



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:23 PM GMT

PDB ID : 3TTF
Title : Crystal structure of E. coli HypF with AMP and carbamoyl phosphate
Authors : Petkun, S.; Shi, R.; Li, Y.; Cygler, M.
Deposited on : 2011-09-14
Resolution : 1.92 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

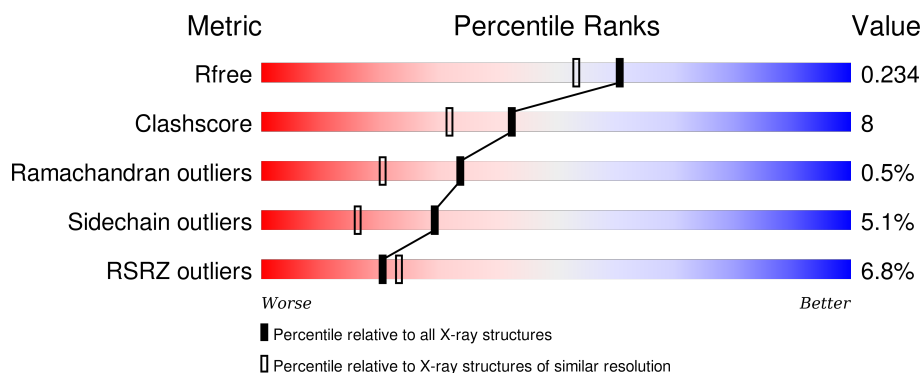
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


The reported resolution of this entry is 1.92 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	5832 (1.94-1.90)
Clashscore	102246	6540 (1.94-1.90)
Ramachandran outliers	100387	6464 (1.94-1.90)
Sidechain outliers	100360	6465 (1.94-1.90)
RSRZ outliers	91569	5846 (1.94-1.90)

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	657	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5380 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Transcriptional regulatory protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	646	Total	C	N	O	S	0	1	0
			4930	3114	886	893	37			

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	90	GLY	-	EXPRESSION TAG	UNP Q7ABC4
A	91	SER	-	EXPRESSION TAG	UNP Q7ABC4
A	571	ALA	GLN	CONFLICT	UNP Q7ABC4
A	572	ALA	GLN	CONFLICT	UNP Q7ABC4
A	573	ALA	GLN	CONFLICT	UNP Q7ABC4

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	3	Total	Zn	0	0
			3	3		

- 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 MONOPHOSPHATE (three-letter code: AMP) (formula: C₁₀H₁₄N₅O₇P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			23	10	5	7	1		

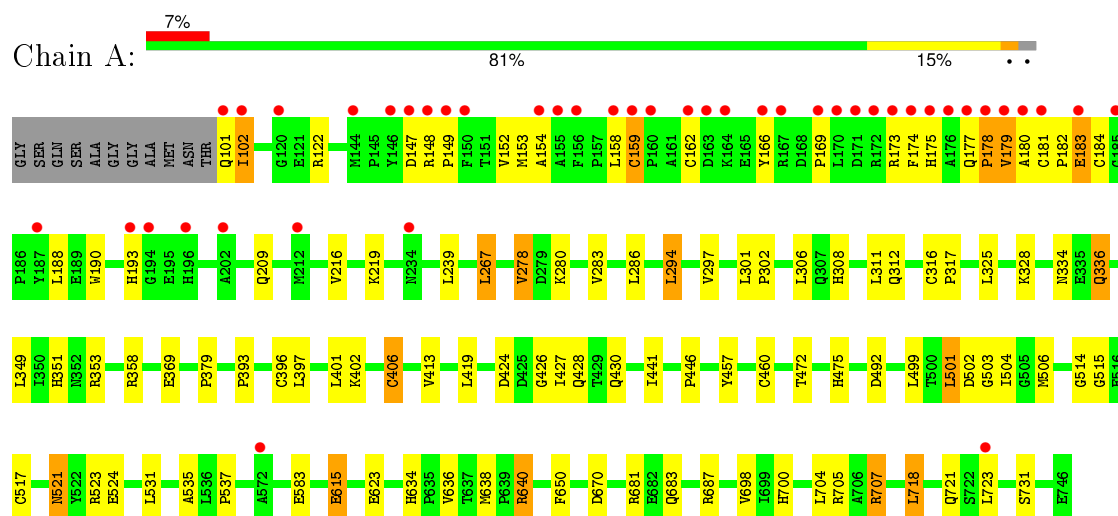
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	423	Total	O	0	0
			423	423		

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 ($\text{RSRZ} > 2$). 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: Transcriptional regulatory protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	46.36 Å 75.63 Å 200.67 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	100.33 – 1.92 42.09 – 1.92	Depositor EDS
% Data completeness (in resolution range)	90.5 (100.33-1.92) 90.5 (42.09-1.92)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.47 (at 1.92 Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.177 , 0.225 0.185 , 0.234	Depositor DCC
R_{free} test set	2532 reflections (5.36%)	DCC
Wilson B-factor (Å ²)	23.6	Xtriage
Anisotropy	0.662	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 56.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 49809 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5380	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.65% 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: AMP, ZN, MG

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	0.75	1/5055 (0.0%)	0.82	7/6902 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	406	CYS	CB-SG	-9.88	1.65	1.82

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	707	ARG	NE-CZ-NH2	-12.64	113.98	120.30
1	A	707	ARG	NE-CZ-NH1	10.92	125.76	120.30
1	A	278	VAL	CG1-CB-CG2	7.34	122.65	110.90
1	A	718	LEU	CB-CG-CD1	6.29	121.70	111.00
1	A	640	ARG	NE-CZ-NH2	-6.14	117.23	120.30
1	A	267	LEU	CA-CB-CG	6.10	129.32	115.30
1	A	681	ARG	NE-CZ-NH2	-5.45	117.58	120.30

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	4930	0	4842	81	0
2	A	3	0	0	0	0
3	A	1	0	0	0	0
4	A	23	0	12	2	0
5	A	423	0	0	19	0
All	All	5380	0	4854	81	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 8.

All (81) 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:297:VAL:HB	5:A:1067:HOH:O	0.97	1.13
1:A:683:GLN:HG3	5:A:790:HOH:O	1.44	1.13
1:A:517:CYS:HB2	5:A:886:HOH:O	0.89	1.05
1:A:506:MET:HE3	5:A:953:HOH:O	1.78	0.82
1:A:475:HIS:HE1	5:A:1069:HOH:O	1.68	0.75
1:A:152:VAL:HG12	1:A:153:MET:HE3	1.68	0.75
1:A:521:ASN:HD22	1:A:523:ARG:H	1.34	0.74
1:A:402:LYS:NZ	1:A:615:GLU:OE1	2.22	0.70
1:A:683:GLN:OE1	1:A:687:ARG:NH2	2.25	0.70
1:A:158:LEU:HD21	1:A:179:VAL:HG13	1.73	0.70
1:A:406:CYS:SG	1:A:413:VAL:CG2	2.80	0.69
1:A:153:MET:HA	1:A:153:MET:HE2	1.75	0.68
1:A:634:HIS:HD2	1:A:636:VAL:H	1.39	0.68
1:A:475:HIS:CE1	5:A:1069:HOH:O	2.42	0.68
1:A:351:HIS:HD2	1:A:353:ARG:H	1.42	0.65
1:A:506:MET:CE	5:A:953:HOH:O	2.39	0.63
1:A:517:CYS:SG	5:A:790:HOH:O	2.49	0.63
1:A:308:HIS:HD2	5:A:961:HOH:O	1.82	0.62
1:A:178:PRO:HD2	5:A:977:HOH:O	2.00	0.61
1:A:521:ASN:ND2	1:A:523:ARG:H	1.99	0.59
1:A:153:MET:HE3	1:A:353:ARG:HG3	1.85	0.59
1:A:148:ARG:NH1	1:A:177:GLN:O	2.36	0.58
1:A:302:PRO:HG3	1:A:311:LEU:HG	1.85	0.58
1:A:358:ARG:HD2	1:A:430:GLN:NE2	2.19	0.58
1:A:169:PRO:HA	1:A:174:PHE:CD1	2.39	0.58
1:A:492:ASP:HB3	5:A:939:HOH:O	2.04	0.57
1:A:504:ILE:HD13	5:A:807:HOH:O	2.04	0.57
1:A:102:ILE:HG12	1:A:441:ILE:HD13	1.88	0.56
1:A:153:MET:CE	1:A:353:ARG:HG3	2.37	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:122:ARG:NH2	1:A:162:CYS:SG	2.75	0.55
1:A:501:LEU:HB3	1:A:698:VAL:HB	1.88	0.55
1:A:308:HIS:HE1	5:A:87:HOH:O	1.90	0.55
1:A:280:LYS:HE3	1:A:294:LEU:O	2.06	0.54
1:A:521:ASN:ND2	1:A:524:GLU:H	2.05	0.54
1:A:325:LEU:HD23	1:A:328:LYS:HD2	1.90	0.53
1:A:670:ASP:OD1	1:A:707:ARG:HD3	2.08	0.53
1:A:159:CYS:HB3	1:A:184:CYS:HB3	1.91	0.53
1:A:369:GLU:OE1	1:A:705:ARG:NH2	2.43	0.52
1:A:396:CYS:HB2	1:A:406:CYS:HB3	1.92	0.51
1:A:179:VAL:HG22	1:A:180:ALA:H	1.73	0.51
1:A:166:TYR:CE1	1:A:175:HIS:HA	2.46	0.51
1:A:169:PRO:HA	1:A:174:PHE:HD1	1.75	0.51
1:A:153:MET:HG3	1:A:178:PRO:HA	1.92	0.50
1:A:173:ARG:HD3	1:A:179:VAL:HG23	1.92	0.50
1:A:638:MET:HG2	1:A:650:PHE:HB2	1.92	0.49
1:A:306:LEU:HD11	1:A:349:LEU:HD21	1.95	0.48
1:A:181:CYS:SG	1:A:183:GLU:HB2	2.54	0.48
1:A:698:VAL:HG22	4:A:747:AMP:N3	2.30	0.47
1:A:499:LEU:HG	1:A:501:LEU:HD13	1.97	0.47
1:A:535:ALA:O	1:A:537:PRO:HD3	2.14	0.46
1:A:406:CYS:SG	1:A:413:VAL:HG21	2.54	0.46
1:A:393:PRO:HB2	1:A:446:PRO:HA	1.98	0.46
1:A:670:ASP:OD1	1:A:707:ARG:CD	2.62	0.46
1:A:336:GLN:HA	1:A:336:GLN:NE2	2.29	0.46
1:A:316:CYS:HB2	1:A:317:PRO:HD2	1.98	0.45
1:A:193:HIS:HE1	5:A:945:HOH:O	1.99	0.45
1:A:634:HIS:HE1	1:A:670:ASP:OD2	1.99	0.45
1:A:308:HIS:O	1:A:312:GLN:HG3	2.16	0.45
1:A:369:GLU:OE2	1:A:700:HIS:HD2	2.00	0.45
1:A:283:VAL:HG12	1:A:286:LEU:HG	1.99	0.44
1:A:190:TRP:CZ2	1:A:209:GLN:HG3	2.53	0.44
1:A:181:CYS:HB2	1:A:182:PRO:CD	2.49	0.43
1:A:402:LYS:NZ	5:A:905:HOH:O	2.50	0.43
1:A:406:CYS:HB2	1:A:731:SER:HB2	2.01	0.43
1:A:501:LEU:HD12	1:A:515:GLY:HA2	2.00	0.43
1:A:179:VAL:HG22	1:A:180:ALA:N	2.33	0.43
1:A:147:ASP:CG	1:A:149:PRO:HD2	2.39	0.43
1:A:426:GLY:N	5:A:1072:HOH:O	2.49	0.42
1:A:424:ASP:O	1:A:427:ILE:HG22	2.18	0.42
1:A:457:TYR:HB2	1:A:460:CYS:SG	2.59	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:503:GLY:C	1:A:504:ILE:HD12	2.40	0.42
1:A:502:ASP:HA	1:A:698:VAL:HG23	2.02	0.42
1:A:634:HIS:CD2	1:A:636:VAL:H	2.28	0.42
1:A:623:GLU:OE1	4:A:747:AMP:N6	2.49	0.42
1:A:148:ARG:O	1:A:154:ALA:HB2	2.20	0.41
1:A:179:VAL:O	1:A:180:ALA:HB3	2.20	0.41
1:A:426:GLY:CA	5:A:1072:HOH:O	2.69	0.41
1:A:721:GLN:HB2	5:A:1035:HOH:O	2.21	0.41
1:A:219:LYS:NZ	1:A:334:ASN:HD21	2.17	0.41
1:A:148:ARG:HB3	1:A:149:PRO:HD3	2.03	0.40
1:A:428:GLN:NE2	5:A:1074:HOH:O	2.54	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	645/657 (98%)	616 (96%)	26 (4%)	3 (0%)	34 20

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	179	VAL
1	A	514	GLY
1	A	178	PRO

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	508/513 (99%)	482 (95%)	26 (5%)	29	16

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	101	GLN
1	A	102	ILE
1	A	159	CYS
1	A	183	GLU
1	A	188	LEU
1	A	216	VAL
1	A	239	LEU
1	A	267	LEU
1	A	278	VAL
1	A	294	LEU
1	A	301	LEU
1	A	336	GLN
1	A	379	PRO
1	A	397	LEU
1	A	401	LEU
1	A	419	LEU
1	A	472	THR
1	A	501	LEU
1	A	521	ASN
1	A	531	LEU
1	A	583	GLU
1	A	615	GLU
1	A	640	ARG
1	A	704	LEU
1	A	718	LEU
1	A	723	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	204	GLN
1	A	295	ASN
1	A	308	HIS
1	A	312	GLN

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Mol	Chain	Res	Type
1	A	334	ASN
1	A	336	GLN
1	A	351	HIS
1	A	430	GLN
1	A	451	HIS
1	A	521	ASN
1	A	587	ASN
1	A	634	HIS
1	A	700	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](#)

Of 5 ligands modelled in this entry, 4 are 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	AMP	A	747	2	20,25,25	1.46	2 (10%)	22,38,38	2.18	3 (13%)

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	AMP	A	747	2	-	0/6/26/26	0/3/3/3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	747	AMP	C5-C4	3.49	1.48	1.40
4	A	747	AMP	O4'-C1'	4.29	1.46	1.41

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	747	AMP	N3-C2-N1	-7.59	123.08	128.89
4	A	747	AMP	C4-C5-N7	-4.24	105.58	109.48
4	A	747	AMP	O4'-C1'-N9	2.73	113.82	108.10

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	747	AMP	2	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	646/657 (98%)	0.39	44 (6%) 20 23	14, 26, 60, 89	0

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	170	LEU	13.1
1	A	149	PRO	12.9
1	A	146	TYR	9.0
1	A	150	PHE	7.8
1	A	175	HIS	7.0
1	A	174	PHE	6.8
1	A	147	ASP	5.9
1	A	178	PRO	5.8
1	A	166	TYR	5.7
1	A	179	VAL	5.7
1	A	167	ARG	5.4
1	A	169	PRO	5.3
1	A	160	PRO	5.2
1	A	148	ARG	4.7
1	A	144	MET	4.6
1	A	156	PHE	4.2
1	A	183	GLU	3.9
1	A	176	ALA	3.9
1	A	159	CYS	3.6
1	A	101	GLN	3.1
1	A	172	ARG	3.0
1	A	177	GLN	2.8
1	A	162	CYS	2.7
1	A	154	ALA	2.7
1	A	193	HIS	2.6
1	A	181	CYS	2.6
1	A	171	ASP	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	180	ALA	2.5
1	A	155	ALA	2.5
1	A	158	LEU	2.5
1	A	202	ALA	2.4
1	A	173	ARG	2.3
1	A	164	LYS	2.2
1	A	723	LEU	2.2
1	A	212	MET	2.2
1	A	234	ASN	2.2
1	A	102	ILE	2.2
1	A	185	GLY	2.1
1	A	194	GLY	2.1
1	A	572	ALA	2.1
1	A	196	HIS	2.1
1	A	120	GLY	2.0
1	A	187	TYR	2.0
1	A	163	ASP	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(Å ²)	Q<0.9
4	AMP	A	747	23/23	0.88	0.16	0.97	31,45,46,48	0
2	ZN	A	1	1/1	0.99	0.05	-1.11	47,47,47,47	0
2	ZN	A	2	1/1	0.97	0.07	-2.14	70,70,70,70	0
2	ZN	A	3	1/1	0.99	0.05	-2.76	39,39,39,39	0
3	MG	A	4	1/1	0.99	0.04	-4.51	20,20,20,20	0

6.5 Other polymers

There are no such residues in this entry.