



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:34 PM GMT

PDB ID : 4IR1
Title : Polymerase-DNA Complex
Authors : Sharma, A.; Nair, D.T.
Deposited on : 2013-01-14
Resolution : 2.38 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 ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) 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

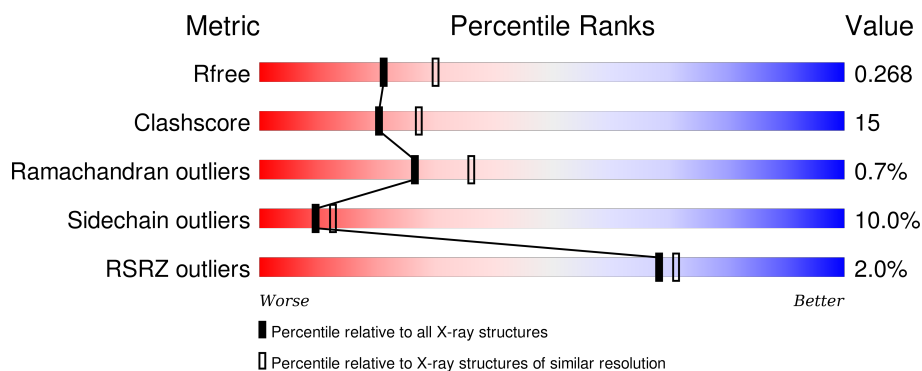
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.38 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}	91344	4019 (2.40-2.36)
Clashscore	102246	4595 (2.40-2.36)
Ramachandran outliers	100387	4520 (2.40-2.36)
Sidechain outliers	100360	4522 (2.40-2.36)
RSRZ outliers	91569	4034 (2.40-2.36)

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 $\leq 5\%$. 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	352	<div> <div>4%</div> <div>61%</div> <div>32%</div> <div>• •</div> </div>
1	F	352	<div> <div>72%</div> <div>22%</div> <div>• •</div> </div>
2	B	18	<div> <div>44%</div> <div>56%</div> </div>
2	G	18	<div> <div>6%</div> <div>67%</div> <div>33%</div> </div>
3	H	14	<div> <div>57%</div> <div>43%</div> </div>

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Mol	Chain	Length	Quality of chain
4	C	17	<div><div></div><div>59%</div><div>41%</div></div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
6	MG	F	903	-	-	-	X

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 7080 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 DNA polymerase IV.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	F	342	Total	C	N	O	S	0	0	0
			2687	1695	494	484	14			
1	A	342	Total	C	N	O	S	0	0	0
			2687	1695	494	484	14			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
F	0	GLY	-	EXPRESSION TAG	UNP Q47155
F	1	SER	-	EXPRESSION TAG	UNP Q47155
A	0	GLY	-	EXPRESSION TAG	UNP Q47155
A	1	SER	-	EXPRESSION TAG	UNP Q47155

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	G	18	Total	C	N	O	P	0	0	0
			364	174	66	107	17			
2	B	18	Total	C	N	O	P	0	0	0
			364	174	66	107	17			

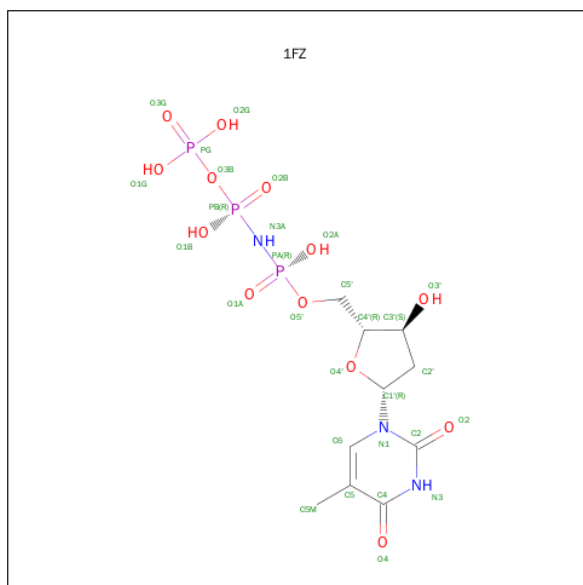
- Molecule 3 is a DNA chain called DNA (5'-D(*GP*GP*GP*TP*CP*CP*TP*AP*GP*GP*AP*CP*CP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	H	14	Total	C	N	O	P	0	0	0
			284	135	54	82	13			

- Molecule 4 is a DNA chain called DNA (5'-D(*CP*TP*AP*GP*GP*GP*TP*CP*CP*TP*AP*GP*GP*AP*CP*CP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	C	17	Total	C	N	O	P	0	0	0
			344	164	64	100	16			

- Molecule 5 is 5'-O-[(R)-HYDROXY{[(R)-HYDROXY(PHOSPHONOOXY)PHOSPHORYL]AMINO}PHOSPHORYL]THYMIDINE (three-letter code: 1FZ) (formula: C₁₀H₁₈N₃O₁₃P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	F	1	Total	C	N	O	P	0	0
			29	10	3	13	3		
5	A	1	Total	C	N	O	P	0	0
			29	10	3	13	3		

- Molecule 6 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	2	Total	Mg	0	0
			2	2		
6	F	2	Total	Mg	0	0
			2	2		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	F	138	Total	O	0	0
			138	138		

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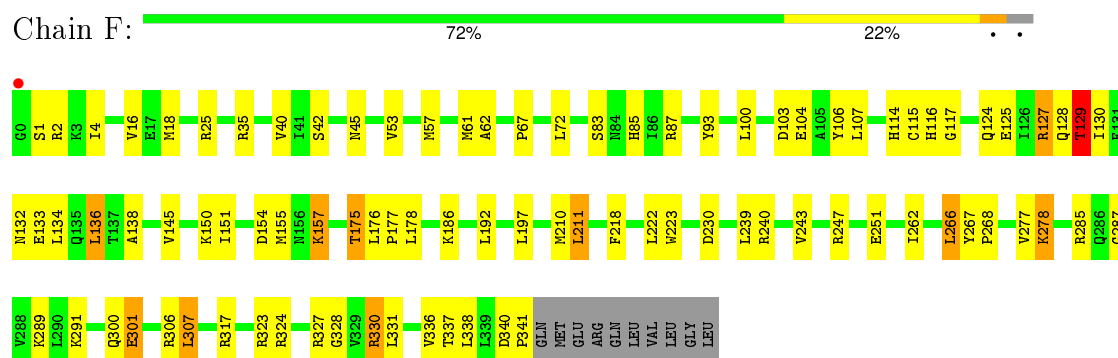
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	54	Total 54	O 54	0	0
7	G	35	Total 35	O 35	0	0
7	H	27	Total 27	O 27	0	0
7	B	20	Total 20	O 20	0	0
7	C	14	Total 14	O 14	0	0

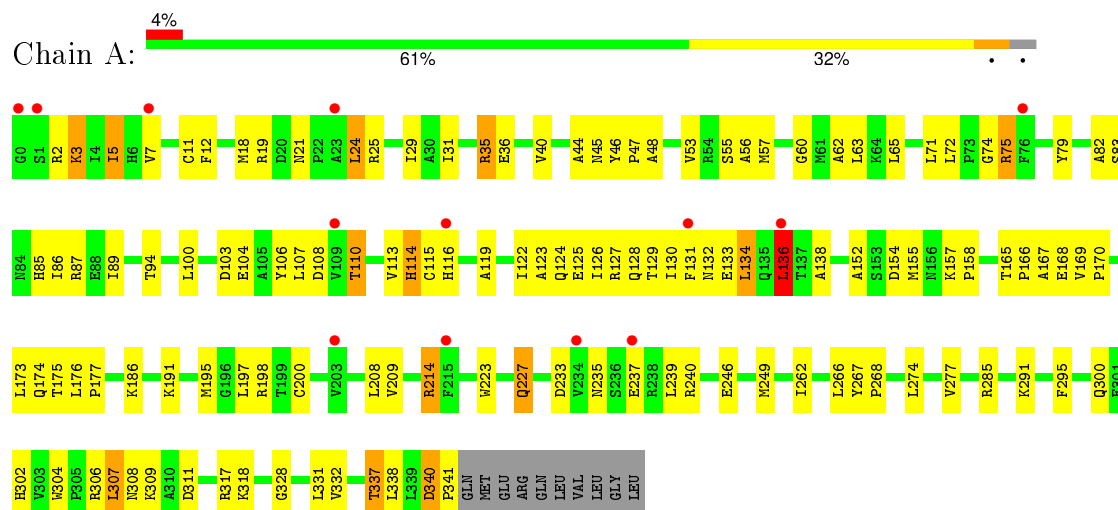
3 Residue-property plots

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.

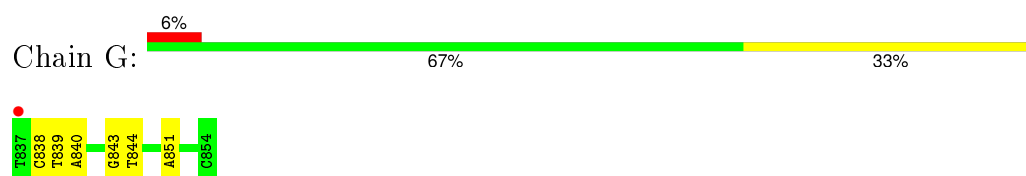
• Molecule 1: DNA polymerase IV



• Molecule 1: DNA polymerase IV



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



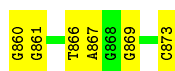
- Molecule 2: DNA (5'-D(*TP*CP*TP*AP*GP*GP*GP*TP*CP*CP*TP*AP*GP*GP*AP*CP*CP*C)-3')

Chain B:  44% 56%



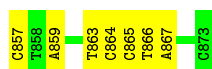
- Molecule 3: DNA (5'-D(*GP*GP*GP*TP*CP*CP*TP*AP*GP*GP*AP*CP*CP*C)-3')

Chain H:  57% 43%



- Molecule 4: DNA (5'-D(*CP*TP*AP*GP*GP*GP*TP*CP*CP*TP*AP*GP*GP*AP*CP*CP*C)-3')

Chain C:  59% 41%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	85.75Å 57.25Å 111.30Å 90.00° 90.44° 90.00°	Depositor
Resolution (Å)	47.66 – 2.38 47.61 – 2.38	Depositor EDS
% Data completeness (in resolution range)	99.8 (47.66-2.38) 99.9 (47.61-2.38)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.11 (at 2.37Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.214 , 0.269 0.214 , 0.268	Depositor DCC
R_{free} test set	2199 reflections (5.30%)	DCC
Wilson B-factor (Å ²)	49.0	Xtriage
Anisotropy	0.209	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 31.6	EDS
Estimated twinning fraction	0.023 for h,-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 43673 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7080	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.10% 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, 0.375 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: MG, 1FZ

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	0.55	0/2738	0.74	1/3699 (0.0%)
1	F	0.63	1/2738 (0.0%)	0.83	3/3699 (0.1%)
2	B	0.43	0/407	0.90	1/626 (0.2%)
2	G	0.53	0/407	0.75	0/626
3	H	0.57	0/318	0.74	0/489
4	C	0.46	0/385	0.94	2/592 (0.3%)
All	All	0.57	1/6993 (0.0%)	0.80	7/9731 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	301	GLU	CD-OE1	5.80	1.32	1.25

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	859	DA	O5'-P-OP1	-9.30	97.33	105.70
2	B	839	DT	O5'-P-OP1	-7.62	98.84	105.70
4	C	859	DA	O5'-P-OP2	6.66	118.70	110.70
1	F	127	ARG	NE-CZ-NH2	-6.09	117.26	120.30
1	F	127	ARG	NE-CZ-NH1	5.80	123.20	120.30
1	A	136	LEU	CA-CB-CG	5.17	127.19	115.30
1	F	129	THR	N-CA-CB	5.06	119.92	110.30

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	2687	0	2739	97	0
1	F	2687	0	2739	70	0
2	B	364	0	204	8	0
2	G	364	0	204	7	0
3	H	284	0	158	6	0
4	C	344	0	192	11	0
5	A	29	0	14	3	0
5	F	29	0	14	0	0
6	A	2	0	0	0	0
6	F	2	0	0	0	0
7	A	54	0	0	4	0
7	B	20	0	0	0	0
7	C	14	0	0	1	0
7	F	138	0	0	12	0
7	G	35	0	0	3	0
7	H	27	0	0	1	0
All	All	7080	0	6264	191	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 15.

All (191) 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:340:ASP:HB2	1:A:341:PRO:HD3	1.29	1.08
1:F:18:MET:HE3	1:F:45:ASN:HD22	1.20	1.03
1:F:18:MET:CE	1:F:45:ASN:HD22	1.70	1.02
1:A:317:ARG:HD3	7:A:515:HOH:O	1.57	1.02
1:F:285:ARG:NH1	3:H:866:DT:OP2	1.97	0.96
1:F:57:MET:HE3	1:F:62:ALA:HA	1.48	0.95
1:A:285:ARG:HB2	1:A:337:THR:HG23	1.52	0.90
1:A:285:ARG:HB2	1:A:337:THR:CG2	2.02	0.90
4:C:864:DC:H2"	4:C:865:DC:H5"	1.57	0.87
1:A:18:MET:HE3	1:A:45:ASN:HD22	1.37	0.86
1:F:57:MET:HG3	1:F:61:MET:HE2	1.58	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:125:GLU:O	1:F:128:GLN:O	1.94	0.85
1:F:57:MET:HE3	1:F:62:ALA:CA	2.05	0.85
1:A:18:MET:CE	1:A:45:ASN:HD22	1.91	0.83
1:A:340:ASP:CB	1:A:341:PRO:HD3	2.10	0.82
1:F:128:GLN:HA	7:F:501:HOH:O	1.82	0.78
1:F:150:LYS:NZ	3:H:873:DC:OP1	2.17	0.78
1:F:57:MET:CE	1:F:62:ALA:HA	2.15	0.77
1:F:129:THR:O	1:F:132:ASN:O	2.03	0.76
1:A:130:ILE:O	1:A:134:LEU:HB2	1.88	0.73
1:F:18:MET:HE2	1:F:25:ARG:O	1.88	0.72
1:A:173:LEU:O	1:A:200:CYS:HB2	1.89	0.72
1:A:132:ASN:O	1:A:133:GLU:HB2	1.89	0.72
1:A:57:MET:HE3	1:A:62:ALA:HA	1.71	0.71
1:A:57:MET:HE3	1:A:62:ALA:CA	2.21	0.70
1:A:340:ASP:HB2	1:A:341:PRO:CD	2.17	0.70
1:F:18:MET:HE3	1:F:45:ASN:ND2	2.02	0.70
1:A:128:GLN:O	1:A:129:THR:HB	1.91	0.68
1:A:57:MET:CE	1:A:62:ALA:HA	2.24	0.68
1:F:150:LYS:HE3	7:F:583:HOH:O	1.92	0.68
1:A:166:PRO:O	1:A:169:VAL:HG12	1.95	0.67
1:F:129:THR:N	7:F:501:HOH:O	1.90	0.66
1:F:157:LYS:HD3	7:F:579:HOH:O	1.94	0.66
1:F:300:GLN:HB2	7:F:553:HOH:O	1.95	0.66
1:F:18:MET:HE1	1:F:45:ASN:HD22	1.59	0.65
4:C:857:DC:H5'	4:C:857:DC:H6	1.61	0.65
1:A:5:ILE:HG23	1:A:107:LEU:HB2	1.77	0.65
1:F:287:GLY:HA3	1:F:301:GLU:HG2	1.80	0.63
1:A:331:LEU:C	1:A:331:LEU:HD23	2.19	0.62
1:A:308:ASN:ND2	1:A:311:ASP:H	1.97	0.62
1:A:308:ASN:HD22	1:A:311:ASP:H	1.48	0.62
1:A:157:LYS:HE2	5:A:875:1FZ:PG	2.40	0.61
1:A:19:ARG:HD3	1:A:136:LEU:HD13	1.83	0.61
1:A:291:LYS:HB3	1:A:331:LEU:HB3	1.83	0.60
1:A:195:MET:HE1	1:A:214:ARG:HG2	1.81	0.60
1:A:29:ILE:HG12	1:A:45:ASN:ND2	2.16	0.60
1:F:104:GLU:OE2	1:F:150:LYS:NZ	2.35	0.59
1:F:340:ASP:O	1:F:341:PRO:C	2.40	0.59
1:F:128:GLN:CA	7:F:501:HOH:O	2.47	0.58
1:F:40:VAL:HG13	2:G:840:DA:H5'	1.86	0.58
1:A:3:LYS:H	1:A:110:THR:HG22	1.68	0.58
1:F:145:VAL:HB	1:F:230:ASP:HB3	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:127:ARG:HD2	1:F:138:ALA:O	2.03	0.57
1:A:82:ALA:O	1:A:86:ILE:HG23	2.04	0.56
1:A:195:MET:CE	1:A:214:ARG:HG2	2.35	0.56
1:A:304:TRP:HD1	1:A:306:ARG:O	1.89	0.56
1:F:240:ARG:NH1	2:G:843:DG:OP1	2.39	0.55
1:A:249:MET:SD	1:A:262:ILE:HD13	2.46	0.55
1:F:128:GLN:O	1:F:129:THR:OG1	2.24	0.55
1:A:18:MET:HE3	1:A:45:ASN:ND2	2.15	0.55
1:A:267:TYR:OH	1:A:307:LEU:HD13	2.06	0.54
1:F:124:GLN:OE1	1:F:127:ARG:NH2	2.40	0.54
1:A:169:VAL:HG13	1:A:170:PRO:HD3	1.90	0.54
1:F:278:LYS:HE2	7:F:572:HOH:O	2.06	0.54
1:F:18:MET:CE	1:F:45:ASN:ND2	2.55	0.54
1:F:40:VAL:HG21	2:G:840:DA:C5	2.43	0.54
4:C:865:DC:H2'	4:C:866:DT:C6	2.43	0.53
1:A:85:HIS:O	1:A:89:ILE:HG13	2.08	0.53
1:F:285:ARG:HB2	1:F:337:THR:HB	1.91	0.53
4:C:865:DC:C2'	4:C:866:DT:C6	2.91	0.53
1:A:74:GLY:HA2	1:A:79:TYR:OH	2.09	0.53
1:A:227:GLN:NE2	7:A:526:HOH:O	2.40	0.53
1:A:75:ARG:HB2	1:A:75:ARG:HH11	1.74	0.53
1:F:85:HIS:CD2	1:F:134:LEU:HD11	2.44	0.52
1:A:167:ALA:O	1:A:170:PRO:HD2	2.10	0.52
1:A:291:LYS:HE2	1:A:295:PHE:O	2.10	0.52
1:F:85:HIS:CG	1:F:134:LEU:HD11	2.45	0.52
4:C:857:DC:C6	4:C:857:DC:H5'	2.44	0.52
4:C:867:DA:N3	7:C:902:HOH:O	2.34	0.51
1:A:55:SER:O	1:A:56:ALA:HB3	2.10	0.51
1:F:93:TYR:OH	1:F:125:GLU:HG2	2.10	0.51
1:A:94:THR:CG2	1:A:107:LEU:HD23	2.40	0.51
1:A:208:LEU:HD23	1:A:223:TRP:CD1	2.46	0.51
3:H:860:DG:N7	7:H:915:HOH:O	2.33	0.51
1:A:21:ASN:HB3	1:A:24:LEU:HD22	1.93	0.51
1:A:46:TYR:CE1	1:A:158:PRO:HG3	2.46	0.51
1:F:18:MET:HE2	1:F:25:ARG:HA	1.93	0.50
1:A:53:VAL:HA	1:A:57:MET:HE2	1.93	0.50
1:A:246:GLU:HG2	2:B:842:DG:OP2	2.12	0.50
1:A:57:MET:HE3	1:A:62:ALA:CB	2.42	0.49
1:F:240:ARG:HH21	1:F:277:VAL:HG21	1.77	0.49
1:A:57:MET:HE3	1:A:62:ALA:HB2	1.94	0.49
1:F:328:GLY:HA3	7:F:507:HOH:O	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:C:865:DC:H2'	4:C:866:DT:C7	2.43	0.49
1:A:31:ILE:HD12	1:A:31:ILE:N	2.28	0.49
1:A:123:ALA:O	1:A:127:ARG:HG3	2.14	0.48
1:F:330:ARG:HD3	7:G:917:HOH:O	2.14	0.48
1:A:152:ALA:HA	1:A:155:MET:HG3	1.96	0.48
1:A:53:VAL:HA	1:A:57:MET:CE	2.43	0.48
1:A:128:GLN:HA	7:A:531:HOH:O	2.14	0.48
1:F:178:LEU:HD13	1:F:192:LEU:HD13	1.96	0.47
1:F:132:ASN:O	1:F:133:GLU:HB2	2.14	0.47
1:F:324:ARG:HD2	1:F:327:ARG:O	2.14	0.47
1:A:128:GLN:O	1:A:129:THR:CB	2.62	0.47
1:F:2:ARG:NH2	1:F:4:ILE:HD11	2.30	0.47
1:A:40:VAL:HG13	2:B:840:DA:H5'	1.97	0.46
1:F:291:LYS:HB3	1:F:331:LEU:HB3	1.98	0.46
1:A:89:ILE:HD13	1:A:129:THR:HG22	1.97	0.46
1:A:317:ARG:NH1	3:H:869:DG:OP1	2.49	0.46
1:A:122:ILE:HD12	1:A:123:ALA:N	2.31	0.46
1:A:127:ARG:HD2	1:A:138:ALA:O	2.15	0.46
1:F:106:TYR:O	1:F:107:LEU:HD12	2.15	0.46
1:A:267:TYR:CG	1:A:309:LYS:HG3	2.51	0.46
1:A:274:LEU:HD22	1:A:307:LEU:HG	1.98	0.45
1:A:302:HIS:NE2	1:A:311:ASP:OD2	2.41	0.45
1:A:94:THR:HG22	1:A:107:LEU:HD23	1.98	0.45
1:A:122:ILE:O	1:A:126:ILE:HG13	2.16	0.45
1:F:154:ASP:HA	1:F:157:LYS:HD2	1.99	0.45
1:F:218:PHE:CE2	1:F:222:LEU:HD22	2.52	0.45
1:F:67:PRO:HD2	7:F:544:HOH:O	2.16	0.45
1:F:57:MET:HG3	1:F:61:MET:CE	2.39	0.45
2:B:851:DA:H2''	2:B:852:DC:H5'	1.98	0.45
2:B:851:DA:H1'	2:B:852:DC:H5'	1.99	0.45
1:A:166:PRO:O	1:A:169:VAL:CG1	2.62	0.45
2:G:839:DT:H1'	7:G:925:HOH:O	2.16	0.44
1:A:328:GLY:HA3	7:A:507:HOH:O	2.17	0.44
1:F:57:MET:SD	1:F:61:MET:HE3	2.57	0.44
1:A:157:LYS:HE3	1:A:157:LYS:HB3	1.82	0.44
1:F:289:LYS:NZ	7:F:547:HOH:O	2.50	0.44
1:A:44:ALA:HB1	1:A:48:ALA:HB3	1.99	0.44
1:A:175:THR:O	1:A:177:PRO:HD3	2.17	0.44
1:A:267:TYR:HB3	1:A:268:PRO:HD3	1.99	0.44
1:F:175:THR:O	1:F:177:PRO:HD3	2.18	0.44
1:A:300:GLN:HE22	1:A:318:LYS:NZ	2.15	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:C:865:DC:H2'	4:C:866:DT:C5	2.53	0.44
1:A:71:LEU:C	1:A:72:LEU:HD12	2.39	0.44
2:B:848:DA:H2''	2:B:849:DG:C8	2.53	0.44
1:A:108:ASP:OD2	1:A:110:THR:HG23	2.17	0.43
1:A:267:TYR:CZ	1:A:307:LEU:HD13	2.53	0.43
1:A:100:LEU:HD11	1:A:106:TYR:CE1	2.53	0.43
1:A:2:ARG:HG3	1:A:2:ARG:HH11	1.83	0.43
1:F:53:VAL:HG13	1:F:57:MET:HE2	2.00	0.43
1:A:5:ILE:HD11	1:A:7:VAL:HG22	1.99	0.43
1:A:304:TRP:CD1	1:A:306:ARG:O	2.70	0.43
1:F:267:TYR:HB3	1:F:268:PRO:HD3	1.99	0.43
1:A:35:ARG:HD2	1:A:60:GLY:HA2	1.99	0.43
1:A:11:CYS:O	1:A:12:PHE:C	2.57	0.43
2:B:844:DT:H2''	2:B:845:DC:C5'	2.48	0.43
1:A:18:MET:HG2	1:A:25:ARG:HA	2.00	0.43
1:A:7:VAL:O	1:A:104:GLU:HA	2.19	0.43
2:G:843:DG:H2''	2:G:844:DT:H5'	2.01	0.43
2:G:838:DC:H4'	2:G:839:DT:O5'	2.18	0.43
1:A:235:ASN:C	1:A:237:GLU:H	2.22	0.43
4:C:865:DC:H2''	4:C:866:DT:C6	2.53	0.43
1:A:18:MET:CE	1:A:45:ASN:ND2	2.70	0.43
1:A:114:HIS:HB3	1:A:115:CYS:H	1.71	0.43
1:F:211:LEU:HD23	1:F:223:TRP:HB2	2.00	0.43
1:A:18:MET:HE3	1:A:45:ASN:CB	2.49	0.42
4:C:863:DT:H2''	4:C:864:DC:C6	2.54	0.42
1:F:300:GLN:HA	3:H:867:DA:OP2	2.18	0.42
1:F:57:MET:HE3	1:F:62:ALA:CB	2.48	0.42
3:H:860:DG:H2''	3:H:861:DG:OP2	2.20	0.42
1:F:18:MET:HE2	1:F:25:ARG:C	2.39	0.42
1:A:165:THR:OG1	1:A:168:GLU:HG3	2.19	0.42
1:F:306:ARG:O	1:F:307:LEU:C	2.58	0.42
1:F:115:CYS:C	1:F:117:GLY:N	2.72	0.42
1:F:130:ILE:O	1:F:134:LEU:HB2	2.20	0.42
1:F:83:SER:O	1:F:87:ARG:HG3	2.19	0.42
1:A:308:ASN:ND2	1:A:311:ASP:N	2.66	0.42
1:F:157:LYS:NZ	7:F:609:HOH:O	2.53	0.41
1:A:63:LEU:HA	1:A:63:LEU:HD23	1.85	0.41
1:F:129:THR:O	1:F:133:GLU:HB2	2.20	0.41
1:A:119:ALA:O	1:A:122:ILE:HG13	2.19	0.41
1:F:262:ILE:HG22	1:F:266:LEU:HD22	2.01	0.41
1:A:125:GLU:O	1:A:128:GLN:O	2.38	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:323:ARG:HD3	7:F:635:HOH:O	2.21	0.41
1:F:100:LEU:HD11	1:F:106:TYR:CE1	2.55	0.41
1:A:46:TYR:N	1:A:47:PRO:CD	2.83	0.41
1:A:308:ASN:HD22	1:A:311:ASP:N	2.16	0.41
4:C:865:DC:H2'	4:C:866:DT:H72	2.02	0.41
1:A:169:VAL:O	1:A:173:LEU:HB2	2.20	0.41
1:F:151:ILE:O	1:F:155:MET:HG3	2.21	0.41
1:F:16:VAL:HG22	1:F:136:LEU:HD21	2.03	0.41
2:B:838:DC:H2'	2:B:838:DC:H6	1.62	0.41
1:A:83:SER:O	1:A:87:ARG:HD3	2.21	0.41
1:A:45:ASN:CG	1:A:47:PRO:HD2	2.41	0.40
2:B:844:DT:H2''	2:B:845:DC:H5'	2.02	0.40
2:G:851:DA:H2'	7:G:921:HOH:O	2.21	0.40
1:F:72:LEU:N	1:F:72:LEU:HD12	2.37	0.40
1:A:11:CYS:HA	5:A:875:1FZ:O1B	2.22	0.40
5:A:875:1FZ:H1	5:A:875:1FZ:O5'	2.21	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	340/352 (97%)	314 (92%)	24 (7%)	2 (1%)	30	40
1	F	340/352 (97%)	321 (94%)	16 (5%)	3 (1%)	21	28
All	All	680/704 (97%)	635 (93%)	40 (6%)	5 (1%)	26	36

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	114	HIS
1	F	116	HIS

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Mol	Chain	Res	Type
1	F	129	THR
1	A	36	GLU
1	A	340	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	286/297 (96%)	252 (88%)	34 (12%)	6	7
1	F	286/297 (96%)	263 (92%)	23 (8%)	15	21
All	All	572/594 (96%)	515 (90%)	57 (10%)	9	12

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

Mol	Chain	Res	Type
1	F	1	SER
1	F	35	ARG
1	F	42	SER
1	F	103	ASP
1	F	136	LEU
1	F	157	LYS
1	F	175	THR
1	F	176	LEU
1	F	186	LYS
1	F	197	LEU
1	F	210	MET
1	F	211	LEU
1	F	239	LEU
1	F	243	VAL
1	F	247	ARG
1	F	251	GLU
1	F	266	LEU
1	F	278	LYS
1	F	307	LEU
1	F	317	ARG

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Mol	Chain	Res	Type
1	F	330	ARG
1	F	336	VAL
1	F	338	LEU
1	A	3	LYS
1	A	5	ILE
1	A	24	LEU
1	A	35	ARG
1	A	65	LEU
1	A	75	ARG
1	A	103	ASP
1	A	110	THR
1	A	113	VAL
1	A	114	HIS
1	A	116	HIS
1	A	124	GLN
1	A	131	PHE
1	A	134	LEU
1	A	136	LEU
1	A	154	ASP
1	A	174	GLN
1	A	176	LEU
1	A	186	LYS
1	A	191	LYS
1	A	197	LEU
1	A	198	ARG
1	A	209	VAL
1	A	214	ARG
1	A	227	GLN
1	A	233	ASP
1	A	239	LEU
1	A	240	ARG
1	A	266	LEU
1	A	277	VAL
1	A	307	LEU
1	A	332	VAL
1	A	337	THR
1	A	338	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	F	45	ASN

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Mol	Chain	Res	Type
1	F	128	GLN
1	F	132	ASN
1	F	300	GLN
1	A	45	ASN
1	A	84	ASN
1	A	255	HIS
1	A	300	GLN
1	A	308	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	1FZ	A	875	6	24,30,30	2.45	7 (29%)	31,47,47	2.31	7 (22%)
5	1FZ	F	876	6	24,30,30	2.22	5 (20%)	31,47,47	2.11	7 (22%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical

component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	1FZ	A	875	6	-	0/13/34/34	0/2/2/2
5	1FZ	F	876	6	-	0/13/34/34	0/2/2/2

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	F	876	1FZ	PB-O1B	-2.78	1.48	1.56
5	A	875	1FZ	PB-O1B	-2.00	1.51	1.56
5	A	875	1FZ	PB-O3B	2.45	1.62	1.59
5	A	875	1FZ	PA-O5'	2.49	1.64	1.57
5	F	876	1FZ	O4-C4	4.31	1.35	1.24
5	A	875	1FZ	C6-N1	4.58	1.41	1.35
5	F	876	1FZ	PB-O2B	4.68	1.51	1.46
5	F	876	1FZ	C6-N1	4.83	1.42	1.35
5	A	875	1FZ	O4-C4	4.94	1.36	1.24
5	A	875	1FZ	PB-O2B	5.00	1.51	1.46
5	F	876	1FZ	PA-O1A	5.55	1.52	1.46
5	A	875	1FZ	PA-O1A	7.02	1.54	1.46

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	875	1FZ	C5-C4-N3	-5.69	118.80	125.14
5	F	876	1FZ	C5-C4-N3	-5.55	118.96	125.14
5	A	875	1FZ	O1A-PA-N3A	-3.04	107.24	111.90
5	A	875	1FZ	PG-O3B-PB	-2.28	125.02	132.67
5	F	876	1FZ	PG-O3B-PB	-2.23	125.18	132.67
5	A	875	1FZ	O3B-PB-N3A	-2.18	100.44	106.44
5	F	876	1FZ	O3B-PB-N3A	-2.05	100.78	106.44
5	F	876	1FZ	PA-O5'-C5'	2.03	126.99	120.79
5	A	875	1FZ	O2G-PG-O1G	3.16	119.41	107.38
5	F	876	1FZ	O1B-PB-O2B	3.53	117.37	110.00
5	A	875	1FZ	O2A-PA-O1A	3.87	118.08	110.00
5	F	876	1FZ	O2A-PA-O1A	3.97	118.28	110.00
5	F	876	1FZ	C4-N3-C2	6.84	121.16	115.25
5	A	875	1FZ	C4-N3-C2	8.22	122.35	115.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	875	1FZ	3	0

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	342/352 (97%)	0.31	13 (3%) 44 49	38, 63, 93, 130	0
1	F	342/352 (97%)	0.06	1 (0%) 94 95	32, 45, 63, 80	0
2	B	18/18 (100%)	-0.54	0 100 100	45, 61, 76, 85	0
2	G	18/18 (100%)	-0.06	1 (5%) 28 32	43, 49, 87, 117	0
3	H	14/14 (100%)	-0.06	0 100 100	42, 45, 81, 84	0
4	C	17/17 (100%)	-0.60	0 100 100	47, 51, 76, 81	0
All	All	751/771 (97%)	0.14	15 (1%) 68 71	32, 50, 87, 130	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	0	GLY	17.1
1	F	0	GLY	5.6
1	A	116	HIS	5.2
1	A	1	SER	4.2
2	G	837	DT	3.2
1	A	76	PHE	3.1
1	A	203	VAL	3.1
1	A	23	ALA	2.6
1	A	136	LEU	2.5
1	A	131	PHE	2.5
1	A	109	VAL	2.2
1	A	215	PHE	2.2
1	A	237	GLU	2.1
1	A	7	VAL	2.1
1	A	234	VAL	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 carbohydrates in this entry.

6.4 Ligands [i](#)

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, 95th 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(\AA^2)	Q<0.9
6	MG	F	903	1/1	0.96	0.33	8.39	40,40,40,40	0
5	1FZ	F	876	29/29	0.99	0.16	0.54	40,41,41,42	0
6	MG	A	901	1/1	0.89	0.13	0.31	49,49,49,49	0
5	1FZ	A	875	29/29	0.96	0.11	-0.47	48,49,52,52	0
6	MG	F	904	1/1	0.89	0.14	-0.47	44,44,44,44	0
6	MG	A	902	1/1	0.95	0.08	-	76,76,76,76	0

6.5 Other polymers [i](#)

There are no such residues in this entry.