



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:02 PM UTC

PDB ID : 4JL3 / pdb_00004jl3
Title : Crystal structure of ms6564-dna complex
Authors : Yang, S.F.; Gao, Z.Q.; He, Z.G.; Dong, Y.H.
Deposited on : 2013-03-12
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 ⓘ 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 ⓘ](#)) were used in the production of this report:

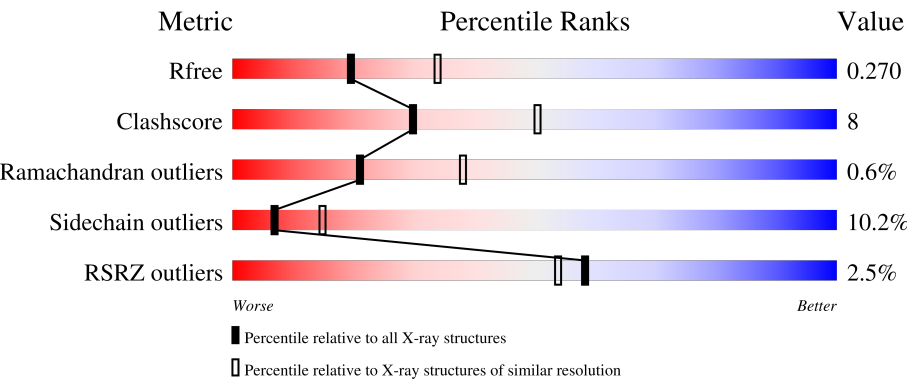
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

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}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (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 $\leq 5\%$. 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	196	3% 69% 21% • 9%
1	B	196	3% 73% 16% • 8%
1	C	196	2% 73% 17% • 7%
1	D	196	2% 65% 22% • 8%
2	E	31	3% 48% 45% 6%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
3	F	31	 A horizontal bar chart showing the quality of chain F. The bar is divided into three segments: green (42%), yellow (55%), and orange (3%). The green segment is labeled '42%' and the yellow segment is labeled '55%'. The orange segment is at the far right and contains a small black dot.

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6858 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 Transcriptional regulator, TetR family.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	179	Total	C	N	O	Se	0	1	0
			1369	846	258	262	3			
1	B	181	Total	C	N	O	Se	0	1	0
			1391	858	266	264	3			
1	C	182	Total	C	N	O	Se	0	0	0
			1386	854	266	263	3			
1	D	180	Total	C	N	O	Se	0	0	0
			1376	849	264	260	3			

There are 60 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-6	HIS	-	expression tag	UNP A0R6I8
A	-5	MSE	-	expression tag	UNP A0R6I8
A	-4	ALA	-	expression tag	UNP A0R6I8
A	-3	SER	-	expression tag	UNP A0R6I8
A	-2	MSE	-	expression tag	UNP A0R6I8
A	-1	THR	-	expression tag	UNP A0R6I8
A	0	GLY	-	expression tag	UNP A0R6I8
A	1	GLY	-	expression tag	UNP A0R6I8
A	2	GLN	-	expression tag	UNP A0R6I8
A	3	GLN	-	expression tag	UNP A0R6I8
A	4	MSE	-	expression tag	UNP A0R6I8
A	5	GLY	-	expression tag	UNP A0R6I8
A	6	ARG	-	expression tag	UNP A0R6I8
A	7	GLY	-	expression tag	UNP A0R6I8
A	8	SER	-	expression tag	UNP A0R6I8
B	-6	HIS	-	expression tag	UNP A0R6I8
B	-5	MSE	-	expression tag	UNP A0R6I8
B	-4	ALA	-	expression tag	UNP A0R6I8
B	-3	SER	-	expression tag	UNP A0R6I8
B	-2	MSE	-	expression tag	UNP A0R6I8
B	-1	THR	-	expression tag	UNP A0R6I8

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
B	0	GLY	-	expression tag	UNP A0R6I8
B	1	GLY	-	expression tag	UNP A0R6I8
B	2	GLN	-	expression tag	UNP A0R6I8
B	3	GLN	-	expression tag	UNP A0R6I8
B	4	MSE	-	expression tag	UNP A0R6I8
B	5	GLY	-	expression tag	UNP A0R6I8
B	6	ARG	-	expression tag	UNP A0R6I8
B	7	GLY	-	expression tag	UNP A0R6I8
B	8	SER	-	expression tag	UNP A0R6I8
C	-6	HIS	-	expression tag	UNP A0R6I8
C	-5	MSE	-	expression tag	UNP A0R6I8
C	-4	ALA	-	expression tag	UNP A0R6I8
C	-3	SER	-	expression tag	UNP A0R6I8
C	-2	MSE	-	expression tag	UNP A0R6I8
C	-1	THR	-	expression tag	UNP A0R6I8
C	0	GLY	-	expression tag	UNP A0R6I8
C	1	GLY	-	expression tag	UNP A0R6I8
C	2	GLN	-	expression tag	UNP A0R6I8
C	3	GLN	-	expression tag	UNP A0R6I8
C	4	MSE	-	expression tag	UNP A0R6I8
C	5	GLY	-	expression tag	UNP A0R6I8
C	6	ARG	-	expression tag	UNP A0R6I8
C	7	GLY	-	expression tag	UNP A0R6I8
C	8	SER	-	expression tag	UNP A0R6I8
D	-6	HIS	-	expression tag	UNP A0R6I8
D	-5	MSE	-	expression tag	UNP A0R6I8
D	-4	ALA	-	expression tag	UNP A0R6I8
D	-3	SER	-	expression tag	UNP A0R6I8
D	-2	MSE	-	expression tag	UNP A0R6I8
D	-1	THR	-	expression tag	UNP A0R6I8
D	0	GLY	-	expression tag	UNP A0R6I8
D	1	GLY	-	expression tag	UNP A0R6I8
D	2	GLN	-	expression tag	UNP A0R6I8
D	3	GLN	-	expression tag	UNP A0R6I8
D	4	MSE	-	expression tag	UNP A0R6I8
D	5	GLY	-	expression tag	UNP A0R6I8
D	6	ARG	-	expression tag	UNP A0R6I8
D	7	GLY	-	expression tag	UNP A0R6I8
D	8	SER	-	expression tag	UNP A0R6I8

- Molecule 2 is a DNA chain called DNA (31-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	31	Total	C	N	O	P	0	0	0
			636	303	114	188	31			

- Molecule 3 is a DNA chain called DNA (31-MER).

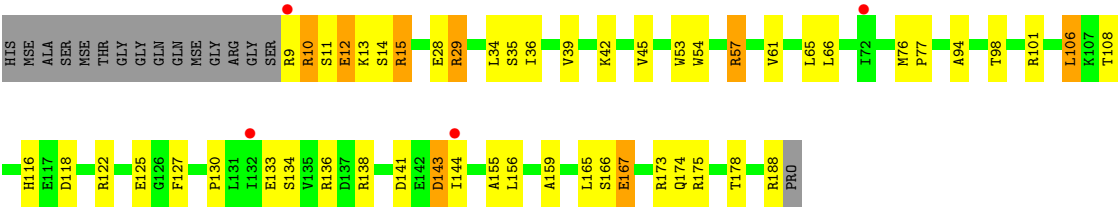
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	F	31	Total	C	N	O	P	0	0	0
			635	302	118	184	31			

- Molecule 4 is water.

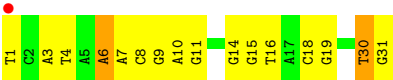
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	12	Total	O	0	0
			12	12		
4	B	10	Total	O	0	0
			10	10		
4	C	12	Total	O	0	0
			12	12		
4	D	11	Total	O	0	0
			11	11		
4	E	9	Total	O	0	0
			9	9		
4	F	11	Total	O	0	0
			11	11		

- Molecule 1: Transcriptional regulator, TetR family





● Molecule 2: DNA (31-MER)



● Molecule 3: DNA (31-MER)



4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, α , β , γ	100.95Å 100.95Å 99.86Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	35.50 – 2.50 35.50 – 2.50	Depositor EDS
% Data completeness (in resolution range)	98.6 (35.50-2.50) 98.6 (35.50-2.50)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.41 (at 2.51Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, R_{free}	0.224 , 0.280 0.214 , 0.270	Depositor DCC
R_{free} test set	1952 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	54.2	Xtriage
Anisotropy	0.635	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 44.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.026 for -h,-k,l 0.059 for h,-h-k,-l 0.026 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6858	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.52% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.56	0/1383	0.93	0/1867
1	B	0.55	0/1405	0.92	1/1895 (0.1%)
1	C	0.54	0/1399	0.87	0/1885
1	D	0.58	0/1389	1.01	5/1872 (0.3%)
2	E	0.34	0/712	1.06	3/1097 (0.3%)
3	F	0.37	0/712	1.11	4/1096 (0.4%)
All	All	0.52	0/7000	0.97	13/9712 (0.1%)

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	29	ARG	N-CA-C	9.21	125.15	113.55
1	D	167	GLU	N-CA-C	-7.81	99.26	110.59
3	F	4	DA	C4'-C3'-O3'	7.33	121.00	110.00
2	E	11	DG	O4'-C1'-N9	6.58	118.26	108.40
2	E	30	DT	C4'-C3'-O3'	5.97	118.95	110.00
3	F	25	DT	C2'-C3'-O3'	5.77	120.16	111.50
3	F	4	DA	N9-C1'-C2'	-5.54	105.19	113.50
1	D	12	GLU	N-CA-C	-5.48	103.67	110.41
1	B	142	GLU	N-CA-C	-5.16	101.99	109.31
3	F	21	DC	C2'-C3'-O3'	5.12	119.17	111.50
1	D	54	TRP	CA-C-N	5.11	125.19	119.87
1	D	54	TRP	C-N-CA	5.11	125.19	119.87
2	E	6	DA	C4'-C3'-O3'	5.09	117.64	110.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1369	0	1386	28	0
1	B	1391	0	1412	21	0
1	C	1386	0	1410	17	0
1	D	1376	0	1402	23	0
2	E	636	0	351	14	0
3	F	635	0	349	14	0
4	A	12	0	0	0	0
4	B	10	0	0	0	0
4	C	12	0	0	0	0
4	D	11	0	0	1	0
4	E	9	0	0	1	0
4	F	11	0	0	1	0
All	All	6858	0	6310	110	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 8.

All (110) 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:76:MSE:HE3	1:A:77:PRO:HD2	1.60	0.81
1:C:58:HIS:HB3	1:C:119:THR:HG21	1.73	0.71
3:F:23:DC:H2'	3:F:24:DG:C8	2.27	0.70
1:C:28:GLU:HG2	1:C:29:ARG:HG3	1.74	0.69
3:F:11:DG:N3	4:F:101:HOH:O	2.27	0.67
1:C:136:ARG:HH21	1:C:150:GLN:HG3	1.59	0.66
1:D:116:HIS:ND1	4:D:211:HOH:O	2.29	0.66
3:F:13:DC:H2'	3:F:14:DG:C8	2.32	0.64
1:A:109:LEU:HD12	1:A:123:LEU:HD13	1.81	0.63
1:C:23:ARG:NH2	1:C:67:GLU:OE1	2.32	0.62
1:A:34:LEU:HB3	1:A:57:ARG:HD2	1.82	0.61
1:A:99:THR:C	1:A:169:ARG:HH12	2.08	0.61
1:C:98:THR:O	1:C:169:ARG:HD2	2.01	0.60
3:F:29:DT:H2''	3:F:30:DG:C8	2.36	0.60
1:B:12:GLU:OE2	1:B:15:ARG:NH1	2.34	0.60

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:24:DG:H2''	3:F:25:DT:H5'	1.84	0.60
1:C:72:ILE:HG22	1:C:73:LEU:HD23	1.84	0.59
1:A:110:MSE:HE3	1:A:165:LEU:HD11	1.84	0.59
3:F:16:DA:H2''	3:F:17:DC:H5''	1.84	0.59
2:E:9:DG:H2''	2:E:10:DA:O5'	2.02	0.58
2:E:14:DG:H2''	2:E:15:DG:H5''	1.84	0.58
1:A:154:ASP:CG	1:D:175:ARG:HH21	2.13	0.57
1:A:155:ALA:HA	1:D:159:ALA:HB2	1.87	0.56
1:D:116:HIS:CE1	1:D:118:ASP:HB2	2.41	0.56
1:A:99:THR:O	1:A:169:ARG:NH1	2.33	0.55
1:A:36:ILE:HG21	1:A:51:TYR:CE2	2.42	0.55
1:D:167:GLU:OE2	1:D:175:ARG:NH1	2.40	0.55
1:D:174:GLN:NE2	1:D:178:THR:OG1	2.40	0.54
1:D:10:ARG:HG3	1:D:53:TRP:CE2	2.42	0.54
1:A:65:LEU:HD13	1:A:109:LEU:HD13	1.88	0.54
1:D:10:ARG:HG3	1:D:53:TRP:CZ2	2.42	0.54
1:D:66:LEU:HD11	1:D:122:ARG:HG2	1.88	0.54
1:B:23:ARG:NH1	1:B:68:ASP:OD1	2.42	0.53
1:C:136:ARG:HE	1:C:150:GLN:HA	1.72	0.53
1:A:22:THR:HG23	1:A:34:LEU:HD21	1.92	0.52
1:B:10:ARG:HD3	1:B:53:TRP:CE2	2.45	0.51
1:C:145:ASP:OD2	1:C:148:HIS:ND1	2.43	0.51
1:D:15:ARG:HB3	1:D:15:ARG:HH11	1.76	0.51
1:D:36:ILE:HD11	1:D:57:ARG:HG3	1.93	0.50
3:F:27:DT:H2''	3:F:28:DA:C8	2.46	0.50
1:C:65:LEU:HD11	1:C:108:THR:HG22	1.93	0.49
1:D:76:MSE:SE	1:D:77:PRO:HD2	2.62	0.49
1:A:72:ILE:HD12	1:A:97:LEU:HD23	1.95	0.49
2:E:6:DA:H2'	2:E:7:DA:C8	2.47	0.49
2:E:30:DT:H2''	2:E:31:DG:C8	2.48	0.49
1:B:110:MSE:HE3	1:B:165:LEU:HD11	1.94	0.49
1:D:35:SER:O	1:D:39:VAL:HG23	2.13	0.49
1:D:34:LEU:O	1:D:57:ARG:NH1	2.46	0.49
1:D:94:ALA:O	1:D:98:THR:OG1	2.24	0.49
1:C:27:LEU:HD21	1:C:101:ARG:HD3	1.94	0.49
1:B:49:THR:HG21	2:E:8:DC:H3'	1.96	0.48
1:A:125:GLU:O	1:A:129:ARG:NH1	2.47	0.48
1:B:49:THR:HG22	1:B:52:ARG:HH12	1.79	0.48
1:A:127:PHE:O	1:A:130:PRO:HD2	2.14	0.48
2:E:1:DT:H2'	2:E:1:DT:OP2	2.14	0.47
1:A:160:VAL:O	1:A:164:VAL:HG23	2.15	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:18:DC:H2'	3:F:19:DG:C8	2.50	0.47
1:A:136:ARG:NH1	1:A:150:GLN:OE1	2.45	0.47
1:B:9:ARG:O	2:E:7:DA:H5''	2.15	0.46
1:D:28:GLU:HG3	1:D:29:ARG:HG3	1.96	0.46
1:B:26:LEU:HD11	1:B:61:VAL:HG13	1.97	0.46
2:E:6:DA:H2''	2:E:7:DA:O5'	2.15	0.46
3:F:26:DT:H2''	3:F:27:DT:H5''	1.98	0.46
2:E:15:DG:H2'	2:E:16:DT:H72	1.98	0.45
1:A:76:MSE:CE	1:A:89:TRP:HB2	2.47	0.45
3:F:28:DA:H1'	3:F:29:DT:C6	2.52	0.45
1:B:45:VAL:HG22	2:E:9:DG:OP2	2.17	0.44
1:A:37:GLU:HG3	4:E:107:HOH:O	2.18	0.44
1:A:76:MSE:HE1	1:A:89:TRP:HB2	1.99	0.44
1:B:151:ALA:O	1:B:154:ASP:HB2	2.17	0.44
1:A:127:PHE:C	1:A:130:PRO:HD2	2.43	0.44
1:C:10:ARG:HB2	1:C:53:TRP:CZ2	2.53	0.44
1:C:101:ARG:O	1:C:105:MSE:HB2	2.17	0.44
1:A:56:SER:OG	1:A:57:ARG:N	2.50	0.44
1:B:116:HIS:HB3	1:B:119:THR:HB	1.99	0.44
1:D:65:LEU:HD11	1:D:108:THR:HG22	1.99	0.44
1:B:62:ALA:HB2	1:B:119:THR:HG23	1.99	0.44
1:C:100:ARG:HG2	1:C:169:ARG:NH2	2.32	0.44
3:F:24:DG:C8	3:F:25:DT:H72	2.53	0.43
1:C:132:ILE:HG13	1:C:157:LEU:HD22	2.01	0.43
1:D:106:LEU:HD13	1:D:165:LEU:HG	2.00	0.43
1:B:45:VAL:HG13	1:B:46:GLY:O	2.19	0.43
1:C:79:THR:HG22	1:C:80:ASP:H	1.83	0.43
2:E:18:DC:H2'	2:E:19:DG:C8	2.53	0.43
1:B:70:ASP:OD2	1:B:70:ASP:N	2.52	0.43
1:A:65:LEU:HD11	1:A:108:THR:HG22	2.01	0.43
1:A:159:ALA:HB2	1:D:155:ALA:HA	2.01	0.42
2:E:6:DA:H2''	2:E:7:DA:C5'	2.49	0.42
1:B:187:LEU:HD23	1:B:187:LEU:HA	1.94	0.42
1:C:87:ALA:O	1:C:177:GLU:HG2	2.20	0.42
1:C:86:LEU:HD21	1:C:138:ARG:HG2	2.02	0.42
1:A:57:ARG:HE	1:A:57:ARG:HB3	1.67	0.42
1:B:13:LYS:HE2	1:B:13:LYS:HB3	1.84	0.42
2:E:3:DA:H4'	2:E:4:DT:OP1	2.19	0.42
1:A:123:LEU:HD23	1:A:123:LEU:HA	1.87	0.42
1:B:83:THR:OG1	1:B:184:VAL:HG11	2.19	0.42
1:B:116:HIS:ND1	1:B:119:THR:OG1	2.33	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:30:DG:OP2	3:F:30:DG:H2'	2.20	0.41
1:B:135:VAL:HG11	1:B:153:ALA:HA	2.02	0.41
1:A:154:ASP:OD1	1:D:175:ARG:NH2	2.52	0.41
1:B:106:LEU:HD13	1:B:165:LEU:HG	2.02	0.41
1:B:123:LEU:HD23	1:B:123:LEU:HA	1.82	0.41
1:D:138:ARG:O	1:D:141:ASP:HB2	2.21	0.41
2:E:8:DC:H2''	2:E:9:DG:C8	2.56	0.41
3:F:31:DA:H5'	3:F:31:DA:C8	2.56	0.41
1:A:22:THR:OG1	1:A:60:LEU:HD21	2.21	0.40
1:D:127:PHE:C	1:D:130:PRO:HD2	2.46	0.40
1:A:63:ASP:OD1	1:A:122:ARG:NH2	2.48	0.40
1:D:11:SER:OG	1:D:12:GLU:O	2.33	0.40
3:F:25:DT:H2''	3:F:26:DT:O5'	2.22	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	178/196 (91%)	171 (96%)	5 (3%)	2 (1%)	11	22
1	B	180/196 (92%)	173 (96%)	7 (4%)	0	100	100
1	C	180/196 (92%)	170 (94%)	9 (5%)	1 (1%)	21	38
1	D	178/196 (91%)	168 (94%)	9 (5%)	1 (1%)	21	38
All	All	716/784 (91%)	682 (95%)	30 (4%)	4 (1%)	21	38

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	100	ARG
1	C	173	ARG

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	99	THR
1	D	143	ASP

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	137/142 (96%)	127 (93%)	10 (7%)	13	27
1	B	139/142 (98%)	125 (90%)	14 (10%)	7	15
1	C	138/142 (97%)	126 (91%)	12 (9%)	9	21
1	D	137/142 (96%)	116 (85%)	21 (15%)	3	5
All	All	551/568 (97%)	494 (90%)	57 (10%)	7	14

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

Mol	Chain	Res	Type
1	A	13	LYS
1	A	15	ARG
1	A	32	ASP
1	A	45	VAL
1	A	73	LEU
1	A	82	VAL
1	A	101	ARG
1	A	117	GLU
1	A	140	ARG
1	A	169	ARG
1	B	9	ARG
1	B	10	ARG
1	B	27	LEU
1	B	34	LEU
1	B	45	VAL
1	B	70	ASP
1	B	98	THR
1	B	106	LEU
1	B	118[A]	ASP

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	118[B]	ASP
1	B	136	ARG
1	B	139	LEU
1	B	142	GLU
1	B	167	GLU
1	C	9	ARG
1	C	15	ARG
1	C	28	GLU
1	C	36	ILE
1	C	72	ILE
1	C	78	LYS
1	C	80	ASP
1	C	119	THR
1	C	125	GLU
1	C	128	SER
1	C	134	SER
1	C	169	ARG
1	D	9	ARG
1	D	10	ARG
1	D	13	LYS
1	D	14	SER
1	D	15	ARG
1	D	42	LYS
1	D	45	VAL
1	D	57	ARG
1	D	61	VAL
1	D	101	ARG
1	D	106	LEU
1	D	125	GLU
1	D	133	GLU
1	D	134	SER
1	D	136	ARG
1	D	143	ASP
1	D	144	ILE
1	D	156	LEU
1	D	166	SER
1	D	173	ARG
1	D	188	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	148	HIS
1	D	116	HIS
1	D	148	HIS

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 oligosaccharides 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

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, 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	176/196 (89%)	0.29	6 (3%)	48	43	32, 56, 89, 110	1 (0%)
1	B	178/196 (90%)	0.28	5 (2%)	55	50	24, 59, 83, 108	1 (0%)
1	C	179/196 (91%)	0.29	3 (1%)	69	65	34, 57, 93, 109	0
1	D	177/196 (90%)	0.28	4 (2%)	61	57	37, 56, 86, 108	0
2	E	31/31 (100%)	0.07	1 (3%)	50	46	39, 59, 109, 128	0
3	F	31/31 (100%)	0.02	0	100	100	40, 58, 104, 110	0
All	All	772/846 (91%)	0.26	19 (2%)	58	54	24, 58, 91, 128	2 (0%)

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	101	ARG	4.6
1	D	72	ILE	4.0
1	D	9	ARG	3.6
1	B	189	PRO	3.2
1	A	72	ILE	3.0
1	A	11	SER	3.0
1	A	189	PRO	2.6
1	A	99	THR	2.4
1	B	185	ALA	2.4
2	E	1	DT	2.4
1	B	9	ARG	2.4
1	B	188	ARG	2.3
1	D	132	ILE	2.3
1	B	118[A]	ASP	2.3
1	A	100	ARG	2.2
1	D	144	ILE	2.1
1	C	73	LEU	2.1
1	C	74	ALA	2.0
1	A	168	GLY	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 oligosaccharides 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.