Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:21 PM GMT

PDB ID	:	1OKJ
Title	:	crystal structure of the essential E. coli YeaZ protein by MAD method using the gadolinium complex "DOTMA"
Authors	:	Abergel, C.; Jeudy, S.; Claverie, J.M.
Deposited on	:	2003-07-26
Resolution	:	2.28 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:

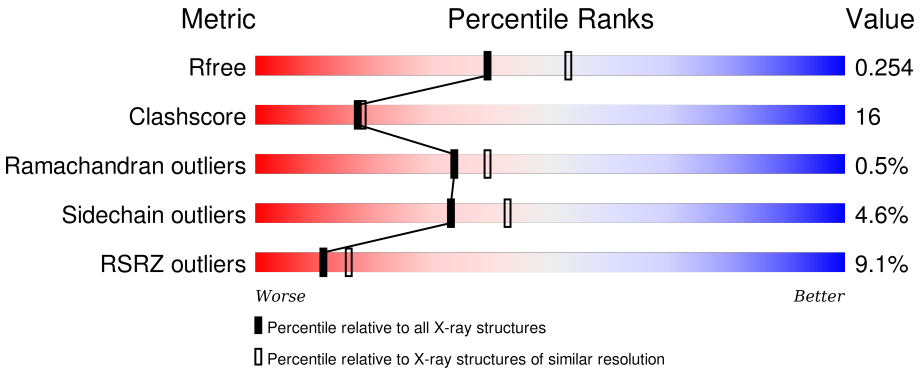
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	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 2.28 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	5193 (2.30-2.26)
Clashscore	102246	5929 (2.30-2.26)
Ramachandran outliers	100387	5851 (2.30-2.26)
Sidechain outliers	100360	5850 (2.30-2.26)
RSRZ outliers	91569	5204 (2.30-2.26)

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 <=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	251	14% 59% 25% • 12%
1	B	251	4% 73% 14% • 11%
1	C	251	9% 63% 22% • 13%
1	D	251	4% 58% 27% • 12%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6983 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 TRNA THREONYLCARBAMOYLADENOSINE BIOSYNTHESIS PROTEIN TSAB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	220	Total	C	N	O	S	0	0	1
			1663	1048	288	316	11			
1	B	223	Total	C	N	O	S	0	0	1
			1686	1063	293	319	11			
1	C	219	Total	C	N	O	S	0	0	1
			1655	1042	287	315	11			
1	D	221	Total	C	N	O	S	0	0	1
			1668	1051	289	317	11			

- Molecule 2 is GADOLINIUM ION (three-letter code: GD3) (formula: Gd).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	8	Total	Gd	0	0
			8	8		

- Molecule 3 is water.

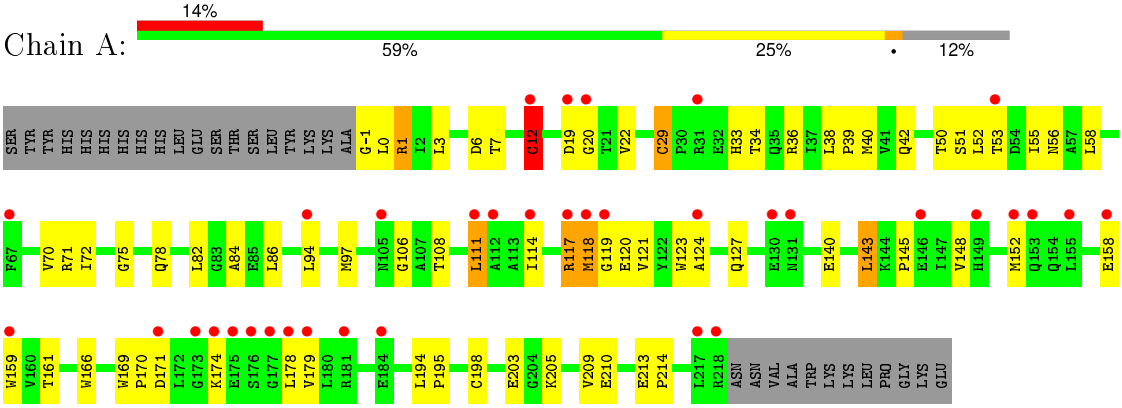
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	64	Total	O	0	0
			64	64		
3	B	95	Total	O	0	0
			95	95		
3	C	66	Total	O	0	0
			66	66		
3	D	78	Total	O	0	0
			78	78		

3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

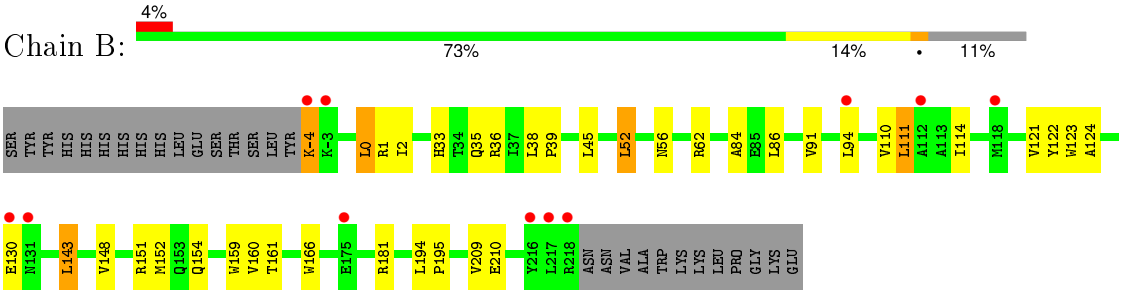
• Molecule 1: TRNA THREONYLCARBAMOYLADENOSINE BIOSYNTHESIS PROTEIN TSAB

Chain A:



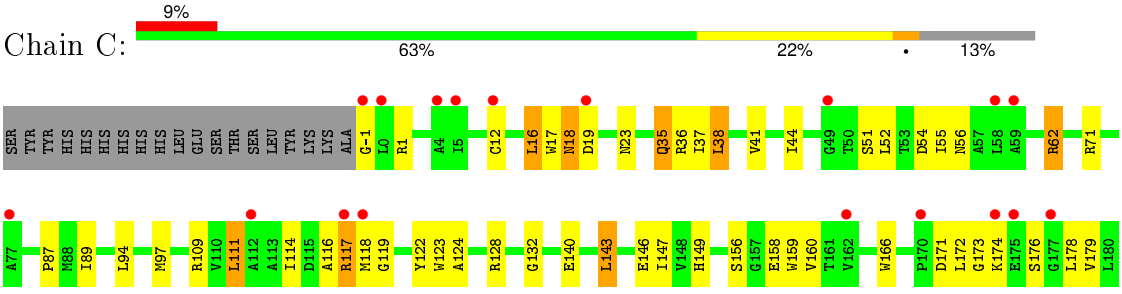
• Molecule 1: TRNA THREONYLCARBAMOYLADENOSINE BIOSYNTHESIS PROTEIN TSAB

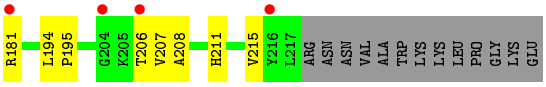
Chain B:



• Molecule 1: TRNA THREONYLCARBAMOYLADENOSINE BIOSYNTHESIS PROTEIN TSAB

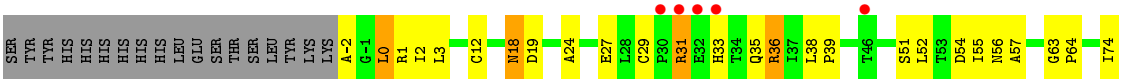
Chain C:

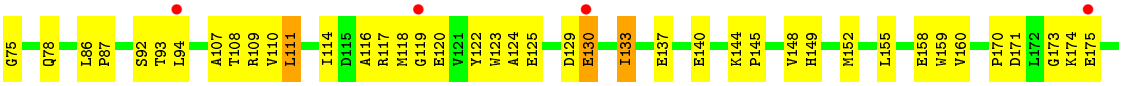


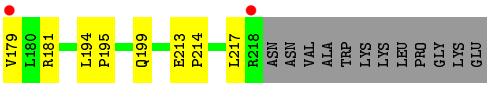


• Molecule 1: TRNA THREONYLCARBAMOYLADENOSINE BIOSYNTHESIS PROTEIN TSAB

Chain D:







4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	76.34Å 97.60Å 141.94Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.28 24.62 – 2.28	Depositor EDS
% Data completeness (in resolution range)	98.3 (20.00-2.28) 98.4 (24.62-2.28)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	4.13 (at 2.28Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.216 , 0.257 0.215 , 0.254	Depositor DCC
R_{free} test set	4871 reflections (11.23%)	DCC
Wilson B-factor (Å^2)	36.1	Xtriage
Anisotropy	0.118	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 45.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Outliers	1 of 48274 reflections (0.002%)	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	6983	wwPDB-VP
Average B, all atoms (Å^2)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.66% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GD3

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	0.34	0/1697	0.62	2/2311 (0.1%)
1	B	0.38	0/1720	0.63	0/2340
1	C	0.34	0/1689	0.63	2/2300 (0.1%)
1	D	0.36	0/1702	0.63	0/2318
All	All	0.36	0/6808	0.63	4/9269 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	12	CYS	CA-CB-SG	-6.77	101.82	114.00
1	C	38	LEU	CA-CB-CG	-5.67	102.26	115.30
1	A	29	CYS	CA-CB-SG	-5.52	104.06	114.00
1	C	119	GLY	N-CA-C	-5.04	100.51	113.10

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	1663	0	1638	64	0
1	B	1686	0	1669	34	0
1	C	1655	0	1627	51	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	1668	0	1645	71	0
2	A	8	0	0	0	0
3	A	64	0	0	1	0
3	B	95	0	0	0	0
3	C	66	0	0	2	0
3	D	78	0	0	1	0
All	All	6983	0	6579	213	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 16.

All (213) 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:1:ARG:H	1:B:56:ASN:HD22	1.08	1.00
1:A:1:ARG:HB2	1:A:1:ARG:HH11	1.26	0.97
1:A:1:ARG:H	1:A:56:ASN:HD22	1.13	0.95
1:D:1:ARG:H	1:D:56:ASN:HD22	1.14	0.95
1:D:64:PRO:O	1:D:93:THR:HG21	1.69	0.92
1:D:174:LYS:O	1:D:175:GLU:HG2	1.71	0.88
1:A:117:ARG:HH11	1:A:120:GLU:HG3	1.37	0.87
1:C:62:ARG:HD3	1:C:215:VAL:HB	1.60	0.84
1:C:147:ILE:HD12	3:C:2045:HOH:O	1.83	0.79
1:A:171:ASP:HA	1:A:174:LYS:HD3	1.64	0.79
1:A:51:SER:OG	1:A:53:THR:HG22	1.85	0.75
1:A:42:GLN:HE22	1:B:210:GLU:HB3	1.50	0.75
1:D:1:ARG:H	1:D:56:ASN:ND2	1.84	0.75
1:B:111:LEU:HD13	1:B:159:TRP:CE3	2.21	0.75
1:C:123:TRP:HB3	1:C:143:LEU:HD21	1.68	0.75
1:B:1:ARG:H	1:B:56:ASN:ND2	1.86	0.74
1:D:31:ARG:HG2	1:D:33:HIS:H	1.52	0.73
1:C:143:LEU:HD23	1:C:143:LEU:N	2.02	0.73
1:D:158:GLU:HB3	1:D:181:ARG:NH2	2.04	0.72
1:C:171:ASP:HA	1:C:174:LYS:HD3	1.71	0.72
1:A:38:LEU:O	1:A:42:GLN:HG3	1.91	0.71
1:D:63:GLY:C	1:D:93:THR:HG22	2.11	0.71
1:A:1:ARG:HH12	1:A:56:ASN:HD21	1.39	0.70
1:D:111:LEU:HD13	1:D:159:TRP:CE3	2.25	0.70
1:A:198:CYS:HB3	3:A:2055:HOH:O	1.92	0.69
1:D:31:ARG:HD2	1:D:31:ARG:H	1.56	0.68
1:A:170:PRO:HG3	1:D:24:ALA:O	1.93	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:18:ASN:C	1:D:18:ASN:HD22	1.97	0.67
1:D:31:ARG:HG3	1:D:33:HIS:CG	2.30	0.66
1:B:151:ARG:HH11	1:B:154:GLN:NE2	1.94	0.66
1:B:62:ARG:HG2	1:B:91:VAL:O	1.96	0.65
1:D:27:GLU:OE1	1:D:36:ARG:NH1	2.30	0.65
1:A:1:ARG:HH12	1:A:56:ASN:ND2	1.94	0.65
1:D:108:THR:HG23	1:D:109:ARG:HG3	1.79	0.65
1:A:117:ARG:NH1	1:A:120:GLU:HG3	2.09	0.64
1:D:152:MET:HE3	1:D:173:GLY:HA2	1.79	0.64
1:A:111:LEU:HD13	1:A:159:TRP:CE3	2.32	0.64
1:D:120:GLU:HG2	1:D:144:LYS:HD2	1.79	0.64
1:D:129:ASP:OD2	1:D:133:ILE:HG23	1.98	0.64
1:D:38:LEU:HB2	1:D:39:PRO:HD3	1.80	0.63
1:D:117:ARG:HG2	1:D:117:ARG:HH11	1.63	0.62
1:D:12:CYS:CB	1:D:29:CYS:SG	2.88	0.61
1:C:62:ARG:CD	1:C:215:VAL:HB	2.28	0.61
1:C:111:LEU:HD13	1:C:159:TRP:CE3	2.35	0.61
1:A:36:ARG:O	1:A:40:MET:HG3	2.00	0.61
1:A:1:ARG:H	1:A:56:ASN:ND2	1.93	0.61
1:A:29:CYS:HB3	1:A:33:HIS:O	2.01	0.61
1:A:94:LEU:HD13	1:A:124:ALA:HB2	1.82	0.60
1:B:160:VAL:HA	1:B:181:ARG:O	2.02	0.60
1:A:84:ALA:O	1:A:86:LEU:HD13	2.01	0.60
1:C:117:ARG:CA	1:C:117:ARG:HH11	2.14	0.60
1:A:170:PRO:O	1:A:174:LYS:HE2	2.01	0.59
1:A:0:LEU:HD23	1:A:1:ARG:N	2.17	0.59
1:D:63:GLY:O	1:D:93:THR:HG22	2.03	0.59
1:A:194:LEU:HB2	1:A:195:PRO:HD3	1.84	0.59
1:B:111:LEU:HG	1:B:123:TRP:CH2	2.38	0.58
1:D:194:LEU:HB2	1:D:195:PRO:HD3	1.83	0.58
1:A:210:GLU:HB2	1:B:39:PRO:HG3	1.84	0.58
1:A:1:ARG:HH11	1:A:1:ARG:CB	2.08	0.58
1:D:31:ARG:HG3	1:D:33:HIS:ND1	2.19	0.58
1:D:75:GLY:HA2	1:D:78:GLN:HE21	1.68	0.58
1:D:148:VAL:HG12	1:D:152:MET:HE2	1.86	0.58
1:A:148:VAL:O	1:A:152:MET:HG3	2.04	0.58
1:D:0:LEU:HD22	1:D:2:ILE:HG13	1.85	0.58
1:D:130:GLU:H	1:D:130:GLU:CD	2.07	0.58
1:B:151:ARG:HH11	1:B:154:GLN:HE21	1.50	0.57
1:C:62:ARG:HG2	1:C:62:ARG:O	2.05	0.57
1:A:7:THR:O	1:A:33:HIS:HE1	1.89	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:152:MET:O	1:A:178:LEU:HD11	2.06	0.56
1:C:117:ARG:HH11	1:C:117:ARG:CB	2.20	0.55
1:B:1:ARG:N	1:B:56:ASN:HD22	1.92	0.55
1:A:1:ARG:HB2	1:A:1:ARG:NH1	2.10	0.54
1:D:107:ALA:HB1	1:D:160:VAL:HG21	1.90	0.54
1:A:171:ASP:HA	1:A:174:LYS:CD	2.36	0.54
1:D:152:MET:CE	1:D:173:GLY:HA2	2.38	0.54
1:A:161:THR:HB	1:A:166:TRP:CE2	2.43	0.54
1:D:-2:ALA:N	3:D:2001:HOH:O	2.40	0.54
1:B:143:LEU:HD13	1:B:148:VAL:CG2	2.38	0.54
1:A:1:ARG:NH1	1:A:56:ASN:ND2	2.55	0.53
1:C:38:LEU:HD13	1:D:78:GLN:HB3	1.89	0.53
1:A:70:VAL:HG13	1:A:71:ARG:N	2.23	0.53
1:A:75:GLY:HA2	1:A:78:GLN:HE21	1.73	0.53
1:B:143:LEU:HD13	1:B:148:VAL:HG22	1.91	0.53
1:C:143:LEU:CD2	1:C:143:LEU:N	2.71	0.53
1:C:123:TRP:O	1:C:140:GLU:HA	2.09	0.53
1:A:121:VAL:HG12	1:A:143:LEU:O	2.09	0.52
1:D:114:ILE:HB	1:D:122:TYR:HB2	1.92	0.52
1:A:82:LEU:C	1:A:82:LEU:HD23	2.30	0.52
1:A:42:GLN:NE2	1:B:210:GLU:HB3	2.24	0.52
1:C:38:LEU:CD1	1:D:78:GLN:HB3	2.40	0.52
1:C:194:LEU:HB2	1:C:195:PRO:HD3	1.92	0.52
1:D:1:ARG:NH1	1:D:18:ASN:HD21	2.09	0.51
1:B:110:VAL:HG22	1:B:160:VAL:CG2	2.41	0.51
1:D:33:HIS:HB3	1:D:35:GLN:HG2	1.92	0.51
1:C:117:ARG:NH1	1:C:117:ARG:HB3	2.24	0.51
1:D:213:GLU:HB2	1:D:214:PRO:HD2	1.90	0.51
1:D:31:ARG:HG2	1:D:33:HIS:N	2.24	0.51
1:A:203:GLU:OE1	1:A:205:LYS:HE2	2.10	0.51
1:C:160:VAL:HA	1:C:181:ARG:O	2.10	0.51
1:B:148:VAL:HG12	1:B:152:MET:HE2	1.92	0.51
1:D:137:GLU:O	1:D:140:GLU:HG3	2.10	0.51
1:A:3:LEU:HB2	1:A:55:ILE:HD13	1.91	0.51
1:C:37:ILE:O	1:C:41:VAL:HG23	2.11	0.51
1:D:52:LEU:HD12	1:D:55:ILE:HD12	1.93	0.50
1:C:52:LEU:HD23	1:C:55:ILE:HD12	1.93	0.50
1:C:207:VAL:CG2	1:C:211:HIS:HB2	2.40	0.50
1:C:158:GLU:HG2	1:C:179:VAL:HB	1.91	0.50
1:B:148:VAL:HG12	1:B:152:MET:CE	2.41	0.50
1:C:97:MET:HE2	1:C:114:ILE:HD11	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:1:ARG:NH1	1:D:18:ASN:ND2	2.60	0.50
1:D:119:GLY:O	1:D:145:PRO:HD3	2.12	0.50
1:C:117:ARG:HH11	1:C:117:ARG:HA	1.76	0.49
1:D:31:ARG:CG	1:D:33:HIS:CG	2.94	0.49
1:D:74:ILE:O	1:D:78:GLN:HG3	2.12	0.49
1:C:51:SER:HB3	1:C:54:ASP:OD2	2.12	0.49
1:C:-1:GLY:HA3	1:C:19:ASP:OD1	2.13	0.49
1:A:6:ASP:O	1:A:12:CYS:HA	2.13	0.48
1:D:158:GLU:HB3	1:D:181:ARG:HH21	1.76	0.48
1:C:71:ARG:HH11	1:C:71:ARG:HG3	1.78	0.48
1:A:117:ARG:HG3	1:A:118:MET:H	1.77	0.48
1:A:36:ARG:C	1:A:39:PRO:HD2	2.33	0.48
1:C:146:GLU:O	1:C:149:HIS:HB3	2.13	0.48
1:D:1:ARG:HH11	1:D:1:ARG:HG2	1.78	0.48
1:B:84:ALA:CB	1:B:86:LEU:HD13	2.43	0.48
1:A:117:ARG:HG3	1:A:120:GLU:HB2	1.96	0.48
1:D:31:ARG:HH11	1:D:31:ARG:HG3	1.79	0.48
1:D:160:VAL:HA	1:D:181:ARG:O	2.14	0.48
1:B:84:ALA:HB3	1:B:86:LEU:HD13	1.94	0.48
1:D:158:GLU:HG2	1:D:179:VAL:HB	1.96	0.48
1:B:0:LEU:HD22	1:B:2:ILE:HG13	1.96	0.47
1:A:210:GLU:CB	1:B:39:PRO:HG3	2.44	0.47
1:A:111:LEU:HG	1:A:123:TRP:CH2	2.50	0.47
1:A:97:MET:HE2	1:A:114:ILE:HD11	1.96	0.47
1:B:94:LEU:HD13	1:B:124:ALA:HB2	1.97	0.47
1:D:18:ASN:C	1:D:18:ASN:ND2	2.64	0.47
1:D:116:ALA:O	1:D:118:MET:HG3	2.14	0.47
1:D:195:PRO:O	1:D:199:GLN:HG3	2.15	0.47
1:D:51:SER:HB3	1:D:54:ASP:OD2	2.14	0.47
1:D:170:PRO:O	1:D:171:ASP:HB2	2.15	0.46
1:D:110:VAL:HG22	1:D:160:VAL:CG2	2.45	0.46
1:A:94:LEU:HD12	1:A:140:GLU:HG2	1.96	0.46
1:A:-1:GLY:HA3	1:A:19:ASP:OD1	2.15	0.46
1:B:-4:LYS:NZ	1:B:-4:LYS:HB3	2.30	0.46
1:C:117:ARG:HH11	1:C:117:ARG:HB3	1.80	0.46
1:B:114:ILE:HB	1:B:122:TYR:HB2	1.97	0.46
1:A:158:GLU:HB3	1:A:179:VAL:HB	1.97	0.46
1:D:12:CYS:SG	1:D:29:CYS:SG	3.13	0.46
1:D:12:CYS:HB3	1:D:29:CYS:SG	2.57	0.46
1:C:116:ALA:C	1:C:118:MET:H	2.20	0.46
1:C:35:GLN:H	1:C:35:GLN:HG3	1.39	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:119:GLY:O	1:A:145:PRO:HD3	2.15	0.45
1:C:166:TRP:CE3	1:C:172:LEU:HD23	2.51	0.45
1:B:38:LEU:N	1:B:38:LEU:HD22	2.32	0.45
1:A:12:CYS:SG	1:A:29:CYS:N	2.89	0.45
1:C:23:ASN:ND2	1:C:44:ILE:HG23	2.32	0.45
1:A:82:LEU:O	1:A:82:LEU:HD23	2.17	0.45
1:D:111:LEU:HD22	1:D:159:TRP:CB	2.47	0.45
1:C:89:ILE:HG12	1:C:206:THR:HG22	1.99	0.45
1:C:156:SER:HA	1:C:178:LEU:CD2	2.47	0.45
1:C:52:LEU:HA	1:C:55:ILE:HD12	1.99	0.45
1:C:94:LEU:HD13	1:C:124:ALA:HB2	2.00	0.44
1:D:149:HIS:HA	1:D:152:MET:CE	2.47	0.44
1:D:1:ARG:NH1	1:D:19:ASP:OD1	2.50	0.44
1:C:114:ILE:HB	1:C:122:TYR:HB2	1.98	0.44
1:C:16:LEU:HB2	1:C:44:ILE:HG21	1.99	0.44
1:B:161:THR:HG22	1:B:166:TRP:CZ2	2.52	0.44
1:D:123:TRP:O	1:D:140:GLU:HA	2.18	0.44
1:C:158:GLU:HG2	1:C:179:VAL:CB	2.48	0.44
1:C:18:ASN:O	1:C:19:ASP:HB2	2.18	0.44
1:D:1:ARG:HH11	1:D:18:ASN:ND2	2.15	0.44
1:A:169:TRP:C	1:A:171:ASP:H	2.21	0.44
1:A:209:VAL:HG23	1:A:210:GLU:N	2.33	0.44
1:A:108:THR:O	1:A:127:GLN:HA	2.17	0.44
1:B:62:ARG:HG2	1:B:91:VAL:C	2.37	0.43
1:D:94:LEU:HD13	1:D:124:ALA:HB2	1.99	0.43
1:C:1:ARG:HA	1:C:17:TRP:O	2.19	0.43
1:A:34:THR:HG21	1:A:72:ILE:CG2	2.49	0.43
1:D:31:ARG:HG3	1:D:31:ARG:NH1	2.34	0.43
1:B:194:LEU:HB2	1:B:195:PRO:HD3	2.00	0.43
1:C:36:ARG:HG3	3:C:2015:HOH:O	2.17	0.43
1:A:1:ARG:N	1:A:56:ASN:HD22	1.96	0.43
1:A:7:THR:HG21	1:A:34:THR:HG22	2.01	0.43
1:B:36:ARG:O	1:B:39:PRO:HD2	2.18	0.43
1:D:125:GLU:OE2	1:D:155:LEU:HD11	2.18	0.43
1:D:86:LEU:HA	1:D:87:PRO:HD3	1.85	0.43
1:A:52:LEU:HD23	1:A:55:ILE:HD12	2.01	0.43
1:C:207:VAL:HG22	1:C:208:ALA:N	2.33	0.43
1:B:45:LEU:HD11	1:B:52:LEU:HD13	2.00	0.43
1:D:18:ASN:HD22	1:D:19:ASP:N	2.16	0.42
1:C:56:ASN:O	1:C:87:PRO:HD2	2.18	0.42
1:D:55:ILE:HG22	1:D:57:ALA:H	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:62:ARG:CG	1:C:215:VAL:HB	2.49	0.42
1:A:7:THR:O	1:A:33:HIS:CE1	2.71	0.42
1:A:22:VAL:HG23	1:A:22:VAL:O	2.20	0.42
1:B:36:ARG:C	1:B:39:PRO:HD2	2.39	0.42
1:B:130:GLU:HA	1:B:130:GLU:OE2	2.19	0.42
1:D:111:LEU:HG	1:D:123:TRP:CH2	2.55	0.41
1:A:117:ARG:O	1:A:118:MET:O	2.37	0.41
1:A:55:ILE:O	1:A:86:LEU:HD21	2.20	0.41
1:B:143:LEU:CD1	1:B:148:VAL:HG22	2.50	0.41
1:C:173:GLY:O	1:C:176:SER:HB3	2.20	0.41
1:A:50:THR:HG22	1:A:51:SER:N	2.35	0.41
1:C:109:ARG:HD3	1:C:158:GLU:O	2.21	0.41
1:A:117:ARG:HG3	1:A:118:MET:N	2.35	0.41
1:C:128:ARG:NH1	1:C:132:GLY:HA2	2.35	0.41
1:B:52:LEU:HD12	1:B:52:LEU:HA	1.88	0.41
1:C:124:ALA:HB2	1:C:140:GLU:HG2	2.03	0.41
1:D:149:HIS:HA	1:D:152:MET:HE2	2.02	0.41
1:C:143:LEU:HD23	1:C:143:LEU:H	1.83	0.40
1:D:117:ARG:HG2	1:D:117:ARG:NH1	2.33	0.40
1:A:213:GLU:HB2	1:A:214:PRO:CD	2.50	0.40
1:D:3:LEU:HB2	1:D:55:ILE:HD13	2.03	0.40
1:C:111:LEU:HG	1:C:123:TRP:CH2	2.56	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	218/251 (87%)	200 (92%)	14 (6%)	4 (2%)	11 8
1	B	221/251 (88%)	213 (96%)	8 (4%)	0	100 100
1	C	217/251 (86%)	206 (95%)	11 (5%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	D	219/251 (87%)	215 (98%)	4 (2%)	0	100 100
All	All	875/1004 (87%)	834 (95%)	37 (4%)	4 (0%)	34 39

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	117	ARG
1	A	118	MET
1	A	20	GLY
1	A	106	GLY

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	170/199 (85%)	165 (97%)	5 (3%)	50 65
1	B	172/199 (86%)	163 (95%)	9 (5%)	29 36
1	C	169/199 (85%)	161 (95%)	8 (5%)	32 42
1	D	170/199 (85%)	161 (95%)	9 (5%)	28 35
All	All	681/796 (86%)	650 (95%)	31 (5%)	33 43

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	ARG
1	A	12	CYS
1	A	58	LEU
1	A	111	LEU
1	A	143	LEU
1	B	-4	LYS
1	B	0	LEU
1	B	33	HIS
1	B	35	GLN
1	B	52	LEU

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Mol	Chain	Res	Type
1	B	111	LEU
1	B	121	VAL
1	B	143	LEU
1	B	209	VAL
1	C	12	CYS
1	C	16	LEU
1	C	18	ASN
1	C	35	GLN
1	C	62	ARG
1	C	111	LEU
1	C	117	ARG
1	C	143	LEU
1	D	0	LEU
1	D	18	ASN
1	D	31	ARG
1	D	36	ARG
1	D	92	SER
1	D	111	LEU
1	D	130	GLU
1	D	133	ILE
1	D	217	LEU

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

Mol	Chain	Res	Type
1	A	23	ASN
1	A	33	HIS
1	A	42	GLN
1	A	56	ASN
1	A	78	GLN
1	A	153	GLN
1	A	199	GLN
1	B	18	ASN
1	B	35	GLN
1	B	56	ASN
1	B	78	GLN
1	B	131	ASN
1	B	149	HIS
1	B	154	GLN
1	B	199	GLN
1	C	18	ASN
1	C	23	ASN

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Mol	Chain	Res	Type
1	C	35	GLN
1	C	78	GLN
1	C	105	ASN
1	C	149	HIS
1	D	18	ASN
1	D	25	HIS
1	D	35	GLN
1	D	42	GLN
1	D	56	ASN
1	D	78	GLN
1	D	105	ASN
1	D	153	GLN
1	D	211	HIS

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 carbohydrates in this entry.

5.6 Ligand geometry

Of 8 ligands modelled in this entry, 8 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	220/251 (87%)	0.84	36 (16%) 2 3	21, 43, 69, 78	0
1	B	223/251 (88%)	0.30	11 (4%) 33 41	19, 31, 51, 66	0
1	C	219/251 (87%)	0.66	22 (10%) 9 13	25, 40, 60, 69	0
1	D	221/251 (88%)	0.33	11 (4%) 32 40	22, 35, 54, 69	0
All	All	883/1004 (87%)	0.53	80 (9%) 11 16	19, 37, 62, 78	0

All (80) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	118	MET	7.3
1	A	119	GLY	6.1
1	A	19	ASP	5.2
1	B	-4	LYS	4.6
1	C	-1	GLY	4.5
1	C	118	MET	4.4
1	B	118	MET	4.3
1	D	31	ARG	3.8
1	A	177	GLY	3.8
1	A	179	VAL	3.7
1	A	176	SER	3.7
1	C	174	LYS	3.5
1	A	153	GLN	3.4
1	A	130	GLU	3.4
1	A	174	LYS	3.4
1	C	206	THR	3.3
1	C	204	GLY	3.3
1	C	117	ARG	3.2
1	D	175	GLU	3.2
1	C	49	GLY	3.1
1	A	181	ARG	3.1

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Mol	Chain	Res	Type	RSRZ
1	B	-3	LYS	3.1
1	C	5	ILE	3.0
1	D	218	ARG	3.0
1	D	33	HIS	3.0
1	A	175	GLU	3.0
1	C	175	GLU	3.0
1	A	184	GLU	2.9
1	A	31	ARG	2.9
1	A	146	GLU	2.9
1	D	30	PRO	2.8
1	A	53	THR	2.8
1	B	130	GLU	2.8
1	C	19	ASP	2.8
1	C	59	ALA	2.8
1	A	124	ALA	2.8
1	A	117	ARG	2.8
1	A	173	GLY	2.7
1	B	217	LEU	2.7
1	A	149	HIS	2.7
1	A	178	LEU	2.6
1	A	152	MET	2.6
1	C	216	TYR	2.6
1	C	177	GLY	2.6
1	D	119	GLY	2.6
1	A	112	ALA	2.5
1	C	77	ALA	2.5
1	A	67	PHE	2.5
1	A	131	ASN	2.4
1	A	218	ARG	2.4
1	B	94	LEU	2.4
1	B	216	TYR	2.3
1	C	170	PRO	2.3
1	D	32	GLU	2.3
1	C	0	LEU	2.3
1	C	4	ALA	2.3
1	A	159	TRP	2.3
1	A	217	LEU	2.3
1	B	218	ARG	2.3
1	A	20	GLY	2.2
1	C	162	VAL	2.2
1	A	12	CYS	2.2
1	C	112	ALA	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	130	GLU	2.2
1	B	175	GLU	2.2
1	A	105	ASN	2.2
1	A	111	LEU	2.1
1	A	171	ASP	2.1
1	B	131	ASN	2.1
1	A	94	LEU	2.1
1	D	179	VAL	2.1
1	B	112	ALA	2.1
1	A	158	GLU	2.1
1	C	181	ARG	2.1
1	C	58	LEU	2.1
1	A	114	ILE	2.1
1	D	46	THR	2.0
1	D	94	LEU	2.0
1	A	155	LEU	2.0
1	C	12	CYS	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 carbohydrates in this entry.

6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å^2)	Q<0.9
2	GD3	A	1218	1/1	0.95	0.14	-	57,57,57,57	1
2	GD3	A	1217	1/1	0.85	0.08	-	34,34,34,34	1
2	GD3	A	1214	1/1	0.98	0.13	-	16,16,16,16	1
2	GD3	A	1213	1/1	0.89	0.07	-	41,41,41,41	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å^2)	Q<0.9
2	GD3	A	1219	1/1	0.98	0.04	-	35,35,35,35	1
2	GD3	A	1215	1/1	0.92	0.10	-	44,44,44,44	1
2	GD3	A	1216	1/1	0.94	0.07	-	38,38,38,38	1
2	GD3	A	1220	1/1	0.91	0.07	-	62,62,62,62	1

6.5 Other polymers (i)

There are no such residues in this entry.