



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 09:06 AM GMT

PDB ID : 3H84
Title : Crystal structure of GET3
Authors : Hu, J.; Li, J.; Qian, X.; Sha, B.
Deposited on : 2009-04-28
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

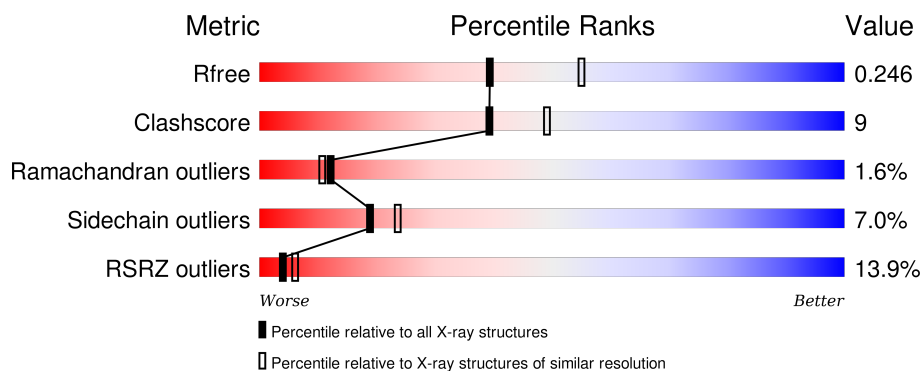
1 Overall quality at a glance

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	354	<div> <div>9%</div> <div>70%</div> <div>18%</div> <div>•</div> <div>8%</div> </div>
1	B	354	<div> <div>16%</div> <div>71%</div> <div>18%</div> <div>•</div> <div>9%</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
4	NA	A	359	-	-	-	X

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 5405 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 ATPase GET3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	326	Total	C	N	O	S	0	0	0
			2547	1600	428	502	17			
1	B	323	Total	C	N	O	S	0	0	0
			2526	1598	420	491	17			

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		

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

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

- Molecule 4 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	3	Total	Na	0	0
			3	3		

- Molecule 5 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	2	Total	Cl	0	0
			2	2		

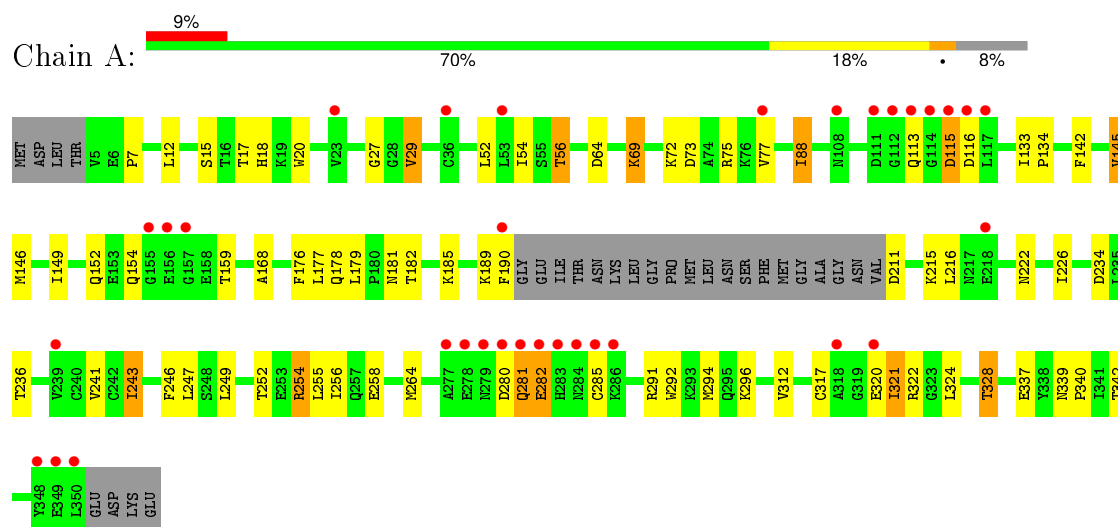
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	202	Total 202	O 202	0	0
6	B	122	Total 122	O 122	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: ATPase GET3



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 2	Depositor
Cell constants a, b, c, α , β , γ	219.40 Å 113.74 Å 48.37 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.30 49.41 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.3 (50.00-2.30) 99.3 (49.41-2.30)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.24 (at 2.29 Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.222 , 0.247 0.220 , 0.246	Depositor DCC
R_{free} test set	2763 reflections (5.33%)	DCC
Wilson B-factor (Å ²)	42.3	Xtriage
Anisotropy	0.258	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 48.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 54558 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5405	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.59% 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: NA, ZN, MG, CL

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	0.39	0/2588	0.58	0/3491
1	B	0.38	0/2569	0.55	0/3466
All	All	0.38	0/5157	0.57	0/6957

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	2547	0	2514	46	0
1	B	2526	0	2512	47	0
2	A	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	3	0	0	0	0
5	B	2	0	0	0	0
6	A	202	0	0	0	0
6	B	122	0	0	1	0
All	All	5405	0	5026	90	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 9.

All (90) 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:B:157:GLY:H	1:B:158:GLU:HA	1.26	0.97
1:A:17:THR:HG22	1:A:234:ASP:O	1.69	0.92
1:A:324:LEU:O	1:A:328:THR:HG23	1.71	0.90
1:A:189:LYS:CB	1:A:190:PHE:HA	2.02	0.88
1:B:10:HIS:HD2	1:B:335:ASN:HD21	1.22	0.88
1:A:189:LYS:HB2	1:A:190:PHE:HA	1.56	0.87
1:B:324:LEU:O	1:B:328:THR:HG23	1.84	0.77
1:A:291:ARG:O	1:A:294:MET:HG3	1.88	0.74
1:B:157:GLY:N	1:B:158:GLU:HA	1.94	0.69
1:B:196:LYS:HG2	1:B:197:LEU:H	1.65	0.62
1:A:29:VAL:HG13	1:A:241:VAL:HG12	1.82	0.62
1:B:15:SER:OG	1:B:18:HIS:HD2	1.84	0.61
1:B:7:PRO:HG2	1:B:337:GLU:HG2	1.83	0.61
1:B:10:HIS:CD2	1:B:335:ASN:HD21	2.11	0.60
1:B:146:MET:HE2	1:B:146:MET:HA	1.83	0.59
1:B:196:LYS:CG	1:B:197:LEU:H	2.16	0.58
1:B:317:CYS:SG	1:B:321:ILE:HD11	2.43	0.58
1:A:320:GLU:HG2	1:A:321:ILE:N	2.18	0.57
1:A:7:PRO:HG2	1:A:337:GLU:HG2	1.85	0.57
1:A:56:THR:CG2	1:A:168:ALA:HB2	2.34	0.57
1:B:259:LEU:HD13	1:B:266:VAL:HG11	1.89	0.55
1:A:142:PHE:HA	1:A:145:VAL:HG13	1.89	0.55
1:B:281:GLN:O	1:B:282:GLU:HB3	2.06	0.55
1:B:196:LYS:CG	1:B:197:LEU:N	2.70	0.54
1:A:189:LYS:CB	1:A:190:PHE:CA	2.83	0.54
1:B:299:LEU:HD22	1:B:313:LYS:HE2	1.89	0.54
1:A:317:CYS:SG	1:A:321:ILE:HD11	2.48	0.54
1:A:320:GLU:HG3	1:A:322:ARG:HG3	1.90	0.53
1:A:246:PHE:HE2	1:B:26:LYS:HE2	1.73	0.53
1:A:15:SER:OG	1:A:18:HIS:HD2	1.92	0.53
1:A:252:THR:O	1:A:256:ILE:HG12	2.10	0.52
1:A:29:VAL:HA	1:A:243:ILE:HD13	1.90	0.52
1:B:281:GLN:NE2	1:B:281:GLN:HA	2.25	0.52
1:A:20:TRP:HB2	1:A:236:THR:HG23	1.91	0.52
1:B:145:VAL:O	1:B:149:ILE:HG12	2.10	0.52
1:A:324:LEU:O	1:A:328:THR:CG2	2.54	0.51
1:A:146:MET:HA	1:A:146:MET:HE2	1.92	0.51
1:A:281:GLN:O	1:A:282:GLU:HB2	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:49:LYS:O	1:B:82:ASN:OD1	2.29	0.51
1:B:284:ASN:N	1:B:284:ASN:HD22	2.09	0.51
1:A:152:GLN:C	1:A:154:GLN:H	2.14	0.51
1:B:122:GLN:HE21	1:B:215:LYS:HB2	1.75	0.51
1:B:281:GLN:HE21	1:B:281:GLN:HA	1.76	0.50
1:B:179:LEU:N	1:B:180:PRO:HD2	2.26	0.50
1:A:246:PHE:CD2	1:B:27:GLY:HA2	2.47	0.50
1:B:157:GLY:N	1:B:158:GLU:CA	2.72	0.50
1:A:69:LYS:CD	1:A:69:LYS:H	2.24	0.49
1:B:238:PHE:HB3	1:B:266:VAL:HG12	1.95	0.48
1:B:29:VAL:HG13	1:B:241:VAL:HG12	1.94	0.48
1:A:292:TRP:NE1	1:A:296:LYS:HD2	2.29	0.48
1:B:281:GLN:O	1:B:282:GLU:CB	2.61	0.48
1:A:27:GLY:HA2	1:B:246:PHE:CD2	2.49	0.48
1:A:69:LYS:HD3	1:A:69:LYS:H	1.80	0.47
1:B:20:TRP:HB2	1:B:236:THR:HG23	1.95	0.47
1:A:254:ARG:NH2	1:A:258:GLU:OE1	2.48	0.47
1:B:77:VAL:HG22	1:B:80:MET:HG3	1.96	0.47
1:A:176:PHE:O	1:A:179:LEU:HB2	2.15	0.47
1:A:64:ASP:OD1	1:A:322:ARG:NH1	2.48	0.47
1:B:324:LEU:O	1:B:328:THR:CG2	2.58	0.46
1:B:345:LYS:O	1:B:349:GLU:HB3	2.16	0.45
1:A:222:ASN:O	1:A:226:ILE:HG12	2.16	0.45
1:A:56:THR:HG22	1:A:168:ALA:HB2	1.98	0.45
1:B:10:HIS:HD2	1:B:335:ASN:ND2	2.03	0.45
1:B:278:GLU:OE1	1:B:292:TRP:HZ3	2.00	0.44
1:A:146:MET:CE	1:A:149:ILE:HD12	2.48	0.44
1:B:201:LEU:O	1:B:205:MET:HG2	2.18	0.44
1:B:6:GLU:O	1:B:310:HIS:HD2	2.01	0.44
1:B:238:PHE:HB3	1:B:266:VAL:CG1	2.48	0.43
1:A:72:LYS:HA	1:A:88:ILE:HG22	2.01	0.43
1:A:178:GLN:O	1:A:182:THR:HG23	2.19	0.43
1:B:43:ALA:O	1:B:82:ASN:ND2	2.51	0.42
1:B:347:ILE:C	1:B:349:GLU:H	2.23	0.42
1:A:189:LYS:HB3	1:A:190:PHE:HA	1.96	0.42
1:A:339:ASN:HA	1:A:340:PRO:HD2	1.93	0.42
1:A:54:ILE:HD11	1:A:88:ILE:CD1	2.50	0.42
1:B:196:LYS:HE2	1:B:201:LEU:HD23	2.02	0.42
1:B:50:GLN:NE2	6:B:379:HOH:O	2.53	0.41
1:B:275:LEU:HD11	1:B:291:ARG:HD3	2.01	0.41
1:A:73:ASP:O	1:A:75:ARG:HD3	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:259:LEU:CD1	1:B:266:VAL:HG11	2.50	0.41
1:A:115:ASP:HA	1:A:116:ASP:HA	1.80	0.41
1:B:306:TYR:HB3	1:B:309:PHE:HB2	2.02	0.41
1:A:292:TRP:CE2	1:A:296:LYS:HD2	2.56	0.41
1:B:292:TRP:CH2	1:B:296:LYS:HD3	2.56	0.41
1:A:181:ASN:O	1:A:185:LYS:HG2	2.20	0.41
1:A:146:MET:HE1	1:A:149:ILE:HD12	2.02	0.40
1:A:133:ILE:HA	1:A:134:PRO:HD2	1.95	0.40
1:A:211:ASP:O	1:A:215:LYS:HG2	2.21	0.40
1:A:320:GLU:CG	1:A:322:ARG:HG3	2.51	0.40
1:B:344:GLY:O	1:B:348:TYR:HD2	2.03	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	322/354 (91%)	297 (92%)	19 (6%)	6 (2%)	10	8
1	B	319/354 (90%)	303 (95%)	12 (4%)	4 (1%)	15	15
All	All	641/708 (90%)	600 (94%)	31 (5%)	10 (2%)	12	11

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	281	GLN
1	A	282	GLU
1	B	283	HIS
1	A	113	GLN
1	A	280	ASP
1	B	282	GLU
1	B	285	CYS

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Mol	Chain	Res	Type
1	B	212	ILE
1	A	285	CYS
1	A	321	ILE

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	286/309 (93%)	265 (93%)	21 (7%)	17	22
1	B	284/309 (92%)	265 (93%)	19 (7%)	20	26
All	All	570/618 (92%)	530 (93%)	40 (7%)	19	23

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	LEU
1	A	29	VAL
1	A	52	LEU
1	A	56	THR
1	A	69	LYS
1	A	77	VAL
1	A	88	ILE
1	A	115	ASP
1	A	145	VAL
1	A	159	THR
1	A	177	LEU
1	A	216	LEU
1	A	243	ILE
1	A	247	LEU
1	A	249	LEU
1	A	254	ARG
1	A	255	LEU
1	A	264	MET
1	A	312	VAL
1	A	328	THR

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Mol	Chain	Res	Type
1	A	342	THR
1	B	12	LEU
1	B	29	VAL
1	B	32	THR
1	B	77	VAL
1	B	129	LEU
1	B	158	GLU
1	B	162	THR
1	B	183	LEU
1	B	218	GLU
1	B	255	LEU
1	B	264	MET
1	B	281	GLN
1	B	284	ASN
1	B	287	ARG
1	B	305	LEU
1	B	312	VAL
1	B	320	GLU
1	B	324	LEU
1	B	328	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	18	HIS
1	A	46	GLN
1	A	60	HIS
1	A	148	HIS
1	A	257	GLN
1	A	332	GLN
1	B	10	HIS
1	B	18	HIS
1	B	46	GLN
1	B	50	GLN
1	B	122	GLN
1	B	154	GLN
1	B	202	ASN
1	B	281	GLN
1	B	284	ASN
1	B	301	GLN
1	B	310	HIS
1	B	332	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](#)

Of 8 ligands modelled in this entry, 8 are monoatomic - leaving 0 for Mogul analysis.

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	326/354 (92%)	0.86	33 (10%) 9 13	39, 47, 68, 76	0
1	B	323/354 (91%)	1.07	57 (17%) 2 3	40, 49, 66, 70	0
All	All	649/708 (91%)	0.97	90 (13%) 4 6	39, 48, 67, 76	0

All (90) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	281	GLN	11.1
1	A	113	GLN	9.3
1	B	120	LEU	8.8
1	B	280	ASP	7.8
1	A	190	PHE	7.7
1	A	283	HIS	7.7
1	B	283	HIS	7.6
1	A	280	ASP	7.5
1	A	115	ASP	7.4
1	B	284	ASN	7.2
1	B	197	LEU	5.7
1	A	114	GLY	5.6
1	A	282	GLU	5.6
1	B	282	GLU	5.6
1	B	205	MET	5.4
1	A	156	GLU	5.3
1	B	204	PHE	5.3
1	A	284	ASN	5.2
1	B	195	ASN	5.2
1	A	157	GLY	4.9
1	B	126	LEU	4.9
1	B	190	PHE	4.7
1	B	121	LEU	4.5
1	B	192	GLU	4.5

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Mol	Chain	Res	Type	RSRZ
1	A	155	GLY	4.4
1	B	214	GLY	4.1
1	B	318	ALA	4.1
1	B	200	MET	4.1
1	A	108	ASN	4.1
1	B	207	ALA	4.1
1	B	158	GLU	4.1
1	A	279	ASN	4.0
1	B	286	LYS	3.9
1	A	318	ALA	3.9
1	A	112	GLY	3.7
1	B	91	SER	3.7
1	B	154	GLN	3.7
1	B	196	LYS	3.7
1	B	187	LEU	3.6
1	A	116	ASP	3.6
1	B	157	GLY	3.4
1	B	228	GLN	3.4
1	B	279	ASN	3.3
1	A	117	LEU	3.3
1	B	208	GLY	3.3
1	A	320	GLU	3.2
1	B	156	GLU	3.2
1	B	119	SER	3.2
1	B	191	GLY	3.1
1	B	155	GLY	2.9
1	A	277	ALA	2.9
1	B	117	LEU	2.9
1	B	183	LEU	2.9
1	B	186	LEU	2.9
1	B	159	THR	2.9
1	A	111	ASP	2.8
1	B	218	GLU	2.8
1	A	286	LYS	2.7
1	B	150	LYS	2.7
1	B	125	ALA	2.6
1	A	278	GLU	2.6
1	A	349	GLU	2.5
1	B	122	GLN	2.5
1	B	291	ARG	2.5
1	B	320	GLU	2.5
1	B	226	ILE	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	133	ILE	2.4
1	A	281	GLN	2.4
1	B	199	PRO	2.3
1	B	194	THR	2.3
1	B	193	ILE	2.3
1	B	201	LEU	2.3
1	B	209	ASN	2.3
1	A	350	LEU	2.3
1	B	216	LEU	2.3
1	B	225	THR	2.3
1	A	218	GLU	2.3
1	A	348	TYR	2.2
1	B	212	ILE	2.2
1	A	23	VAL	2.2
1	A	36	CYS	2.2
1	A	77	VAL	2.2
1	A	285	CYS	2.2
1	B	293	LYS	2.1
1	B	348	TYR	2.1
1	B	129	LEU	2.1
1	B	285	CYS	2.1
1	A	239	VAL	2.1
1	A	53	LEU	2.0
1	B	143	MET	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	NA	A	359	1/1	0.95	0.22	2.90	53,53,53,53	0
5	CL	B	357	1/1	0.89	0.16	0.83	67,67,67,67	0
4	NA	A	357	1/1	0.90	0.21	0.03	58,58,58,58	0
4	NA	A	358	1/1	0.87	0.17	-0.34	46,46,46,46	0
5	CL	B	356	1/1	0.93	0.08	-2.82	73,73,73,73	0
2	ZN	A	355	1/1	0.98	0.03	-4.39	54,54,54,54	0
3	MG	B	355	1/1	0.93	0.39	-	40,40,40,40	0
3	MG	A	356	1/1	0.98	0.41	-	30,30,30,30	0

6.5 Other polymers

There are no such residues in this entry.