



Full wwPDB X-ray Structure Validation Report ⓘ

May 23, 2016 – 11:42 AM EDT

PDB ID : 5JJ6
Title : Fic-1 (aa134 - 508) from *C. elegans*
Authors : Cruz, V.E.; Schwartz, T.U.
Deposited on : 2016-04-22
Resolution : 2.91 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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20027674
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) : rb-20027674

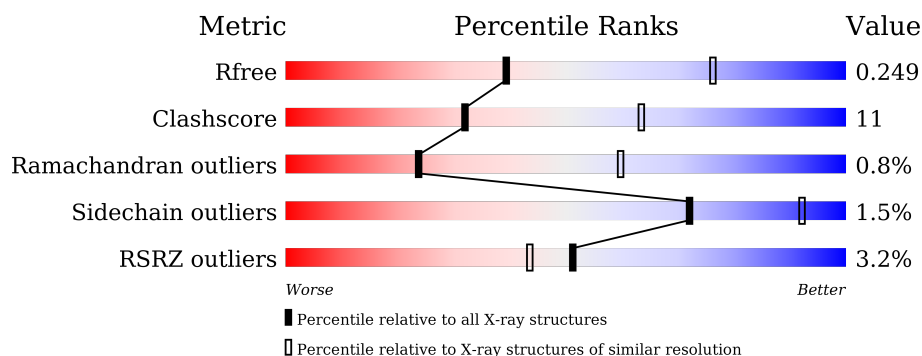
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.91 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	1451 (2.90-2.90)
Clashscore	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (2.90-2.90)
RSRZ outliers	91569	1456 (2.90-2.90)

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	382	<div> <div>3%</div> <div> <div></div> <div>65%</div> <div>20%</div> <div>•</div> <div>14%</div> </div> </div>
1	B	382	<div> <div>2%</div> <div> <div></div> <div>73%</div> <div>14%</div> <div>13%</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5100 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 Adenosine monophosphate-protein transferase FICD homolog.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	330	Total	C	N	O	S	0	0	0
			2503	1573	454	462	14			
1	B	333	Total	C	N	O	S	0	0	0
			2581	1625	466	473	17			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	127	MET	-	initiating methionine	UNP Q23544
A	128	HIS	-	expression tag	UNP Q23544
A	129	HIS	-	expression tag	UNP Q23544
A	130	HIS	-	expression tag	UNP Q23544
A	131	HIS	-	expression tag	UNP Q23544
A	132	HIS	-	expression tag	UNP Q23544
A	133	HIS	-	expression tag	UNP Q23544
B	127	MET	-	initiating methionine	UNP Q23544
B	128	HIS	-	expression tag	UNP Q23544
B	129	HIS	-	expression tag	UNP Q23544
B	130	HIS	-	expression tag	UNP Q23544
B	131	HIS	-	expression tag	UNP Q23544
B	132	HIS	-	expression tag	UNP Q23544
B	133	HIS	-	expression tag	UNP Q23544

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

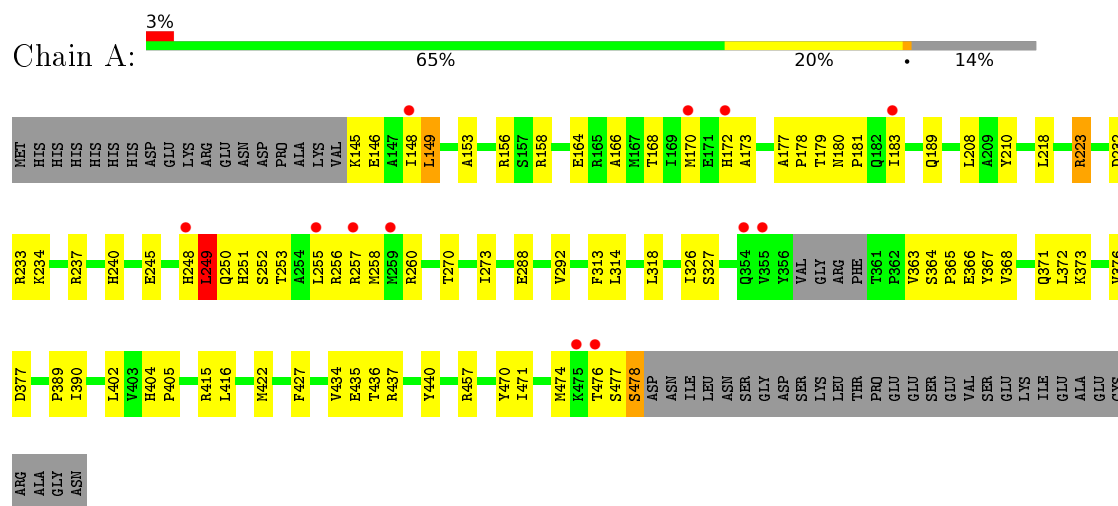
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	O	0	0
			2	2		
3	B	4	Total	O	0	0
			4	4		

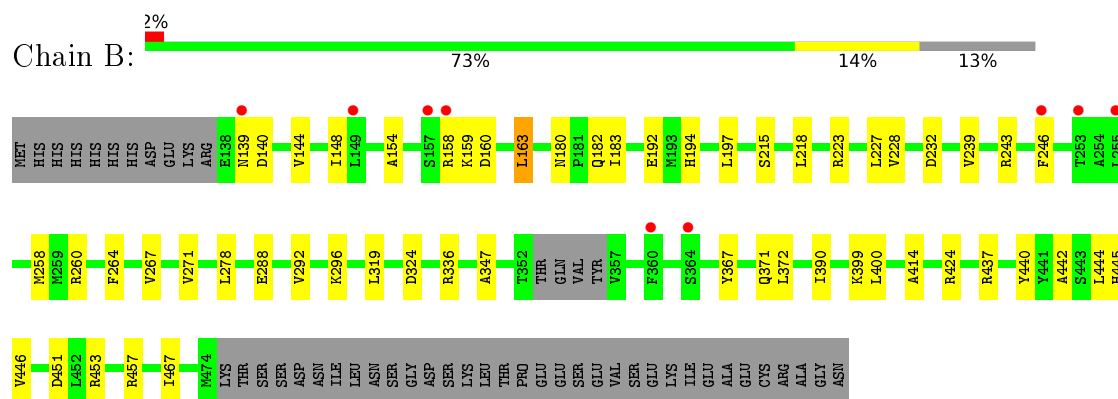
3 Residue-property plots

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: Adenosine monophosphate-protein transferase FICD homolog



- Molecule 1: Adenosine monophosphate-protein transferase FICD homolog



4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	192.15Å 192.15Å 182.52Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	122.97 – 2.91 122.97 – 2.91	Depositor EDS
% Data completeness (in resolution range)	97.5 (122.97-2.91) 92.1 (122.97-2.91)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.14 (at 2.91Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.206 , 0.255 0.211 , 0.249	Depositor DCC
R_{free} test set	1872 reflections (7.11%)	DCC
Wilson B-factor (Å ²)	58.5	Xtrriage
Anisotropy	0.445	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 55.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5100	wwPDB-VP
Average B, all atoms (Å ²)	66.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.02% 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: SO4

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	0.46	0/2545	0.68	2/3457 (0.1%)
1	B	0.44	0/2626	0.61	0/3557
All	All	0.45	0/5171	0.65	2/7014 (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
1	A	0	1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	249	LEU	CA-CB-CG	9.36	136.82	115.30
1	A	149	LEU	CA-CB-CG	6.25	129.69	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	367	TYR	Peptide

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	2503	0	2470	75	0
1	B	2581	0	2575	33	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
3	A	2	0	0	0	0
3	B	4	0	0	0	0
All	All	5100	0	5045	107	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 11.

All (107) 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:257:ARG:NH2	1:A:474:MET:SD	2.29	1.03
1:A:258:MET:SD	1:A:470:TYR:CE2	2.58	0.96
1:A:234:LYS:HB2	1:A:237:ARG:HD2	1.56	0.88
1:A:223:ARG:CB	1:A:223:ARG:HH11	1.89	0.85
1:A:148:ILE:HD13	1:A:172:HIS:HE1	1.41	0.85
1:A:258:MET:SD	1:A:470:TYR:CD2	2.70	0.84
1:A:258:MET:HE3	1:A:470:TYR:HD2	1.40	0.84
1:A:258:MET:HE3	1:A:470:TYR:CD2	2.15	0.81
1:A:258:MET:CE	1:A:470:TYR:CD2	2.64	0.81
1:A:363:VAL:HG21	1:A:404:HIS:HB3	1.61	0.80
1:A:148:ILE:HG12	1:A:173:ALA:HA	1.67	0.75
1:A:177:ALA:HB1	1:A:183:ILE:HD13	1.69	0.74
1:A:148:ILE:HD13	1:A:172:HIS:CE1	2.23	0.72
1:A:232:ASP:OD2	1:A:457:ARG:NH1	2.23	0.72
1:A:248:HIS:O	1:A:249:LEU:HB3	1.92	0.69
1:A:223:ARG:CG	1:A:223:ARG:HH11	2.06	0.68
1:B:192:GLU:OE1	1:B:223:ARG:NH2	2.26	0.68
1:B:215:SER:HA	1:B:218:LEU:HD12	1.76	0.67
1:B:144:VAL:HG13	1:B:183:ILE:HD11	1.77	0.67
1:A:223:ARG:HB3	1:A:223:ARG:HH11	1.60	0.64
1:B:163:LEU:CB	1:B:194:HIS:NE2	2.62	0.62
1:A:234:LYS:HG3	1:A:234:LYS:O	2.00	0.62
1:A:253:THR:HA	1:A:256:ARG:HB3	1.82	0.62
1:A:270:THR:HG23	1:A:415:ARG:HH11	1.64	0.61
1:B:158:ARG:HG3	1:B:159:LYS:H	1.66	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:234:LYS:CG	1:A:237:ARG:HB2	2.32	0.60
1:B:446:VAL:HG12	1:B:451:ASP:HB3	1.83	0.59
1:A:245:GLU:HG2	1:A:248:HIS:CE1	2.38	0.58
1:A:292:VAL:HG11	1:B:292:VAL:HG13	1.84	0.58
1:A:368:VAL:HA	1:A:371:GLN:OE1	2.04	0.58
1:A:234:LYS:HA	1:A:237:ARG:H	1.69	0.58
1:A:245:GLU:HA	1:A:248:HIS:CE1	2.40	0.57
1:A:252:SER:HA	1:A:255:LEU:HD13	1.87	0.56
1:B:246:PHE:HB2	1:B:467:ILE:HG23	1.88	0.55
1:A:245:GLU:HG2	1:A:248:HIS:HE1	1.73	0.54
1:A:470:TYR:N	1:A:470:TYR:CD1	2.73	0.54
1:A:257:ARG:HH11	1:A:470:TYR:HB3	1.72	0.54
1:A:148:ILE:HD11	1:A:173:ALA:HB2	1.91	0.53
1:A:164:GLU:O	1:A:168:THR:HG23	2.09	0.52
1:A:146:GLU:HA	1:A:149:LEU:HG	1.92	0.51
1:A:234:LYS:HE2	1:A:237:ARG:HD2	1.93	0.50
1:A:314:LEU:HA	1:A:318:LEU:HD13	1.93	0.50
1:A:326:ILE:O	1:A:376:VAL:HG13	2.11	0.50
1:A:234:LYS:CB	1:A:237:ARG:HB2	2.42	0.49
1:A:233:ARG:O	1:A:237:ARG:HG3	2.11	0.49
1:A:252:SER:O	1:A:255:LEU:HB2	2.12	0.49
1:A:158:ARG:O	1:A:166:ALA:HB2	2.12	0.49
1:A:260:ARG:NH2	1:A:288:GLU:OE2	2.42	0.49
1:A:273:ILE:HA	1:A:434:VAL:HG23	1.95	0.49
1:B:197:LEU:HD21	1:B:227:LEU:HD13	1.94	0.49
1:A:422:MET:O	1:A:427:PHE:HB2	2.13	0.48
1:A:372:LEU:HD11	1:A:405:PRO:HB3	1.94	0.48
1:A:234:LYS:HG3	1:A:237:ARG:HB2	1.95	0.48
1:A:470:TYR:N	1:A:470:TYR:HD1	2.09	0.48
1:A:363:VAL:HG23	1:A:402:LEU:O	2.13	0.48
1:A:148:ILE:O	1:A:148:ILE:HG13	2.13	0.47
1:A:390:ILE:HG12	1:A:422:MET:HE3	1.96	0.47
1:B:239:VAL:HG11	1:B:390:ILE:HD11	1.96	0.47
1:B:336:ARG:HA	1:B:347:ALA:HB1	1.96	0.47
1:A:146:GLU:CB	1:A:149:LEU:HG	2.45	0.47
1:B:148:ILE:HG21	1:B:182:GLN:OE1	2.13	0.47
1:B:264:PHE:CZ	1:B:288:GLU:HB2	2.50	0.46
1:A:223:ARG:NH1	1:A:223:ARG:HB3	2.29	0.46
1:B:437:ARG:O	1:B:440:TYR:HB3	2.15	0.46
1:A:208:LEU:HD21	1:A:218:LEU:HG	1.98	0.45
1:A:476:THR:C	1:A:478:SER:H	2.19	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:476:THR:C	1:A:478:SER:N	2.69	0.45
1:A:390:ILE:HG12	1:A:422:MET:CE	2.47	0.45
1:B:267:VAL:O	1:B:271:VAL:HG22	2.17	0.45
1:B:278:LEU:HD23	1:B:296:LYS:HE3	1.98	0.44
1:B:367:TYR:O	1:B:371:GLN:HG3	2.18	0.44
1:A:153:ALA:C	1:A:156:ARG:H	2.20	0.44
1:A:240:HIS:CE1	1:A:389:PRO:HD2	2.52	0.44
1:B:399:LYS:HB3	1:B:399:LYS:HE2	1.86	0.44
1:B:180:ASN:HB3	1:B:183:ILE:HD12	2.00	0.44
1:A:234:LYS:CG	1:A:234:LYS:O	2.64	0.43
1:B:158:ARG:HG3	1:B:159:LYS:N	2.31	0.43
1:A:180:ASN:HA	1:A:181:PRO:HD3	1.82	0.43
1:B:442:ALA:O	1:B:445:HIS:HB3	2.19	0.43
1:A:179:THR:HG22	1:A:210:TYR:CE2	2.54	0.43
1:B:243:ARG:HA	1:B:467:ILE:HD13	2.01	0.43
1:B:400:LEU:HD23	1:B:414:ALA:HA	2.01	0.42
1:B:319:LEU:HD22	1:B:424:ARG:HG2	2.02	0.42
1:A:170:MET:O	1:A:173:ALA:HB3	2.18	0.42
1:A:177:ALA:CB	1:A:183:ILE:HD13	2.43	0.42
1:A:363:VAL:HG21	1:A:404:HIS:CB	2.43	0.42
1:A:364:SER:HA	1:A:365:PRO:HD3	1.97	0.41
1:A:416:LEU:HD23	1:A:416:LEU:HA	1.92	0.41
1:B:260:ARG:HA	1:B:260:ARG:HD3	1.57	0.41
1:A:148:ILE:O	1:A:148:ILE:CG1	2.69	0.41
1:A:437:ARG:O	1:A:440:TYR:HB3	2.21	0.41
1:A:470:TYR:O	1:A:471:ILE:C	2.56	0.41
1:B:159:LYS:HE2	1:B:159:LYS:HB3	1.81	0.41
1:B:444:LEU:HD23	1:B:444:LEU:HA	1.82	0.41
1:B:232:ASP:OD2	1:B:457:ARG:NH1	2.54	0.41
1:B:228:VAL:HG21	1:B:453:ARG:NH1	2.36	0.41
1:A:145:LYS:HD3	1:A:145:LYS:HA	1.69	0.41
1:B:139:ASN:OD1	1:B:140:ASP:O	2.38	0.41
1:A:249:LEU:HD12	1:A:251:HIS:H	1.86	0.41
1:B:372:LEU:HA	1:B:372:LEU:HD23	1.74	0.41
1:A:435:GLU:HG2	1:A:436:THR:HG23	2.03	0.40
1:A:146:GLU:CA	1:A:149:LEU:HG	2.50	0.40
1:B:215:SER:HA	1:B:218:LEU:HB2	2.03	0.40
1:A:313:PHE:CE2	1:A:318:LEU:HD11	2.56	0.40
1:A:373:LYS:HG2	1:A:377:ASP:OD2	2.21	0.40
1:A:223:ARG:HG3	1:A:223:ARG:HH11	1.86	0.40
1:B:160:ASP:O	1:B:160:ASP:OD1	2.39	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	326/382 (85%)	308 (94%)	16 (5%)	2 (1%)	30	67
1	B	329/382 (86%)	313 (95%)	13 (4%)	3 (1%)	21	57
All	All	655/764 (86%)	621 (95%)	29 (4%)	5 (1%)	24	60

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	163	LEU
1	A	178	PRO
1	B	154	ALA
1	B	324	ASP
1	A	366	GLU

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	257/328 (78%)	250 (97%)	7 (3%)	52	84
1	B	269/328 (82%)	268 (100%)	1 (0%)	93	98
All	All	526/656 (80%)	518 (98%)	8 (2%)	72	92

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	189	GLN
1	A	223	ARG
1	A	249	LEU
1	A	250	GLN
1	A	327	SER
1	A	477	SER
1	A	478	SER
1	B	258	MET

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

Mol	Chain	Res	Type
1	A	172	HIS
1	A	251	HIS
1	B	445	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

2 ligands are 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	$\# Z > 2$
2	SO4	A	601	-	4,4,4	0.11	0	6,6,6	0.24	0
2	SO4	B	601	-	4,4,4	0.23	0	6,6,6	0.52	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
2	SO4	A	601	-	-	0/0/0/0	0/0/0/0
2	SO4	B	601	-	-	0/0/0/0	0/0/0/0

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	330/382 (86%)	0.51	12 (3%) 46 38	27, 71, 134, 144	0
1	B	333/382 (87%)	0.53	9 (2%) 58 52	30, 52, 105, 138	0
All	All	663/764 (86%)	0.52	21 (3%) 51 43	27, 61, 127, 144	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	259	MET	5.9
1	A	148	ILE	3.2
1	B	149	LEU	3.0
1	B	139	ASN	2.8
1	A	475	LYS	2.8
1	A	248	HIS	2.7
1	A	172	HIS	2.7
1	A	257	ARG	2.7
1	B	360	PHE	2.6
1	B	246	PHE	2.5
1	B	364	SER	2.5
1	A	476	THR	2.4
1	B	158	ARG	2.4
1	A	183	ILE	2.2
1	B	157	SER	2.2
1	A	355	VAL	2.1
1	B	255	LEU	2.1
1	A	170	MET	2.1
1	B	253	THR	2.0
1	A	255	LEU	2.0
1	A	354	GLN	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
2	SO4	B	601	5/5	0.96	0.22	0.71	40,44,49,57	0
2	SO4	A	601	5/5	0.97	0.13	-2.20	54,57,80,80	0

6.5 Other polymers [i](#)

There are no such residues in this entry.