Full wwPDB X-ray Structure Validation Report (i)

Feb 18, 2024 – 10:12 PM EST

PDB ID	:	4GCK
Title	:	structure of no-dna complex
Authors	:	Schumacher, M.A.
Deposited on	:	2012-07-30
Resolution	:	2.05 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:

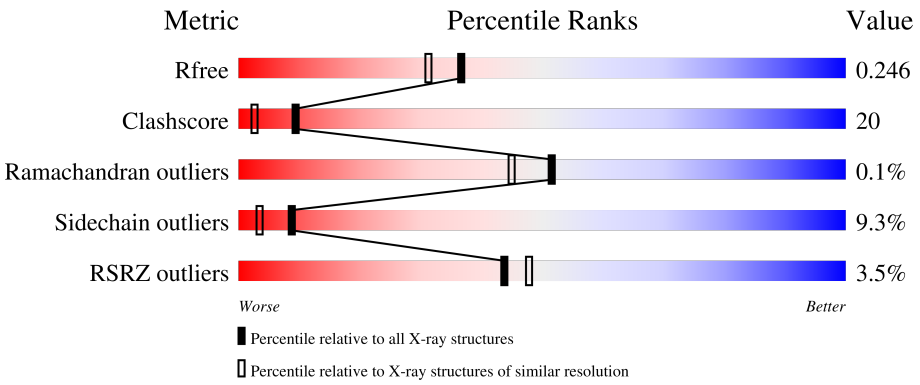
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 2.05 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	1692 (2.04-2.04)
Clashscore	141614	1773 (2.04-2.04)
Ramachandran outliers	138981	1752 (2.04-2.04)
Sidechain outliers	138945	1752 (2.04-2.04)
RSRZ outliers	127900	1672 (2.04-2.04)

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	212	2%, 56%, 29%, 5%, 10%
1	B	212	5%, 60%, 23%, 6%, 11%
1	C	212	2%, 62%, 25%, •, 10%
1	D	212	4%, 59%, 25%, 5%, 11%
2	W	12	17%, 75%, 8%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
2	Z	12	42% 50% 8%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 7088 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 Nucleoid occlusion factor SlmA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	190	1544	973	279	285	7	0	0	0
1	B	189	1533	967	275	284	7	0	0	0
1	C	190	1544	973	279	285	7	0	0	0
1	D	189	1533	967	275	284	7	0	0	0

There are 56 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-13	MET	-	expression tag	UNP B5XTG2
A	-12	PRO	-	expression tag	UNP B5XTG2
A	-11	PRO	-	expression tag	UNP B5XTG2
A	-10	GLY	-	expression tag	UNP B5XTG2
A	-9	LYS	-	expression tag	UNP B5XTG2
A	-8	CYS	-	expression tag	UNP B5XTG2
A	-7	LEU	-	expression tag	UNP B5XTG2
A	-6	PHE	-	expression tag	UNP B5XTG2
A	-5	SER	-	expression tag	UNP B5XTG2
A	-4	GLY	-	expression tag	UNP B5XTG2
A	-3	VAL	-	expression tag	UNP B5XTG2
A	-2	PHE	-	expression tag	UNP B5XTG2
A	-1	CYS	-	expression tag	UNP B5XTG2
A	0	ASN	-	expression tag	UNP B5XTG2
B	-13	MET	-	expression tag	UNP B5XTG2
B	-12	PRO	-	expression tag	UNP B5XTG2
B	-11	PRO	-	expression tag	UNP B5XTG2
B	-10	GLY	-	expression tag	UNP B5XTG2
B	-9	LYS	-	expression tag	UNP B5XTG2
B	-8	CYS	-	expression tag	UNP B5XTG2
B	-7	LEU	-	expression tag	UNP B5XTG2

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	-6	PHE	-	expression tag	UNP B5XTG2
B	-5	SER	-	expression tag	UNP B5XTG2
B	-4	GLY	-	expression tag	UNP B5XTG2
B	-3	VAL	-	expression tag	UNP B5XTG2
B	-2	PHE	-	expression tag	UNP B5XTG2
B	-1	CYS	-	expression tag	UNP B5XTG2
B	0	ASN	-	expression tag	UNP B5XTG2
C	-13	MET	-	expression tag	UNP B5XTG2
C	-12	PRO	-	expression tag	UNP B5XTG2
C	-11	PRO	-	expression tag	UNP B5XTG2
C	-10	GLY	-	expression tag	UNP B5XTG2
C	-9	LYS	-	expression tag	UNP B5XTG2
C	-8	CYS	-	expression tag	UNP B5XTG2
C	-7	LEU	-	expression tag	UNP B5XTG2
C	-6	PHE	-	expression tag	UNP B5XTG2
C	-5	SER	-	expression tag	UNP B5XTG2
C	-4	GLY	-	expression tag	UNP B5XTG2
C	-3	VAL	-	expression tag	UNP B5XTG2
C	-2	PHE	-	expression tag	UNP B5XTG2
C	-1	CYS	-	expression tag	UNP B5XTG2
C	0	ASN	-	expression tag	UNP B5XTG2
D	-13	MET	-	expression tag	UNP B5XTG2
D	-12	PRO	-	expression tag	UNP B5XTG2
D	-11	PRO	-	expression tag	UNP B5XTG2
D	-10	GLY	-	expression tag	UNP B5XTG2
D	-9	LYS	-	expression tag	UNP B5XTG2
D	-8	CYS	-	expression tag	UNP B5XTG2
D	-7	LEU	-	expression tag	UNP B5XTG2
D	-6	PHE	-	expression tag	UNP B5XTG2
D	-5	SER	-	expression tag	UNP B5XTG2
D	-4	GLY	-	expression tag	UNP B5XTG2
D	-3	VAL	-	expression tag	UNP B5XTG2
D	-2	PHE	-	expression tag	UNP B5XTG2
D	-1	CYS	-	expression tag	UNP B5XTG2
D	0	ASN	-	expression tag	UNP B5XTG2

- Molecule 2 is a DNA chain called DNA (5'-D(*GP*TP*GP*AP*GP*TP*AP*CP*TP*CP*AP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	W	12	243	117	45	70	11	0	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	Z	12	243	117	45	70	11	0	0	0

- Molecule 3 is water.

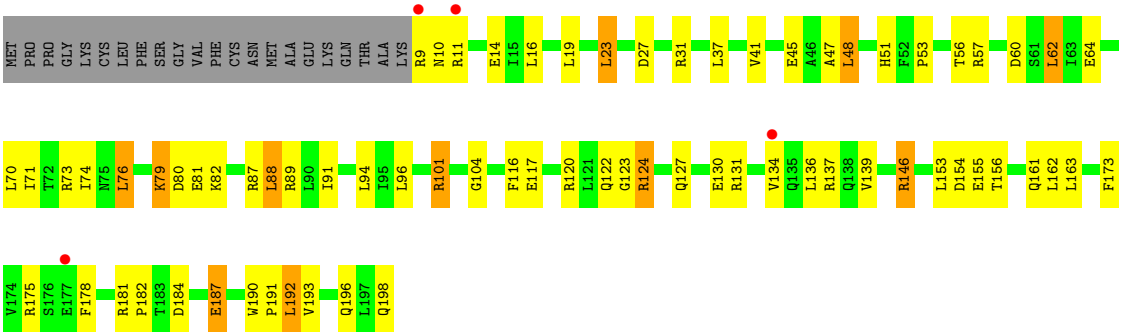
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	102	102	102	0	0
3	B	80	80	80	0	0
3	C	116	116	116	0	0
3	D	108	108	108	0	0
3	W	21	21	21	0	0
3	Z	21	21	21	0	0

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.

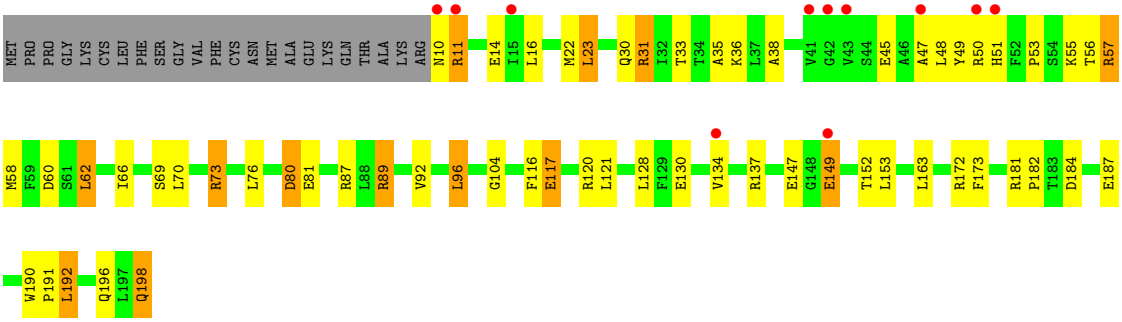
• Molecule 1: Nucleoid occlusion factor SlmA

Chain A: 2% 56% 29% 5% 10%



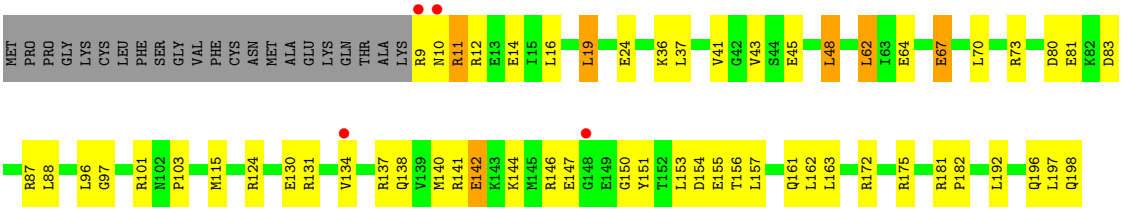
• Molecule 1: Nucleoid occlusion factor SlmA

Chain B: 5% 60% 23% 6% 11%



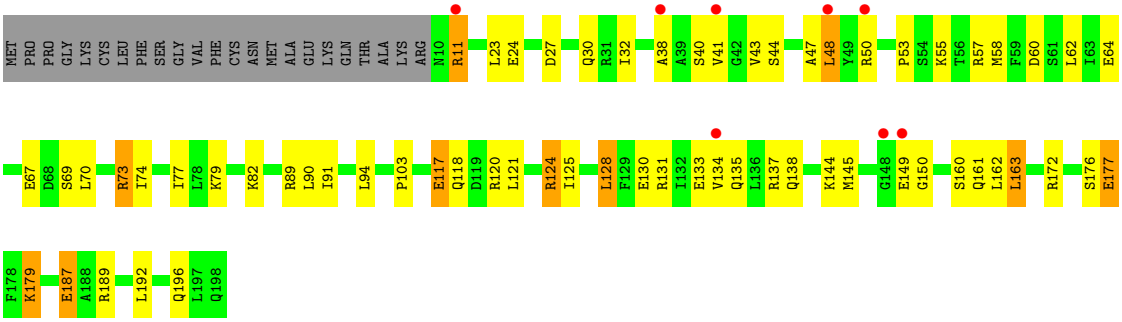
• Molecule 1: Nucleoid occlusion factor SlmA

Chain C: 2% 62% 25% • 10%



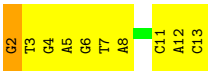
• Molecule 1: Nucleoid occlusion factor SlmA

Chain D:



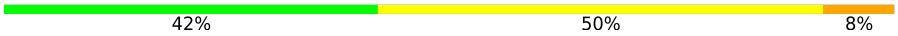
• Molecule 2: DNA (5'-D(*GP*TP*GP*AP*GP*TP*AP*CP*TP*CP*AP*C)-3')

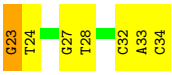
Chain W:



• Molecule 2: DNA (5'-D(*GP*TP*GP*AP*GP*TP*AP*CP*TP*CP*AP*C)-3')

Chain Z:





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	71.49Å 80.75Å 201.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	74.00 – 2.05 74.94 – 2.05	Depositor EDS
% Data completeness (in resolution range)	94.8 (74.00-2.05) 94.9 (74.94-2.05)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.21 (at 2.05Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.218 , 0.243 0.221 , 0.246	Depositor DCC
R_{free} test set	7056 reflections (10.02%)	wwPDB-VP
Wilson B-factor (Å^2)	27.6	Xtriage
Anisotropy	0.812	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.38 , 55.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	7088	wwPDB-VP
Average B, all atoms (Å^2)	37.0	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 44.95 % 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 1.4170e-04. 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	0.54	0/1564	0.73	1/2098 (0.0%)
1	B	0.54	0/1553	0.70	1/2084 (0.0%)
1	C	0.59	0/1564	0.76	2/2098 (0.1%)
1	D	0.55	0/1553	0.72	2/2084 (0.1%)
2	W	0.61	0/272	0.88	0/418
2	Z	0.60	0/272	0.95	1/418 (0.2%)
All	All	0.56	0/6778	0.75	7/9200 (0.1%)

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
2	W	0	1
2	Z	0	1
All	All	0	2

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	80	ASP	CB-CG-OD1	8.70	126.13	118.30
1	D	89	ARG	NE-CZ-NH2	5.98	123.29	120.30
1	A	89	ARG	NE-CZ-NH1	-5.91	117.35	120.30
2	Z	32	DC	OP2-P-O3'	5.70	117.73	105.20
1	B	80	ASP	CB-CG-OD1	5.51	123.26	118.30
1	C	175	ARG	CB-CA-C	-5.43	99.54	110.40
1	D	89	ARG	NE-CZ-NH1	-5.27	117.67	120.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	W	2	DG	Sidechain
2	Z	23	DG	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	1544	0	1585	72	0
1	B	1533	0	1572	56	0
1	C	1544	0	1585	70	1
1	D	1533	0	1572	60	1
2	W	243	0	137	15	0
2	Z	243	0	137	5	0
3	A	102	0	0	10	0
3	B	80	0	0	1	0
3	C	116	0	0	16	0
3	D	108	0	0	7	0
3	W	21	0	0	0	0
3	Z	21	0	0	0	0
All	All	7088	0	6588	260	1

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 20.

All (260) 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:55:LYS:HA	1:B:58:MET:HE3	1.33	1.09
1:D:55:LYS:HA	1:D:58:MET:HE3	1.27	1.08
1:C:161:GLN:HE21	1:C:196:GLN:NE2	1.53	1.03
1:C:161:GLN:NE2	1:C:196:GLN:HE22	1.67	0.92
1:A:161:GLN:HE21	1:A:196:GLN:HE22	1.17	0.88
1:B:81:GLU:O	1:B:87:ARG:HD2	1.74	0.88
1:A:130:GLU:O	1:A:134:VAL:HG23	1.76	0.86
1:C:134:VAL:HG22	1:C:137:ARG:NH2	1.91	0.86
1:D:73:ARG:HG2	1:D:73:ARG:HH11	1.42	0.85
1:A:101:ARG:HG2	1:A:101:ARG:HH11	1.42	0.83

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:10:ASN:O	1:A:14:GLU:HB2	1.77	0.83
1:A:81:GLU:O	1:A:87:ARG:HD2	1.80	0.82
1:D:134:VAL:HG22	1:D:137:ARG:NH2	1.95	0.81
1:C:161:GLN:HE21	1:C:196:GLN:HE22	0.86	0.81
1:B:11:ARG:HA	1:B:14:GLU:HG3	1.63	0.81
1:C:81:GLU:O	1:C:87:ARG:HD2	1.80	0.81
1:C:147:GLU:OE2	3:C:312:HOH:O	1.99	0.80
1:C:142:GLU:HG2	1:C:146:ARG:HE	1.46	0.80
1:C:67:GLU:HG3	3:C:227:HOH:O	1.81	0.80
1:B:53:PRO:HD2	1:B:57:ARG:HG3	1.63	0.78
1:C:24:GLU:HG3	1:C:103:PRO:HB2	1.63	0.78
1:C:196:GLN:HE21	1:D:192:LEU:HB3	1.50	0.77
1:B:117:GLU:OE2	3:B:267:HOH:O	2.03	0.77
1:C:115:MET:HA	1:C:115:MET:CE	2.15	0.77
1:C:64:GLU:HG3	1:C:124:ARG:HH22	1.47	0.76
1:C:67:GLU:OE1	3:C:314:HOH:O	2.03	0.76
1:C:163:LEU:HD23	1:D:172:ARG:HH12	1.50	0.76
1:B:55:LYS:HA	1:B:58:MET:CE	2.15	0.75
1:A:161:GLN:HE21	1:A:196:GLN:NE2	1.85	0.75
1:D:138:GLN:HG2	3:D:266:HOH:O	1.86	0.74
1:A:11:ARG:NH1	2:W:2:DG:H5"	2.03	0.73
1:C:134:VAL:HG13	3:C:265:HOH:O	1.88	0.73
1:D:118:GLN:OE1	1:D:120:ARG:HG2	1.89	0.73
2:W:2:DG:H2'	2:W:3:DT:C7	2.19	0.72
1:B:130:GLU:O	1:B:134:VAL:HG23	1.88	0.72
1:D:161:GLN:HE21	1:D:196:GLN:NE2	1.88	0.72
2:W:5:DA:H1'	2:W:6:DG:H5'	1.71	0.72
1:A:101:ARG:HH11	1:A:101:ARG:CG	2.03	0.72
1:A:73:ARG:HH12	1:A:101:ARG:HH22	1.38	0.71
1:B:134:VAL:HG22	1:B:137:ARG:NH2	2.05	0.71
1:D:53:PRO:HD2	1:D:57:ARG:HG3	1.72	0.71
1:A:187:GLU:HB2	3:A:260:HOH:O	1.90	0.71
1:A:9:ARG:HG3	1:A:9:ARG:HH11	1.57	0.69
1:D:43:VAL:HG23	1:D:44:SER:O	1.93	0.69
1:A:134:VAL:HG21	3:A:246:HOH:O	1.91	0.69
1:D:64:GLU:CG	1:D:124:ARG:HH12	2.06	0.69
1:C:12:ARG:NH1	3:C:275:HOH:O	2.03	0.69
1:D:70:LEU:O	1:D:74:ILE:HG12	1.93	0.69
1:A:146:ARG:HG3	1:A:146:ARG:HH11	1.57	0.68
1:A:70:LEU:O	1:A:74:ILE:HG12	1.92	0.68
1:D:74:ILE:HD12	1:D:91:ILE:HG12	1.76	0.68

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:153:LEU:HD11	1:A:198:GLN:HB2	1.75	0.68
1:C:64:GLU:HG3	1:C:124:ARG:NH2	2.08	0.68
1:D:64:GLU:HG2	1:D:124:ARG:HH12	1.58	0.68
1:D:11:ARG:HG2	1:D:11:ARG:HH21	1.58	0.67
1:B:57:ARG:HH11	1:B:57:ARG:HA	1.60	0.67
1:A:73:ARG:HH12	1:A:101:ARG:NH2	1.91	0.66
1:C:101:ARG:HB3	3:C:308:HOH:O	1.96	0.65
1:C:130:GLU:O	1:C:134:VAL:HG23	1.96	0.65
1:A:153:LEU:CD1	1:A:198:GLN:HB2	2.26	0.65
1:A:101:ARG:HG2	3:A:229:HOH:O	1.98	0.63
1:C:19:LEU:HD13	1:C:62:LEU:HD22	1.80	0.63
1:D:38:ALA:HB2	1:D:48:LEU:HD22	1.80	0.63
1:B:11:ARG:HG2	1:B:11:ARG:HH21	1.64	0.62
1:C:73:ARG:NH1	1:C:101:ARG:HH22	1.97	0.62
1:C:134:VAL:HG23	3:C:306:HOH:O	1.99	0.62
1:B:181:ARG:HB2	1:B:184:ASP:CG	2.21	0.61
1:C:161:GLN:NE2	1:C:196:GLN:NE2	2.35	0.61
1:C:64:GLU:CG	1:C:124:ARG:HH22	2.15	0.60
1:C:19:LEU:HD13	1:C:62:LEU:CD2	2.32	0.60
1:C:134:VAL:CG2	3:C:306:HOH:O	2.49	0.59
1:D:135:GLN:NE2	3:D:233:HOH:O	2.27	0.59
1:C:115:MET:HA	1:C:115:MET:HE2	1.83	0.59
1:C:138:GLN:HG2	3:C:279:HOH:O	2.01	0.59
1:A:88:LEU:HD13	1:A:139:VAL:HG11	1.84	0.59
1:D:118:GLN:NE2	3:D:251:HOH:O	2.36	0.59
1:D:73:ARG:HG2	1:D:73:ARG:NH1	2.16	0.59
1:D:177:GLU:OE2	1:D:179:LYS:NZ	2.36	0.58
1:B:38:ALA:HB2	1:B:48:LEU:HD22	1.85	0.58
1:C:144:LYS:HE3	1:C:150:GLY:HA3	1.85	0.58
2:W:2:DG:H2'	2:W:3:DT:H72	1.85	0.58
1:A:193:VAL:HA	1:B:196:GLN:HE21	1.69	0.57
2:Z:27:DG:H2"	2:Z:28:DT:H5"	1.84	0.57
1:D:130:GLU:O	1:D:134:VAL:HG23	2.05	0.57
1:A:64:GLU:HG3	1:A:124:ARG:HH22	1.70	0.56
1:A:123:GLY:O	1:A:127:GLN:HG3	2.05	0.56
1:B:69:SER:O	1:B:73:ARG:HG3	2.04	0.56
1:C:154:ASP:OD1	1:C:156:THR:N	2.36	0.55
1:A:74:ILE:CD1	1:A:94:LEU:HD23	2.37	0.55
1:C:134:VAL:HG22	1:C:137:ARG:HH22	1.70	0.55
1:C:10:ASN:O	1:C:14:GLU:HB2	2.06	0.54
1:C:16:LEU:HD22	1:C:62:LEU:HD13	1.88	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:45:GLU:O	1:C:48:LEU:HB2	2.07	0.54
1:A:31:ARG:HB3	1:A:116:PHE:CD2	2.43	0.54
1:C:73:ARG:HH12	1:C:101:ARG:NH2	2.05	0.53
2:W:2:DG:H2'	2:W:3:DT:H73	1.90	0.53
1:A:31:ARG:HD2	3:A:276:HOH:O	2.07	0.53
1:A:134:VAL:HG22	1:A:137:ARG:NH2	2.22	0.53
1:B:57:ARG:NH1	1:B:57:ARG:HG2	2.24	0.53
1:D:117:GLU:HB3	1:D:121:LEU:HD12	1.90	0.53
1:A:74:ILE:HD12	1:A:91:ILE:HG23	1.91	0.53
1:C:196:GLN:HG2	1:D:192:LEU:O	2.07	0.53
1:C:115:MET:HA	1:C:115:MET:HE3	1.91	0.53
1:B:190:TRP:HB3	1:B:191:PRO:HD3	1.91	0.53
1:C:144:LYS:CD	1:C:150:GLY:HA3	2.38	0.53
1:C:153:LEU:HD12	1:C:198:GLN:HB2	1.90	0.53
1:D:30:GLN:O	1:D:32:ILE:HD12	2.09	0.53
1:A:60:ASP:OD2	1:A:120:ARG:NH2	2.43	0.52
1:A:153:LEU:HD12	1:A:198:GLN:O	2.09	0.52
1:A:37:LEU:O	1:A:41:VAL:HG23	2.10	0.52
1:B:60:ASP:OD1	1:B:120:ARG:NH2	2.43	0.52
1:B:11:ARG:HB2	1:B:51:HIS:NE2	2.25	0.52
1:C:9:ARG:HG3	1:C:9:ARG:HH11	1.75	0.52
1:A:101:ARG:CG	1:A:101:ARG:NH1	2.69	0.52
1:A:45:GLU:O	1:A:48:LEU:HB2	2.10	0.51
2:Z:23:DG:H2'	2:Z:24:DT:C7	2.40	0.51
1:A:74:ILE:CD1	1:A:91:ILE:HG23	2.40	0.51
1:D:57:ARG:NH2	1:D:60:ASP:OD2	2.43	0.51
1:B:49:TYR:CE1	2:W:13:DC:H2'	2.45	0.51
1:D:60:ASP:CG	1:D:120:ARG:HH22	2.14	0.51
1:D:134:VAL:HG22	1:D:137:ARG:HH21	1.73	0.51
1:A:193:VAL:HA	1:B:196:GLN:NE2	2.25	0.51
1:B:30:GLN:HA	1:B:116:PHE:CZ	2.46	0.51
1:C:73:ARG:NH1	1:C:101:ARG:NH2	2.59	0.51
1:D:149:GLU:OE2	1:D:149:GLU:N	2.44	0.51
1:D:47:ALA:O	1:D:50:ARG:HG2	2.11	0.50
1:D:144:LYS:HG2	1:D:150:GLY:HA3	1.93	0.50
2:Z:23:DG:H2'	2:Z:24:DT:H72	1.92	0.50
1:B:173:PHE:CD1	1:B:182:PRO:HD3	2.46	0.50
1:C:144:LYS:CE	1:C:150:GLY:HA3	2.41	0.50
1:C:172:ARG:HH12	1:D:160:SER:HB2	1.77	0.50
1:D:67:GLU:HG3	1:D:128:LEU:HG	1.94	0.50
1:B:11:ARG:HH21	1:B:11:ARG:CG	2.22	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:115:MET:CE	3:C:281:HOH:O	2.59	0.50
1:C:131:ARG:HD2	3:C:216:HOH:O	2.11	0.50
1:B:60:ASP:CG	1:B:120:ARG:HH22	2.15	0.50
1:D:73:ARG:HH11	1:D:73:ARG:CG	2.19	0.50
1:A:154:ASP:OD1	1:A:156:THR:N	2.44	0.49
1:D:79:LYS:HG3	3:D:224:HOH:O	2.12	0.49
1:D:24:GLU:HG3	1:D:103:PRO:HB2	1.94	0.49
1:A:79:LYS:HD2	1:A:80:ASP:OD1	2.13	0.49
1:B:35:ALA:HB2	1:B:45:GLU:OE1	2.13	0.49
1:C:153:LEU:CD1	1:C:198:GLN:HB2	2.42	0.49
1:D:134:VAL:HG13	3:D:253:HOH:O	2.12	0.49
1:A:192:LEU:HD11	1:B:153:LEU:CD1	2.42	0.49
1:D:77:ILE:HD12	1:D:90:LEU:HB3	1.94	0.49
1:B:45:GLU:OE2	2:W:13:DC:N4	2.37	0.48
1:A:74:ILE:HD13	1:A:94:LEU:HD23	1.95	0.48
1:A:161:GLN:NE2	1:A:196:GLN:HE22	1.98	0.48
1:D:133:GLU:OE1	1:D:163:LEU:HD12	2.13	0.48
1:A:146:ARG:HH11	1:A:146:ARG:CG	2.25	0.48
1:B:11:ARG:HA	1:B:14:GLU:CG	2.41	0.48
1:B:57:ARG:HH11	1:B:57:ARG:HG2	1.79	0.47
1:B:89:ARG:HG3	1:B:190:TRP:CD2	2.50	0.47
1:D:176:SER:O	1:D:179:LYS:HD3	2.14	0.47
1:A:134:VAL:HG13	3:A:302:HOH:O	2.14	0.47
2:Z:23:DG:H8	2:Z:23:DG:HO5'	1.62	0.47
1:A:146:ARG:HB3	3:A:299:HOH:O	2.14	0.47
1:A:11:ARG:HD3	1:A:51:HIS:NE2	2.30	0.47
1:D:11:ARG:HG2	1:D:11:ARG:NH2	2.27	0.46
1:A:9:ARG:HH11	1:A:9:ARG:CG	2.25	0.46
1:B:10:ASN:O	1:B:14:GLU:HG3	2.15	0.46
1:C:73:ARG:HH12	1:C:101:ARG:CZ	2.29	0.46
1:D:74:ILE:CD1	1:D:94:LEU:HD23	2.46	0.46
1:D:64:GLU:HG3	1:D:124:ARG:HH12	1.80	0.46
1:C:192:LEU:O	1:D:196:GLN:HG2	2.16	0.46
1:A:173:PHE:CD1	1:A:178:PHE:HA	2.52	0.45
1:A:196:GLN:HE21	1:B:192:LEU:HB3	1.82	0.45
1:D:73:ARG:NH1	1:D:73:ARG:CG	2.78	0.45
1:D:30:GLN:HB2	3:D:279:HOH:O	2.17	0.45
1:A:73:ARG:NH1	1:A:101:ARG:HH22	2.11	0.45
1:C:151:TYR:CE1	1:C:197:LEU:HB3	2.52	0.45
1:C:172:ARG:NH1	1:D:160:SER:HB2	2.31	0.45
1:B:57:ARG:HH11	1:B:57:ARG:CA	2.28	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:11:ARG:NH1	1:C:43:VAL:CG1	2.80	0.45
1:A:56:THR:CG2	1:A:120:ARG:HH22	2.29	0.44
1:B:11:ARG:CG	1:B:11:ARG:NH2	2.80	0.44
1:B:181:ARG:HB2	1:B:184:ASP:OD1	2.17	0.44
1:C:36:LYS:HE2	3:C:256:HOH:O	2.17	0.44
1:D:125:ILE:HD13	1:D:125:ILE:HA	1.85	0.44
2:Z:33:DA:C5	2:Z:34:DC:C4	3.04	0.44
1:B:22:MET:HE3	1:B:36:LYS:HG2	1.99	0.44
1:B:47:ALA:O	1:B:50:ARG:HG2	2.17	0.44
1:B:57:ARG:HH11	1:B:57:ARG:CG	2.30	0.44
1:C:146:ARG:NH1	3:C:288:HOH:O	2.28	0.44
1:A:64:GLU:CG	1:A:124:ARG:HH22	2.30	0.44
1:D:64:GLU:HG3	1:D:124:ARG:HH22	1.83	0.44
1:D:53:PRO:CD	1:D:57:ARG:HG3	2.45	0.44
1:A:23:LEU:HB3	1:A:104:GLY:CA	2.48	0.44
2:W:7:DT:H6	2:W:7:DT:H2'	1.67	0.44
1:A:79:LYS:O	1:A:82:LYS:HE2	2.18	0.43
1:D:74:ILE:CD1	1:D:91:ILE:HG12	2.45	0.43
1:B:56:THR:HG23	1:B:121:LEU:HD11	1.99	0.43
1:D:176:SER:C	1:D:177:GLU:HG3	2.38	0.43
1:A:53:PRO:CG	1:A:57:ARG:HG3	2.49	0.43
1:C:115:MET:CE	1:C:115:MET:CA	2.93	0.43
1:C:141:ARG:HG3	1:C:155:GLU:HG2	2.00	0.43
1:D:60:ASP:OD1	1:D:120:ARG:NH2	2.47	0.43
2:W:6:DG:H4'	2:W:7:DT:OP1	2.17	0.43
1:D:43:VAL:CG2	1:D:47:ALA:HB3	2.48	0.43
2:W:7:DT:H2"	2:W:8:DA:C8	2.54	0.43
1:C:11:ARG:NH1	1:C:43:VAL:HG11	2.34	0.43
1:D:69:SER:O	1:D:73:ARG:HG3	2.19	0.43
1:B:62:LEU:HD12	1:B:62:LEU:HA	1.82	0.43
1:A:101:ARG:NH1	3:A:244:HOH:O	2.45	0.43
1:B:31:ARG:HH12	2:W:12:DA:H5"	1.84	0.43
1:D:74:ILE:HD11	1:D:94:LEU:HD23	2.00	0.43
1:D:134:VAL:HG23	3:D:232:HOH:O	2.18	0.43
1:A:79:LYS:CE	1:A:80:ASP:OD1	2.67	0.43
1:B:66:ILE:O	1:B:70:LEU:HD23	2.19	0.43
2:W:3:DT:H2"	2:W:4:DG:N7	2.34	0.42
1:A:190:TRP:HB3	1:A:191:PRO:HD3	2.02	0.42
1:B:62:LEU:O	1:B:66:ILE:HG13	2.19	0.42
1:A:9:ARG:HG3	1:A:9:ARG:NH1	2.30	0.42
1:A:9:ARG:CG	1:A:9:ARG:NH1	2.81	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:19:LEU:HD23	1:A:62:LEU:HD22	2.01	0.42
1:B:76:LEU:O	1:B:80:ASP:HB2	2.19	0.42
1:C:197:LEU:HB2	3:C:287:HOH:O	2.20	0.42
1:D:69:SER:O	1:D:73:ARG:CG	2.68	0.42
1:B:23:LEU:HB3	1:B:104:GLY:CA	2.49	0.42
1:A:134:VAL:CG2	3:A:246:HOH:O	2.61	0.42
1:A:134:VAL:CG2	3:A:277:HOH:O	2.68	0.42
1:C:97:GLY:O	1:C:101:ARG:HG2	2.20	0.42
1:C:144:LYS:HG3	1:C:150:GLY:N	2.34	0.42
1:C:83:ASP:O	1:C:87:ARG:HD3	2.19	0.42
1:A:181:ARG:HB2	1:A:184:ASP:OD2	2.19	0.41
1:B:48:LEU:HD12	1:B:48:LEU:HA	1.93	0.41
1:C:172:ARG:HH12	1:D:160:SER:CB	2.33	0.41
2:W:5:DA:C1'	2:W:6:DG:H5'	2.45	0.41
1:A:74:ILE:HD12	1:A:91:ILE:HG12	2.00	0.41
1:B:33:THR:HG22	2:W:11:DC:H3'	2.02	0.41
1:A:181:ARG:HA	1:A:182:PRO:HD3	1.88	0.41
1:A:154:ASP:OD1	1:A:156:THR:HB	2.20	0.41
1:A:74:ILE:HD11	1:A:94:LEU:HD23	2.02	0.41
1:B:152:THR:HB	1:B:198:GLN:OXT	2.20	0.41
1:C:140:MET:CE	1:C:162:LEU:HD12	2.50	0.41
1:A:163:LEU:HD23	1:B:172:ARG:HH12	1.85	0.41
1:C:115:MET:HE1	3:C:281:HOH:O	2.19	0.41
1:B:134:VAL:HG22	1:B:137:ARG:HH22	1.84	0.41
1:B:149:GLU:OE1	1:B:149:GLU:O	2.39	0.41
2:W:12:DA:C5	2:W:13:DC:C4	3.09	0.41
1:A:76:LEU:HD13	1:A:76:LEU:HA	1.75	0.41
1:C:37:LEU:O	1:C:41:VAL:HG23	2.21	0.41
1:C:115:MET:HE3	3:C:281:HOH:O	2.19	0.41
1:C:9:ARG:HG3	1:C:9:ARG:NH1	2.36	0.40
1:C:157:LEU:HD11	1:D:189:ARG:HG2	2.02	0.40
1:C:181:ARG:HA	1:C:182:PRO:HD3	1.88	0.40
1:C:11:ARG:HH11	1:C:43:VAL:HG11	1.86	0.40
1:A:11:ARG:HH12	1:A:47:ALA:HB1	1.85	0.40
1:A:16:LEU:HD22	1:A:62:LEU:HD13	2.02	0.40
1:B:16:LEU:HD22	1:B:62:LEU:HD13	2.01	0.40
1:B:73:ARG:HH11	1:B:73:ARG:HG2	1.86	0.40
1:B:92:VAL:HG12	1:B:96:LEU:HD22	2.03	0.40
1:C:24:GLU:OE1	1:C:103:PRO:HG2	2.21	0.40
1:D:48:LEU:HD12	1:D:48:LEU:HA	1.95	0.40
1:B:50:ARG:HG3	1:B:51:HIS:ND1	2.36	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:163:LEU:HD12	1:B:163:LEU:HA	1.80	0.40
1:A:56:THR:HG22	1:A:120:ARG:HH22	1.87	0.40
1:A:117:GLU:HB3	3:A:204:HOH:O	2.22	0.40
1:A:173:PHE:CD1	1:A:182:PRO:HD3	2.57	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:146:ARG:NH2	1:D:187:GLU:OE2[4_555]	2.07	0.13

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	188/212 (89%)	186 (99%)	2 (1%)	0	100	100
1	B	187/212 (88%)	181 (97%)	6 (3%)	0	100	100
1	C	188/212 (89%)	184 (98%)	4 (2%)	0	100	100
1	D	187/212 (88%)	181 (97%)	5 (3%)	1 (0%)	29	18
All	All	750/848 (88%)	732 (98%)	17 (2%)	1 (0%)	51	45

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	41	VAL

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	167/185 (90%)	147 (88%)	20 (12%)	5 1
1	B	166/185 (90%)	151 (91%)	15 (9%)	9 4
1	C	167/185 (90%)	158 (95%)	9 (5%)	22 13
1	D	166/185 (90%)	148 (89%)	18 (11%)	6 2
All	All	666/740 (90%)	604 (91%)	62 (9%)	9 3

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	23	LEU
1	A	27	ASP
1	A	48	LEU
1	A	62	LEU
1	A	71	ILE
1	A	76	LEU
1	A	79	LYS
1	A	88	LEU
1	A	96	LEU
1	A	101	ARG
1	A	122	GLN
1	A	124	ARG
1	A	131	ARG
1	A	136	LEU
1	A	146	ARG
1	A	155	GLU
1	A	162	LEU
1	A	175	ARG
1	A	187	GLU
1	A	192	LEU
1	B	11	ARG
1	B	23	LEU
1	B	31	ARG
1	B	57	ARG
1	B	62	LEU
1	B	73	ARG
1	B	89	ARG
1	B	96	LEU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	117	GLU
1	B	128	LEU
1	B	147	GLU
1	B	149	GLU
1	B	187	GLU
1	B	192	LEU
1	B	198	GLN
1	C	11	ARG
1	C	19	LEU
1	C	48	LEU
1	C	62	LEU
1	C	67	GLU
1	C	70	LEU
1	C	88	LEU
1	C	96	LEU
1	C	142	GLU
1	D	11	ARG
1	D	23	LEU
1	D	27	ASP
1	D	40	SER
1	D	48	LEU
1	D	62	LEU
1	D	73	ARG
1	D	82	LYS
1	D	117	GLU
1	D	124	ARG
1	D	128	LEU
1	D	131	ARG
1	D	145	MET
1	D	162	LEU
1	D	163	LEU
1	D	177	GLU
1	D	179	LYS
1	D	187	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	GLN
1	A	196	GLN
1	B	10	ASN
1	B	118	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	196	GLN
1	C	122	GLN
1	C	196	GLN
1	D	17	GLN
1	D	122	GLN
1	D	196	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	190/212 (89%)	0.14	4 (2%)	63	67	22, 33, 54, 88	0
1	B	189/212 (89%)	0.43	11 (5%)	23	25	23, 34, 68, 90	0
1	C	190/212 (89%)	0.22	4 (2%)	63	67	17, 29, 48, 95	0
1	D	189/212 (89%)	0.37	8 (4%)	36	39	21, 32, 65, 76	0
2	W	12/12 (100%)	-0.12	0	100	100	33, 44, 52, 53	0
2	Z	12/12 (100%)	-0.29	0	100	100	34, 40, 50, 50	0
All	All	782/872 (89%)	0.27	27 (3%)	44	48	17, 33, 60, 95	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	41	VAL	5.2
1	A	134	VAL	5.1
1	B	11	ARG	4.5
1	D	41	VAL	4.4
1	C	9	ARG	4.1
1	A	11	ARG	3.9
1	D	134	VAL	3.9
1	C	134	VAL	3.7
1	B	134	VAL	3.6
1	C	10	ASN	3.0
1	B	43	VAL	3.0
1	B	50	ARG	2.9
1	B	51	HIS	2.7
1	B	149	GLU	2.7
1	D	38	ALA	2.6
1	A	9	ARG	2.5
1	B	47	ALA	2.5
1	D	50	ARG	2.4
1	D	11	ARG	2.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	42	GLY	2.3
1	B	10	ASN	2.2
1	B	15	ILE	2.2
1	D	149	GLU	2.1
1	D	148	GLY	2.1
1	A	177	GLU	2.0
1	C	148	GLY	2.0
1	D	48	LEU	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.