GLN
2	F	361	ILE
2	F	363	SER
2	F	367	THR
2	P	2	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	18	ASN
1	A	46	HIS
1	A	250	GLN
1	A	275	HIS
1	A	328	ASN
1	A	596	HIS
1	A	630	ASN
1	A	661	HIS
2	D	360	GLN
1	B	18	ASN
1	B	46	HIS
1	B	250	GLN
1	B	275	HIS
1	B	328	ASN
1	B	596	HIS
1	B	630	ASN
1	B	661	HIS
1	C	18	ASN
1	C	46	HIS
1	C	250	GLN
1	C	275	HIS
1	C	328	ASN
1	C	596	HIS
1	C	630	ASN

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Mol	Chain	Res	Type
1	C	661	HIS

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 carbohydrates 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	734/761 (96%)	-0.41	29 (3%) 42 35	10, 24, 77, 108	0
1	B	734/761 (96%)	-0.00	30 (4%) 41 34	10, 24, 77, 108	0
1	C	734/761 (96%)	-0.07	24 (3%) 50 42	10, 24, 77, 108	0
2	D	16/20 (80%)	0.80	3 (18%) 2 1	45, 82, 91, 97	0
2	E	16/20 (80%)	0.88	2 (12%) 5 3	45, 82, 91, 97	0
2	F	16/20 (80%)	0.95	1 (6%) 23 17	45, 82, 91, 97	0
2	P	3/20 (15%)	0.37	0 100 100	30, 30, 37, 43	0
All	All	2253/2363 (95%)	-0.14	89 (3%) 42 35	10, 25, 83, 108	0

All (89) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	737	GLY	6.9
1	A	297	VAL	6.7
1	A	296	GLY	6.5
1	B	294	GLN	6.0
1	C	4	ASN	6.0
1	C	297	VAL	5.5
1	C	737	GLY	5.3
1	A	12	GLY	5.2
1	A	270	GLY	5.1
1	A	11	ASP	5.1
1	B	296	GLY	5.0
1	B	297	VAL	5.0
1	C	296	GLY	4.7
2	F	360	GLN	4.7
1	B	269	ARG	4.7
1	B	10	ARG	4.7
1	C	298	ARG	4.5

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Mol	Chain	Res	Type	RSRZ
1	B	737	GLY	4.4
1	A	4	ASN	4.4
1	A	299	GLY	4.3
1	C	294	GLN	4.3
1	C	736	ASP	4.2
1	A	273	ALA	4.2
1	A	271	GLY	4.0
1	A	323	ARG	3.9
2	E	360	GLN	3.9
1	C	269	ARG	3.9
1	B	11	ASP	3.8
1	B	273	ALA	3.8
1	A	294	GLN	3.7
1	A	298	ARG	3.7
1	A	13	SER	3.7
1	C	10	ARG	3.6
1	B	4	ASN	3.6
1	C	274	PHE	3.6
1	A	322	ASN	3.6
1	C	270	GLY	3.6
1	C	11	ASP	3.5
1	A	269	ARG	3.5
1	B	12	GLY	3.4
1	B	322	ASN	3.4
1	B	323	ARG	3.4
1	C	273	ALA	3.3
1	A	736	ASP	3.3
1	C	271	GLY	3.2
1	C	299	GLY	3.1
1	B	736	ASP	3.1
1	B	270	GLY	3.1
2	D	360	GLN	3.1
1	B	271	GLY	3.1
1	B	268	ILE	3.0
1	B	326	GLU	2.9
1	A	10	ARG	2.9
1	C	16	ARG	2.9
2	D	372	ASN	2.9
1	C	13	SER	2.8
1	A	6	LEU	2.8
1	A	274	PHE	2.8
1	A	24	ARG	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	592	ASN	2.7
1	B	267	PRO	2.7
1	C	275	HIS	2.6
1	A	17	ILE	2.6
1	A	623	GLU	2.6
1	B	298	ARG	2.6
2	E	368	ASP	2.5
1	A	160	ARG	2.5
1	C	326	GLU	2.4
1	C	323	ARG	2.4
1	B	299	GLY	2.4
1	B	623	GLU	2.4
1	C	325	VAL	2.4
1	B	266	SER	2.4
1	C	623	GLU	2.4
1	A	295	GLY	2.4
1	C	160	ARG	2.3
1	B	13	SER	2.3
1	A	268	ILE	2.3
1	B	18	ASN	2.3
1	B	160	ARG	2.2
1	B	16	ARG	2.2
2	D	362	ASP	2.1
1	B	272	GLU	2.1
1	A	272	GLU	2.1
1	B	187	GLU	2.1
1	C	272	GLU	2.1
1	B	648	LYS	2.1
1	B	5	LEU	2.1
1	A	161	VAL	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

There are no carbohydrates 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.