Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 08:16 AM GMT

PDB ID	:	3E0Y
Title	:	The crystal structure of a conserved domain from a protein of Geobacter sulfurreducens PCA
Authors	:	Tan, K.; Bigelow, L.; Clancy, S.; Joachimiak, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on	:	2008-08-01
Resolution	:	3.10 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:

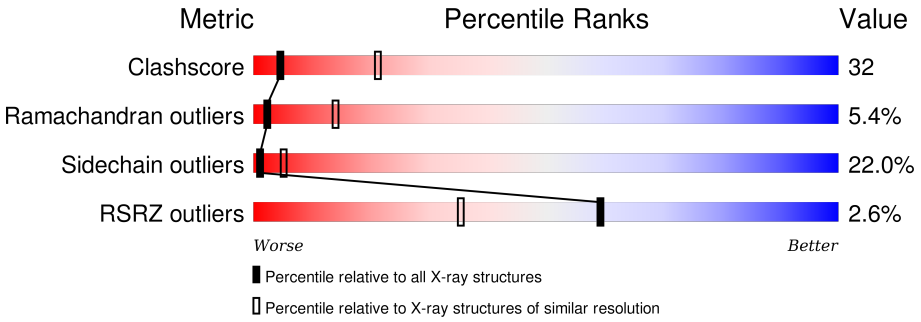
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	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 3.10 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	102246	1222 (3.14-3.06)
Ramachandran outliers	100387	1174 (3.14-3.06)
Sidechain outliers	100360	1174 (3.14-3.06)
RSRZ outliers	91569	1119 (3.14-3.06)

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	181	3% 41% 31% 13% • 14%
1	B	181	% 33% 39% 10% • 16%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 2350 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 Conserved domain protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	155	Total	C	N	O	S	Se	0	0	0
			1177	747	200	227	1	2			
1	B	152	Total	C	N	O	S	Se	0	0	0
			1171	746	197	225	1	2			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	65	SER	-	expression tag	UNP Q74E48
A	66	ASN	-	expression tag	UNP Q74E48
A	67	ALA	-	expression tag	UNP Q74E48
B	65	SER	-	expression tag	UNP Q74E48
B	66	ASN	-	expression tag	UNP Q74E48
B	67	ALA	-	expression tag	UNP Q74E48

- Molecule 2 is SODIUM ION (three-letter code: NA) (formula: Na).

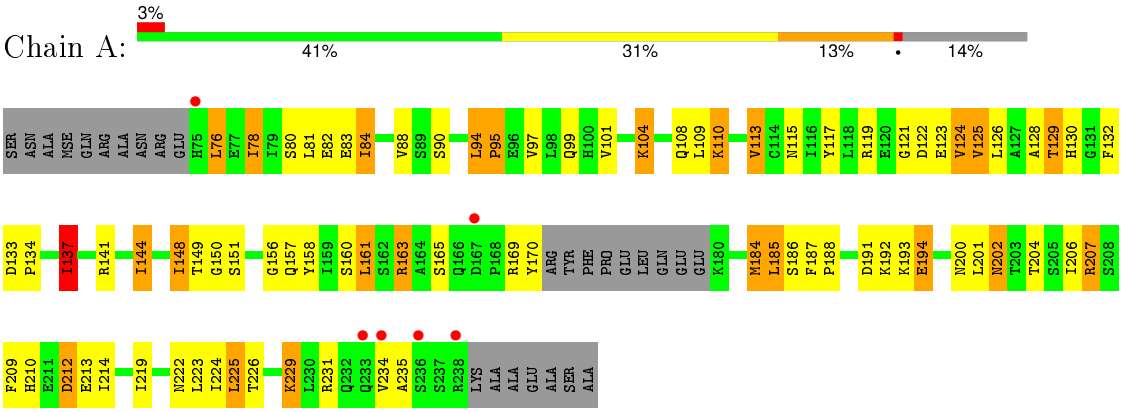
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Na	0	0
			1	1		
2	A	1	Total	Na	0	0
			1	1		

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.

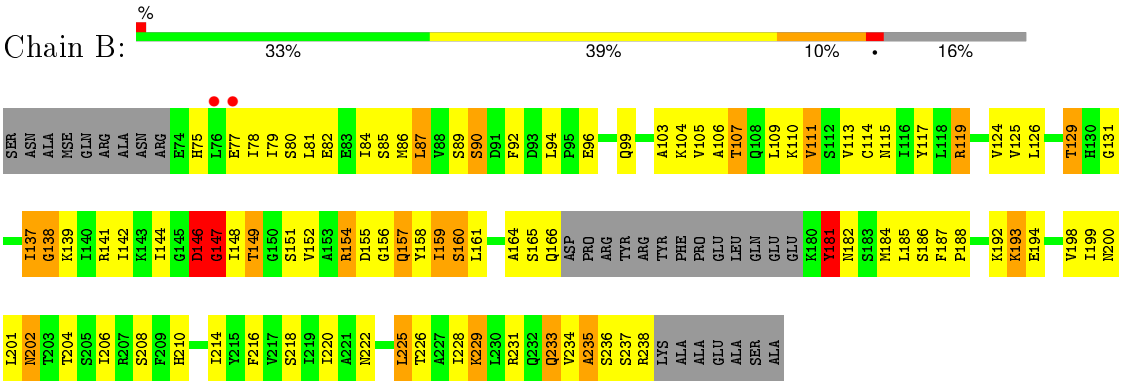
- Molecule 1: Conserved domain protein

Chain A:



- Molecule 1: Conserved domain protein

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α, β, γ	59.42Å 59.42Å 223.18Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	42.30 – 3.10 42.32 – 3.10	Depositor EDS
% Data completeness (in resolution range)	99.5 (42.30-3.10) 99.5 (42.32-3.10)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.95 (at 3.12Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.216 , 0.291 0.217 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å^2)	90.3	Xtriage
Anisotropy	0.003	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 89.2	EDS
Estimated twinning fraction	0.054 for -h,-k,l	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Outliers	2 of 8837 reflections (0.023%)	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	2350	wwPDB-VP
Average B, all atoms (Å^2)	69.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.76% 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: NA

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	0.81	1/1190 (0.1%)	0.88	1/1607 (0.1%)
1	B	0.78	0/1184	0.92	0/1597
All	All	0.80	1/2374 (0.0%)	0.90	1/3204 (0.0%)

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	B	0	3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	170	TYR	C-O	9.37	1.41	1.23

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	137	ILE	CG1-CB-CG2	-5.34	99.65	111.40

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	129	THR	Peptide
1	B	138	GLY	Peptide
1	B	147	GLY	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	1177	0	1182	81	0
1	B	1171	0	1188	76	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
All	All	2350	0	2370	152	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 32.

All (152) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:115:ASN:ND2	1:A:129:THR:HG22	1.40	1.31
1:A:80:SER:HB3	1:A:83:GLU:HG3	1.32	1.11
1:A:115:ASN:HD22	1:A:129:THR:HG22	1.32	0.94
1:B:154:ARG:HG3	1:B:155:ASP:N	1.84	0.92
1:A:148:ILE:C	1:A:148:ILE:HD13	1.91	0.90
1:B:138:GLY:HA3	1:B:139:LYS:HE3	1.54	0.87
1:A:222:ASN:O	1:A:226:THR:HG22	1.74	0.86
1:A:117:TYR:HB3	1:A:124:VAL:HG13	1.55	0.85
1:B:158:TYR:HB3	1:B:188:PRO:HD3	1.58	0.85
1:A:115:ASN:HD21	1:A:129:THR:HG22	1.45	0.80
1:A:115:ASN:ND2	1:A:129:THR:CG2	2.36	0.77
1:A:187:PHE:CD2	1:A:188:PRO:HD2	2.26	0.71
1:A:109:LEU:O	1:A:110:LYS:HG2	1.90	0.71
1:B:158:TYR:HB3	1:B:188:PRO:CD	2.21	0.70
1:A:231:ARG:O	1:A:235:ALA:HB3	1.92	0.69
1:B:81:LEU:HA	1:B:84:ILE:HD12	1.73	0.69
1:A:117:TYR:CB	1:A:124:VAL:HG13	2.23	0.68
1:A:126:LEU:HD23	1:A:137:ILE:HA	1.77	0.67
1:A:113:VAL:HA	1:A:130:HIS:O	1.95	0.67
1:A:78:ILE:O	1:A:78:ILE:HG23	1.95	0.66
1:B:206:ILE:O	1:B:206:ILE:HG23	1.95	0.66
1:B:186:SER:HA	1:B:200:ASN:HD22	1.61	0.66
1:B:159:ILE:HG13	1:B:160:SER:N	2.11	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:158:TYR:CB	1:B:188:PRO:HD3	2.27	0.65
1:B:106:ALA:HA	1:B:111:VAL:HG12	1.78	0.65
1:B:156:GLY:O	1:B:188:PRO:HG3	1.96	0.65
1:A:184:MSE:HE2	1:A:202:ASN:ND2	2.13	0.64
1:B:126:LEU:HD23	1:B:137:ILE:HA	1.78	0.64
1:B:144:ILE:HG22	1:B:154:ARG:HB3	1.81	0.64
1:A:117:TYR:HB3	1:A:124:VAL:CG1	2.27	0.63
1:A:161:LEU:HD23	1:A:163:ARG:HG3	1.80	0.63
1:B:78:ILE:HG22	1:B:78:ILE:O	1.97	0.63
1:A:129:THR:HG21	1:A:132:PHE:HB2	1.80	0.62
1:A:80:SER:CB	1:A:83:GLU:HG3	2.20	0.62
1:A:133:ASP:OD1	1:A:134:PRO:HD2	2.00	0.62
1:A:81:LEU:HA	1:A:84:ILE:HD11	1.81	0.62
1:B:186:SER:HA	1:B:200:ASN:ND2	2.14	0.62
1:B:157:GLN:HA	1:B:157:GLN:HE21	1.63	0.62
1:A:223:LEU:HA	1:A:226:THR:CG2	2.32	0.60
1:A:122:ASP:O	1:A:144:ILE:HB	2.02	0.60
1:A:88:VAL:HG13	1:B:226:THR:HG21	1.84	0.59
1:A:148:ILE:O	1:A:148:ILE:HD13	2.03	0.59
1:B:138:GLY:CA	1:B:139:LYS:HE3	2.31	0.59
1:A:210:HIS:O	1:A:214:ILE:HD12	2.02	0.59
1:A:113:VAL:HG12	1:A:202:ASN:O	2.03	0.58
1:A:78:ILE:HG12	1:A:108:GLN:HE22	1.68	0.58
1:A:184:MSE:HE2	1:A:202:ASN:HD21	1.68	0.57
1:B:235:ALA:O	1:B:238:ARG:HD2	2.05	0.57
1:A:104:LYS:HE2	1:A:108:GLN:CG	2.35	0.56
1:B:105:VAL:HG13	1:B:109:LEU:HD23	1.87	0.56
1:B:115:ASN:ND2	1:B:129:THR:HG22	2.21	0.56
1:A:225:LEU:HD22	1:A:229:LYS:HD2	1.87	0.56
1:B:84:ILE:O	1:B:87:LEU:HB2	2.06	0.56
1:A:158:TYR:CD2	1:A:158:TYR:C	2.79	0.56
1:B:210:HIS:O	1:B:214:ILE:HD12	2.06	0.56
1:B:158:TYR:HB2	1:B:187:PHE:HA	1.87	0.56
1:A:184:MSE:HB2	1:A:202:ASN:HD22	1.69	0.55
1:A:126:LEU:HD11	1:A:129:THR:HG23	1.88	0.55
1:A:80:SER:HB3	1:A:83:GLU:CG	2.22	0.55
1:B:222:ASN:N	1:B:222:ASN:HD22	2.04	0.55
1:A:101:VAL:HG11	1:A:224:ILE:HD11	1.87	0.55
1:B:161:LEU:C	1:B:161:LEU:HD12	2.26	0.55
1:A:184:MSE:CB	1:A:202:ASN:HD22	2.21	0.54
1:A:157:GLN:HA	1:A:157:GLN:OE1	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:99:GLN:HE22	1:A:128:ALA:HA	1.74	0.53
1:B:78:ILE:O	1:B:78:ILE:CG2	2.57	0.53
1:B:142:ILE:HD13	1:B:149:THR:HG21	1.91	0.53
1:B:164:ALA:HA	1:B:181:TYR:HB3	1.90	0.52
1:A:148:ILE:CD1	1:A:148:ILE:C	2.67	0.52
1:A:185:LEU:HB2	1:A:209:PHE:CE1	2.45	0.52
1:A:231:ARG:O	1:A:235:ALA:CB	2.57	0.52
1:B:165:SER:O	1:B:166:GLN:HB2	2.10	0.52
1:A:144:ILE:O	1:A:150:GLY:CA	2.59	0.51
1:B:113:VAL:HB	1:B:202:ASN:HD22	1.75	0.51
1:B:142:ILE:HG21	1:B:149:THR:HG22	1.93	0.50
1:B:164:ALA:HA	1:B:181:TYR:CB	2.42	0.50
1:B:82:GLU:HA	1:B:85:SER:HB3	1.94	0.50
1:B:161:LEU:O	1:B:161:LEU:HD12	2.11	0.50
1:A:117:TYR:CB	1:A:124:VAL:CG1	2.89	0.50
1:B:99:GLN:HA	1:B:99:GLN:OE1	2.11	0.50
1:B:114:CYS:SG	1:B:199:ILE:HD11	2.53	0.49
1:B:113:VAL:HG22	1:B:131:GLY:HA3	1.94	0.49
1:A:223:LEU:HA	1:A:226:THR:HG22	1.94	0.49
1:B:161:LEU:HD11	1:B:184:MSE:HE3	1.95	0.49
1:B:117:TYR:HA	1:B:125:VAL:O	2.13	0.48
1:A:219:ILE:HD12	1:B:81:LEU:HB3	1.95	0.48
1:B:152:VAL:HG12	1:B:198:VAL:HG22	1.93	0.48
1:A:200:ASN:C	1:A:201:LEU:HD23	2.34	0.48
1:A:125:VAL:HB	1:A:141:ARG:HG2	1.96	0.48
1:A:225:LEU:CD2	1:A:229:LYS:HD2	2.43	0.48
1:B:146:ASP:O	1:B:147:GLY:O	2.32	0.48
1:A:115:ASN:HD22	1:A:129:THR:CG2	2.15	0.48
1:A:104:LYS:HE2	1:A:108:GLN:HG2	1.95	0.47
1:B:234:VAL:HG13	1:B:235:ALA:N	2.29	0.47
1:B:225:LEU:O	1:B:228:ILE:HB	2.14	0.47
1:B:184:MSE:CG	1:B:185:LEU:N	2.78	0.47
1:B:79:ILE:HD12	1:B:104:LYS:HE2	1.96	0.47
1:A:148:ILE:HD13	1:A:149:THR:N	2.28	0.47
1:B:103:ALA:O	1:B:107:THR:HB	2.16	0.46
1:B:80:SER:O	1:B:84:ILE:HG13	2.16	0.46
1:A:186:SER:HA	1:A:200:ASN:HD22	1.81	0.45
1:A:161:LEU:CD2	1:A:163:ARG:HG3	2.45	0.45
1:A:210:HIS:CD2	1:A:212:ASP:OD2	2.69	0.45
1:B:125:VAL:HG11	1:B:141:ARG:NH1	2.31	0.45
1:B:193:LYS:HG3	1:B:194:GLU:HB2	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:160:SER:HB3	1:A:185:LEU:HD23	1.98	0.45
1:B:154:ARG:HG2	1:B:154:ARG:NH1	2.32	0.45
1:A:84:ILE:HD13	1:B:79:ILE:HG21	1.98	0.45
1:A:104:LYS:HE2	1:A:108:GLN:HG3	1.98	0.45
1:A:94:LEU:HB2	1:A:95:PRO:HD2	1.99	0.45
1:A:123:GLU:OE2	1:A:141:ARG:HD3	2.17	0.45
1:A:200:ASN:O	1:A:201:LEU:HD23	2.18	0.44
1:B:81:LEU:O	1:B:84:ILE:HB	2.16	0.44
1:A:134:PRO:HA	1:A:137:ILE:CD1	2.48	0.44
1:B:220:ILE:HD13	1:B:220:ILE:HA	1.80	0.44
1:B:109:LEU:HD21	1:B:216:PHE:CD2	2.53	0.44
1:B:184:MSE:HG2	1:B:185:LEU:N	2.32	0.44
1:B:234:VAL:CG1	1:B:235:ALA:N	2.80	0.44
1:A:78:ILE:CG2	1:A:78:ILE:O	2.65	0.44
1:B:206:ILE:O	1:B:206:ILE:CG2	2.63	0.44
1:A:191:ASP:OD1	1:A:194:GLU:N	2.51	0.43
1:B:144:ILE:HG22	1:B:154:ARG:CB	2.47	0.43
1:A:95:PRO:O	1:A:99:GLN:HG2	2.18	0.43
1:A:76:LEU:O	1:B:81:LEU:HG	2.19	0.43
1:A:234:VAL:HG22	1:B:233:GLN:OE1	2.18	0.43
1:A:109:LEU:HD12	1:A:201:LEU:HD11	2.01	0.43
1:B:158:TYR:CD1	1:B:158:TYR:C	2.90	0.42
1:B:87:LEU:HD22	1:B:87:LEU:HA	1.79	0.42
1:B:117:TYR:HB3	1:B:124:VAL:CG1	2.49	0.42
1:B:149:THR:HG23	1:B:149:THR:O	2.17	0.42
1:A:94:LEU:CB	1:A:95:PRO:CD	2.98	0.42
1:A:134:PRO:HA	1:A:137:ILE:HD12	2.02	0.42
1:B:187:PHE:CE2	1:B:218:SER:HA	2.55	0.42
1:A:156:GLY:O	1:A:188:PRO:HG3	2.20	0.42
1:B:126:LEU:HD21	1:B:129:THR:HG23	2.01	0.42
1:B:94:LEU:HD12	1:B:94:LEU:HA	1.88	0.42
1:A:161:LEU:O	1:A:161:LEU:HD22	2.20	0.41
1:A:119:ARG:NE	1:A:121:GLY:O	2.53	0.41
1:B:154:ARG:HH11	1:B:154:ARG:HG2	1.85	0.41
1:A:187:PHE:CD2	1:A:188:PRO:CD	3.01	0.41
1:B:119:ARG:HH21	1:B:144:ILE:HG13	1.84	0.41
1:A:94:LEU:O	1:A:95:PRO:C	2.59	0.41
1:B:231:ARG:O	1:B:234:VAL:HG12	2.20	0.41
1:B:225:LEU:HD22	1:B:229:LYS:HD3	2.03	0.41
1:A:184:MSE:HB2	1:A:202:ASN:ND2	2.35	0.41
1:B:200:ASN:HA	1:B:200:ASN:HD22	1.63	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:122:ASP:OD2	1:A:122:ASP:C	2.58	0.41
1:B:222:ASN:N	1:B:222:ASN:ND2	2.68	0.41
1:B:117:TYR:CB	1:B:124:VAL:CG1	2.99	0.41
1:A:207:ARG:HG3	1:A:207:ARG:O	2.21	0.41
1:A:97:VAL:O	1:A:101:VAL:HG23	2.22	0.40
1:A:158:TYR:HB2	1:A:187:PHE:HA	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	151/181 (83%)	129 (85%)	16 (11%)	6 (4%)	4 21
1	B	148/181 (82%)	125 (84%)	13 (9%)	10 (7%)	1 8
All	All	299/362 (83%)	254 (85%)	29 (10%)	16 (5%)	2 14

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	76	LEU
1	B	75	HIS
1	B	89	SER
1	A	169	ARG
1	B	90	SER
1	B	147	GLY
1	B	146	ASP
1	B	181	TYR
1	A	144	ILE
1	B	235	ALA
1	B	77	GLU
1	B	236	SER

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Mol	Chain	Res	Type
1	A	192	LYS
1	B	237	SER
1	A	95	PRO
1	A	113	VAL

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	129/153 (84%)	101 (78%)	28 (22%)	1 5
1	B	130/153 (85%)	101 (78%)	29 (22%)	1 5
All	All	259/306 (85%)	202 (78%)	57 (22%)	1 5

All (57) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	78	ILE
1	A	82	GLU
1	A	84	ILE
1	A	90	SER
1	A	94	LEU
1	A	104	LYS
1	A	110	LYS
1	A	124	VAL
1	A	125	VAL
1	A	129	THR
1	A	137	ILE
1	A	148	ILE
1	A	151	SER
1	A	161	LEU
1	A	163	ARG
1	A	165	SER
1	A	184	MSE
1	A	185	LEU
1	A	193	LYS

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Mol	Chain	Res	Type
1	A	194	GLU
1	A	202	ASN
1	A	204	THR
1	A	206	ILE
1	A	207	ARG
1	A	212	ASP
1	A	213	GLU
1	A	225	LEU
1	A	229	LYS
1	B	86	MSE
1	B	87	LEU
1	B	90	SER
1	B	92	PHE
1	B	96	GLU
1	B	107	THR
1	B	110	LYS
1	B	111	VAL
1	B	119	ARG
1	B	137	ILE
1	B	146	ASP
1	B	148	ILE
1	B	149	THR
1	B	151	SER
1	B	154	ARG
1	B	157	GLN
1	B	159	ILE
1	B	160	SER
1	B	181	TYR
1	B	182	ASN
1	B	192	LYS
1	B	193	LYS
1	B	201	LEU
1	B	202	ASN
1	B	204	THR
1	B	208	SER
1	B	225	LEU
1	B	229	LYS
1	B	233	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	99	GLN
1	A	108	GLN
1	A	115	ASN
1	A	182	ASN
1	A	200	ASN
1	A	202	ASN
1	A	233	GLN
1	B	115	ASN
1	B	130	HIS
1	B	157	GLN
1	B	200	ASN
1	B	202	ASN
1	B	210	HIS
1	B	222	ASN

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 2 ligands modelled in this entry, 2 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 (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	153/181 (84%)	0.01	6 (3%)	43	21	51, 68, 87, 98	0
1	B	150/181 (82%)	-0.03	2 (1%)	79	62	53, 67, 84, 97	0
All	All	303/362 (83%)	-0.01	8 (2%)	59	35	51, 68, 85, 98	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	77	GLU	4.5
1	A	233	GLN	4.3
1	A	236	SER	3.8
1	A	167	ASP	2.6
1	A	75	HIS	2.6
1	A	238	ARG	2.6
1	B	76	LEU	2.3
1	A	234	VAL	2.2

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

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	NA	B	2	1/1	0.94	0.17	-1.84	49,49,49,49	0
2	NA	A	1	1/1	0.97	0.11	-5.03	46,46,46,46	0

6.5 Other polymers (i)

There are no such residues in this entry.