



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 12:39 AM GMT

PDB ID : 2B9J
Title : Crystal structure of Fus3 with a docking motif from Far1
Authors : Remenyi, A.; Good, M.C.; Bhattacharyya, R.P.; Lim, W.A.
Deposited on : 2005-10-11
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
<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

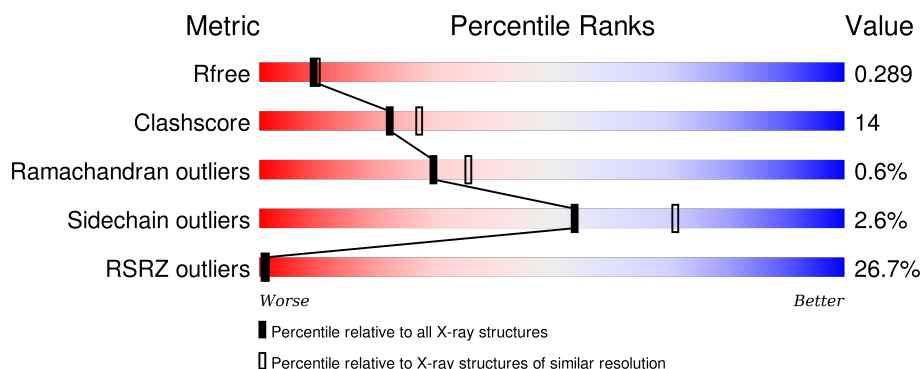
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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}	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (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. 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	353	
2	C	13	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3093 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 Mitogen-activated protein kinase FUS3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	337	Total	C	N	O	S	0	0	0
			2721	1763	460	484	14			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	180	VAL	THR	ENGINEERED	UNP P16892
A	182	PHE	TYR	ENGINEERED	UNP P16892

- Molecule 2 is a protein called Cyclin-dependent kinase inhibitor FAR1.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	11	Total	C	N	O	0	0	0
			80	52	16	12			

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		

- Molecule 4 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: C₁₀H₁₅N₅O₁₀P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

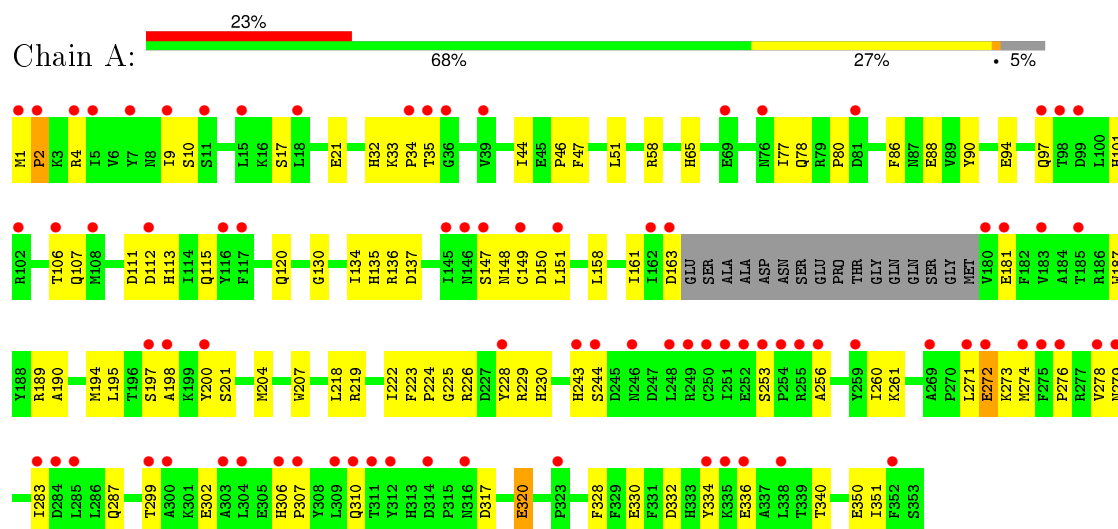
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	255	Total	O	0	0
			255	255		
5	C	9	Total	O	0	0
			9	9		

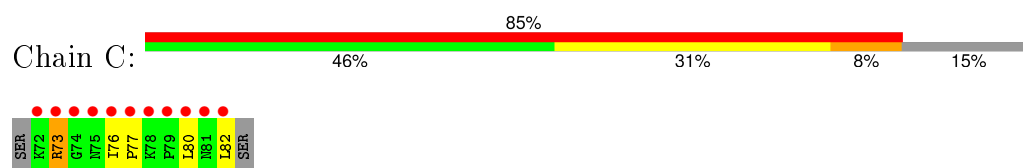
3 Residue-property plots [i](#)

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: Mitogen-activated protein kinase FUS3



• Molecule 2: Cyclin-dependent kinase inhibitor FAR1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	58.30 Å 63.24 Å 100.85 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.30 42.86 – 2.49	Depositor EDS
% Data completeness (in resolution range)	97.6 (20.00-2.30) 96.4 (42.86-2.49)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.59 (at 2.48 Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.196 , 0.263 0.286 , 0.289	Depositor DCC
R_{free} test set	1350 reflections (10.29%)	DCC
Wilson B-factor (Å ²)	31.2	Xtriage
Anisotropy	0.412	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 49.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 13125 reflections	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	3093	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.03% 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, ADP

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	0.52	0/2792	0.70	0/3788
2	C	0.47	0/81	0.73	0/109
All	All	0.52	0/2873	0.70	0/3897

There are no bond length outliers.

There are no bond angle outliers.

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	2721	0	2708	74	0
2	C	80	0	85	7	0
3	A	1	0	0	0	0
4	A	27	0	12	0	0
5	A	255	0	0	7	0
5	C	9	0	0	0	0
All	All	3093	0	2805	78	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 14.

All (78) 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:134:ILE:HD11	1:A:200:TYR:HB3	1.38	1.05
1:A:10:SER:H	1:A:78:GLN:HE22	1.09	0.95
1:A:1:MET:HG3	1:A:17:SER:HA	1.53	0.89
1:A:134:ILE:CD1	1:A:200:TYR:HB3	2.10	0.82
2:C:73:ARG:HG3	2:C:73:ARG:O	1.80	0.79
1:A:330:GLU:OE2	1:A:334:TYR:HE1	1.65	0.78
1:A:32:HIS:HD2	1:A:35:THR:OG1	1.70	0.75
1:A:189:ARG:HD2	1:A:194:MET:HG2	1.69	0.74
1:A:33:LYS:HB2	1:A:34:PRO:HD3	1.68	0.74
1:A:226:ARG:H	1:A:230:HIS:HD2	1.37	0.72
2:C:80:LEU:HD12	2:C:82:LEU:HD11	1.69	0.72
1:A:271:LEU:HD23	1:A:274:MET:CE	2.21	0.70
1:A:278:VAL:HB	1:A:283:ILE:HD11	1.73	0.69
1:A:271:LEU:HA	1:A:274:MET:HE3	1.75	0.68
2:C:73:ARG:O	2:C:73:ARG:CG	2.40	0.68
1:A:195:LEU:HD13	1:A:260:ILE:HD11	1.76	0.67
1:A:9:ILE:HA	1:A:78:GLN:OE1	1.94	0.67
1:A:330:GLU:OE2	1:A:334:TYR:CE1	2.46	0.66
1:A:10:SER:H	1:A:78:GLN:NE2	1.89	0.65
1:A:226:ARG:H	1:A:230:HIS:CD2	2.13	0.65
1:A:120:GLN:NE2	1:A:151:LEU:H	1.97	0.63
1:A:201:SER:H	1:A:204:MET:HE3	1.62	0.62
1:A:101:HIS:HD2	5:A:764:HOH:O	1.82	0.61
1:A:271:LEU:HD23	1:A:274:MET:HE1	1.84	0.59
1:A:273:LYS:O	1:A:276:PRO:HD3	2.02	0.59
1:A:226:ARG:N	1:A:230:HIS:HD2	2.01	0.57
1:A:4:ARG:HH12	1:A:21:GLU:CD	2.09	0.56
1:A:44:ILE:O	1:A:46:PRO:HD3	2.07	0.54
1:A:107:GLN:O	1:A:219:ARG:NH1	2.40	0.54
1:A:130:GLY:HA3	1:A:320:GLU:O	2.08	0.53
1:A:271:LEU:HD23	1:A:274:MET:HE3	1.90	0.52
1:A:317:ASP:OD2	2:C:73:ARG:NH1	2.43	0.52
1:A:218:LEU:O	1:A:219:ARG:HB2	2.10	0.52
1:A:271:LEU:HD12	1:A:287:GLN:CG	2.41	0.50
1:A:272:GLU:CD	1:A:272:GLU:H	2.15	0.50
1:A:225:GLY:HA2	1:A:230:HIS:CD2	2.47	0.50
1:A:161:ILE:CD1	1:A:328:PHE:HB3	2.44	0.48
1:A:148:ASN:O	1:A:149:CYS:HB2	2.13	0.48
1:A:113:HIS:CD2	2:C:77:PRO:HB2	2.48	0.48
1:A:197:SER:HA	5:A:733:HOH:O	2.14	0.48
1:A:243:HIS:ND1	1:A:244:SER:N	2.61	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:9:ILE:HG22	1:A:78:GLN:OE1	2.15	0.47
1:A:120:GLN:HE22	1:A:151:LEU:H	1.62	0.47
1:A:201:SER:O	1:A:204:MET:HG2	2.15	0.47
1:A:10:SER:N	1:A:78:GLN:HE22	1.92	0.46
1:A:334:TYR:HB3	1:A:336:GLU:O	2.16	0.46
2:C:76:ILE:HG23	2:C:77:PRO:HD2	1.98	0.45
1:A:135:HIS:CE1	1:A:137:ASP:HB3	2.52	0.45
1:A:65:HIS:HE1	5:A:687:HOH:O	1.99	0.45
1:A:307:PRO:O	1:A:310:GLN:HG2	2.17	0.45
1:A:120:GLN:HE22	1:A:150:ASP:HA	1.82	0.44
1:A:80:PRO:HG2	1:A:86:PHE:CD1	2.53	0.44
1:A:106:THR:O	1:A:107:GLN:HG3	2.18	0.44
1:A:299:THR:OG1	1:A:302:GLU:HG3	2.18	0.44
1:A:97:GLN:NE2	1:A:147:SER:H	2.15	0.43
1:A:51:LEU:HD11	1:A:181:GLU:HA	2.00	0.43
1:A:78:GLN:NE2	5:A:635:HOH:O	2.50	0.43
1:A:279:ASN:O	1:A:283:ILE:HD12	2.17	0.43
1:A:77:ILE:HG21	1:A:351:ILE:HG12	2.00	0.43
1:A:187:TRP:CD1	1:A:224:PRO:HA	2.53	0.43
1:A:317:ASP:HB3	2:C:73:ARG:HH12	1.84	0.43
1:A:306:HIS:CG	1:A:307:PRO:HD2	2.54	0.43
1:A:88:GLU:HB2	1:A:90:TYR:CE2	2.54	0.42
1:A:261:LYS:HD3	5:A:846:HOH:O	2.18	0.42
1:A:58:ARG:HD3	1:A:332:ASP:OD2	2.20	0.42
1:A:136:ARG:HD3	1:A:158:LEU:O	2.20	0.41
1:A:111:ASP:OD2	1:A:115:GLN:NE2	2.49	0.41
1:A:253:SER:O	1:A:256:ALA:HB3	2.20	0.41
1:A:222:ILE:HG23	1:A:223:PHE:CD2	2.55	0.41
1:A:271:LEU:HD12	1:A:287:GLN:HG3	2.03	0.41
1:A:340:THR:HG23	5:A:832:HOH:O	2.21	0.41
1:A:1:MET:O	1:A:2:PRO:C	2.59	0.41
1:A:32:HIS:CD2	1:A:35:THR:OG1	2.60	0.41
1:A:46:PRO:HG2	1:A:47:PHE:CE2	2.56	0.41
1:A:306:HIS:HA	1:A:307:PRO:HD3	1.96	0.41
1:A:190:ALA:HA	1:A:207:TRP:CD1	2.56	0.41
1:A:112:ASP:HB2	5:A:672:HOH:O	2.20	0.41
1:A:1:MET:HG2	1:A:2:PRO:HD2	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	333/353 (94%)	309 (93%)	22 (7%)	2 (1%)	30	36
2	C	9/13 (69%)	7 (78%)	2 (22%)	0	100	100
All	All	342/366 (93%)	316 (92%)	24 (7%)	2 (1%)	30	36

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	198	ALA
1	A	2	PRO

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	297/319 (93%)	290 (98%)	7 (2%)	57	74
2	C	8/12 (67%)	7 (88%)	1 (12%)	6	6
All	All	305/331 (92%)	297 (97%)	8 (3%)	54	71

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

Mol	Chain	Res	Type
1	A	94	GLU
1	A	163	ASP
1	A	228	TYR
1	A	229	ARG

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Mol	Chain	Res	Type
1	A	272	GLU
1	A	320	GLU
1	A	350	GLU
2	C	73	ARG

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

Mol	Chain	Res	Type
1	A	8	ASN
1	A	32	HIS
1	A	65	HIS
1	A	85	ASN
1	A	97	GLN
1	A	101	HIS
1	A	120	GLN
1	A	230	HIS
1	A	246	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 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	$\# Z > 2$
4	ADP	A	500	3	22,29,29	0.94	0	27,45,45	1.80	4 (14%)

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
4	ADP	A	500	3	-	0/12/32/32	0/3/3/3

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
4	A	500	ADP	N3-C2-N1	-6.96	123.57	128.89
4	A	500	ADP	C2'-C1'-N9	-2.09	111.10	114.29
4	A	500	ADP	C2-N1-C6	2.12	122.55	118.77
4	A	500	ADP	C4'-O4'-C1'	3.20	113.24	109.72

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	337/353 (95%)	1.46	82 (24%) 1 1	10, 21, 47, 61	0
2	C	11/13 (84%)	7.28	11 (100%) 0 0	36, 44, 63, 69	0
All	All	348/366 (95%)	1.64	93 (26%) 1 1	10, 21, 48, 69	0

All (93) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	76	ILE	12.7
2	C	77	PRO	9.6
2	C	82	LEU	8.9
2	C	79	PRO	8.2
2	C	78	LYS	8.1
2	C	72	LYS	7.5
2	C	73	ARG	6.8
1	A	253	SER	6.1
2	C	74	GLY	6.0
2	C	75	ASN	5.8
1	A	112	ASP	5.3
1	A	312	TYR	4.9
1	A	197	SER	4.7
1	A	5	ILE	4.7
1	A	9	ILE	4.5
1	A	248	LEU	4.4
1	A	275	PHE	4.3
1	A	18	LEU	4.2
1	A	336	GLU	4.1
1	A	198	ALA	3.9
1	A	149	CYS	3.9
1	A	304	LEU	3.9
1	A	35	THR	3.9
1	A	246	ASN	3.8

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Mol	Chain	Res	Type	RSRZ
1	A	250	CYS	3.8
1	A	36	GLY	3.8
1	A	274	MET	3.7
1	A	252	GLU	3.7
1	A	256	ALA	3.7
1	A	334	TYR	3.6
1	A	11	SER	3.5
1	A	310	GLN	3.4
1	A	272	GLU	3.4
1	A	146	ASN	3.4
2	C	80	LEU	3.4
1	A	147	SER	3.4
1	A	228	TYR	3.3
1	A	285	LEU	3.2
1	A	34	PRO	3.1
1	A	314	ASP	3.1
1	A	338	LEU	3.1
1	A	180	VAL	3.1
1	A	303	ALA	3.0
1	A	145	ILE	3.0
2	C	81	ASN	3.0
1	A	4	ARG	3.0
1	A	311	THR	3.0
1	A	200	TYR	2.9
1	A	271	LEU	2.9
1	A	284	ASP	2.8
1	A	335	LYS	2.8
1	A	183	VAL	2.8
1	A	251	ILE	2.8
1	A	116	TYR	2.8
1	A	99	ASP	2.7
1	A	316	ASN	2.7
1	A	1	MET	2.7
1	A	254	PRO	2.7
1	A	163	ASP	2.7
1	A	98	THR	2.7
1	A	309	LEU	2.7
1	A	76	ASN	2.6
1	A	299	THR	2.6
1	A	269	ALA	2.6
1	A	108	MET	2.6
1	A	249	ARG	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	181	GLU	2.6
1	A	276	PRO	2.6
1	A	352	PHE	2.5
1	A	106	THR	2.5
1	A	244	SER	2.5
1	A	81	ASP	2.5
1	A	278	VAL	2.4
1	A	306	HIS	2.4
1	A	283	ILE	2.3
1	A	162	ILE	2.3
1	A	97	GLN	2.3
1	A	102	ARG	2.2
1	A	39	VAL	2.2
1	A	259	TYR	2.2
1	A	279	ASN	2.2
1	A	7	TYR	2.1
1	A	117	PHE	2.1
1	A	69	GLU	2.1
1	A	2	PRO	2.1
1	A	243	HIS	2.1
1	A	185	THR	2.1
1	A	15	LEU	2.0
1	A	323	PRO	2.0
1	A	255	ARG	2.0
1	A	307	PRO	2.0
1	A	151	LEU	2.0
1	A	300	ALA	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
4	ADP	A	500	27/27	0.90	0.20	0.08	11,15,18,21	0
3	MG	A	600	1/1	0.91	0.18	-	21,21,21,21	0

6.5 Other polymers [i](#)

There are no such residues in this entry.