Full wwPDB X-ray Structure Validation Report (i)

Jun 15, 2024 – 08:13 PM EDT

PDB ID	:	4PGM
Title	:	SACCHAROMYCES CEREVISIAE PHOSPHOGLYCERATE MUTASE
Authors	:	Rigden, D.J.; Alexeev, D.; Phillips, S.E.V.; Fothergill-Gilmore, L.A.
Deposited on	:	1997-04-25
Resolution	:	2.30 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:

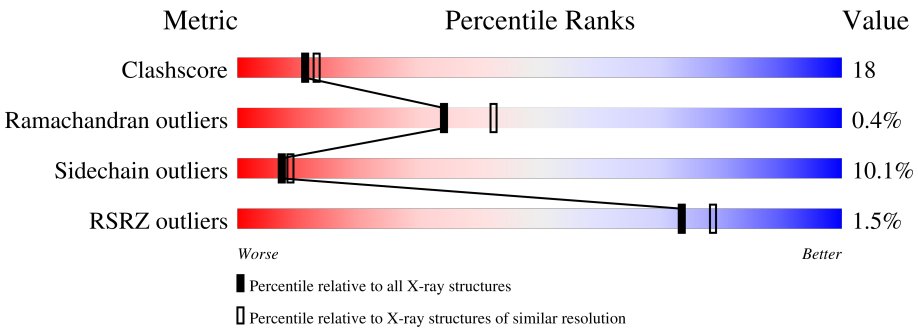
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

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 2.30 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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	246	2%, 59%, 33%, • •
1	B	246	2%, 59%, 31%, 6%, •
1	C	246	55%, 33%, 7%, 5%
1	D	246	2%, 59%, 34%, • •

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 8322 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 PHOSPHOGLYCERATE MUTASE 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	235	1873	1199	323	350	1	0	0	0
1	B	236	1877	1201	324	351	1	0	0	0
1	C	234	1868	1196	322	349	1	0	0	0
1	D	236	1877	1201	324	351	1	0	0	0

- Molecule 2 is water.

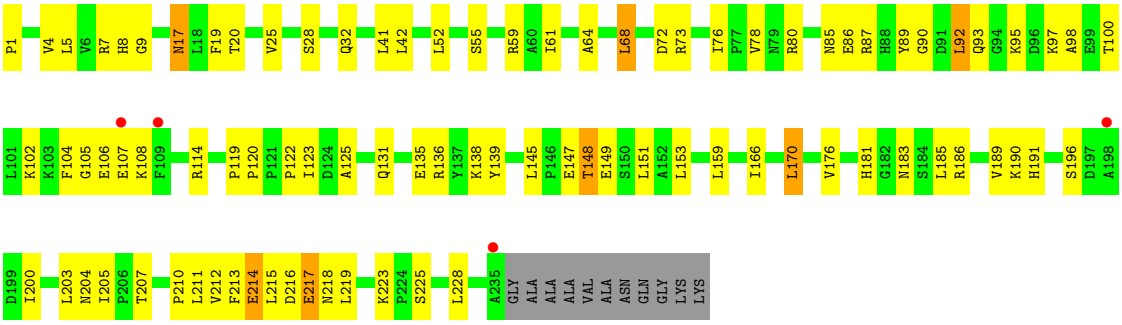
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	201	201	201	0	0
2	B	236	236	236	0	0
2	C	223	223	223	0	0
2	D	167	167	167	0	0

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.

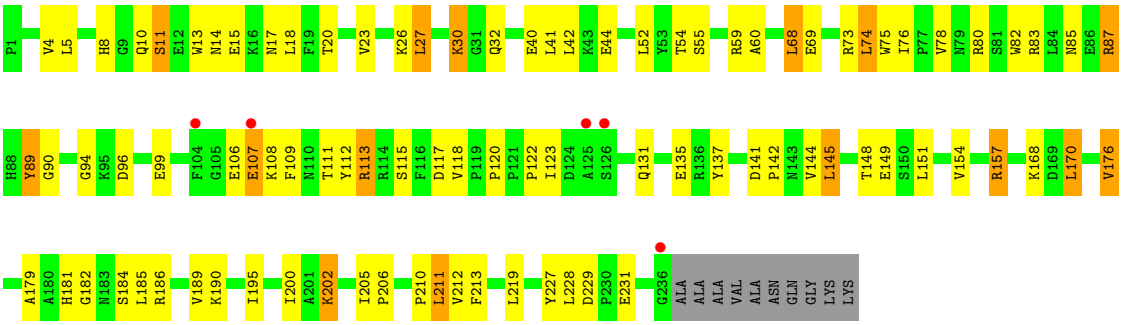
• Molecule 1: PHOSPHOGLYCERATE MUTASE 1

Chain A:



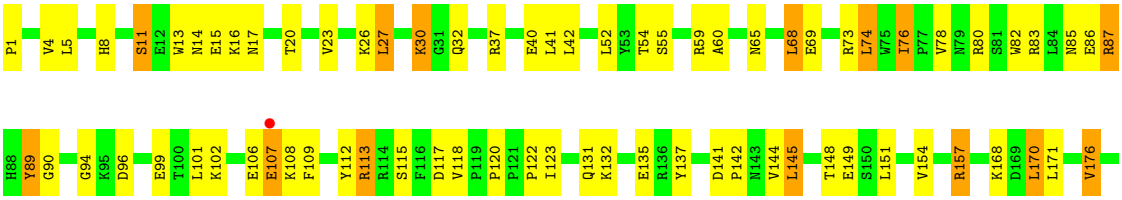
• Molecule 1: PHOSPHOGLYCERATE MUTASE 1

Chain B:



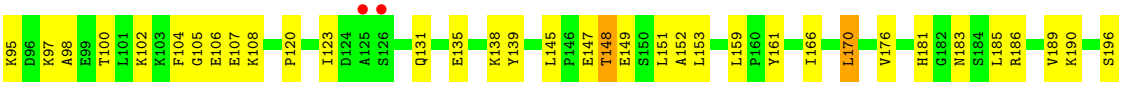
• Molecule 1: PHOSPHOGLYCERATE MUTASE 1

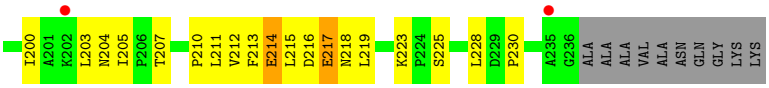
Chain C:



• Molecule 1: PHOSPHOGLYCERATE MUTASE 1

Chain D:





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	81.47Å 84.56Å 88.88Å 90.00° 111.72° 90.00°	Depositor
Resolution (Å)	25.00 – 2.30 24.68 – 2.30	Depositor EDS
% Data completeness (in resolution range)	88.7 (25.00-2.30) 73.4 (24.68-2.30)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.24 (at 2.31Å)	Xtriage
Refinement program	X-PLOR 3.843	Depositor
R, R_{free}	0.192 , 0.286 0.207 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å^2)	16.9	Xtriage
Anisotropy	0.327	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.30 , 49.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.44$, $< L^2 > = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	8322	wwPDB-VP
Average B, all atoms (Å^2)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 93.95 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.9158e-09. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[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	0.50	0/1917	0.76	0/2599
1	B	0.50	0/1921	0.78	1/2604 (0.0%)
1	C	0.50	0/1912	0.79	1/2592 (0.0%)
1	D	0.50	0/1921	0.76	0/2604
All	All	0.50	0/7671	0.77	2/10399 (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	C	27	LEU	CA-CB-CG	5.34	127.58	115.30
1	B	27	LEU	CA-CB-CG	5.33	127.56	115.30

There are no chirality outliers.

There are no planarity outliers.

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	1873	0	1894	65	0
1	B	1877	0	1897	73	0
1	C	1868	0	1889	80	0
1	D	1877	0	1897	66	0
2	A	201	0	0	3	0
2	B	236	0	0	5	0
2	C	223	0	0	13	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	167	0	0	6	0
All	All	8322	0	7577	270	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 18.

All (270) 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:D:170:LEU:HD13	1:D:176:VAL:HG13	1.54	0.90
1:A:170:LEU:HD13	1:A:176:VAL:HG13	1.61	0.81
1:B:120:PRO:O	1:B:148:THR:HG21	1.82	0.80
1:C:120:PRO:O	1:C:148:THR:HG21	1.82	0.78
1:B:189:VAL:HG21	1:B:205:ILE:HD11	1.65	0.77
1:D:61:ILE:HG23	1:D:80:ARG:NH2	2.00	0.77
1:A:120:PRO:O	1:A:148:THR:HG21	1.85	0.76
1:D:120:PRO:O	1:D:148:THR:HG21	1.85	0.76
1:A:61:ILE:HG23	1:A:80:ARG:NH2	2.02	0.75
1:C:189:VAL:HG21	1:C:205:ILE:HD11	1.68	0.74
1:B:210:PRO:HB2	1:B:228:LEU:HD12	1.70	0.74
1:C:122:PRO:HA	1:C:148:THR:OG1	1.88	0.74
1:B:122:PRO:HA	1:B:148:THR:OG1	1.88	0.73
1:B:69:GLU:HA	1:B:74:LEU:HD22	1.70	0.72
1:C:210:PRO:HB2	1:C:228:LEU:HD12	1.71	0.71
1:D:55:SER:HB2	1:D:85:ASN:OD1	1.89	0.71
1:A:55:SER:HB2	1:A:85:ASN:OD1	1.91	0.69
1:C:1:PRO:HA	2:C:253:HOH:O	1.95	0.66
1:A:213:PHE:CE1	1:A:225:SER:HB3	2.31	0.65
1:C:69:GLU:HA	1:C:74:LEU:HD22	1.77	0.65
1:A:104:PHE:HB3	1:A:108:LYS:HB2	1.79	0.64
1:D:213:PHE:CE1	1:D:225:SER:HB3	2.31	0.64
1:A:125:ALA:HB2	2:A:422:HOH:O	1.97	0.64
1:D:104:PHE:HB3	1:D:108:LYS:HB2	1.80	0.64
1:C:220:LYS:HG2	2:C:337:HOH:O	1.99	0.62
1:D:40:GLU:HG2	2:D:333:HOH:O	1.99	0.62
1:C:96:ASP:HB3	1:C:99:GLU:HB2	1.83	0.61
1:D:190:LYS:HB2	1:D:200:ILE:HD13	1.83	0.61
1:C:52:LEU:HD23	1:C:78:VAL:HG22	1.82	0.61
1:B:13:TRP:CZ3	1:B:94:GLY:HA2	2.35	0.61
1:D:90:GLY:HA3	1:D:148:THR:HG23	1.82	0.60
1:B:96:ASP:HB3	1:B:99:GLU:HB2	1.82	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1:PRO:HB2	1:A:215:LEU:HB2	1.82	0.60
1:A:73:ARG:HH11	1:A:76:ILE:HD11	1.66	0.60
1:A:90:GLY:HA3	1:A:148:THR:HG23	1.83	0.60
1:C:211:LEU:HD13	1:C:213:PHE:HE1	1.66	0.60
1:A:189:VAL:HG21	1:A:205:ILE:HD11	1.84	0.60
1:B:185:LEU:O	1:B:189:VAL:HG23	2.02	0.60
1:D:189:VAL:HG21	1:D:205:ILE:HD11	1.84	0.60
1:B:211:LEU:HD13	1:B:213:PHE:HE1	1.67	0.59
1:C:5:LEU:HG	1:C:185:LEU:HD22	1.84	0.59
1:A:190:LYS:HB2	1:A:200:ILE:HD13	1.84	0.58
1:B:5:LEU:HG	1:B:185:LEU:HD22	1.85	0.58
1:B:69:GLU:HA	1:B:74:LEU:CD2	2.33	0.58
1:B:211:LEU:HD13	1:B:213:PHE:CE1	2.39	0.58
1:C:211:LEU:HD13	1:C:213:PHE:CE1	2.38	0.58
1:C:80:ARG:NH1	1:D:76:ILE:O	2.37	0.57
1:D:152:ALA:HB1	2:D:307:HOH:O	2.04	0.57
1:C:13:TRP:CZ3	1:C:94:GLY:HA2	2.39	0.57
1:D:52:LEU:HD23	1:D:78:VAL:HG13	1.87	0.57
1:D:73:ARG:HH11	1:D:76:ILE:HD11	1.69	0.57
1:A:7:ARG:HD3	1:A:207:THR:HG22	1.87	0.57
1:C:185:LEU:O	1:C:189:VAL:HG23	2.04	0.56
1:A:135:GLU:O	1:A:138:LYS:HG2	2.05	0.56
1:D:1:PRO:HB2	1:D:215:LEU:HB2	1.86	0.56
1:B:23:VAL:O	1:B:59:ARG:NH2	2.38	0.56
1:C:23:VAL:O	1:C:59:ARG:NH2	2.38	0.56
1:A:210:PRO:HB2	1:A:228:LEU:HD12	1.86	0.56
1:C:135:GLU:OE1	1:D:73:ARG:NH2	2.38	0.55
1:A:139:TYR:HB3	1:C:168:LYS:HG3	1.88	0.55
1:B:190:LYS:HB2	1:B:200:ILE:CD1	2.36	0.55
1:C:55:SER:HB2	1:C:85:ASN:OD1	2.07	0.55
1:C:82:TRP:CZ3	1:C:157:ARG:HD2	2.41	0.55
1:B:15:GLU:OE2	1:B:30:LYS:HD3	2.07	0.55
1:C:85:ASN:O	1:C:184:SER:HB3	2.07	0.55
1:C:15:GLU:OE2	1:C:30:LYS:HD3	2.07	0.55
1:C:69:GLU:HA	1:C:74:LEU:CD2	2.37	0.55
1:C:190:LYS:HB2	1:C:200:ILE:CD1	2.37	0.55
1:A:217:GLU:CD	1:A:217:GLU:H	2.10	0.54
1:B:108:LYS:HE2	1:B:112:TYR:CZ	2.42	0.54
1:B:170:LEU:CD1	1:B:176:VAL:HG13	2.37	0.54
1:D:186:ARG:HD3	1:D:204:ASN:HA	1.88	0.54
1:A:76:ILE:O	1:B:80:ARG:NH1	2.40	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:210:PRO:HG3	2:C:381:HOH:O	2.08	0.54
1:D:217:GLU:CD	1:D:217:GLU:H	2.09	0.54
1:D:5:LEU:HB3	1:D:185:LEU:HD22	1.90	0.54
1:D:210:PRO:HB2	1:D:228:LEU:HD12	1.89	0.54
1:A:5:LEU:HB3	1:A:185:LEU:HD22	1.90	0.54
1:C:204:ASN:HB2	2:C:397:HOH:O	2.07	0.54
1:D:61:ILE:HG23	1:D:80:ARG:HH22	1.72	0.53
1:C:108:LYS:HE3	2:C:306:HOH:O	2.07	0.53
1:C:108:LYS:HE2	1:C:112:TYR:CZ	2.44	0.53
1:D:135:GLU:O	1:D:138:LYS:HG2	2.07	0.53
1:B:52:LEU:HD23	1:B:78:VAL:HG22	1.89	0.53
1:A:186:ARG:HD3	1:A:204:ASN:HA	1.89	0.53
1:A:52:LEU:HD23	1:A:78:VAL:HG13	1.91	0.52
1:B:55:SER:HB2	1:B:85:ASN:OD1	2.09	0.52
1:C:90:GLY:HA3	1:C:148:THR:HG23	1.91	0.52
1:C:149:GLU:OE2	1:C:157:ARG:HD3	2.10	0.52
1:B:122:PRO:HA	1:B:148:THR:HG1	1.75	0.52
1:B:149:GLU:OE2	1:B:157:ARG:HD3	2.10	0.52
1:C:182:GLY:O	1:C:186:ARG:HG3	2.10	0.52
1:C:83:ARG:HD3	1:C:157:ARG:O	2.10	0.52
1:D:7:ARG:HD3	1:D:207:THR:HG22	1.91	0.52
1:C:122:PRO:HA	1:C:148:THR:HG1	1.75	0.52
1:B:182:GLY:O	1:B:186:ARG:HG3	2.10	0.52
1:D:98:ALA:O	1:D:102:LYS:HG3	2.10	0.52
1:B:210:PRO:HB2	1:B:228:LEU:CD1	2.39	0.52
1:A:98:ALA:O	1:A:102:LYS:HG3	2.10	0.51
1:B:85:ASN:O	1:B:184:SER:HB3	2.10	0.51
1:C:37:ARG:NH1	2:C:392:HOH:O	2.43	0.51
1:B:13:TRP:HB2	1:B:20:THR:HG22	1.91	0.51
1:D:20:THR:HG21	1:D:25:VAL:HG12	1.92	0.51
1:A:61:ILE:HG23	1:A:80:ARG:HH22	1.74	0.51
1:A:8:HIS:CE1	1:A:59:ARG:HH11	2.29	0.51
1:A:210:PRO:HB2	1:A:228:LEU:CD1	2.40	0.51
1:B:82:TRP:CZ3	1:B:157:ARG:HD2	2.46	0.50
1:C:170:LEU:HD13	1:C:176:VAL:HG13	1.93	0.50
1:C:170:LEU:CD1	1:C:176:VAL:HG13	2.41	0.50
1:C:210:PRO:CB	1:C:228:LEU:HD12	2.39	0.50
1:A:105:GLY:O	1:A:107:GLU:N	2.44	0.50
1:C:108:LYS:HE2	1:C:112:TYR:OH	2.11	0.50
1:A:20:THR:HG21	1:A:25:VAL:HG12	1.92	0.50
1:B:170:LEU:HD13	1:B:176:VAL:HG13	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:206:PRO:HD2	1:C:227:TYR:OH	2.12	0.50
1:C:210:PRO:HB2	1:C:228:LEU:CD1	2.39	0.50
1:D:105:GLY:O	1:D:107:GLU:N	2.45	0.50
1:C:99:GLU:HG2	2:C:451:HOH:O	2.11	0.50
1:A:131:GLN:HA	2:A:281:HOH:O	2.11	0.50
1:B:90:GLY:HA3	1:B:148:THR:HG23	1.93	0.50
1:D:92:LEU:HA	1:D:95:LYS:HD2	1.93	0.50
1:B:108:LYS:HE2	1:B:112:TYR:OH	2.12	0.49
1:D:120:PRO:C	1:D:148:THR:HG21	2.32	0.49
1:B:83:ARG:HD3	1:B:157:ARG:O	2.12	0.49
1:D:131:GLN:HA	2:D:274:HOH:O	2.11	0.49
1:D:210:PRO:HB2	1:D:228:LEU:CD1	2.43	0.49
1:C:13:TRP:HB2	1:C:20:THR:HG22	1.94	0.49
1:C:41:LEU:HD21	1:C:229:ASP:HB2	1.95	0.49
1:A:5:LEU:N	1:A:5:LEU:HD22	2.28	0.49
1:A:120:PRO:C	1:A:148:THR:HG21	2.32	0.49
1:B:41:LEU:HD21	1:B:229:ASP:HB2	1.95	0.49
1:D:5:LEU:N	1:D:5:LEU:HD22	2.28	0.49
1:C:107:GLU:HG3	1:C:108:LYS:H	1.77	0.48
1:D:8:HIS:CE1	1:D:59:ARG:HH11	2.31	0.48
1:D:28:SER:O	1:D:32:GLN:HG3	2.13	0.48
1:D:17:ASN:ND2	1:D:97:LYS:HB2	2.28	0.48
1:B:168:LYS:HG3	1:D:139:TYR:HB3	1.94	0.48
1:A:135:GLU:OE1	1:B:73:ARG:NH2	2.46	0.48
1:A:28:SER:O	1:A:32:GLN:HG3	2.13	0.48
1:D:59:ARG:HG2	2:D:275:HOH:O	2.13	0.48
1:A:87:ARG:HB3	1:A:183:ASN:HD22	1.79	0.48
1:C:120:PRO:HD2	1:C:148:THR:HG22	1.94	0.48
1:D:87:ARG:HB3	1:D:183:ASN:HD22	1.79	0.48
1:B:107:GLU:HG3	1:B:108:LYS:H	1.78	0.47
1:B:210:PRO:CB	1:B:228:LEU:HD12	2.40	0.47
1:A:17:ASN:ND2	1:A:97:LYS:HB2	2.29	0.47
1:B:54:THR:O	1:B:80:ARG:HA	2.15	0.47
1:C:65:ASN:ND2	2:C:307:HOH:O	2.46	0.47
1:A:92:LEU:HA	1:A:95:LYS:HD2	1.96	0.47
1:B:142:PRO:HG2	1:D:218:ASN:HA	1.97	0.47
1:A:87:ARG:HG2	1:A:149:GLU:HB2	1.97	0.47
1:A:122:PRO:HD3	2:A:301:HOH:O	2.15	0.47
1:C:54:THR:O	1:C:80:ARG:HA	2.15	0.47
1:B:154:VAL:O	1:B:157:ARG:HG3	2.15	0.46
1:D:166:ILE:HG23	1:D:176:VAL:HG21	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:8:HIS:CE1	1:B:59:ARG:HD2	2.49	0.46
1:A:214:GLU:OE2	1:A:223:LYS:HD3	2.15	0.46
1:C:68:LEU:HB3	1:C:74:LEU:HD13	1.98	0.46
1:D:90:GLY:HA2	1:D:147:GLU:O	2.16	0.46
1:C:141:ASP:O	1:C:144:VAL:HG12	2.16	0.46
1:B:30:LYS:HD2	2:B:333:HOH:O	2.16	0.45
1:D:216:ASP:HB2	1:D:217:GLU:OE2	2.15	0.45
1:B:44:GLU:HB2	2:B:379:HOH:O	2.17	0.45
1:B:202:LYS:HD2	1:B:202:LYS:N	2.31	0.45
1:D:17:ASN:HD21	1:D:97:LYS:HB2	1.82	0.45
1:D:190:LYS:HB2	1:D:200:ILE:CD1	2.46	0.45
1:D:214:GLU:OE2	1:D:223:LYS:HD3	2.16	0.45
1:C:200:ILE:HD12	1:C:200:ILE:HA	1.80	0.45
1:A:73:ARG:NH2	1:B:135:GLU:OE1	2.50	0.45
1:A:86:GLU:OE1	1:A:181:HIS:HB3	2.17	0.45
1:D:87:ARG:HG2	1:D:149:GLU:HB2	1.98	0.45
1:A:7:ARG:CD	1:A:207:THR:HG22	2.46	0.45
1:C:76:ILE:HG22	2:C:270:HOH:O	2.17	0.45
1:A:218:ASN:HA	1:C:142:PRO:HG2	1.99	0.44
1:B:137:TYR:HB3	1:B:145:LEU:HD21	1.99	0.44
1:C:60:ALA:HB2	1:C:181:HIS:CE1	2.52	0.44
1:C:73:ARG:NH2	1:D:135:GLU:OE1	2.51	0.44
1:B:141:ASP:O	1:B:144:VAL:HG12	2.17	0.44
1:C:102:LYS:HD2	2:C:451:HOH:O	2.16	0.44
1:A:135:GLU:CD	1:B:73:ARG:HH22	2.20	0.44
1:C:4:VAL:HG13	1:C:212:VAL:HG22	2.00	0.44
1:C:202:LYS:HD2	1:C:202:LYS:N	2.31	0.44
1:B:120:PRO:HD2	1:B:148:THR:HG22	2.00	0.44
1:B:206:PRO:HD2	1:B:227:TYR:OH	2.18	0.44
1:C:16:LYS:HB2	2:C:347:HOH:O	2.16	0.44
1:B:4:VAL:HG13	1:B:212:VAL:HG22	2.00	0.44
1:D:131:GLN:HG2	2:D:268:HOH:O	2.16	0.44
1:A:159:LEU:HD23	1:A:159:LEU:HA	1.85	0.44
1:C:154:VAL:O	1:C:157:ARG:HG3	2.17	0.44
1:D:159:LEU:HD23	1:D:159:LEU:HA	1.85	0.44
1:D:203:LEU:HG	1:D:204:ASN:N	2.33	0.44
1:A:119:PRO:HA	1:A:120:PRO:HD3	1.86	0.43
1:B:109:PHE:O	1:B:113:ARG:HB2	2.17	0.43
1:A:90:GLY:C	1:A:123:ILE:HB	2.39	0.43
1:C:132:LYS:HE3	2:C:351:HOH:O	2.17	0.43
1:A:166:ILE:HG23	1:A:176:VAL:HG21	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:8:HIS:CE1	1:B:59:ARG:CD	3.01	0.43
1:B:115:SER:OG	1:B:118:VAL:HG23	2.18	0.43
1:C:90:GLY:C	1:C:123:ILE:HB	2.39	0.43
1:C:190:LYS:HD2	1:C:195:ILE:O	2.19	0.43
1:A:20:THR:O	1:A:93:GLN:HG2	2.17	0.43
1:B:108:LYS:HD2	2:B:393:HOH:O	2.18	0.43
1:C:73:ARG:HH22	1:D:135:GLU:CD	2.22	0.43
1:B:190:LYS:HD2	1:B:195:ILE:O	2.19	0.43
1:C:30:LYS:HE3	1:C:30:LYS:HB3	1.70	0.43
1:D:19:PHE:HB2	1:D:92:LEU:O	2.18	0.43
1:A:17:ASN:HD21	1:A:97:LYS:HB2	1.82	0.43
1:B:170:LEU:HD13	1:B:176:VAL:CG1	2.49	0.43
1:C:115:SER:OG	1:C:118:VAL:HG23	2.19	0.43
1:D:86:GLU:OE1	1:D:181:HIS:HB3	2.18	0.43
1:A:19:PHE:HB2	1:A:92:LEU:O	2.18	0.43
1:C:113:ARG:NH1	2:C:464:HOH:O	2.51	0.43
1:B:90:GLY:C	1:B:123:ILE:HB	2.40	0.43
1:B:111:THR:HA	2:B:473:HOH:O	2.18	0.43
1:B:190:LYS:HB2	1:B:200:ILE:HD13	2.00	0.43
1:C:13:TRP:CB	1:C:20:THR:HG22	2.49	0.43
1:B:13:TRP:CB	1:B:20:THR:HG22	2.49	0.43
1:B:82:TRP:O	1:B:85:ASN:HB2	2.19	0.43
1:D:90:GLY:C	1:D:123:ILE:HB	2.39	0.43
1:B:68:LEU:HB3	1:B:74:LEU:HD13	2.01	0.42
1:C:11:SER:OG	1:C:14:ASN:HB2	2.19	0.42
1:C:137:TYR:HB3	1:C:145:LEU:HD21	2.01	0.42
1:A:4:VAL:HG13	1:A:212:VAL:HG22	2.01	0.42
1:A:216:ASP:HB2	1:A:217:GLU:OE2	2.19	0.42
1:C:109:PHE:O	1:C:113:ARG:HB2	2.19	0.42
1:A:90:GLY:HA2	1:A:147:GLU:O	2.19	0.42
1:A:114:ARG:HA	1:A:114:ARG:HD2	1.92	0.42
1:D:90:GLY:HA3	1:D:148:THR:CG2	2.47	0.42
1:D:170:LEU:HD13	1:D:176:VAL:CG1	2.38	0.42
1:A:90:GLY:HA3	1:A:148:THR:CG2	2.50	0.42
1:B:27:LEU:HD23	1:B:32:GLN:HG2	2.01	0.42
1:A:190:LYS:HB2	1:A:200:ILE:CD1	2.48	0.42
1:C:54:THR:HG22	1:C:179:ALA:HB3	2.01	0.42
1:A:64:ALA:O	1:A:68:LEU:HB2	2.20	0.42
1:A:19:PHE:O	1:A:93:GLN:O	2.37	0.42
1:C:8:HIS:CE1	1:C:59:ARG:HD2	2.54	0.42
1:D:7:ARG:CD	1:D:207:THR:HG22	2.48	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:8:HIS:CE1	1:A:59:ARG:HD2	2.55	0.42
1:B:11:SER:OG	1:B:14:ASN:HB2	2.19	0.42
1:D:50:ASP:O	1:D:76:ILE:HG13	2.20	0.42
1:A:9:GLY:HA2	1:A:207:THR:HB	2.02	0.41
1:D:64:ALA:O	1:D:68:LEU:HB2	2.20	0.41
1:D:213:PHE:CD1	1:D:225:SER:HB3	2.54	0.41
1:D:95:LYS:HE2	2:D:352:HOH:O	2.19	0.41
1:A:136:ARG:HB3	1:B:75:TRP:CD2	2.56	0.41
1:C:76:ILE:HA	1:C:76:ILE:HD12	1.85	0.41
1:D:9:GLY:HA2	1:D:207:THR:HB	2.03	0.41
1:D:19:PHE:O	1:D:93:GLN:O	2.38	0.41
1:A:203:LEU:HG	1:A:204:ASN:N	2.36	0.41
1:B:87:ARG:NH2	1:B:89:TYR:OH	2.53	0.41
1:C:87:ARG:NH2	1:C:89:TYR:OH	2.53	0.41
1:B:60:ALA:HB2	1:B:181:HIS:CE1	2.55	0.41
1:B:231:GLU:CD	1:B:231:GLU:H	2.24	0.41
1:C:27:LEU:HD23	1:C:32:GLN:HG2	2.01	0.41
1:A:138:LYS:O	1:C:171:LEU:HD13	2.21	0.41
1:B:30:LYS:HE3	1:B:30:LYS:HB3	1.68	0.41
1:C:109:PHE:CE2	1:C:113:ARG:HD3	2.55	0.41
1:B:18:LEU:HD22	1:B:94:GLY:O	2.21	0.41
1:B:54:THR:HG22	1:B:179:ALA:HB3	2.03	0.41
1:D:4:VAL:HG13	1:D:212:VAL:HG22	2.03	0.41
1:D:83:ARG:HD2	1:D:161:TYR:HB2	2.03	0.41
1:A:159:LEU:HD21	1:A:191:HIS:CD2	2.55	0.41
1:B:109:PHE:CE2	1:B:113:ARG:HD3	2.56	0.41
1:D:228:LEU:C	1:D:230:PRO:HD3	2.41	0.40
1:A:213:PHE:CD1	1:A:225:SER:HB3	2.56	0.40
1:B:10:GLN:HG2	2:B:469:HOH:O	2.21	0.40
1:C:86:GLU:HG3	1:C:87:ARG:N	2.36	0.40
1:C:101:LEU:HD12	1:C:101:LEU:O	2.21	0.40
1:D:53:TYR:CD1	1:D:79:ASN:HB2	2.56	0.40
1:C:231:GLU:H	1:C:231:GLU:CD	2.25	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	233/246 (95%)	222 (95%)	10 (4%)	1 (0%)	34	42
1	B	234/246 (95%)	222 (95%)	11 (5%)	1 (0%)	34	42
1	C	232/246 (94%)	221 (95%)	10 (4%)	1 (0%)	34	42
1	D	234/246 (95%)	222 (95%)	11 (5%)	1 (0%)	34	42
All	All	933/984 (95%)	887 (95%)	42 (4%)	4 (0%)	34	42

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	106	GLU
1	D	106	GLU
1	B	106	GLU
1	C	106	GLU

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	202/207 (98%)	184 (91%)	18 (9%)	9	11
1	B	202/207 (98%)	179 (89%)	23 (11%)	5	6
1	C	202/207 (98%)	179 (89%)	23 (11%)	5	6
1	D	202/207 (98%)	184 (91%)	18 (9%)	9	11
All	All	808/828 (98%)	726 (90%)	82 (10%)	7	9

All (82) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	ASN
1	A	41	LEU
1	A	42	LEU
1	A	68	LEU
1	A	72	ASP
1	A	89	TYR
1	A	92	LEU
1	A	100	THR
1	A	145	LEU
1	A	148	THR
1	A	151	LEU
1	A	153	LEU
1	A	170	LEU
1	A	196	SER
1	A	211	LEU
1	A	214	GLU
1	A	217	GLU
1	A	219	LEU
1	B	11	SER
1	B	17	ASN
1	B	26	LYS
1	B	30	LYS
1	B	40	GLU
1	B	42	LEU
1	B	68	LEU
1	B	74	LEU
1	B	76	ILE
1	B	87	ARG
1	B	89	TYR
1	B	107	GLU
1	B	113	ARG
1	B	117	ASP
1	B	131	GLN
1	B	145	LEU
1	B	151	LEU
1	B	157	ARG
1	B	170	LEU
1	B	176	VAL
1	B	202	LYS
1	B	211	LEU
1	B	219	LEU
1	C	11	SER

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Mol	Chain	Res	Type
1	C	17	ASN
1	C	26	LYS
1	C	30	LYS
1	C	40	GLU
1	C	42	LEU
1	C	68	LEU
1	C	74	LEU
1	C	76	ILE
1	C	87	ARG
1	C	89	TYR
1	C	107	GLU
1	C	113	ARG
1	C	117	ASP
1	C	131	GLN
1	C	145	LEU
1	C	151	LEU
1	C	157	ARG
1	C	170	LEU
1	C	176	VAL
1	C	202	LYS
1	C	211	LEU
1	C	219	LEU
1	D	17	ASN
1	D	41	LEU
1	D	42	LEU
1	D	68	LEU
1	D	72	ASP
1	D	89	TYR
1	D	92	LEU
1	D	100	THR
1	D	145	LEU
1	D	148	THR
1	D	151	LEU
1	D	153	LEU
1	D	170	LEU
1	D	196	SER
1	D	211	LEU
1	D	214	GLU
1	D	217	GLU
1	D	219	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	ASN
1	A	65	ASN
1	B	14	ASN
1	B	17	ASN
1	B	32	GLN
1	B	65	ASN
1	B	191	HIS
1	B	218	ASN
1	C	14	ASN
1	C	17	ASN
1	C	65	ASN
1	C	218	ASN
1	D	10	GLN
1	D	17	ASN
1	D	62	GLN
1	D	65	ASN

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	235/246 (95%)	-0.24	4 (1%)	70	76	3, 19, 43, 84	0
1	B	236/246 (95%)	-0.29	5 (2%)	63	70	3, 17, 50, 75	0
1	C	234/246 (95%)	-0.29	1 (0%)	92	95	3, 18, 43, 82	0
1	D	236/246 (95%)	-0.20	4 (1%)	70	76	4, 19, 45, 75	0
All	All	941/984 (95%)	-0.26	14 (1%)	73	79	3, 19, 46, 84	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	125	ALA	5.5
1	B	236	GLY	3.9
1	A	235	ALA	3.2
1	D	235	ALA	3.1
1	C	107	GLU	3.0
1	B	125	ALA	2.7
1	D	202	LYS	2.6
1	B	104	PHE	2.4
1	A	107	GLU	2.4
1	B	107	GLU	2.2
1	B	126	SER	2.1
1	D	126	SER	2.1
1	A	109	PHE	2.1
1	A	198	ALA	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 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.