Full wwPDB X-ray Structure Validation Report (i)

Jun 12, 2024 – 11:45 AM EDT

PDB ID	:	2JAN
Title	:	TYROSYL-TRNA SYNTHETASE FROM MYCOBACTERIUM TUBERCULOSIS IN UNLIGANDED STATE
Authors	:	Hartmann, M.D.; Shkolnaya, L.A.; Bourenkov, G.P.; Strizhov, N.I.; Bartunik, H.D.
Deposited on	:	2006-11-29
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 (i) 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 (1)) were used in the production of this report:

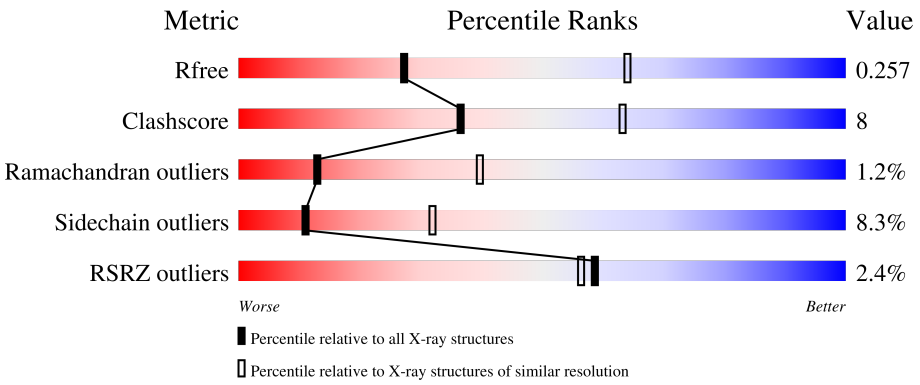
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.20.1
EDS	:	2.36.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

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))
R_{free}	130704	1957 (2.90-2.90)
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	432	3% 75% 20% • •
1	B	432	2% 74% 18% • 6%
1	C	432	3% 76% 15% • 6%
1	D	432	% 75% 16% • 6%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 12591 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 TYROSYL-TRNA SYNTHETASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	419	3159	1993	564	596	6	0	0	0
1	B	405	3095	1949	565	575	6	0	0	0
1	C	407	3116	1965	568	577	6	0	1	0
1	D	407	3116	1964	567	579	6	0	1	0

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	31	31	31	0	0
2	B	24	24	24	0	0
2	C	18	18	18	0	0
2	D	32	32	32	0	0

3 Residue-property plots (i)

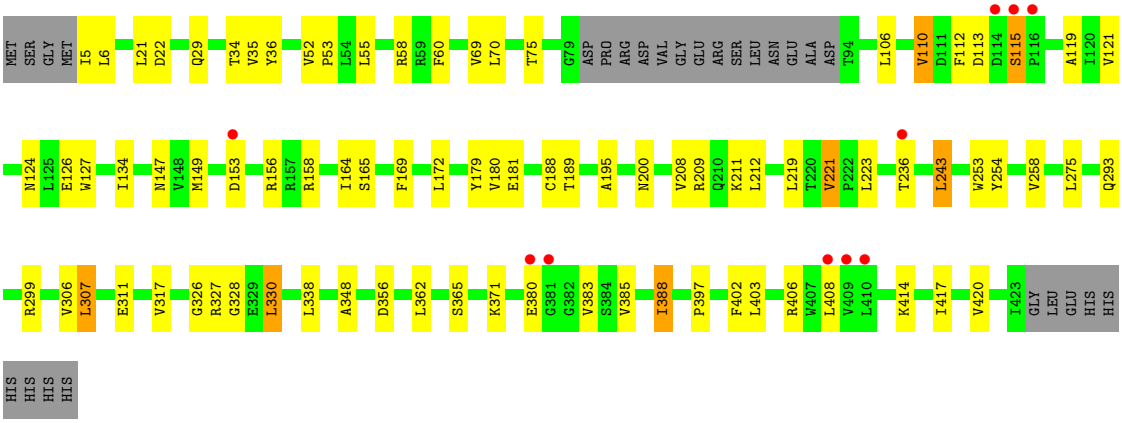
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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

Chain A:

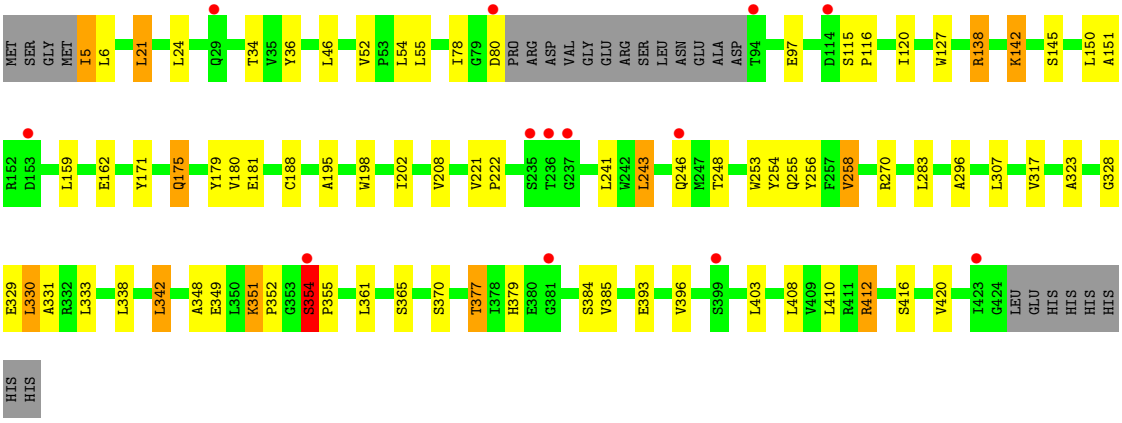
• Molecule 1: TYROSYL-TRNA SYNTHETASE

Chain B:



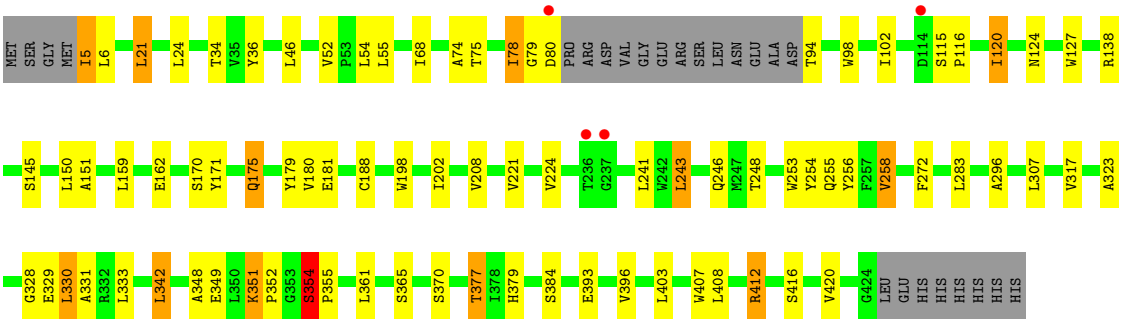
• Molecule 1: TYROSYL-TRNA SYNTHETASE

Chain C:



- Molecule 1: TYROSYL-TRNA SYNTHETASE

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	71.26Å 78.44Å 81.63Å 74.42° 83.19° 82.95°	Depositor
Resolution (Å)	19.89 – 2.90 19.88 – 2.90	Depositor EDS
% Data completeness (in resolution range)	98.4 (19.89-2.90) 98.4 (19.88-2.90)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.00 (at 2.88Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.201 , 0.257 0.206 , 0.257	Depositor DCC
R_{free} test set	1864 reflections (5.10%)	wwPDB-VP
Wilson B-factor (Å^2)	44.9	Xtriage
Anisotropy	0.018	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.30 , 37.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.105 for -h,-l,-k	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	12591	wwPDB-VP
Average B, all atoms (Å^2)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.88% 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 (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	0.47	0/3222	0.61	0/4386
1	B	0.46	0/3156	0.60	0/4288
1	C	0.47	0/3181	0.61	0/4322
1	D	0.47	0/3181	0.61	0/4323
All	All	0.47	0/12740	0.61	0/17319

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	C	0	2
1	D	0	3
All	All	0	6

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	85	GLY	Peptide
1	C	328	GLY	Peptide
1	C	354	SER	Peptide
1	D	328	GLY	Peptide
1	D	354	SER	Peptide
1	D	78	ILE	Peptide

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3159	0	3076	68	0
1	B	3095	0	3052	48	0
1	C	3116	0	3077	40	0
1	D	3116	0	3070	45	0
2	A	31	0	0	0	0
2	B	24	0	0	0	0
2	C	18	0	0	0	0
2	D	32	0	0	0	0
All	All	12591	0	12275	196	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 8.

All (196) 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:B:55:LEU:HD21	1:B:307:LEU:HD11	1.34	1.09
1:A:69:VAL:HG21	1:A:106:LEU:HD22	1.39	1.03
1:A:5:ILE:HG23	1:A:6:LEU:HD12	1.52	0.90
1:A:87:ARG:HH11	1:A:87:ARG:HG2	1.37	0.89
1:A:87:ARG:HB3	1:A:87:ARG:NH1	1.91	0.85
1:B:5:ILE:HG23	1:B:6:LEU:HD12	1.60	0.83
1:C:52:VAL:HG12	1:C:221:VAL:HG21	1.59	0.83
1:D:52:VAL:HG12	1:D:221:VAL:HG21	1.61	0.82
1:A:55:LEU:HD21	1:A:307:LEU:HD11	1.60	0.82
1:A:91:GLU:O	1:A:95:VAL:HG23	1.79	0.82
1:A:87:ARG:HH11	1:A:87:ARG:CG	1.92	0.82
1:D:55:LEU:HD21	1:D:307:LEU:HD11	1.64	0.79
1:B:127:TRP:CG	1:B:181:GLU:HG2	2.23	0.73
1:B:55:LEU:HD21	1:B:307:LEU:CD1	2.15	0.73
1:A:330:LEU:HD13	1:A:403:LEU:HD21	1.69	0.72
1:B:195:ALA:HB3	1:B:223:LEU:HD12	1.73	0.71
1:D:46:LEU:HD13	1:D:54:LEU:HD11	1.74	0.69
1:B:330:LEU:HD13	1:B:403:LEU:HD21	1.75	0.69
1:B:69:VAL:HG21	1:B:106:LEU:HD22	1.73	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:6:LEU:HD23	1:A:21:LEU:HD22	1.73	0.69
1:C:55:LEU:HD21	1:C:307:LEU:HD11	1.74	0.68
1:B:6:LEU:HD23	1:B:21:LEU:HD22	1.76	0.68
1:A:134:ILE:HG23	1:D:78:ILE:HG13	1.76	0.68
1:C:6:LEU:HD21	1:C:24:LEU:HG	1.76	0.67
1:A:82:ARG:O	1:A:87:ARG:HD3	1.95	0.67
1:A:195:ALA:HB3	1:A:223:LEU:HD12	1.77	0.67
1:B:110:VAL:HG13	1:B:119:ALA:HB2	1.78	0.65
1:A:87:ARG:NH1	1:A:87:ARG:CB	2.59	0.65
1:D:6:LEU:HD21	1:D:24:LEU:HG	1.79	0.65
1:A:43:ALA:HB1	1:A:44:PRO:CD	2.28	0.64
1:A:166:TYR:OH	1:D:170:SER:OG	2.04	0.64
1:A:70:LEU:HD11	1:A:124:ASN:HB3	1.78	0.64
1:B:70:LEU:HD11	1:B:124:ASN:HB3	1.81	0.63
1:B:195:ALA:HB3	1:B:223:LEU:CD1	2.27	0.62
1:A:385:VAL:CG1	1:A:408:LEU:HD23	2.30	0.62
1:A:43:ALA:HB1	1:A:44:PRO:HD2	1.81	0.61
1:A:52:VAL:HG12	1:A:221:VAL:HG11	1.81	0.61
1:B:385:VAL:CG1	1:B:408:LEU:HD23	2.31	0.61
1:D:329:GLU:O	1:D:331:ALA:N	2.34	0.60
1:C:150:LEU:HD22	1:C:159:LEU:HD11	1.83	0.60
1:A:55:LEU:HD21	1:A:307:LEU:CD1	2.30	0.60
1:A:69:VAL:HG21	1:A:106:LEU:CD2	2.25	0.59
1:D:241:LEU:HD11	1:D:256:TYR:CG	2.38	0.59
1:B:112:PHE:CE1	1:B:121:VAL:HG23	2.38	0.59
1:A:275:LEU:HD21	1:A:299:ARG:NH2	2.19	0.58
1:C:198:TRP:CZ2	1:C:202:ILE:HD13	2.40	0.57
1:C:241:LEU:HD11	1:C:256:TYR:CG	2.38	0.57
1:D:241:LEU:HD22	1:D:248:THR:HG21	1.85	0.57
1:A:396:VAL:HG22	1:A:397:PRO:HD2	1.86	0.57
1:A:195:ALA:HB3	1:A:223:LEU:CD1	2.35	0.57
1:B:112:PHE:HE1	1:B:121:VAL:HG23	1.70	0.56
1:D:75:THR:O	1:D:79:GLY:N	2.38	0.56
1:D:150:LEU:HD22	1:D:159:LEU:HD11	1.87	0.56
1:C:379:HIS:NE2	1:C:393:GLU:OE1	2.39	0.56
1:D:127:TRP:CD2	1:D:181:GLU:HG2	2.40	0.56
1:A:362:LEU:HD11	1:A:383:VAL:HG11	1.87	0.56
1:C:255:GLN:O	1:C:258:VAL:HG12	2.05	0.56
1:C:329:GLU:O	1:C:331:ALA:N	2.40	0.55
1:A:87:ARG:HB3	1:A:87:ARG:CZ	2.36	0.55
1:D:5:ILE:HD12	1:D:5:ILE:O	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:323:ALA:HB1	1:D:330:LEU:HB3	1.87	0.55
1:A:110:VAL:HG13	1:A:119:ALA:HB2	1.87	0.55
1:C:323:ALA:HB3	1:C:333:LEU:HD11	1.89	0.54
1:B:243:LEU:HD23	1:B:243:LEU:N	2.23	0.54
1:C:323:ALA:HB1	1:C:330:LEU:HB3	1.89	0.54
1:C:5:ILE:HD12	1:C:5:ILE:O	2.07	0.54
1:D:379:HIS:NE2	1:D:393:GLU:OE1	2.41	0.54
1:A:110:VAL:CG1	1:A:119:ALA:HB2	2.38	0.54
1:C:241:LEU:HD22	1:C:248:THR:HG21	1.90	0.54
1:A:84:VAL:O	1:A:88:SER:OG	2.26	0.53
1:C:6:LEU:HD23	1:C:21:LEU:HD13	1.91	0.53
1:D:127:TRP:CG	1:D:181:GLU:HG2	2.43	0.53
1:B:275:LEU:HD21	1:B:299:ARG:NH2	2.23	0.53
1:B:243:LEU:HD22	1:B:253:TRP:CH2	2.44	0.53
1:C:127:TRP:CG	1:C:181:GLU:HG2	2.44	0.53
1:D:6:LEU:HD23	1:D:21:LEU:HD13	1.91	0.53
1:D:198:TRP:CZ2	1:D:202:ILE:HD13	2.43	0.53
1:B:52:VAL:HG12	1:B:221:VAL:HG11	1.91	0.52
1:C:377:THR:HG21	1:C:412:ARG:HD3	1.91	0.52
1:D:255:GLN:O	1:D:258:VAL:HG12	2.10	0.52
1:A:398:GLN:CB	1:A:401:ASP:OD1	2.58	0.52
1:A:385:VAL:HG13	1:A:408:LEU:HD23	1.91	0.52
1:B:164:ILE:HG23	1:B:165:SER:O	2.10	0.52
1:A:84:VAL:C	1:A:86:GLU:H	2.13	0.51
1:A:87:ARG:HH11	1:A:87:ARG:CB	2.23	0.51
1:A:164:ILE:HG23	1:A:165:SER:O	2.10	0.51
1:B:362:LEU:HD11	1:B:383:VAL:HG11	1.92	0.51
1:A:87:ARG:CB	1:A:87:ARG:CZ	2.88	0.51
1:D:283:LEU:HD22	1:D:296:ALA:HA	1.91	0.51
1:A:112:PHE:CE1	1:A:121:VAL:HG23	2.45	0.51
1:C:254:TYR:CB	1:C:317:VAL:HG11	2.40	0.51
1:D:254:TYR:CB	1:D:317:VAL:HG11	2.40	0.51
1:B:147:ASN:OD1	1:C:159:LEU:HD22	2.11	0.51
1:A:88:SER:O	1:A:90:ASN:N	2.44	0.51
1:D:377:THR:HG21	1:D:412:ARG:HD3	1.92	0.51
1:D:348:ALA:HB2	1:D:365:SER:HA	1.93	0.51
1:A:52:VAL:HG12	1:A:221:VAL:CG1	2.41	0.51
1:C:46:LEU:HD13	1:C:54:LEU:HD11	1.92	0.51
1:D:323:ALA:HB3	1:D:333:LEU:HD11	1.93	0.51
1:A:243:LEU:HD22	1:A:253:TRP:CH2	2.47	0.50
1:C:348:ALA:HB2	1:C:365:SER:HA	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:180:VAL:HA	1:A:208:VAL:HG22	1.94	0.48
1:A:89:LEU:O	1:A:89:LEU:HD23	2.13	0.48
1:B:180:VAL:HA	1:B:208:VAL:HG22	1.95	0.48
1:A:356:ASP:HB3	1:A:397:PRO:HD2	1.95	0.48
1:B:110:VAL:CG1	1:B:119:ALA:CB	2.92	0.47
1:B:127:TRP:CD2	1:B:181:GLU:HG2	2.49	0.47
1:C:150:LEU:CD2	1:C:159:LEU:HD11	2.45	0.47
1:C:34:THR:HB	1:C:188:CYS:HA	1.96	0.47
1:B:158:ARG:HG2	1:B:164:ILE:HD13	1.96	0.47
1:C:283:LEU:HD22	1:C:296:ALA:HA	1.97	0.47
1:D:180:VAL:HA	1:D:208:VAL:HG22	1.96	0.47
1:A:243:LEU:N	1:A:243:LEU:HD23	2.30	0.47
1:B:113:ASP:OD1	1:B:115:SER:HB3	2.15	0.47
1:B:134:ILE:HG23	1:C:78:ILE:HG13	1.97	0.47
1:D:243:LEU:HD22	1:D:253:TRP:CH2	2.49	0.47
1:C:115:SER:CB	1:C:116:PRO:CD	2.92	0.47
1:B:385:VAL:HG13	1:B:408:LEU:HD23	1.96	0.47
1:D:68:ILE:HG12	1:D:120:ILE:HG23	1.96	0.47
1:A:112:PHE:CD1	1:A:119:ALA:HB3	2.50	0.47
1:B:149:MET:HE2	1:B:169:PHE:CZ	2.50	0.47
1:C:127:TRP:CD2	1:C:181:GLU:HG2	2.50	0.47
1:D:34:THR:HB	1:D:188:CYS:HA	1.97	0.47
1:A:127:TRP:CG	1:A:181:GLU:HG2	2.50	0.46
1:B:211:LYS:HB3	1:B:212:LEU:HD13	1.97	0.46
1:D:150:LEU:CD2	1:D:159:LEU:HD11	2.45	0.46
1:A:66:ARG:NH1	1:A:120:ILE:HD11	2.30	0.46
1:B:127:TRP:CB	1:B:181:GLU:HG2	2.46	0.46
1:B:110:VAL:CG1	1:B:119:ALA:HB2	2.42	0.46
1:D:323:ALA:CB	1:D:333:LEU:HD11	2.46	0.46
1:D:254:TYR:HB2	1:D:317:VAL:HG11	1.98	0.46
1:A:84:VAL:C	1:A:86:GLU:N	2.69	0.46
1:A:211:LYS:HB3	1:A:212:LEU:HD13	1.97	0.46
1:C:351:LYS:HB2	1:C:352:PRO:HD2	1.98	0.46
1:D:74:ALA:HB3	1:D:124:ASN:HD21	1.80	0.46
1:A:110:VAL:CG1	1:A:119:ALA:CB	2.94	0.45
1:A:147:ASN:OD1	1:D:159:LEU:HD22	2.16	0.45
1:B:219:LEU:C	1:B:219:LEU:HD23	2.36	0.45
1:C:138:ARG:O	1:C:142:LYS:HB2	2.15	0.45
1:D:98:TRP:CE2	1:D:102:ILE:HG13	2.52	0.45
1:B:388:ILE:HG22	1:B:388:ILE:O	2.15	0.45
1:A:396:VAL:CG2	1:A:397:PRO:HD2	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:408:LEU:HB2	1:C:420:VAL:HB	1.98	0.45
1:D:6:LEU:HD23	1:D:21:LEU:CD1	2.45	0.45
1:D:351:LYS:HB2	1:D:352:PRO:HD2	1.97	0.45
1:B:356:ASP:HB3	1:B:397:PRO:HD2	1.99	0.45
1:D:75:THR:HG23	1:D:171[A]:TYR:CZ	2.52	0.45
1:D:115:SER:CB	1:D:116:PRO:CD	2.94	0.45
1:D:52:VAL:CG1	1:D:221:VAL:HG21	2.39	0.45
1:C:323:ALA:CB	1:C:333:LEU:HD11	2.46	0.44
1:C:385:VAL:HG22	1:C:410:LEU:CD2	2.47	0.44
1:A:306:VAL:HG22	1:A:311:GLU:HA	1.99	0.44
1:B:34:THR:HB	1:B:188:CYS:HA	1.99	0.44
1:C:254:TYR:HB2	1:C:317:VAL:HG11	1.98	0.44
1:A:219:LEU:C	1:A:219:LEU:HD23	2.37	0.44
1:B:326:GLY:O	1:B:328:GLY:N	2.49	0.44
1:A:103:ARG:HA	1:A:106:LEU:HD12	2.00	0.44
1:C:52:VAL:CG1	1:C:221:VAL:HG21	2.38	0.43
1:B:52:VAL:HB	1:B:53:PRO:HD3	1.99	0.43
1:C:171[B]:TYR:CE2	1:C:175:GLN:OE1	2.71	0.43
1:C:243:LEU:HD22	1:C:253:TRP:CH2	2.53	0.43
1:A:5:ILE:HG23	1:A:6:LEU:N	2.33	0.43
1:B:402:PHE:CE2	1:B:408:LEU:CD1	3.01	0.43
1:C:180:VAL:HA	1:C:208:VAL:HG22	2.01	0.43
1:D:408:LEU:HB2	1:D:420:VAL:HB	2.00	0.43
1:A:34:THR:HB	1:A:188:CYS:HA	2.01	0.43
1:A:388:ILE:O	1:A:388:ILE:HG22	2.19	0.43
1:B:254:TYR:CB	1:B:317:VAL:HG11	2.49	0.43
1:D:75:THR:O	1:D:79:GLY:CA	2.67	0.42
1:A:275:LEU:HD21	1:A:299:ARG:CZ	2.48	0.42
1:A:69:VAL:HG11	1:A:106:LEU:HD11	2.01	0.42
1:B:172:LEU:C	1:B:172:LEU:HD13	2.40	0.42
1:B:385:VAL:HG11	1:B:408:LEU:HD23	2.01	0.42
1:A:52:VAL:HB	1:A:53:PRO:HD3	2.01	0.42
1:C:6:LEU:HD23	1:C:21:LEU:CD1	2.49	0.42
1:A:41:PRO:HD2	1:A:98:TRP:CZ2	2.55	0.42
1:A:392:ASN:OD1	1:A:395:TRP:N	2.53	0.42
1:C:115:SER:CB	1:C:116:PRO:HD2	2.50	0.42
1:B:408:LEU:HB2	1:B:420:VAL:HB	2.02	0.41
1:A:84:VAL:O	1:A:86:GLU:N	2.54	0.41
1:A:149:MET:HE2	1:A:169:PHE:CZ	2.55	0.41
1:D:68:ILE:HG12	1:D:120:ILE:CG2	2.51	0.41
1:D:171[B]:TYR:CE2	1:D:175:GLN:OE1	2.74	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:254:TYR:CB	1:A:317:VAL:HG11	2.50	0.41
1:B:35:VAL:HG21	1:B:60:PHE:CD1	2.55	0.41
1:D:224:VAL:HG21	1:D:272:PHE:HZ	1.86	0.41
1:D:342:LEU:HD23	1:D:407:TRP:CE3	2.56	0.41
1:B:348:ALA:HB2	1:B:365:SER:HA	2.03	0.41
1:C:338:LEU:CD1	1:C:342:LEU:HD22	2.51	0.41
1:B:306:VAL:HG22	1:B:311:GLU:HA	2.03	0.40
1:A:88:SER:C	1:A:90:ASN:H	2.23	0.40
1:A:348:ALA:HB2	1:A:365:SER:HA	2.03	0.40
1:C:195:ALA:HB2	1:C:222:PRO:HA	2.03	0.40
1:A:402:PHE:CE2	1:A:408:LEU:CD1	3.04	0.40
1:B:5:ILE:HG23	1:B:6:LEU:N	2.36	0.40
1:B:243:LEU:N	1:B:243:LEU:CD2	2.84	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	417/432 (96%)	389 (93%)	22 (5%)	6 (1%)	11	36
1	B	401/432 (93%)	376 (94%)	21 (5%)	4 (1%)	15	45
1	C	404/432 (94%)	379 (94%)	20 (5%)	5 (1%)	13	40
1	D	404/432 (94%)	379 (94%)	20 (5%)	5 (1%)	13	40
All	All	1626/1728 (94%)	1523 (94%)	83 (5%)	20 (1%)	13	40

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	42	THR
1	B	327	ARG

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Mol	Chain	Res	Type
1	C	355	PRO
1	D	355	PRO
1	A	380	GLU
1	B	380	GLU
1	C	120	ILE
1	C	330	LEU
1	D	330	LEU
1	B	236	THR
1	C	354	SER
1	D	120	ILE
1	D	354	SER
1	A	89	LEU
1	A	236	THR
1	A	115	SER
1	B	115	SER
1	C	151	ALA
1	D	151	ALA
1	A	30	ARG

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	312/341 (92%)	286 (92%)	26 (8%)	11 32
1	B	310/341 (91%)	285 (92%)	25 (8%)	11 33
1	C	312/341 (92%)	285 (91%)	27 (9%)	10 30
1	D	312/341 (92%)	287 (92%)	25 (8%)	12 33
All	All	1246/1364 (91%)	1143 (92%)	103 (8%)	11 32

All (103) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	ASP
1	A	29	GLN
1	A	36	TYR

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Mol	Chain	Res	Type
1	A	42	THR
1	A	58	ARG
1	A	75	THR
1	A	83	ASP
1	A	87	ARG
1	A	88	SER
1	A	90	ASN
1	A	93	ASP
1	A	110	VAL
1	A	153	ASP
1	A	179	TYR
1	A	189	THR
1	A	200	ASN
1	A	221	VAL
1	A	243	LEU
1	A	258	VAL
1	A	293	GLN
1	A	307	LEU
1	A	330	LEU
1	A	338	LEU
1	A	371	LYS
1	A	388	ILE
1	A	417	ILE
1	B	22	ASP
1	B	29	GLN
1	B	36	TYR
1	B	58	ARG
1	B	75	THR
1	B	110	VAL
1	B	126	GLU
1	B	153	ASP
1	B	156	ARG
1	B	179	TYR
1	B	189	THR
1	B	200	ASN
1	B	209	ARG
1	B	221	VAL
1	B	243	LEU
1	B	258	VAL
1	B	293	GLN
1	B	307	LEU
1	B	330	LEU

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Mol	Chain	Res	Type
1	B	338	LEU
1	B	371	LYS
1	B	388	ILE
1	B	406	ARG
1	B	414	LYS
1	B	417	ILE
1	C	5	ILE
1	C	21	LEU
1	C	36	TYR
1	C	80	ASP
1	C	97	GLU
1	C	138	ARG
1	C	142	LYS
1	C	145	SER
1	C	162	GLU
1	C	175	GLN
1	C	179	TYR
1	C	243	LEU
1	C	246	GLN
1	C	258	VAL
1	C	270	ARG
1	C	342	LEU
1	C	349	GLU
1	C	351	LYS
1	C	354	SER
1	C	361	LEU
1	C	370	SER
1	C	377	THR
1	C	384	SER
1	C	396	VAL
1	C	403	LEU
1	C	412	ARG
1	C	416	SER
1	D	5	ILE
1	D	21	LEU
1	D	36	TYR
1	D	80	ASP
1	D	94	THR
1	D	138	ARG
1	D	145	SER
1	D	162	GLU
1	D	175	GLN

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Mol	Chain	Res	Type
1	D	179	TYR
1	D	243	LEU
1	D	246	GLN
1	D	258	VAL
1	D	342	LEU
1	D	349	GLU
1	D	351	LYS
1	D	354	SER
1	D	361	LEU
1	D	370	SER
1	D	377	THR
1	D	384	SER
1	D	396	VAL
1	D	403	LEU
1	D	412	ARG
1	D	416	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (30) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	GLN
1	A	90	ASN
1	A	191	GLN
1	A	200	ASN
1	A	289	GLN
1	B	17	GLN
1	B	191	GLN
1	B	200	ASN
1	B	289	GLN
1	B	293	GLN
1	C	17	GLN
1	C	29	GLN
1	C	175	GLN
1	C	191	GLN
1	C	246	GLN
1	C	259	ASN
1	C	285	GLN
1	C	289	GLN
1	C	292	GLN
1	C	398	GLN
1	D	17	GLN
1	D	29	GLN

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Mol	Chain	Res	Type
1	D	175	GLN
1	D	191	GLN
1	D	246	GLN
1	D	259	ASN
1	D	285	GLN
1	D	289	GLN
1	D	292	GLN
1	D	398	GLN

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 monosaccharides 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 (i)

6.1 Protein, DNA and RNA chains (i)

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	419/432 (96%)	-0.09	12 (2%)	51	47	20, 39, 72, 79	0
1	B	405/432 (93%)	-0.12	10 (2%)	57	55	20, 39, 72, 85	0
1	C	407/432 (94%)	-0.11	13 (3%)	47	43	20, 40, 63, 73	0
1	D	407/432 (94%)	-0.22	4 (0%)	82	82	20, 40, 62, 73	0
All	All	1638/1728 (94%)	-0.13	39 (2%)	59	56	20, 39, 67, 85	0

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	114	ASP	4.1
1	C	236	THR	3.8
1	D	114	ASP	3.7
1	A	153	ASP	3.5
1	B	381	GLY	3.2
1	A	114	ASP	3.2
1	C	94	THR	3.1
1	C	114	ASP	2.9
1	B	115	SER	2.9
1	A	380	GLU	2.9
1	C	235	SER	2.9
1	A	399	SER	2.8
1	A	352	PRO	2.8
1	A	381	GLY	2.8
1	B	116	PRO	2.8
1	A	423	ILE	2.8
1	C	399	SER	2.5
1	B	409	VAL	2.5
1	C	246	GLN	2.5
1	C	237	GLY	2.5
1	C	29	GLN	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	410	LEU	2.4
1	B	408	LEU	2.4
1	C	153	ASP	2.3
1	C	80	ASP	2.3
1	A	408	LEU	2.3
1	B	153	ASP	2.3
1	C	354	SER	2.3
1	A	236	THR	2.2
1	A	409	VAL	2.2
1	D	80	ASP	2.2
1	D	237	GLY	2.2
1	D	236	THR	2.2
1	C	381	GLY	2.2
1	A	116	PRO	2.1
1	B	380	GLU	2.1
1	B	236	THR	2.1
1	B	410	LEU	2.1
1	C	423	ILE	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.