Full wwPDB X-ray Structure Validation Report (i)

Jan 31, 2016 – 06:57 PM GMT

PDB ID	:	1DC5
Title	:	STRUCTURAL ANALYSIS OF GLYCERALDEHYDE 3-PHOSPHATE DEHYDROGENASE FROM ESCHERICHIA COLI: DIRECT EVIDENCE FOR SUBSTRATE BINDING AND COFACTOR-INDUCED CONFORMATIONAL CHANGES
Authors	:	Yun, M.; Park, C.-G.; Kim, J.-Y.; Park, H.-W.
Deposited on	:	1999-11-04
Resolution	:	2.00 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

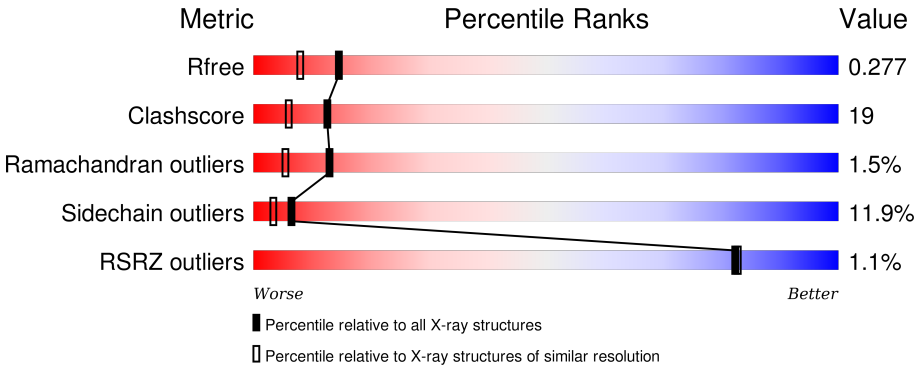
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	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 2.00 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	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

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	330	2% 58% 35% 7%
1	B	330	% 62% 33% 5%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 5206 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 GLYCERALDEHYDE-3-PHOSPHATE DEHYDROGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	330	Total	C	N	O	S	0	0	0
			2488	1563	431	484	10			
1	B	330	Total	C	N	O	S	0	0	0
			2488	1563	431	484	10			

- Molecule 2 is water.

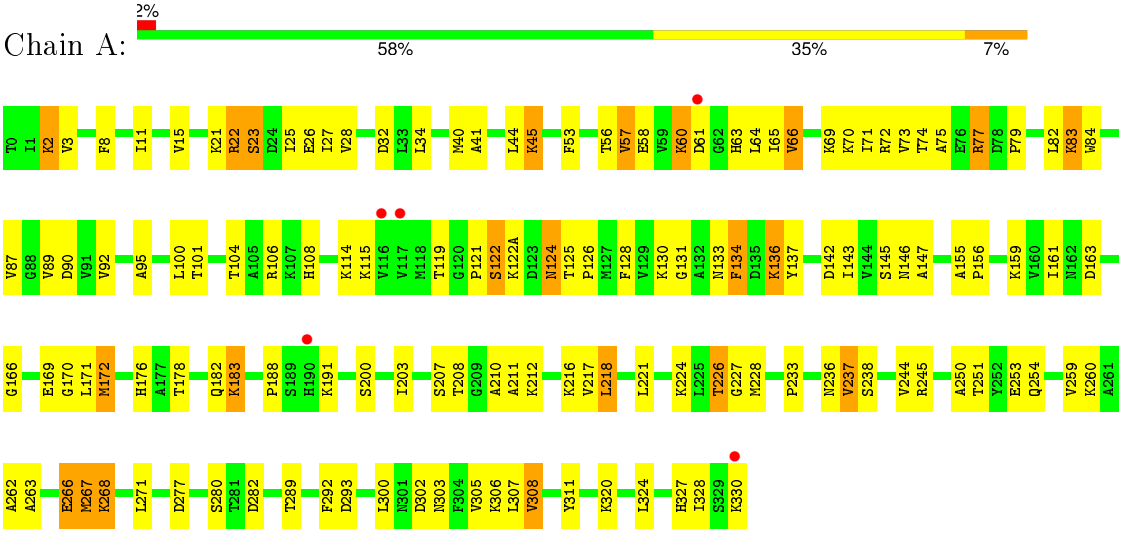
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	118	Total	O	0	0
			118	118		
2	B	112	Total	O	0	0
			112	112		

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.

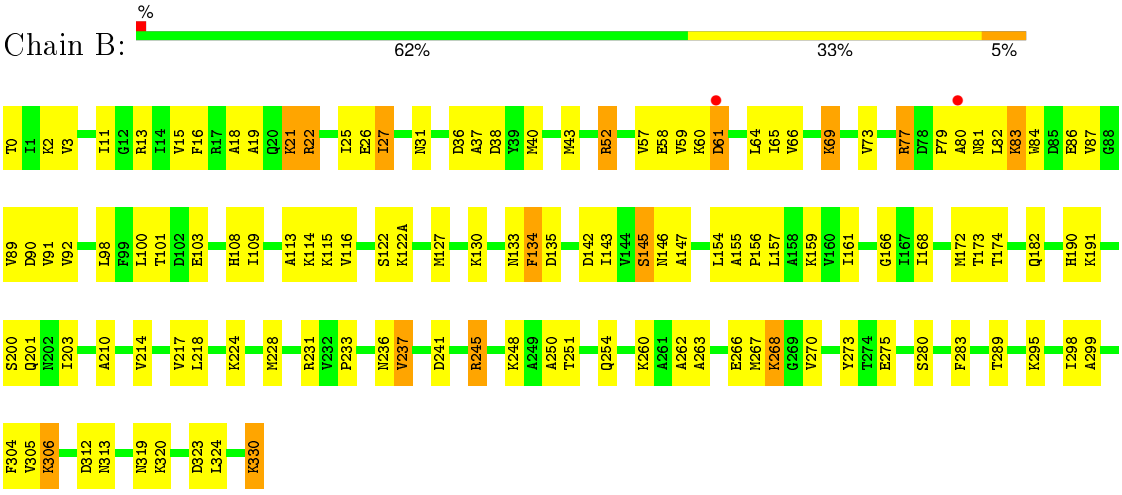
- Molecule 1: GLYCERALDEHYDE-3-PHOSPHATE DEHYDROGENASE

Chain A:



- Molecule 1: GLYCERALDEHYDE-3-PHOSPHATE DEHYDROGENASE

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	81.87Å 134.54Å 68.23Å 90.00° 106.12° 90.00°	Depositor
Resolution (Å)	20.00 – 2.00 19.58 – 2.03	Depositor EDS
% Data completeness (in resolution range)	(Not available) (20.00-2.00) 86.2 (19.58-2.03)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.42 (at 2.02Å)	Xtriage
Refinement program	X-PLOR	Depositor
R, R_{free}	0.206 , 0.281 0.201 , 0.277	Depositor DCC
R_{free} test set	2009 reflections (5.10%)	DCC
Wilson B-factor (Å^2)	25.2	Xtriage
Anisotropy	0.264	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.40 , 72.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Outliers	0 of 42342 reflections	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	5206	wwPDB-VP
Average B, all atoms (Å^2)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.56% 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

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	0.40	0/2527	0.67	1/3421 (0.0%)
1	B	0.39	0/2527	0.67	1/3421 (0.0%)
All	All	0.39	0/5054	0.67	2/6842 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	203	ILE	N-CA-C	-7.12	91.77	111.00
1	A	203	ILE	N-CA-C	-6.09	94.56	111.00

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	2488	0	2506	98	0
1	B	2488	0	2506	97	0
2	A	118	0	0	8	0
2	B	112	0	0	7	0
All	All	5206	0	5012	194	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 19.

All (194) 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:245:ARG:HG3	1:B:245:ARG:HH11	1.26	1.00
1:A:262:ALA:HB1	1:A:267:MET:HG3	1.44	0.96
1:B:77:ARG:HH11	1:B:77:ARG:HG2	1.30	0.95
1:A:2:LYS:HG3	1:A:28:VAL:HG11	1.49	0.94
1:A:124:ASN:H	1:A:124:ASN:HD22	1.07	0.93
1:B:251:THR:H	1:B:254:GLN:HE21	1.09	0.92
1:B:251:THR:H	1:B:254:GLN:NE2	1.65	0.92
1:B:245:ARG:CG	1:B:245:ARG:HH11	1.87	0.87
1:A:251:THR:H	1:A:254:GLN:HE21	1.24	0.84
1:A:251:THR:H	1:A:254:GLN:NE2	1.75	0.83
1:B:18:ALA:O	1:B:21:LYS:HG2	1.78	0.83
1:B:77:ARG:NH1	1:B:77:ARG:HG2	1.98	0.79
1:B:82:LEU:HD13	1:B:84:TRP:CZ2	2.18	0.79
1:B:241:ASP:OD1	1:B:306:LYS:HD2	1.84	0.78
1:B:251:THR:N	1:B:254:GLN:HE21	1.83	0.76
1:A:170:GLY:HA3	1:A:244:VAL:HG12	1.69	0.75
1:B:245:ARG:NH1	1:B:245:ARG:HG3	2.01	0.74
1:B:90:ASP:O	1:B:114:LYS:HB3	1.86	0.74
1:A:124:ASN:H	1:A:124:ASN:ND2	1.86	0.73
1:B:159:LYS:HG3	1:B:218:LEU:HD22	1.72	0.71
1:B:155:ALA:HB3	1:B:156:PRO:HD3	1.71	0.71
1:A:32:ASP:O	1:A:75:ALA:HA	1.91	0.70
1:A:83:LYS:NZ	1:A:83:LYS:HB2	2.07	0.70
1:B:299:ALA:CB	1:B:305:VAL:HG12	2.22	0.70
1:B:319:ASN:HB3	2:B:2164:HOH:O	1.93	0.69
1:B:200:SER:HA	1:B:233:PRO:HB3	1.75	0.69
1:B:83:LYS:HB3	1:B:86:GLU:HG3	1.75	0.68
1:B:115:LYS:HG2	1:B:142:ASP:HA	1.75	0.68
1:B:58:GLU:HG2	1:B:65:ILE:HB	1.74	0.68
1:B:60:LYS:HB3	1:B:65:ILE:CD1	2.24	0.67
1:B:22:ARG:NH1	1:B:25:ILE:HD12	2.09	0.67
1:A:218:LEU:HD23	1:A:221:LEU:HD12	1.75	0.67
1:A:155:ALA:HB3	1:A:156:PRO:HD3	1.76	0.66
1:B:159:LYS:HG3	1:B:218:LEU:CD2	2.25	0.66
1:A:124:ASN:N	1:A:124:ASN:HD22	1.86	0.65
1:A:218:LEU:HD23	1:A:221:LEU:CD1	2.27	0.65
1:A:171:LEU:HD23	1:A:226:THR:HG22	1.78	0.65
1:A:289:THR:HG23	1:A:320:LYS:HE2	1.79	0.64
1:A:115:LYS:HD3	1:A:328:ILE:HD12	1.78	0.64
1:B:52:ARG:HG2	2:B:2038:HOH:O	1.97	0.64
1:B:116:VAL:HG13	1:B:143:ILE:HG23	1.79	0.63

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:210:ALA:O	1:B:214:VAL:HG23	1.99	0.63
1:B:262:ALA:HB1	1:B:267:MET:HG3	1.81	0.63
1:B:11:ILE:O	1:B:15:VAL:HG13	1.98	0.62
1:B:130:LYS:HG2	1:B:270:VAL:HG21	1.82	0.62
1:A:327:HIS:HA	1:A:330:LYS:NZ	2.14	0.62
1:B:130:LYS:NZ	2:B:2140:HOH:O	2.33	0.61
1:B:19:ALA:CB	1:B:27:ILE:HD11	2.30	0.61
1:A:266:GLU:HG2	1:A:267:MET:HG2	1.82	0.61
1:A:56:THR:O	1:A:66:VAL:HA	2.00	0.61
1:B:298:ILE:HD12	1:B:299:ALA:H	1.66	0.60
1:B:100:LEU:HB2	1:B:122(A):LYS:HG3	1.84	0.60
1:B:172:MET:HG2	1:B:173:THR:N	2.17	0.60
1:A:92:VAL:HG11	1:A:108:HIS:CD2	2.37	0.59
1:A:63:HIS:HD2	1:A:73:VAL:H	1.51	0.59
1:B:84:TRP:HB3	1:B:89:VAL:HB	1.85	0.58
1:B:182:GLN:HB3	2:B:2005:HOH:O	2.02	0.58
1:A:2:LYS:HG3	1:A:28:VAL:CG1	2.30	0.58
1:B:100:LEU:HD12	1:B:122(A):LYS:HE3	1.87	0.56
1:A:66:VAL:HG23	1:A:71:ILE:HG13	1.88	0.56
1:A:72:ARG:HD2	2:A:2032:HOH:O	2.05	0.56
1:A:293:ASP:HB3	1:A:308:VAL:HG22	1.87	0.55
1:B:26:GLU:OE2	1:B:26:GLU:HA	2.06	0.55
1:A:266:GLU:HG2	1:A:267:MET:N	2.22	0.55
1:B:22:ARG:NH1	1:B:25:ILE:CD1	2.70	0.55
1:B:245:ARG:CB	1:B:245:ARG:HH11	2.20	0.54
1:A:83:LYS:HZ3	1:A:83:LYS:HB2	1.72	0.54
1:B:330:LYS:HG2	1:B:330:LYS:OXT	2.08	0.54
1:B:79:PRO:HB3	1:B:108:HIS:CE1	2.44	0.53
1:B:22:ARG:HH12	1:B:25:ILE:HD12	1.71	0.53
1:A:72:ARG:HD3	1:A:74:THR:HG23	1.89	0.53
1:B:154:LEU:HD23	1:B:214:VAL:HG21	1.91	0.53
1:A:82:LEU:HD13	1:A:84:TRP:CZ2	2.44	0.53
1:A:172:MET:HG3	1:A:227:GLY:HA3	1.90	0.53
1:A:95:ALA:HA	1:A:119:THR:OG1	2.09	0.53
1:A:2:LYS:HB3	1:A:89:VAL:HA	1.90	0.52
1:B:3:VAL:HG13	1:B:27:ILE:HA	1.91	0.52
1:B:236:ASN:O	1:B:237:VAL:HB	2.10	0.52
1:A:65:ILE:HG22	2:A:2168:HOH:O	2.08	0.52
1:B:109:ILE:HG23	1:B:113:ALA:O	2.09	0.52
1:B:22:ARG:NH2	1:B:323:ASP:OD1	2.43	0.51
1:A:11:ILE:O	1:A:15:VAL:HG23	2.10	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:216:LYS:HE3	2:A:2109:HOH:O	2.10	0.51
1:A:251:THR:N	1:A:254:GLN:HE21	2.00	0.51
1:A:137:TYR:HB3	1:A:327:HIS:HE1	1.75	0.51
1:B:38:ASP:HA	1:B:59:VAL:HG21	1.91	0.51
1:A:100:LEU:HD13	1:A:121:PRO:O	2.10	0.51
1:A:159:LYS:HE3	1:A:163:ASP:OD1	2.10	0.51
1:B:92:VAL:HG11	1:B:108:HIS:CD2	2.46	0.51
1:B:57:VAL:HG23	1:B:66:VAL:HG22	1.92	0.51
1:B:299:ALA:HB1	1:B:305:VAL:HG12	1.91	0.51
1:A:327:HIS:HA	1:A:330:LYS:HZ2	1.76	0.51
1:B:275:GLU:O	1:B:295:LYS:HD2	2.11	0.51
1:A:263:ALA:O	1:A:268:LYS:HA	2.11	0.51
1:A:183:LYS:HE3	1:A:188:PRO:O	2.11	0.51
1:A:131:GLY:CA	1:A:267:MET:HE1	2.41	0.50
1:A:130:LYS:HG3	1:A:134:PHE:CE2	2.45	0.50
1:A:237:VAL:HG11	1:A:280:SER:O	2.11	0.50
1:A:236:ASN:O	1:A:237:VAL:HB	2.11	0.50
1:B:134:PHE:HB2	2:B:2224:HOH:O	2.09	0.50
1:B:174:THR:O	1:B:174:THR:HG23	2.12	0.50
1:A:327:HIS:CD2	1:A:330:LYS:NZ	2.80	0.50
1:A:25:ILE:CG2	1:A:26:GLU:N	2.74	0.50
1:B:133:ASN:HB3	1:B:217:VAL:O	2.12	0.50
1:A:327:HIS:CD2	1:A:330:LYS:HZ2	2.30	0.49
1:A:292:PHE:HE1	1:A:307:LEU:HD22	1.77	0.49
1:B:40:MET:CE	1:B:73:VAL:HG13	2.42	0.49
1:A:200:SER:HA	1:A:233:PRO:HB3	1.94	0.49
1:A:3:VAL:HB	1:A:27:ILE:HD13	1.95	0.49
1:B:31:ASN:CG	1:B:82:LEU:HD21	2.32	0.49
1:B:130:LYS:HG2	1:B:270:VAL:CG2	2.43	0.48
1:B:237:VAL:HG11	1:B:280:SER:O	2.13	0.48
1:B:127:MET:HA	1:B:145:SER:O	2.13	0.48
1:B:3:VAL:CG1	1:B:27:ILE:HD13	2.44	0.48
1:A:133:ASN:N	2:A:2123:HOH:O	2.47	0.48
1:A:136:LYS:HD3	2:A:2095:HOH:O	2.14	0.48
1:B:289:THR:HG23	1:B:320:LYS:HE2	1.94	0.47
1:A:40:MET:CE	1:A:73:VAL:HG13	2.44	0.47
1:B:182:GLN:OE1	1:B:231:ARG:HD2	2.15	0.47
1:A:172:MET:HG2	1:A:211:ALA:HB2	1.96	0.47
1:B:40:MET:HE1	1:B:73:VAL:HG13	1.96	0.47
1:B:250:ALA:HA	1:B:254:GLN:NE2	2.30	0.46
1:A:282:ASP:O	1:B:52:ARG:NH1	2.46	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:18:ALA:HA	1:B:21:LYS:HD3	1.97	0.46
1:A:146:ASN:O	1:A:147:ALA:HB3	2.16	0.46
1:B:16:PHE:C	1:B:16:PHE:CD1	2.89	0.46
1:B:146:ASN:HD22	1:B:324:LEU:HD22	1.80	0.46
1:B:69:LYS:HE2	1:B:69:LYS:HB2	1.69	0.46
1:A:79:PRO:HB3	1:A:108:HIS:CE1	2.50	0.46
1:A:131:GLY:O	1:A:267:MET:HE1	2.17	0.45
1:B:201:GLN:HG3	2:B:2076:HOH:O	2.16	0.45
1:B:245:ARG:HG2	1:B:304:PHE:CD1	2.51	0.45
1:B:22:ARG:HH12	1:B:25:ILE:CD1	2.29	0.45
1:A:106:ARG:NH2	1:A:143:ILE:HD12	2.32	0.45
1:A:159:LYS:HG3	1:A:218:LEU:HD11	1.97	0.45
1:B:79:PRO:C	1:B:81:ASN:H	2.20	0.45
1:A:221:LEU:HA	1:A:224:LYS:HD2	1.99	0.45
1:A:100:LEU:HB3	1:A:122:SER:HA	1.99	0.45
1:B:11:ILE:HG12	2:B:2090:HOH:O	2.16	0.45
1:A:178:THR:HA	1:A:182:GLN:OE1	2.17	0.45
1:B:320:LYS:HD2	1:B:323:ASP:OD2	2.17	0.45
1:A:60:LYS:NZ	2:A:2102:HOH:O	2.47	0.44
1:A:305:VAL:HG12	1:A:306:LYS:N	2.32	0.44
1:B:0:THR:OG1	1:B:2:LYS:NZ	2.49	0.44
1:B:84:TRP:CE3	1:B:89:VAL:HG21	2.52	0.44
1:A:245:ARG:HA	1:A:303:ASN:O	2.17	0.44
1:B:60:LYS:HB3	1:B:65:ILE:HD11	1.98	0.44
1:A:324:LEU:O	1:A:328:ILE:HG12	2.17	0.44
1:B:60:LYS:O	1:B:61:ASP:HB2	2.18	0.44
1:A:45:LYS:HG3	1:A:57:VAL:HG11	2.00	0.44
1:A:11:ILE:HG21	1:A:119:THR:HB	2.00	0.44
1:A:259:VAL:HG13	1:A:271:LEU:HD21	1.99	0.44
1:B:37:ALA:O	1:B:64:LEU:HD21	2.18	0.43
1:A:218:LEU:HB3	1:A:221:LEU:HD12	2.00	0.43
1:B:268:LYS:HB3	1:B:268:LYS:HE2	1.91	0.43
1:A:133:ASN:OD1	1:A:217:VAL:HG13	2.18	0.43
1:A:126:PRO:HB2	1:A:128:PHE:CE2	2.54	0.43
1:A:22:ARG:HB3	1:A:23:SER:H	1.74	0.43
1:B:13:ARG:HD3	1:B:43:MET:HE3	1.99	0.43
1:B:3:VAL:HG11	1:B:27:ILE:HD13	2.01	0.43
1:A:171:LEU:HD23	1:A:226:THR:CG2	2.47	0.43
1:B:168:ILE:O	1:B:224:LYS:HD3	2.19	0.43
1:A:90:ASP:O	1:A:114:LYS:HB2	2.19	0.42
1:B:262:ALA:O	1:B:267:MET:HB2	2.18	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:176:HIS:HA	1:A:238:SER:HB3	2.01	0.42
1:B:260:LYS:HD3	1:B:273:TYR:CE1	2.54	0.42
1:A:44:LEU:O	1:A:53:PHE:HB2	2.19	0.42
1:A:101:THR:OG1	1:A:104:THR:HG23	2.18	0.42
1:B:251:THR:N	1:B:254:GLN:NE2	2.48	0.42
1:A:41:ALA:HB2	1:A:64:LEU:CD2	2.48	0.42
1:A:115:LYS:HB3	1:A:142:ASP:O	2.19	0.42
1:A:238:SER:HB2	1:A:311:TYR:CZ	2.54	0.42
1:A:292:PHE:CE1	1:A:307:LEU:HD22	2.54	0.42
1:A:131:GLY:O	1:A:267:MET:CE	2.68	0.42
1:B:77:ARG:CG	1:B:77:ARG:NH1	2.71	0.42
1:B:91:VAL:HA	1:B:115:LYS:O	2.19	0.42
1:A:169:GLU:OE2	1:A:245:ARG:NH1	2.42	0.41
1:A:124:ASN:O	1:A:125:THR:C	2.59	0.41
1:A:251:THR:N	1:A:254:GLN:NE2	2.57	0.41
1:A:60:LYS:HB2	1:A:65:ILE:CG1	2.51	0.41
1:A:212:LYS:HG3	2:A:2117:HOH:O	2.21	0.41
1:B:283:PHE:O	1:B:312:ASP:HB2	2.20	0.41
1:A:77:ARG:HG2	1:A:77:ARG:H	1.40	0.41
1:A:320:LYS:HD3	1:A:320:LYS:N	2.36	0.41
1:A:327:HIS:HA	1:A:330:LYS:HZ1	1.86	0.41
1:A:57:VAL:HB	1:A:66:VAL:HG13	2.03	0.41
1:A:41:ALA:HB2	1:A:64:LEU:HD21	2.01	0.41
1:A:182:GLN:HB3	2:A:2003:HOH:O	2.20	0.40
1:B:245:ARG:HG2	1:B:304:PHE:HD1	1.86	0.40
1:A:250:ALA:HA	1:A:254:GLN:NE2	2.36	0.40
1:B:146:ASN:O	1:B:147:ALA:HB3	2.20	0.40
1:B:260:LYS:O	1:B:263:ALA:HB3	2.21	0.40
1:B:157:LEU:O	1:B:161:ILE:HG12	2.20	0.40
1:B:101:THR:HG22	1:B:122(A):LYS:HB2	2.02	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	328/330 (99%)	292 (89%)	31 (10%)	5 (2%)	13 5
1	B	328/330 (99%)	291 (89%)	32 (10%)	5 (2%)	13 5
All	All	656/660 (99%)	583 (89%)	63 (10%)	10 (2%)	13 5

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	210	ALA
1	A	23	SER
1	A	302	ASP
1	B	166	GLY
1	B	237	VAL
1	A	237	VAL
1	A	166	GLY
1	B	80	ALA
1	B	313	ASN
1	B	135	ASP

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	265/265 (100%)	226 (85%)	39 (15%)	4 2
1	B	265/265 (100%)	241 (91%)	24 (9%)	12 6
All	All	530/530 (100%)	467 (88%)	63 (12%)	6 3

All (63) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	LYS
1	A	8	PHE
1	A	21	LYS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	22	ARG
1	A	34	LEU
1	A	45	LYS
1	A	57	VAL
1	A	58	GLU
1	A	60	LYS
1	A	61	ASP
1	A	66	VAL
1	A	69	LYS
1	A	70	LYS
1	A	77	ARG
1	A	83	LYS
1	A	87	VAL
1	A	122	SER
1	A	122(A)	LYS
1	A	124	ASN
1	A	134	PHE
1	A	136	LYS
1	A	145	SER
1	A	161	ILE
1	A	172	MET
1	A	183	LYS
1	A	191	LYS
1	A	207	SER
1	A	208	THR
1	A	218	LEU
1	A	226	THR
1	A	228	MET
1	A	253	GLU
1	A	260	LYS
1	A	266	GLU
1	A	267	MET
1	A	268	LYS
1	A	277	ASP
1	A	300	LEU
1	A	308	VAL
1	B	21	LYS
1	B	22	ARG
1	B	27	ILE
1	B	36	ASP
1	B	52	ARG
1	B	61	ASP

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	69	LYS
1	B	77	ARG
1	B	83	LYS
1	B	87	VAL
1	B	98	LEU
1	B	103	GLU
1	B	122	SER
1	B	134	PHE
1	B	145	SER
1	B	190	HIS
1	B	191	LYS
1	B	228	MET
1	B	245	ARG
1	B	248	LYS
1	B	266	GLU
1	B	268	LYS
1	B	306	LYS
1	B	330	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	63	HIS
1	A	124	ASN
1	A	146	ASN
1	A	202	ASN
1	A	254	GLN
1	A	327	HIS
1	B	146	ASN
1	B	222	ASN
1	B	254	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

There are no carbohydrates in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

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	330/330 (100%)	-0.13	5 (1%)	76	77	16, 28, 41, 54	0
1	B	330/330 (100%)	-0.12	2 (0%)	90	90	17, 28, 43, 51	0
All	All	660/660 (100%)	-0.13	7 (1%)	82	83	16, 28, 43, 54	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	61	ASP	2.9
1	A	330	LYS	2.8
1	A	61	ASP	2.5
1	A	190	HIS	2.3
1	A	116	VAL	2.2
1	A	117	VAL	2.1
1	B	80	ALA	2.1

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

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.