



Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 06:56 am BST

PDB ID : 4L18
Title : Crystal structure of Runx1 and Ets1 bound to TCR alpha promoter (crystal form 3)
Authors : Tahirov, T.H.
Deposited on : 2013-06-02
Resolution : 2.30 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
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) : 2.11

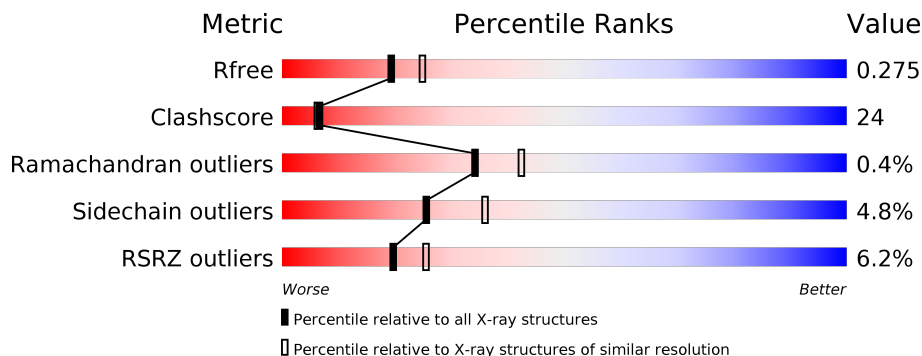
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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.



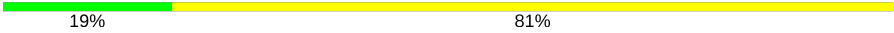

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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 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	167	<div style="display: flex; align-items: center;"> <div style="width: 8%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 53%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 28%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 15%; height: 10px; background-color: grey; margin-right: 5px;"></div> </div> <p style="font-size: small; margin-top: 5px;">8% 53% 28% 15%</p>
1	E	167	<div style="display: flex; align-items: center;"> <div style="width: 7%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 51%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 29%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 16%; height: 10px; background-color: grey; margin-right: 5px;"></div> </div> <p style="font-size: small; margin-top: 5px;">7% 51% 29% 16%</p>
2	B	146	<div style="display: flex; align-items: center;"> <div style="width: 0%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 47%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 27%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 27%; height: 10px; background-color: grey; margin-right: 5px;"></div> </div> <p style="font-size: small; margin-top: 5px;">% 47% 27% 27%</p>
2	F	146	<div style="display: flex; align-items: center;"> <div style="width: 5%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 49%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 23%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 27%; height: 10px; background-color: grey; margin-right: 5px;"></div> </div> <p style="font-size: small; margin-top: 5px;">5% 49% 23% 27%</p>
3	C	16	<div style="display: flex; align-items: center;"> <div style="width: 6%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 38%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 50%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 13%; height: 10px; background-color: orange; margin-right: 5px;"></div> </div> <p style="font-size: small; margin-top: 5px;">6% 38% 50% 13%</p>
3	G	16	<div style="display: flex; align-items: center;"> <div style="width: 13%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 81%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 6%; height: 10px; background-color: orange; margin-right: 5px;"></div> </div> <p style="font-size: small; margin-top: 5px;">13% 81% 6%</p>

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Mol	Chain	Length	Quality of chain
4	D	16	 19% 81%
4	H	16	 13% 88%

2 Entry composition i

There are 6 unique types of molecules in this entry. The entry contains 5431 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 Runt-related transcription factor 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	142	1103	692	201	206	4	0	0	0
1	E	141	1092	686	197	205	4	0	0	0

- Molecule 2 is a protein called Protein C-ets-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	107	897	580	154	159	4	0	0	0
2	F	106	890	576	153	157	4	0	0	0

- Molecule 3 is a DNA chain called 5'-D(*GP*GP*AP*AP*GP*CP*CP*AP*CP*AP*TP*C P*CP*TP*CP*T)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
3	C	16	321	154	59	93	15	0	0	0
3	G	16	321	154	59	93	15	0	0	0

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
4	D	16	329	157	62	95	15	0	0	0
4	H	16	329	157	62	95	15	0	0	0

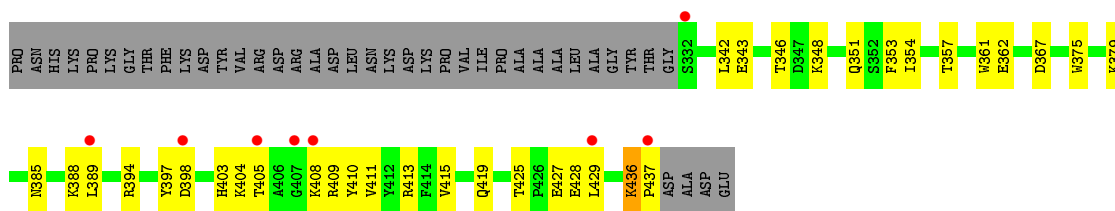
- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	D	1	Total	C	O	0	0
			6	3	3		

- Molecule 6 is water.

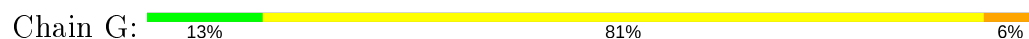
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	27	Total	O	0	0
			27	27		
6	B	25	Total	O	0	0
			25	25		
6	C	13	Total	O	0	0
			13	13		
6	D	18	Total	O	0	0
			18	18		
6	E	32	Total	O	0	0
			32	32		
6	F	14	Total	O	0	0
			14	14		
6	G	8	Total	O	0	0
			8	8		
6	H	6	Total	O	0	0
			6	6		



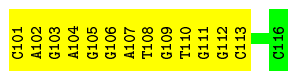
• Molecule 3: 5'-D(*GP*GP*AP*AP*GP*CP*CP*AP*CP*AP*TP*CP*CP*TP*CP*T)-3'



• Molecule 3: 5'-D(*GP*GP*AP*AP*GP*CP*CP*AP*CP*AP*TP*CP*CP*TP*CP*T)-3'



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



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



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	66.15Å 100.03Å 72.06Å 90.00° 110.23° 90.00°	Depositor
Resolution (Å)	29.64 – 2.30 29.64 – 2.31	Depositor EDS
% Data completeness (in resolution range)	87.5 (29.64-2.30) 87.6 (29.64-2.31)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.10 (at 2.31Å)	Xtrriage
Refinement program	CNS	Depositor
R, R_{free}	0.231 , 0.280 0.226 , 0.275	Depositor DCC
R_{free} test set	1753 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	44.0	Xtrriage
Anisotropy	0.137	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 61.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5431	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.94% 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

Bond lengths and bond angles in the following residue types are not validated in this section: GOL

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	0.44	0/1125	0.73	0/1527
1	E	0.46	0/1113	0.75	1/1512 (0.1%)
2	B	0.43	0/922	0.61	0/1243
2	F	0.37	0/915	0.58	0/1233
3	C	0.68	0/359	0.87	0/551
3	G	0.48	0/359	0.78	0/551
4	D	0.57	0/369	0.85	0/569
4	H	0.49	0/369	0.83	0/569
All	All	0.47	0/5531	0.73	1/7755 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
3	C	0	2
3	G	0	1
4	D	0	1
All	All	0	4

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	55	LEU	N-CA-C	-6.91	92.36	111.00

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	C	13	DC	Sidechain
3	C	5	DG	Sidechain
4	D	103	DG	Sidechain
3	G	205	DG	Sidechain

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	1103	0	1104	51	0
1	E	1092	0	1092	49	0
2	B	897	0	889	29	0
2	F	890	0	885	41	0
3	C	321	0	181	15	0
3	G	321	0	181	24	0
4	D	329	0	182	17	0
4	H	329	0	182	27	0
5	D	6	0	8	0	0
6	A	27	0	0	5	0
6	B	25	0	0	1	0
6	C	13	0	0	1	0
6	D	18	0	0	0	0
6	E	32	0	0	5	0
6	F	14	0	0	0	0
6	G	8	0	0	1	0
6	H	6	0	0	1	0
All	All	5431	0	4704	236	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 24.

All (236) 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)
4:H:304:DA:H2''	4:H:305:DG:H5''	1.25	1.10
4:H:307:DA:H2''	4:H:308:DT:H5'	1.34	1.07
4:H:315:DT:H2''	4:H:316:DC:H5''	1.46	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:348:LYS:HA	2:B:351:GLN:HE21	1.34	0.92
1:A:147:THR:HG23	1:A:163:HIS:HA	1.51	0.91
1:E:56:ALA:HB1	1:E:61:GLU:HG2	1.55	0.89
4:D:104:DA:H2''	4:D:105:DG:H5'	1.56	0.86
1:E:56:ALA:HB2	1:E:62:LEU:HG	1.59	0.83
2:F:419:GLN:NE2	2:F:425:THR:HG22	1.94	0.83
1:E:116:GLU:HG2	1:E:137:VAL:HB	1.61	0.83
2:F:348:LYS:HA	2:F:351:GLN:HE21	1.48	0.79
3:C:12:DC:H1'	3:C:13:DC:H5'	1.64	0.79
1:A:56:ALA:HB1	1:A:61:GLU:HG2	1.64	0.79
4:H:307:DA:H2''	4:H:308:DT:C5'	2.12	0.79
2:B:361:TRP:HZ3	2:B:429:LEU:HD23	1.47	0.78
3:C:2:DG:H2''	3:C:3:DA:N7	2.00	0.77
3:G:206:DC:H2''	3:G:207:DC:H5'	1.66	0.77
3:C:2:DG:H2''	3:C:3:DA:C8	2.20	0.76
1:A:192:LEU:HD23	1:A:192:LEU:H	1.50	0.76
3:G:212:DC:H1'	3:G:213:DC:H5'	1.67	0.76
4:H:313:DC:C6	4:H:314:DT:H72	2.20	0.75
2:B:348:LYS:HG3	2:B:351:GLN:NE2	2.02	0.75
1:E:138:GLY:HA3	6:E:332:HOH:O	1.86	0.75
2:F:425:THR:OG1	2:F:428:GLU:HG3	1.88	0.73
1:A:58:HIS:CE1	1:A:61:GLU:HB2	2.24	0.72
4:H:302:DA:H1'	4:H:303:DG:H5'	1.71	0.71
3:G:210:DA:H1'	3:G:211:DT:H5''	1.72	0.70
3:C:6:DC:H1'	3:C:7:DC:H5'	1.72	0.70
1:E:52:VAL:HG22	1:E:55:LEU:HG	1.75	0.68
2:F:405:THR:HG21	2:F:408:LYS:HE2	1.76	0.68
4:H:314:DT:H1'	4:H:315:DT:H5''	1.75	0.68
1:A:56:ALA:HB2	1:A:62:LEU:HG	1.75	0.68
2:F:348:LYS:HA	2:F:351:GLN:NE2	2.10	0.67
3:G:205:DG:H2''	3:G:206:DC:H5'	1.77	0.66
4:H:304:DA:C2'	4:H:305:DG:H5''	2.15	0.66
1:E:192:LEU:H	1:E:192:LEU:HD23	1.61	0.66
1:E:56:ALA:CB	1:E:62:LEU:HG	2.25	0.66
4:D:104:DA:H2''	4:D:105:DG:C5'	2.25	0.65
3:C:11:DT:H1'	3:C:12:DC:H5'	1.79	0.64
1:A:68:PRO:O	1:A:94:LEU:HD12	1.98	0.64
2:F:419:GLN:HE21	2:F:425:THR:HG22	1.61	0.64
1:A:126:ASN:HD22	1:A:126:ASN:N	1.94	0.64
1:A:176:PRO:HB3	6:A:323:HOH:O	1.97	0.63
1:A:99:ASP:O	1:E:192:LEU:HD23	1.99	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:336:GLN:HG3	2:B:339:GLN:OE1	1.99	0.63
1:A:69:ASN:HB3	6:A:325:HOH:O	1.99	0.62
1:A:56:ALA:CB	1:A:61:GLU:HG2	2.31	0.60
1:A:194:PHE:CZ	2:B:429:LEU:HD13	2.37	0.60
2:B:362:GLU:OE1	2:B:411:VAL:HG11	2.02	0.60
1:A:87:ILE:HD12	1:A:87:ILE:O	2.02	0.59
2:F:389:LEU:HD23	2:F:389:LEU:C	2.22	0.59
3:G:212:DC:H1'	3:G:213:DC:C5'	2.31	0.59
1:A:55:LEU:HG	1:A:55:LEU:O	2.02	0.59
2:B:373:ARG:HH21	2:B:373:ARG:HG2	1.67	0.58
2:B:361:TRP:CZ3	2:B:429:LEU:HD23	2.34	0.58
4:D:112:DG:H2''	4:D:113:DC:H5'	1.85	0.58
3:G:204:DA:N7	6:G:303:HOH:O	2.31	0.58
4:H:307:DA:C2'	4:H:308:DT:H5'	2.22	0.58
4:H:308:DT:H2''	4:H:309:DG:OP2	2.03	0.58
3:C:10:DA:H2''	3:C:11:DT:H5'	1.85	0.58
1:A:174:ARG:NH2	4:D:111:DG:N7	2.49	0.58
2:F:419:GLN:NE2	2:F:425:THR:HA	2.19	0.58
4:H:311:DG:H2''	4:H:312:DG:OP2	2.04	0.58
2:F:403:HIS:CG	2:F:415:VAL:HG11	2.39	0.58
2:F:436:LYS:O	2:F:436:LYS:HD2	2.04	0.57
1:E:139:ARG:HD3	1:E:139:ARG:H	1.70	0.57
4:D:108:DT:H2''	4:D:109:DG:C8	2.39	0.57
3:G:201:DG:H2''	3:G:202:DG:C8	2.40	0.57
1:A:139:ARG:HB2	1:A:139:ARG:HH11	1.68	0.57
2:F:405:THR:HG22	2:F:408:LYS:HG3	1.85	0.56
1:E:144:LYS:O	1:E:167:LYS:HG3	2.04	0.56
4:H:312:DG:H2''	4:H:313:DC:H5''	1.87	0.56
1:E:53:GLU:CD	1:E:64:ARG:HG3	2.25	0.56
1:A:79:TRP:CZ2	1:A:86:PRO:HD3	2.40	0.56
1:A:192:LEU:HD23	1:A:192:LEU:N	2.20	0.56
1:E:193:SER:O	1:E:197:ARG:HG3	2.05	0.56
3:G:206:DC:C2'	3:G:207:DC:H5'	2.35	0.56
3:G:211:DT:H2''	3:G:212:DC:H5'	1.88	0.56
1:A:139:ARG:HD2	6:A:307:HOH:O	2.06	0.56
1:E:52:VAL:HG13	1:E:52:VAL:O	2.06	0.55
2:F:427:GLU:H	2:F:427:GLU:CD	2.03	0.55
1:A:139:ARG:CD	6:A:307:HOH:O	2.54	0.55
1:A:199:SER:O	1:A:203:GLN:HG3	2.07	0.55
1:E:87:ILE:O	1:E:87:ILE:HD12	2.07	0.55
2:F:388:LYS:NZ	3:G:211:DT:OP2	2.37	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:302:DA:N3	6:H:401:HOH:O	2.33	0.55
1:A:63:VAL:HG22	1:A:64:ARG:N	2.23	0.54
1:E:58:HIS:CE1	1:E:61:GLU:HB2	2.42	0.54
2:F:348:LYS:HG3	2:F:351:GLN:NE2	2.22	0.54
2:F:409:ARG:HG2	2:F:410:TYR:CD2	2.42	0.54
1:A:203:GLN:O	1:A:204:LEU:C	2.46	0.54
2:F:348:LYS:HG3	2:F:351:GLN:HE22	1.73	0.53
2:B:404:LYS:HG2	4:D:104:DA:OP1	2.08	0.53
4:H:304:DA:H2''	4:H:305:DG:C5'	2.18	0.53
2:B:337:LEU:HD13	2:B:375:TRP:CE2	2.44	0.53
3:C:5:DG:H2''	3:C:6:DC:H5'	1.91	0.53
3:C:10:DA:H1'	3:C:11:DT:H5''	1.91	0.52
2:F:389:LEU:HD23	2:F:389:LEU:O	2.09	0.52
1:A:139:ARG:H	1:A:139:ARG:HD3	1.74	0.52
3:G:210:DA:C2	4:H:309:DG:C2	2.98	0.52
4:H:313:DC:H2'	4:H:314:DT:H72	1.91	0.52
1:A:192:LEU:CD2	1:E:98:PRO:HB2	2.40	0.52
4:H:307:DA:C2'	4:H:308:DT:C5'	2.85	0.52
3:C:9:DC:H2''	3:C:10:DA:C8	2.45	0.52
1:A:70:PHE:CE2	1:A:152:VAL:HG21	2.46	0.51
3:G:209:DC:H2''	3:G:210:DA:C8	2.45	0.51
1:A:192:LEU:HD23	1:E:99:ASP:O	2.10	0.51
4:D:101:DC:H2''	4:D:102:DA:C8	2.45	0.51
4:D:107:DA:H1'	4:D:108:DT:H5''	1.93	0.51
1:A:192:LEU:HD22	1:E:98:PRO:HB2	1.92	0.51
1:E:98:PRO:HG2	1:E:101:THR:OG1	2.10	0.51
1:E:56:ALA:HB2	1:E:62:LEU:CG	2.34	0.50
1:E:68:PRO:O	1:E:94:LEU:HD12	2.12	0.50
2:F:425:THR:OG1	2:F:427:GLU:HG2	2.11	0.50
1:E:147:THR:HG23	1:E:163:HIS:HA	1.93	0.50
1:E:198:LEU:CD1	2:F:343:GLU:HA	2.42	0.50
2:F:436:LYS:C	2:F:436:LYS:HD2	2.32	0.50
1:A:75:LEU:HD11	1:A:148:LEU:HD11	1.94	0.49
3:G:210:DA:H2''	3:G:211:DT:H5'	1.94	0.49
2:B:425:THR:OG1	2:B:428:GLU:HG3	2.11	0.49
4:H:312:DG:H2''	4:H:313:DC:C5'	2.42	0.49
4:H:305:DG:H2''	4:H:306:DG:H8	1.76	0.49
2:B:357:THR:O	2:B:438:ASP:HB2	2.12	0.49
1:A:56:ALA:CB	1:A:62:LEU:HG	2.43	0.49
1:E:132:ASN:C	1:E:132:ASN:ND2	2.65	0.49
2:F:436:LYS:HD2	2:F:437:PRO:O	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:11:DT:H2''	3:C:12:DC:H5'	1.94	0.49
2:B:353:PHE:HB2	2:B:367:ASP:HB3	1.95	0.48
1:E:132:ASN:HD22	1:E:132:ASN:C	2.15	0.48
1:E:61:GLU:HG3	1:E:62:LEU:HD23	1.95	0.48
2:F:409:ARG:HG2	2:F:410:TYR:CE2	2.48	0.48
2:F:404:LYS:NZ	4:H:303:DG:OP1	2.45	0.48
1:A:91:VAL:HG22	1:A:150:ILE:HD13	1.95	0.48
2:F:362:GLU:HB2	2:F:413:ARG:NH1	2.29	0.48
2:F:408:LYS:HE3	2:F:411:VAL:HG11	1.96	0.48
1:A:139:ARG:HD3	1:A:139:ARG:N	2.28	0.48
2:B:426:PRO:HG2	2:B:427:GLU:OE2	2.14	0.47
4:D:104:DA:C8	4:D:104:DA:H5'	2.49	0.47
4:D:104:DA:H1'	4:D:105:DG:H5''	1.96	0.47
1:E:177:ARG:NH1	6:E:327:HOH:O	2.41	0.47
3:G:215:DC:H2''	3:G:216:DT:O5'	2.14	0.47
2:F:362:GLU:HB2	2:F:413:ARG:CZ	2.45	0.47
4:D:111:DG:H1'	4:D:112:DG:C8	2.50	0.47
1:A:75:LEU:CD1	1:A:148:LEU:HD11	2.44	0.47
2:F:404:LYS:HE3	4:H:303:DG:OP1	2.15	0.47
2:B:405:THR:HG21	2:B:408:LYS:HD2	1.97	0.47
2:F:342:LEU:O	2:F:346:THR:HG23	2.14	0.47
3:G:210:DA:H2''	3:G:211:DT:C5'	2.45	0.47
2:B:428:GLU:O	2:B:432:MET:HG3	2.15	0.47
4:H:312:DG:C2'	4:H:313:DC:H5''	2.44	0.47
1:A:126:ASN:ND2	1:A:126:ASN:N	2.63	0.47
1:A:136:PHE:HB2	1:A:168:ILE:HD12	1.96	0.47
1:A:82:ASN:OD1	1:A:118:ARG:NH1	2.47	0.47
4:D:108:DT:H2''	4:D:109:DG:H8	1.79	0.47
2:F:353:PHE:HB2	2:F:367:ASP:HB3	1.97	0.47
2:F:419:GLN:HE22	2:F:425:THR:HG22	1.76	0.46
1:E:171:ASP:HB3	1:E:174:ARG:HD2	1.95	0.46
1:E:70:PHE:CE2	1:E:152:VAL:HG21	2.51	0.46
1:A:75:LEU:HD21	1:A:89:PHE:CD2	2.51	0.46
4:D:111:DG:H2''	4:D:112:DG:OP2	2.16	0.46
1:E:54:VAL:HG12	6:E:321:HOH:O	2.15	0.46
2:F:425:THR:HG23	2:F:428:GLU:OE1	2.16	0.46
1:E:118:ARG:HG2	6:E:328:HOH:O	2.16	0.46
1:E:63:VAL:HG22	1:E:64:ARG:N	2.30	0.46
3:G:211:DT:H1'	3:G:212:DC:H5'	1.98	0.45
2:B:429:LEU:HD12	2:B:433:LEU:HD23	1.97	0.45
4:D:109:DG:H1'	4:D:110:DT:H5'	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:304:DA:H5'	4:H:304:DA:H8	1.81	0.45
1:A:63:VAL:CG2	1:A:64:ARG:N	2.79	0.45
1:A:84:THR:HG23	6:A:301:HOH:O	2.15	0.45
2:B:370:GLU:HA	2:B:370:GLU:OE1	2.17	0.45
2:B:341:LEU:HD11	2:B:393:LEU:HD21	1.98	0.45
1:E:92:VAL:HG13	1:E:128:VAL:HG22	1.98	0.45
4:D:107:DA:H2''	4:D:108:DT:H5'	1.98	0.45
2:F:375:TRP:CH2	2:F:379:LYS:HG3	2.52	0.45
1:A:56:ALA:HB2	1:A:62:LEU:CG	2.45	0.45
1:E:57:ASP:OD2	1:E:57:ASP:N	2.48	0.45
4:H:307:DA:H1'	4:H:308:DT:H5''	1.99	0.45
2:B:373:ARG:HG2	2:B:373:ARG:NH2	2.32	0.44
3:C:11:DT:C2'	3:C:12:DC:H5'	2.47	0.44
2:F:394:ARG:HA	2:F:397:TYR:CE2	2.51	0.44
1:A:78:HIS:NE2	1:A:172:GLY:HA2	2.33	0.44
1:A:193:SER:O	1:A:197:ARG:HG3	2.17	0.44
2:B:378:ARG:HG3	2:B:378:ARG:HH11	1.83	0.44
2:B:385:ASN:OD1	2:B:387:GLU:HB3	2.18	0.44
1:E:116:GLU:HG2	1:E:137:VAL:CB	2.40	0.44
2:F:385:ASN:ND2	2:F:388:LYS:HE2	2.33	0.44
2:B:430:HIS:HB3	2:B:435:VAL:HB	2.00	0.44
4:H:305:DG:H2''	4:H:306:DG:C8	2.53	0.44
3:G:206:DC:H1'	3:G:207:DC:H5'	1.99	0.44
1:A:58:HIS:O	1:A:61:GLU:HB3	2.18	0.43
3:G:212:DC:C1'	3:G:213:DC:H5'	2.43	0.43
1:E:171:ASP:HB3	1:E:174:ARG:CD	2.48	0.43
1:E:175:GLU:HA	1:E:175:GLU:OE2	2.17	0.43
1:E:78:HIS:NE2	1:E:172:GLY:HA2	2.33	0.43
3:G:211:DT:C2'	3:G:212:DC:H5'	2.47	0.43
2:B:361:TRP:HB3	2:B:414:PHE:HB2	2.00	0.43
3:G:206:DC:H1'	3:G:207:DC:C5'	2.49	0.43
1:E:79:TRP:HB2	1:E:166:ILE:HD11	2.01	0.43
2:F:408:LYS:HB2	2:F:411:VAL:HB	2.00	0.43
3:G:206:DC:H2''	3:G:207:DC:C5'	2.45	0.43
2:B:397:TYR:HB2	6:B:523:HOH:O	2.18	0.43
2:F:357:THR:HG21	2:F:362:GLU:HG2	1.99	0.43
2:B:336:GLN:HE21	2:B:336:GLN:HB3	1.68	0.42
1:E:155:ASN:HA	1:E:156:PRO:HA	1.92	0.42
3:G:203:DA:C6	3:G:204:DA:C6	3.07	0.42
3:G:202:DG:H2''	3:G:203:DA:OP2	2.19	0.42
3:C:11:DT:C1'	3:C:12:DC:H5'	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:98:PRO:HB2	1:E:192:LEU:HD22	2.02	0.42
1:E:196:GLU:O	1:E:200:GLU:HG3	2.20	0.42
1:E:63:VAL:CG2	1:E:64:ARG:N	2.83	0.42
1:A:112:ASN:N	1:A:112:ASN:HD22	2.18	0.42
3:C:10:DA:H1'	3:C:11:DT:C5'	2.49	0.42
1:E:139:ARG:HB3	6:E:325:HOH:O	2.19	0.42
1:A:79:TRP:NE1	1:A:83:LYS:HG2	2.34	0.42
1:E:139:ARG:HD3	1:E:139:ARG:N	2.32	0.42
1:A:142:ARG:HH22	3:C:3:DA:H2	1.68	0.42
2:B:390:SER:O	2:B:394:ARG:HG3	2.19	0.42
2:F:357:THR:CG2	2:F:362:GLU:HG2	2.50	0.42
1:A:80:ARG:HG3	1:A:171:ASP:O	2.19	0.41
1:A:79:TRP:HB2	1:A:166:ILE:HD11	2.02	0.41
4:D:104:DA:C2'	4:D:105:DG:C5'	2.98	0.41
2:F:404:LYS:CE	4:H:303:DG:OP1	2.69	0.41
2:B:426:PRO:O	2:B:430:HIS:CG	2.73	0.41
1:A:112:ASN:O	1:A:113:TYR:C	2.58	0.41
1:E:139:ARG:N	1:E:139:ARG:CD	2.84	0.41
4:H:304:DA:C6	4:H:305:DG:C6	3.09	0.41
2:B:391:ARG:NH2	4:D:106:DG:N7	2.53	0.41
2:F:354:ILE:O	2:F:354:ILE:HG23	2.19	0.41
2:F:361:TRP:HZ3	2:F:429:LEU:HD22	1.85	0.41
2:F:388:LYS:HZ1	3:G:211:DT:P	2.40	0.41
1:A:131:PHE:N	1:A:131:PHE:CD2	2.88	0.40
3:C:13:DC:H2'	6:C:110:HOH:O	2.21	0.40
1:E:119:ASN:ND2	1:E:133:ASP:O	2.54	0.40
1:E:60:GLY:O	1:E:74:VAL:HG21	2.22	0.40
1:E:61:GLU:HG3	1:E:62:LEU:N	2.34	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	138/167 (83%)	131 (95%)	5 (4%)	2 (1%)	11	11
1	E	137/167 (82%)	131 (96%)	6 (4%)	0	100	100
2	B	105/146 (72%)	100 (95%)	5 (5%)	0	100	100
2	F	104/146 (71%)	101 (97%)	3 (3%)	0	100	100
All	All	484/626 (77%)	463 (96%)	19 (4%)	2 (0%)	34	42

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	204	LEU
1	A	190	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	122/146 (84%)	114 (93%)	8 (7%)	16	22
1	E	121/146 (83%)	112 (93%)	9 (7%)	13	17
2	B	96/127 (76%)	94 (98%)	2 (2%)	53	70
2	F	96/127 (76%)	94 (98%)	2 (2%)	53	70
All	All	435/546 (80%)	414 (95%)	21 (5%)	25	36

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	53	GLU
1	A	61	GLU
1	A	83	LYS
1	A	112	ASN
1	A	126	ASN
1	A	139	ARG
1	A	162	TYR
1	A	192	LEU
2	B	434	ASP

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Mol	Chain	Res	Type
2	B	436	LYS
1	E	57	ASP
1	E	61	GLU
1	E	77	THR
1	E	114	SER
1	E	130	ARG
1	E	132	ASN
1	E	139	ARG
1	E	162	TYR
1	E	192	LEU
2	F	398	ASP
2	F	436	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	58	HIS
1	A	112	ASN
1	A	126	ASN
2	B	336	GLN
2	B	351	GLN
2	B	380	ASN
1	E	58	HIS
1	E	112	ASN
1	E	132	ASN
2	F	351	GLN
2	F	419	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

1 ligand is modelled in this entry.

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	GOL	D	201	-	5,5,5	0.57	0	5,5,5	0.14	0

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	GOL	D	201	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	142/167 (85%)	0.54	13 (9%) 9 12	18, 50, 82, 101	0
1	E	141/167 (84%)	0.48	11 (7%) 13 17	18, 47, 86, 111	0
2	B	107/146 (73%)	0.27	2 (1%) 66 73	28, 48, 74, 109	0
2	F	106/146 (72%)	0.54	8 (7%) 14 19	39, 59, 80, 111	0
3	C	16/16 (100%)	0.13	1 (6%) 20 25	28, 39, 65, 71	0
3	G	16/16 (100%)	0.08	0 100 100	38, 55, 85, 88	0
4	D	16/16 (100%)	-0.40	0 100 100	34, 46, 56, 62	0
4	H	16/16 (100%)	-0.24	0 100 100	44, 57, 67, 73	0
All	All	560/690 (81%)	0.40	35 (6%) 20 25	18, 51, 82, 111	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	56	ALA	8.7
1	E	56	ALA	5.9
1	E	55	LEU	5.7
1	A	55	LEU	5.7
1	A	54	VAL	4.8
1	E	113	TYR	4.5
2	F	437	PRO	4.4
2	F	332	SER	3.9
1	A	96	ASP	3.7
1	E	57	ASP	3.5
1	A	113	TYR	3.4
1	A	57	ASP	3.4
1	E	175	GLU	3.3
2	B	439	ALA	3.2
1	E	110	ASP	3.1
1	A	177	ARG	3.1

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Mol	Chain	Res	Type	RSRZ
2	F	398	ASP	2.7
1	A	175	GLU	2.7
1	E	111	GLU	2.6
1	A	155	ASN	2.6
2	F	408	LYS	2.6
1	E	53	GLU	2.6
1	E	190	GLY	2.4
3	C	1	DG	2.3
1	E	156	PRO	2.3
2	B	433	LEU	2.3
2	F	407	GLY	2.2
1	A	64	ARG	2.2
2	F	389	LEU	2.2
1	A	122	ALA	2.1
1	E	96	ASP	2.1
1	A	194	PHE	2.1
2	F	429	LEU	2.0
2	F	405	THR	2.0
1	A	102	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 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. 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	B-factors(Å ²)	Q<0.9
5	GOL	D	201	6/6	0.90	0.14	53,54,55,56	0

6.5 Other polymers

There are no such residues in this entry.