Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 02:42 AM GMT

PDB ID	:	2IBI
Title	:	Covalent Ubiquitin-USP2 Complex
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Deposited on	:	2006-09-11
Resolution	:	2.20 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

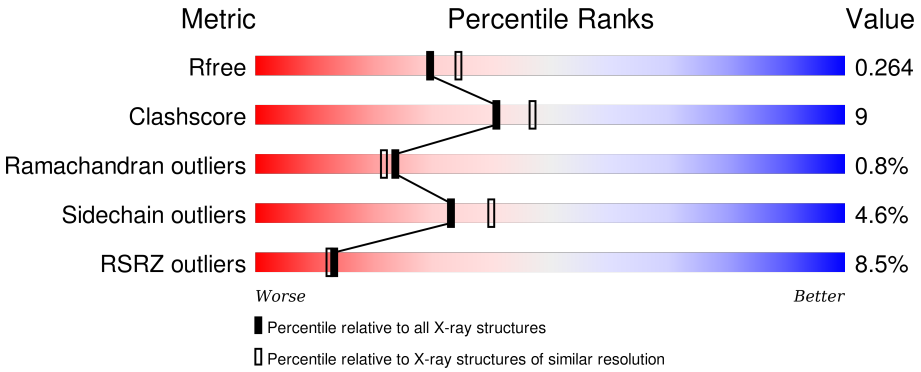
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	3774 (2.20-2.20)
Clashscore	102246	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (2.20-2.20)
RSRZ outliers	91569	3781 (2.20-2.20)

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 <=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	374	9%, 68%, 20%, •, 10%
2	B	75	3%, 85%, 12%, •

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 3422 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 Ubiquitin carboxyl-terminal hydrolase 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	335	2710	1699	487	504	20	0	1	0

There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	232	MET	-	INITIATING METHIONINE	UNP O75604
A	233	GLY	-	CLONING ARTIFACT	UNP O75604
A	234	SER	-	CLONING ARTIFACT	UNP O75604
A	235	SER	-	CLONING ARTIFACT	UNP O75604
A	236	HIS	-	EXPRESSION TAG	UNP O75604
A	237	HIS	-	EXPRESSION TAG	UNP O75604
A	238	HIS	-	EXPRESSION TAG	UNP O75604
A	239	HIS	-	EXPRESSION TAG	UNP O75604
A	240	HIS	-	EXPRESSION TAG	UNP O75604
A	241	HIS	-	EXPRESSION TAG	UNP O75604
A	242	SER	-	CLONING ARTIFACT	UNP O75604
A	243	SER	-	CLONING ARTIFACT	UNP O75604
A	244	GLY	-	CLONING ARTIFACT	UNP O75604
A	245	LEU	-	CLONING ARTIFACT	UNP O75604
A	246	VAL	-	CLONING ARTIFACT	UNP O75604
A	247	PRO	-	CLONING ARTIFACT	UNP O75604
A	248	ARG	-	CLONING ARTIFACT	UNP O75604
A	249	GLY	-	CLONING ARTIFACT	UNP O75604
A	250	SER	-	CLONING ARTIFACT	UNP O75604
A	472	ALA	GLU	ENGINEERED	UNP O75604
A	473	ALA	LYS	ENGINEERED	UNP O75604

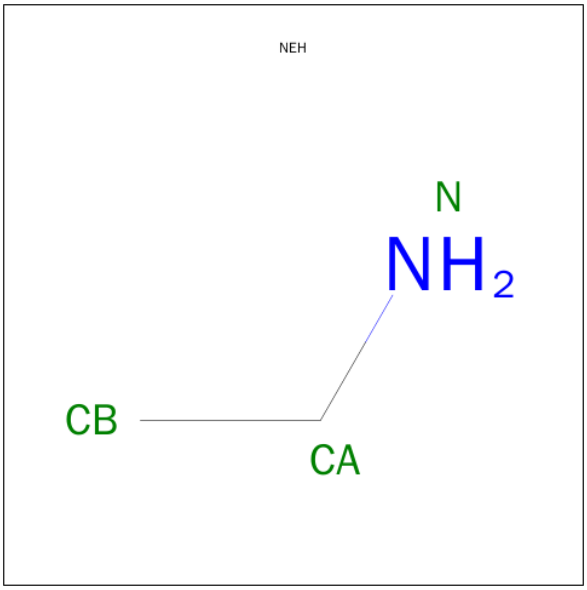
- Molecule 2 is a protein called Ubiquitin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	75	Total	C	N	O	S	0	1	0
			608	382	108	117	1			

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 4 is ETHANAMINE (three-letter code: NEH) (formula: C_2H_7N).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	N	0	0
			3	2	1		

- Molecule 5 is water.

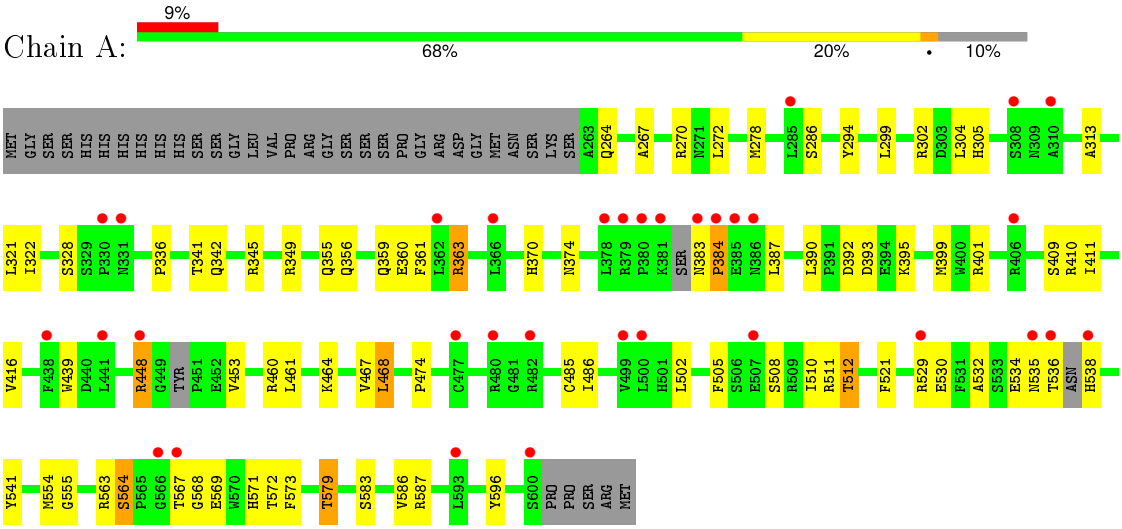
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	80	Total	O	0	0
			80	80		
5	B	20	Total	O	0	0
			20	20		

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.

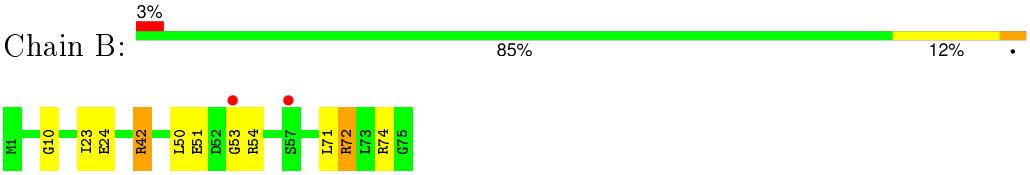
- Molecule 1: Ubiquitin carboxyl-terminal hydrolase 2

Chain A:



- Molecule 2: Ubiquitin

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	102.84Å 54.40Å 75.02Å 90.00° 107.66° 90.00°	Depositor
Resolution (Å)	28.01 – 2.20 28.00 – 2.20	Depositor EDS
% Data completeness (in resolution range)	96.3 (28.01-2.20) 96.3 (28.00-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.16	Depositor
$< I/\sigma(I) >$ [1]	1.76 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.190 , 0.259 0.193 , 0.264	Depositor DCC
R_{free} test set	1002 reflections (5.42%)	DCC
Wilson B-factor (Å^2)	28.8	Xtriage
Anisotropy	0.117	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 52.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Outliers	0 of 19493 reflections	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	3422	wwPDB-VP
Average B, all atoms (Å^2)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.49% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: ZN, NEH

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	0.66	0/2768	0.69	0/3736
2	B	0.64	0/614	0.87	4/825 (0.5%)
All	All	0.65	0/3382	0.73	4/4561 (0.1%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	42	ARG	NE-CZ-NH2	-9.78	115.41	120.30
2	B	72	ARG	NE-CZ-NH1	6.31	123.45	120.30
2	B	42	ARG	NE-CZ-NH1	6.12	123.36	120.30
2	B	72	ARG	NE-CZ-NH2	-5.70	117.45	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	2710	0	2642	46	0
2	B	608	0	638	14	0
3	A	1	0	0	0	0
4	B	3	0	4	0	0
5	A	80	0	0	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	20	0	0	3	0
All	All	3422	0	3284	57	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 (57) 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:521:PHE:O	1:A:587:ARG:HD3	1.74	0.88
1:A:579:THR:HG22	5:A:33:HOH:O	1.77	0.85
2:B:42:ARG:HD2	2:B:72:ARG:CG	2.16	0.75
2:B:42:ARG:HD2	2:B:72:ARG:HG3	1.69	0.73
1:A:572:THR:HB	1:A:579:THR:HG23	1.76	0.67
1:A:387:LEU:HD12	1:A:399:MET:CE	2.25	0.66
2:B:72:ARG:NH2	5:B:95:HOH:O	2.29	0.59
1:A:341:THR:O	1:A:345:ARG:HD3	2.03	0.58
1:A:294:TYR:CD1	1:A:410:ARG:HG2	2.39	0.58
1:A:554[A]:MET:HA	2:B:74[A]:ARG:HD2	1.86	0.57
1:A:387:LEU:HB3	1:A:390:LEU:HD12	1.89	0.55
2:B:42:ARG:HD3	2:B:71:LEU:O	2.07	0.55
1:A:360:GLU:OE1	2:B:72:ARG:HD3	2.06	0.55
1:A:510:ILE:HG22	1:A:511:ARG:HB2	1.89	0.55
2:B:10:GLY:CA	5:B:86:HOH:O	2.55	0.54
2:B:23:ILE:HG12	2:B:50:LEU:HB3	1.90	0.53
1:A:532:ALA:CB	1:A:535:ASN:HB2	2.39	0.53
1:A:541:TYR:HB3	1:A:596:TYR:HB3	1.89	0.53
1:A:363:ARG:NH1	5:A:75:HOH:O	2.40	0.52
1:A:270:ARG:HD2	1:A:272:LEU:HD23	1.92	0.52
1:A:305:HIS:HE1	1:A:374:ASN:O	1.92	0.51
1:A:349:ARG:HD2	5:A:40:HOH:O	2.10	0.51
1:A:286:SER:HA	1:A:322:ILE:HD12	1.92	0.51
1:A:563:ARG:HD2	1:A:568:GLY:HA2	1.92	0.51
1:A:416:VAL:HG22	1:A:439:TRP:CE2	2.45	0.51
2:B:23:ILE:HD12	2:B:54:ARG:O	2.12	0.49
1:A:567:THR:HG22	1:A:569:GLU:OE2	2.12	0.49
1:A:341:THR:O	1:A:345:ARG:CD	2.61	0.49
1:A:355:GLN:OE1	1:A:555:GLY:HA2	2.14	0.48
1:A:532:ALA:HB1	1:A:535:ASN:HB2	1.96	0.48
1:A:564:SER:HB3	1:A:567:THR:HB	1.96	0.47
1:A:321:LEU:HD11	1:A:336:PRO:HB3	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:535:ASN:ND2	1:A:536:THR:H	2.13	0.46
1:A:583:SER:HA	1:A:586:VAL:HG13	1.96	0.46
1:A:278:MET:HG3	1:A:361:PHE:CZ	2.51	0.46
1:A:512:THR:HB	5:A:101:HOH:O	2.15	0.46
1:A:461:LEU:HD12	1:A:464:LYS:HD2	1.98	0.45
1:A:474:PRO:HG2	1:A:485:CYS:SG	2.56	0.45
1:A:313:ALA:HB3	1:A:342:GLN:OE1	2.15	0.45
1:A:267:ALA:HB2	1:A:328:SER:OG	2.17	0.45
1:A:467:VAL:HG22	1:A:486:ILE:HD12	1.99	0.45
1:A:460:ARG:HD3	1:A:530:GLU:OE2	2.16	0.45
1:A:359:GLN:OE1	2:B:42:ARG:NH2	2.42	0.45
1:A:468:LEU:HB3	1:A:474:PRO:HD3	2.00	0.44
2:B:24:GLU:OE2	2:B:53:GLY:N	2.51	0.44
1:A:387:LEU:HD12	1:A:399:MET:HE2	2.00	0.43
1:A:383:ASN:HA	1:A:384:PRO:HD3	1.87	0.43
2:B:42:ARG:HD2	2:B:72:ARG:HG2	1.96	0.43
1:A:374:ASN:HA	1:A:409:SER:HB3	2.00	0.43
2:B:51:GLU:O	2:B:54:ARG:HG2	2.19	0.42
1:A:387:LEU:CD1	1:A:399:MET:CE	2.97	0.41
2:B:10:GLY:N	5:B:86:HOH:O	2.53	0.41
1:A:571:HIS:HB2	1:A:573:PHE:CE2	2.56	0.41
1:A:583:SER:HA	1:A:586:VAL:CG1	2.51	0.41
1:A:370:HIS:HA	1:A:411:ILE:HG13	2.02	0.40
1:A:392:ASP:HA	1:A:395:LYS:HG2	2.02	0.40
1:A:299:LEU:HD23	1:A:302:ARG:HE	1.86	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	328/374 (88%)	312 (95%)	13 (4%)	3 (1%)	21 19

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
2	B	74/75 (99%)	71 (96%)	3 (4%)	0	100 100
All	All	402/449 (90%)	383 (95%)	16 (4%)	3 (1%)	24 25

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	508	SER
1	A	448	ARG
1	A	384	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	302/335 (90%)	285 (94%)	17 (6%)	26 29
2	B	69/68 (102%)	69 (100%)	0	100 100
All	All	371/403 (92%)	354 (95%)	17 (5%)	33 40

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	264	GLN
1	A	304	LEU
1	A	356	GLN
1	A	363	ARG
1	A	393	ASP
1	A	401	ARG
1	A	448	ARG
1	A	453	VAL
1	A	468	LEU
1	A	502	LEU
1	A	505	PHE
1	A	512	THR
1	A	529	ARG

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Mol	Chain	Res	Type
1	A	534	GLU
1	A	538	HIS
1	A	564	SER
1	A	579	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	305	HIS
1	A	535	ASN
1	A	548	ASN
2	B	31	GLN
2	B	60	ASN

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 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	NEH	B	76	1,2	2,2,2	0.41	0	0,1,1	0.00	-

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	NEH	B	76	1,2	-	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, 95^{th} 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	335/374 (89%)	0.42	33 (9%) 9 8	32, 43, 74, 111	0
2	B	75/75 (100%)	0.14	2 (2%) 58 57	32, 42, 56, 61	0
All	All	410/449 (91%)	0.37	35 (8%) 13 12	32, 43, 68, 111	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	384	PRO	6.9
1	A	383	ASN	6.8
1	A	538	HIS	6.7
1	A	380	PRO	6.0
1	A	381	LYS	5.5
1	A	448	ARG	5.0
1	A	477	CYS	4.5
1	A	378	LEU	4.3
1	A	480	ARG	4.2
1	A	536	THR	4.2
2	B	53	GLY	3.9
1	A	535	ASN	3.6
1	A	566	GLY	3.5
1	A	308	SER	3.3
1	A	386	ASN	3.2
1	A	499	VAL	3.1
1	A	441	LEU	3.0
1	A	567	THR	2.8
1	A	366	LEU	2.8
1	A	330	PRO	2.7
1	A	362	LEU	2.7
1	A	593	LEU	2.7
2	B	57	SER	2.5
1	A	310	ALA	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	600	SER	2.5
1	A	385	GLU	2.5
1	A	500	LEU	2.5
1	A	379	ARG	2.4
1	A	331	ASN	2.3
1	A	529	ARG	2.3
1	A	482	ARG	2.2
1	A	507	GLU	2.2
1	A	406	ARG	2.1
1	A	438	PHE	2.1
1	A	285	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. 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, 95^{th} 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	NEH	B	76	3/3	0.94	0.14	-0.02	35,35,36,37	0
3	ZN	A	1	1/1	1.00	0.04	-2.78	45,45,45,45	0

6.5 Other polymers (i)

There are no such residues in this entry.