Full wwPDB X-ray Structure Validation Report (i)

Aug 27, 2023 – 09:45 PM EDT

PDB ID	:	3I6X
Title	:	Crystal structure of the calponin homology domain of IQGAP1
Authors	:	Kurella, V.B.
Deposited on	:	2009-07-07
Resolution	:	2.50 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:

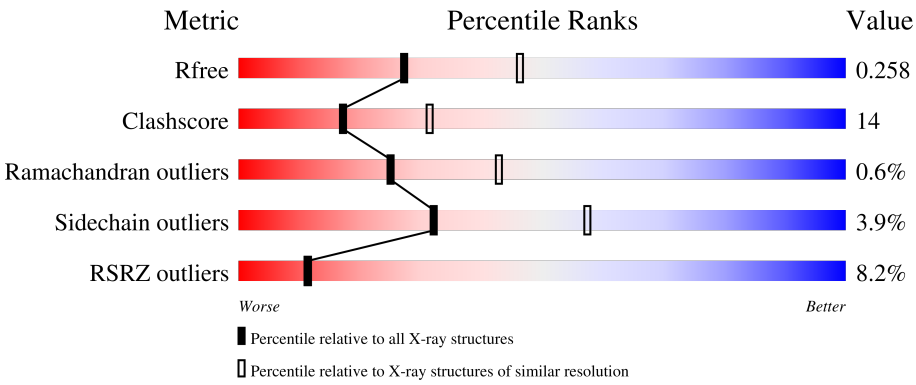
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

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 2.50 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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	193	7%, 63%, 20%, •, 15%
1	B	193	6%, 66%, 17%, •, 15%
1	C	193	7%, 62%, 21%, •, 15%
1	D	193	9%, 60%, 23%, •, 15%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 5374 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 Ras GTPase-activating-like protein IQGAP1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	164	1334	853	225	248	8	0	0	0
1	B	164	1334	853	225	248	8	0	0	0
1	C	164	1334	853	225	248	8	0	0	0
1	D	164	1334	853	225	248	8	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP P46940
A	0	ALA	-	expression tag	UNP P46940
B	-1	GLY	-	expression tag	UNP P46940
B	0	ALA	-	expression tag	UNP P46940
C	-1	GLY	-	expression tag	UNP P46940
C	0	ALA	-	expression tag	UNP P46940
D	-1	GLY	-	expression tag	UNP P46940
D	0	ALA	-	expression tag	UNP P46940

- Molecule 2 is water.

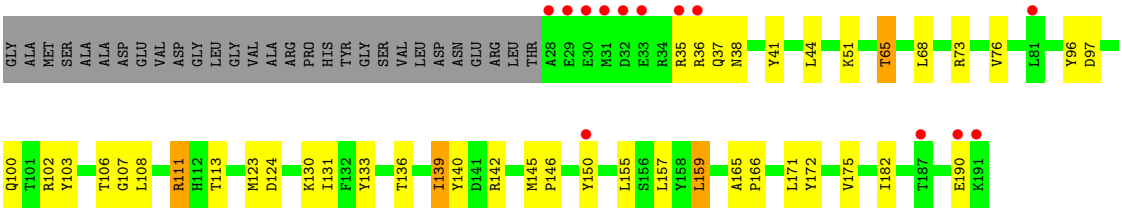
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	17	17	17	0	0
2	B	9	9	9	0	0
2	C	6	6	6	0	0
2	D	6	6	6	0	0

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.

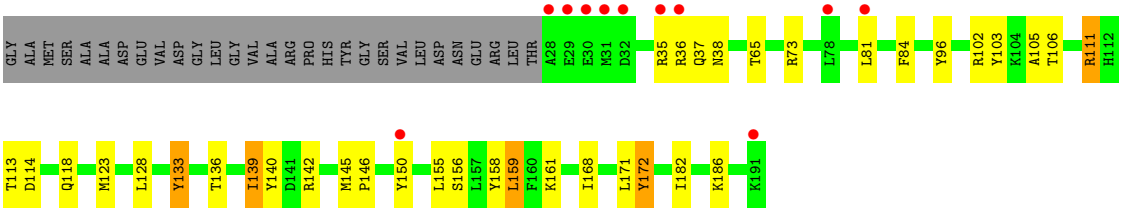
• Molecule 1: Ras GTPase-activating-like protein IQGAP1

Chain A:



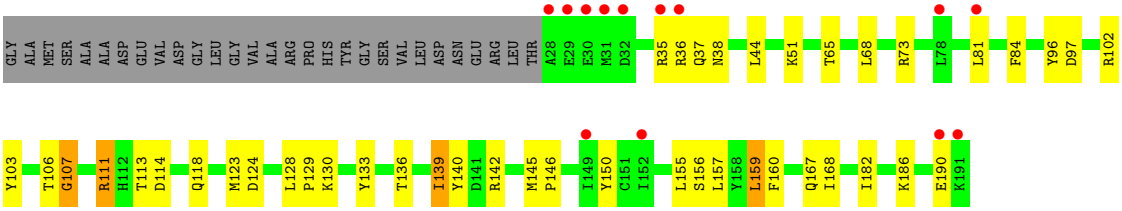
• Molecule 1: Ras GTPase-activating-like protein IQGAP1

Chain B:



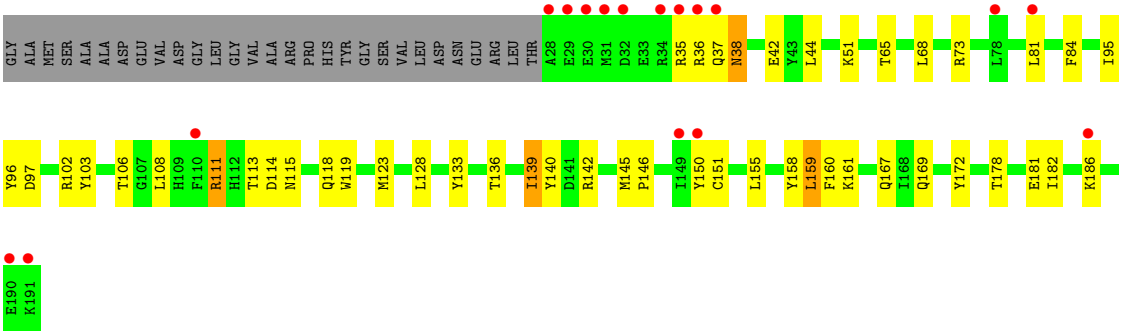
• Molecule 1: Ras GTPase-activating-like protein IQGAP1

Chain C:



• Molecule 1: Ras GTPase-activating-like protein IQGAP1

Chain D:


GLY
ALA
MET
SER
ALA
ALA
ASP
GLU
VAL
ASP
GLY
LEU
GLY
VAL
ALA
ARG
PRO
HIS
TYR
GLY
SER
VAL
LEU
ASP
ASN
GLU
ARG
LEU
THR
A28
E29
E30
M31
D32
E33
R34
R35
R36
Q37
N38
E42
Y43
L44
K51
T65
L68
R73
L78
L81
F84
I95
Y96
D97
R102
Y103
T106
G107
L108
H109
F110
R111
H112
T113
D114
N115
Q118
W119
M123
L128
Y133
T136
I139
Y140
D141
R142
M145
P146
I149
Y150
C151
L155
Y158
L159
F160
K161
Q167
I168
Q169
Y172
T178
E181
I182
K186
E190
K191

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	57.43Å 66.42Å 83.01Å 99.55° 104.07° 108.69°	Depositor
Resolution (Å)	25.00 – 2.50 24.80 – 2.39	Depositor EDS
% Data completeness (in resolution range)	98.5 (25.00-2.50) 96.7 (24.80-2.39)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$< I/\sigma(I) >$ [1]	1.36 (at 2.39Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.234 , 0.263 0.234 , 0.258	Depositor DCC
R_{free} test set	1858 reflections (4.47%)	wwPDB-VP
Wilson B-factor (Å^2)	40.9	Xtriage
Anisotropy	0.155	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 45.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	5374	wwPDB-VP
Average B, all atoms (Å^2)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.83% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\#\|Z\| >5$	RMSZ	$\#\|Z\| >5$
1	A	0.39	0/1364	0.58	0/1844
1	B	0.40	0/1364	0.57	0/1844
1	C	0.37	0/1364	0.56	0/1844
1	D	0.37	0/1364	0.56	0/1844
All	All	0.38	0/5456	0.57	0/7376

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	133	TYR	Sidechain

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	1334	0	1295	39	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1334	0	1295	32	0
1	C	1334	0	1295	41	0
1	D	1334	0	1295	38	0
2	A	17	0	0	3	0
2	B	9	0	0	0	0
2	C	6	0	0	0	0
2	D	6	0	0	1	0
All	All	5374	0	5180	146	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 14.

All (146) 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:139:ILE:HG21	2:A:199:HOH:O	1.51	1.10
1:C:37:GLN:HE21	1:C:182:ILE:HG22	1.41	0.86
1:C:113:THR:HG23	1:C:136:THR:HG21	1.57	0.85
1:D:113:THR:HG23	1:D:136:THR:HG21	1.62	0.81
1:B:73:ARG:HD3	1:B:142:ARG:HB2	1.63	0.80
1:A:73:ARG:HD3	1:A:142:ARG:HB2	1.64	0.80
1:C:73:ARG:HD3	1:C:142:ARG:HB2	1.65	0.78
1:B:113:THR:HG23	1:B:136:THR:HG21	1.67	0.75
1:D:73:ARG:HD3	1:D:142:ARG:HB2	1.67	0.75
1:B:96:TYR:CD1	1:B:111:ARG:HG2	2.23	0.73
1:D:38:ASN:O	1:D:42:GLU:HG3	1.88	0.73
1:D:37:GLN:HE21	1:D:182:ILE:HG22	1.52	0.72
1:D:123:MET:HE2	1:D:155:LEU:HD22	1.73	0.70
1:C:44:LEU:HD12	1:C:182:ILE:HD13	1.73	0.69
1:C:123:MET:HE2	1:C:155:LEU:HD22	1.73	0.69
1:D:103:TYR:HE1	1:D:108:LEU:HD22	1.58	0.69
1:B:123:MET:CE	1:B:155:LEU:HD22	2.24	0.68
1:A:113:THR:HG23	1:A:136:THR:HG21	1.77	0.66
1:D:73:ARG:HA	1:D:139:ILE:HD12	1.76	0.66
1:A:35:ARG:HH11	1:A:35:ARG:HA	1.60	0.66
1:B:35:ARG:HH11	1:B:35:ARG:HA	1.60	0.66
1:C:123:MET:CE	1:C:155:LEU:HD22	2.27	0.64
1:B:123:MET:HE2	1:B:155:LEU:HD13	1.80	0.64
1:B:123:MET:HE2	1:B:155:LEU:HD22	1.80	0.64
1:A:35:ARG:HA	1:A:35:ARG:NH1	2.13	0.64
1:A:37:GLN:HE21	1:A:182:ILE:HG22	1.61	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:35:ARG:HH11	1:D:35:ARG:HA	1.62	0.64
1:C:73:ARG:HA	1:C:139:ILE:HD12	1.78	0.63
1:C:35:ARG:HA	1:C:35:ARG:HH11	1.62	0.62
1:C:96:TYR:CD1	1:C:111:ARG:HG2	2.33	0.62
1:A:123:MET:HE2	1:A:155:LEU:HD22	1.80	0.62
1:B:37:GLN:HE21	1:B:182:ILE:HG22	1.64	0.62
1:A:65:THR:HG22	2:A:201:HOH:O	1.98	0.62
1:C:35:ARG:HA	1:C:35:ARG:NH1	2.15	0.62
1:D:35:ARG:HA	1:D:35:ARG:NH1	2.15	0.62
1:B:73:ARG:HA	1:B:139:ILE:HD12	1.80	0.61
1:B:35:ARG:HA	1:B:35:ARG:NH1	2.14	0.61
1:A:73:ARG:HA	1:A:139:ILE:HD12	1.83	0.61
1:C:113:THR:HG23	1:C:136:THR:CG2	2.31	0.61
1:C:139:ILE:HG13	1:C:140:TYR:N	2.17	0.60
1:D:37:GLN:HE21	1:D:182:ILE:CG2	2.15	0.60
1:A:96:TYR:CD1	1:A:111:ARG:HG2	2.37	0.60
1:A:123:MET:CE	1:A:155:LEU:HD22	2.33	0.58
1:D:123:MET:CE	1:D:155:LEU:HD22	2.33	0.58
1:C:37:GLN:HE21	1:C:182:ILE:CG2	2.15	0.57
1:C:103:TYR:O	1:C:106:THR:O	2.22	0.57
1:A:106:THR:HB	2:A:200:HOH:O	2.04	0.57
1:D:139:ILE:HG13	1:D:140:TYR:N	2.19	0.57
1:A:97:ASP:OD2	1:A:102:ARG:HB3	2.05	0.57
1:C:146:PRO:O	1:C:150:TYR:HD1	1.89	0.56
1:D:113:THR:HG23	1:D:136:THR:CG2	2.35	0.56
1:D:44:LEU:HD12	1:D:182:ILE:HD13	1.88	0.56
1:A:146:PRO:O	1:A:150:TYR:HD1	1.89	0.55
1:A:157:LEU:HD11	1:D:158:TYR:HD1	1.73	0.54
1:C:96:TYR:CE2	1:C:111:ARG:HB3	2.42	0.54
1:A:155:LEU:HG	1:A:159:LEU:HD22	1.89	0.54
1:C:106:THR:HG22	1:C:107:GLY:N	2.23	0.53
1:A:139:ILE:HG13	1:A:140:TYR:N	2.22	0.53
1:A:96:TYR:CE2	1:A:111:ARG:HB3	2.43	0.53
1:C:97:ASP:CG	1:C:102:ARG:HB3	2.30	0.52
1:B:139:ILE:HG13	1:B:140:TYR:N	2.25	0.52
1:A:44:LEU:HD12	1:A:182:ILE:HD13	1.92	0.51
1:B:37:GLN:HE21	1:B:182:ILE:CG2	2.23	0.51
1:A:113:THR:HG23	1:A:136:THR:CG2	2.40	0.51
1:A:124:ASP:OD2	1:A:130:LYS:HE3	2.11	0.50
1:A:123:MET:HE2	1:A:155:LEU:HD13	1.94	0.50
1:B:113:THR:HG23	1:B:136:THR:CG2	2.37	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:97:ASP:CG	1:D:102:ARG:HB3	2.33	0.50
1:B:133:TYR:HD1	1:B:133:TYR:N	2.10	0.49
1:C:145:MET:N	1:C:146:PRO:CD	2.75	0.49
1:A:171:LEU:O	1:A:172:TYR:C	2.51	0.49
1:B:133:TYR:N	1:B:133:TYR:CD1	2.79	0.48
1:B:145:MET:N	1:B:146:PRO:CD	2.76	0.48
1:A:123:MET:CE	1:A:155:LEU:HB2	2.44	0.48
1:B:123:MET:HE3	1:B:155:LEU:HD22	1.95	0.48
1:A:76:VAL:HG21	1:A:100:GLN:HG2	1.97	0.47
1:A:165:ALA:HB1	1:A:166:PRO:HD2	1.96	0.47
1:C:81:LEU:O	1:C:84:PHE:HB3	2.15	0.47
1:B:155:LEU:HG	1:B:159:LEU:HD22	1.96	0.47
1:C:146:PRO:O	1:C:150:TYR:CD1	2.67	0.47
1:C:44:LEU:CD1	1:C:182:ILE:HD13	2.42	0.47
1:A:103:TYR:HE1	1:A:108:LEU:HD22	1.80	0.46
1:D:106:THR:HG22	1:D:106:THR:O	2.15	0.46
1:B:161:LYS:HA	1:B:161:LYS:HD3	1.75	0.46
1:D:123:MET:CE	1:D:155:LEU:HB2	2.46	0.46
1:A:36:ARG:HD2	1:A:190:GLU:CD	2.35	0.46
1:A:37:GLN:HE21	1:A:182:ILE:CG2	2.25	0.46
1:B:96:TYR:CG	1:B:111:ARG:HG2	2.50	0.46
1:C:124:ASP:OD2	1:C:130:LYS:HE3	2.16	0.46
1:D:145:MET:N	1:D:146:PRO:CD	2.80	0.45
1:A:103:TYR:O	1:A:106:THR:O	2.34	0.45
1:D:178:THR:OG1	1:D:181:GLU:HG3	2.16	0.45
1:A:157:LEU:HD11	1:D:158:TYR:CD1	2.51	0.45
1:B:36:ARG:HG2	1:B:186:LYS:HG3	1.99	0.45
1:B:73:ARG:HA	1:B:139:ILE:CD1	2.47	0.45
1:C:128:LEU:HD12	1:C:129:PRO:HD2	1.99	0.44
1:B:123:MET:HE1	1:B:128:LEU:CD2	2.48	0.44
1:C:36:ARG:HD2	1:C:190:GLU:CD	2.37	0.44
1:C:73:ARG:HA	1:C:139:ILE:CD1	2.47	0.43
1:C:142:ARG:O	1:C:142:ARG:HG2	2.17	0.43
1:D:96:TYR:CD1	1:D:111:ARG:HG2	2.52	0.43
1:D:142:ARG:O	1:D:142:ARG:HG2	2.18	0.43
1:A:131:ILE:HD11	1:D:150:TYR:HD1	1.81	0.43
1:B:123:MET:CE	1:B:155:LEU:HB2	2.48	0.43
1:C:114:ASP:O	1:C:118:GLN:HG3	2.19	0.43
1:D:81:LEU:O	1:D:84:PHE:HB3	2.19	0.43
1:B:146:PRO:O	1:B:150:TYR:CD1	2.72	0.43
1:D:169:GLN:HG3	2:D:193:HOH:O	2.17	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:36:ARG:HG2	1:D:186:LYS:HG3	2.01	0.43
1:C:37:GLN:NE2	1:C:182:ILE:HG22	2.22	0.43
1:C:159:LEU:HD12	1:C:159:LEU:HA	1.91	0.43
1:D:161:LYS:HD3	1:D:161:LYS:HA	1.81	0.42
1:B:156:SER:OG	1:B:168:ILE:HB	2.19	0.42
1:D:133:TYR:N	1:D:133:TYR:CD1	2.88	0.42
1:D:155:LEU:HG	1:D:159:LEU:HD22	2.00	0.42
1:A:51:LYS:HB2	1:A:68:LEU:HD22	2.02	0.42
1:B:81:LEU:O	1:B:84:PHE:HB3	2.20	0.42
1:B:103:TYR:O	1:B:106:THR:O	2.36	0.42
1:C:123:MET:CE	1:C:155:LEU:HB2	2.50	0.42
1:D:123:MET:HE1	1:D:151:CYS:SG	2.59	0.42
1:C:51:LYS:HB2	1:C:68:LEU:HD22	2.02	0.42
1:C:133:TYR:N	1:C:133:TYR:CD1	2.88	0.42
1:C:160:PHE:CE1	1:C:167:GLN:HB2	2.56	0.41
1:B:114:ASP:O	1:B:118:GLN:HG3	2.21	0.41
1:D:160:PHE:CE1	1:D:167:GLN:HB2	2.54	0.41
1:A:146:PRO:O	1:A:150:TYR:CD1	2.73	0.41
1:C:156:SER:OG	1:C:168:ILE:HB	2.20	0.41
1:C:36:ARG:HG2	1:C:186:LYS:HG3	2.02	0.41
1:D:51:LYS:HB2	1:D:68:LEU:HD22	2.02	0.41
1:D:119:TRP:CZ2	1:D:123:MET:HE3	2.55	0.41
1:B:146:PRO:O	1:B:150:TYR:HD1	2.03	0.41
1:C:44:LEU:HD12	1:C:182:ILE:CD1	2.46	0.41
1:B:171:LEU:O	1:B:172:TYR:C	2.57	0.41
1:C:123:MET:HE1	1:C:128:LEU:CD2	2.51	0.41
1:A:145:MET:N	1:A:146:PRO:CD	2.84	0.41
1:C:113:THR:CG2	1:C:136:THR:HG21	2.39	0.41
1:D:95:ILE:HG23	1:D:115:ASN:OD1	2.19	0.41
1:D:114:ASP:O	1:D:118:GLN:HG3	2.21	0.41
1:A:41:TYR:OH	1:A:175:VAL:O	2.39	0.40
1:A:97:ASP:CG	1:A:102:ARG:HB3	2.40	0.40
1:B:158:TYR:HD1	1:C:157:LEU:HD11	1.86	0.40
1:D:123:MET:HE1	1:D:128:LEU:CD2	2.51	0.40
1:A:36:ARG:HD2	1:A:190:GLU:OE2	2.22	0.40
1:A:73:ARG:HA	1:A:139:ILE:CD1	2.49	0.40
1:C:96:TYR:CG	1:C:111:ARG:HG2	2.56	0.40
1:D:42:GLU:HG2	1:D:172:TYR:HE1	1.85	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	162/193 (84%)	157 (97%)	4 (2%)	1 (1%)	25	43
1	B	162/193 (84%)	156 (96%)	4 (2%)	2 (1%)	13	24
1	C	162/193 (84%)	158 (98%)	3 (2%)	1 (1%)	25	43
1	D	162/193 (84%)	159 (98%)	3 (2%)	0	100	100
All	All	648/772 (84%)	630 (97%)	14 (2%)	4 (1%)	25	43

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	172	TYR
1	A	107	GLY
1	B	105	ALA
1	C	107	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	140/168 (83%)	135 (96%)	5 (4%)	35	61
1	B	140/168 (83%)	133 (95%)	7 (5%)	24	46
1	C	140/168 (83%)	135 (96%)	5 (4%)	35	61
1	D	140/168 (83%)	135 (96%)	5 (4%)	35	61
All	All	560/672 (83%)	538 (96%)	22 (4%)	32	57

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	38	ASN
1	A	65	THR
1	A	111	ARG
1	A	139	ILE
1	A	159	LEU
1	B	38	ASN
1	B	65	THR
1	B	102	ARG
1	B	111	ARG
1	B	133	TYR
1	B	139	ILE
1	B	159	LEU
1	C	38	ASN
1	C	65	THR
1	C	111	ARG
1	C	139	ILE
1	C	159	LEU
1	D	38	ASN
1	D	65	THR
1	D	111	ARG
1	D	139	ILE
1	D	159	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	118	GLN
1	B	118	GLN
1	C	118	GLN
1	D	118	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains (i)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

There are no ligands in this entry.

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	164/193 (84%)	0.38	13 (7%)	12	12	18, 36, 92, 115	0
1	B	164/193 (84%)	0.37	11 (6%)	17	18	18, 36, 90, 117	0
1	C	164/193 (84%)	0.59	13 (7%)	12	12	22, 45, 87, 118	0
1	D	164/193 (84%)	0.48	17 (10%)	6	6	25, 44, 87, 115	0
All	All	656/772 (84%)	0.46	54 (8%)	11	11	18, 41, 93, 118	0

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	191	LYS	7.5
1	C	28	ALA	7.2
1	D	28	ALA	6.6
1	A	28	ALA	6.3
1	C	29	GLU	5.9
1	B	28	ALA	5.8
1	A	191	LYS	5.8
1	B	191	LYS	5.6
1	A	31	MET	5.6
1	D	191	LYS	5.1
1	C	31	MET	4.6
1	B	31	MET	4.5
1	B	32	ASP	4.3
1	D	31	MET	4.3
1	D	32	ASP	4.3
1	B	29	GLU	4.0
1	C	35	ARG	3.8
1	A	32	ASP	3.7
1	D	35	ARG	3.7
1	D	29	GLU	3.6
1	A	35	ARG	3.5

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Mol	Chain	Res	Type	RSRZ
1	A	29	GLU	3.5
1	C	32	ASP	3.4
1	C	78	LEU	3.4
1	C	30	GLU	3.4
1	B	35	ARG	3.1
1	B	78	LEU	3.1
1	B	81	LEU	3.1
1	D	150	TYR	3.0
1	D	186	LYS	3.0
1	D	81	LEU	3.0
1	A	81	LEU	2.9
1	D	30	GLU	2.9
1	D	190	GLU	2.9
1	C	149	ILE	2.8
1	B	30	GLU	2.7
1	B	150	TYR	2.7
1	A	33	GLU	2.6
1	C	190	GLU	2.6
1	D	110	PHE	2.5
1	A	150	TYR	2.5
1	C	81	LEU	2.5
1	A	30	GLU	2.5
1	A	187	THR	2.4
1	C	152	ILE	2.3
1	D	37	GLN	2.3
1	D	36	ARG	2.2
1	D	34	ARG	2.2
1	A	36	ARG	2.2
1	D	149	ILE	2.2
1	C	36	ARG	2.1
1	D	78	LEU	2.1
1	B	36	ARG	2.1
1	A	190	GLU	2.1

6.2 Non-standard residues in protein, DNA, RNA chains (i)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.