



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 19, 2026 – 12:26 PM UTC

PDB ID : 8AIF / pdb_00008aif
Title : TapA acts as specific chaperone in TasA non-amyloid filament formation
Authors : Roske, Y.; Heinemann, U.
Deposited on : 2022-07-26
Resolution : 1.07 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

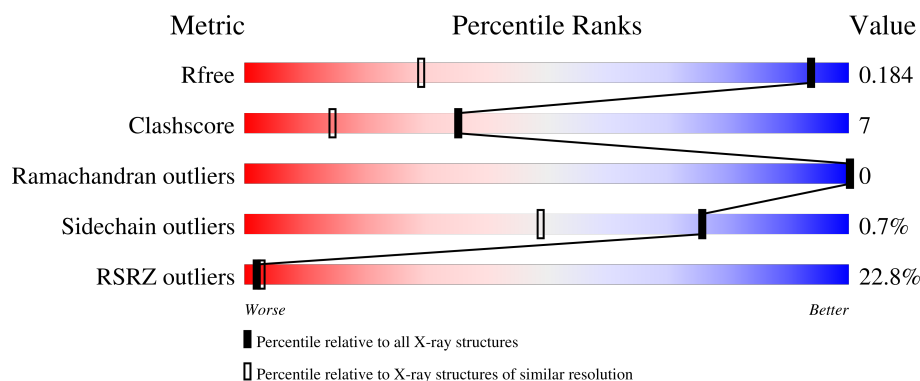
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.07 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}	180053	1687 (1.10-1.06)
Clashscore	190562	1727 (1.10-1.06)
Ramachandran outliers	187476	1679 (1.10-1.06)
Sidechain outliers	187428	1676 (1.10-1.06)
RSRZ outliers	180081	1687 (1.10-1.06)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	120	19% 87% 12% .
1	B	120	22% 94% 6%
1	C	120	26% 86% 12% ..

2 Entry composition [i](#)

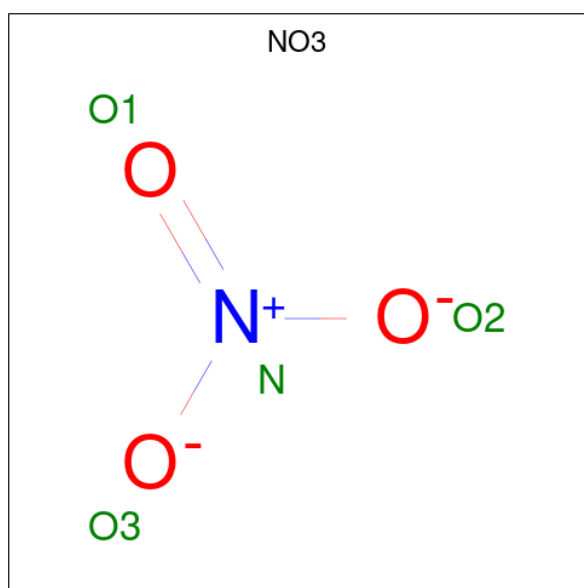
There are 4 unique types of molecules in this entry. The entry contains 3719 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 YqxM protein required for localization of TasA to extracellular matrix.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	118	Total	C	N	O	S	0	7	0
			984	642	164	174	4			
1	B	120	Total	C	N	O	S	0	5	0
			992	636	166	186	4			
1	C	118	Total	C	N	O	S	0	8	0
			972	632	159	177	4			

- Molecule 2 is NITRATE ION (CCD ID: NO3) (formula: NO₃).



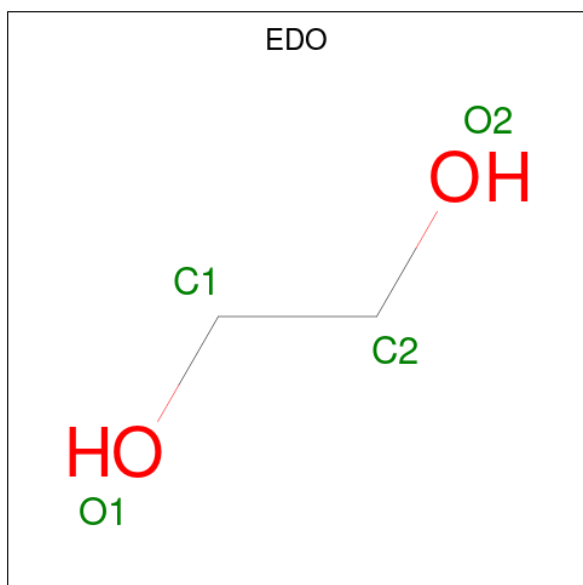
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	N	O	0	0
			4	1	3		
2	B	1	Total	N	O	0	0
			4	1	3		
2	B	1	Total	N	O	0	0
			4	1	3		

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

- Molecule 3 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



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

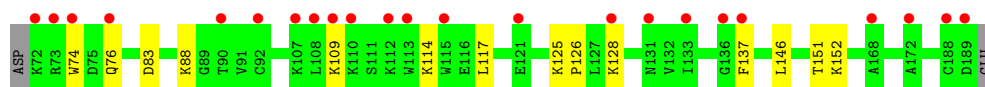
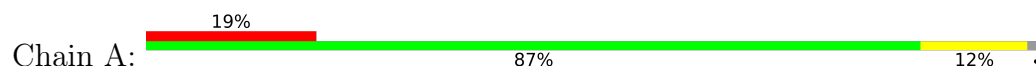
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	257	Total	O	0	0
			257	257		
4	B	269	Total	O	0	0
			269	269		
4	C	217	Total	O	0	0
			217	217		

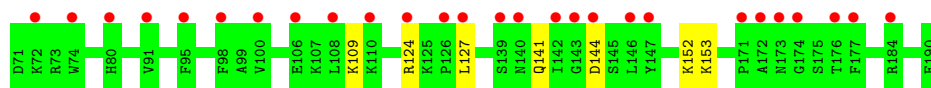
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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.

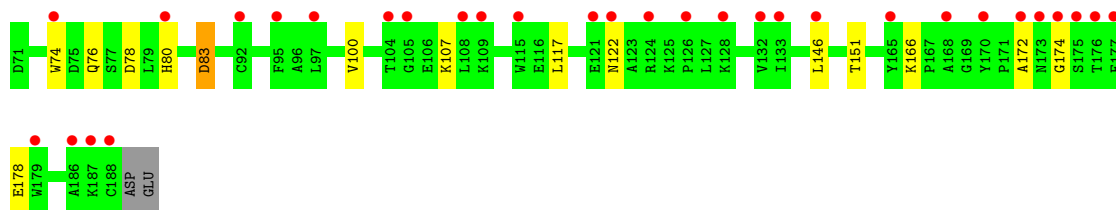
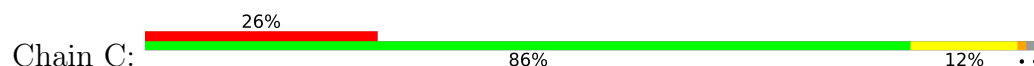
- Molecule 1: YqxM protein required for localization of TasA to extracellular matrix



- Molecule 1: YqxM protein required for localization of TasA to extracellular matrix



- Molecule 1: YqxM protein required for localization of TasA to extracellular matrix



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	39.20Å 52.94Å 167.76Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.44 – 1.07 38.44 – 1.07	Depositor EDS
% Data completeness (in resolution range)	98.8 (38.44-1.07) 98.9 (38.44-1.07)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.08 (at 1.07Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.148 , 0.177 0.158 , 0.184	Depositor DCC
R_{free} test set	2100 reflections (1.37%)	wwPDB-VP
Wilson B-factor (Å ²)	13.1	Xtriage
Anisotropy	0.173	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 100.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3719	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.54% 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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NO3, EDO

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	0.85	0/1032	0.91	0/1384
1	B	0.86	0/1031	0.95	0/1384
1	C	0.94	0/1023	1.03	1/1379 (0.1%)
All	All	0.88	0/3086	0.96	1/4147 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	83	ASP	CA-CB-CG	8.22	120.82	112.60

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	984	0	1025	21	0
1	B	992	0	1001	7	1
1	C	972	0	984	18	0
2	A	4	0	0	0	0
2	B	12	0	0	2	0
2	C	4	0	0	0	0
3	B	4	0	6	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	4	0	6	2	0
4	A	257	0	0	7	0
4	B	269	0	0	4	5
4	C	217	0	0	10	2
All	All	3719	0	3022	44	5

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 7.

All (44) 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:114[A]:LYS:HG2	1:A:137[A]:PHE:CE1	1.73	1.22
1:C:80:HIS:CE1	4:C:303:HOH:O	1.95	1.19
1:A:128:LYS:NZ	4:A:501:HOH:O	1.69	1.18
1:C:83:ASP:OD1	4:C:301:HOH:O	1.73	1.06
1:A:83:ASP:OD2	4:A:502:HOH:O	1.80	0.98
1:A:114[A]:LYS:CG	1:A:137[A]:PHE:CE1	2.48	0.96
1:C:80:HIS:HE1	4:C:303:HOH:O	1.39	0.93
1:B:124:ARG:HB2	4:B:328:HOH:O	1.69	0.92
1:A:114[A]:LYS:CG	1:A:137[A]:PHE:HE1	1.91	0.78
1:C:100[A]:VAL:CG1	1:C:146:LEU:HD11	2.17	0.73
1:C:78:ASP:OD2	4:C:303:HOH:O	2.07	0.73
1:B:152:LYS:NZ	4:B:301:HOH:O	2.22	0.73
1:A:88:LYS:HG3	4:A:532:HOH:O	1.90	0.70
1:B:124:ARG:HD3	2:B:204:NO3:O1	1.91	0.69
1:A:146[B]:LEU:HD11	1:C:146:LEU:HD13	1.75	0.68
1:A:114[B]:LYS:HG3	4:A:511:HOH:O	1.99	0.62
1:C:83:ASP:HA	4:C:301:HOH:O	2.01	0.61
1:C:83:ASP:HB3	1:C:100[B]:VAL:HG22	1.82	0.61
1:B:141:GLN:NE2	1:B:144:ASP:O	2.33	0.61
1:B:124:ARG:CB	4:B:328:HOH:O	2.40	0.60
1:A:88:LYS:CG	4:A:532:HOH:O	2.46	0.59
1:A:152:LYS:HE3	4:A:596:HOH:O	2.05	0.56
1:A:114[A]:LYS:HZ2	1:A:137[A]:PHE:HE1	1.48	0.56
1:A:114[A]:LYS:HG3	1:A:137[A]:PHE:HE1	1.73	0.53
1:C:172:ALA:HA	3:C:202:EDO:H22	1.91	0.52
1:B:127:LEU:HG	2:B:203:NO3:O1	2.10	0.51
1:A:88:LYS:HE3	4:C:458:HOH:O	2.13	0.49
1:A:114[A]:LYS:NZ	4:A:504:HOH:O	2.46	0.49
1:A:114[A]:LYS:NZ	1:A:137[A]:PHE:CE1	2.76	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:125:LYS:HA	1:A:126:PRO:C	2.39	0.48
1:C:122:ASN:HB2	4:C:400:HOH:O	2.14	0.48
1:A:117:LEU:HD22	1:A:151:THR:CG2	2.45	0.47
1:C:74:TRP:CE2	1:C:76:GLN:HA	2.51	0.46
1:C:166[A]:LYS:NZ	4:C:311:HOH:O	2.48	0.46
1:C:117:LEU:HD22	1:C:151:THR:CG2	2.46	0.45
1:C:174:GLY:N	3:C:202:EDO:O2	2.49	0.45
1:A:74:TRP:CE2	1:A:76:GLN:HA	2.52	0.45
1:A:146[B]:LEU:CD1	1:C:146:LEU:HD13	2.46	0.44
1:A:114[A]:LYS:NZ	1:A:137[A]:PHE:HE1	2.13	0.43
1:C:83:ASP:HB3	1:C:100[B]:VAL:CG2	2.46	0.43
1:C:178:GLU:OE2	4:C:304:HOH:O	2.21	0.43
1:A:126:PRO:O	1:A:128:LYS:HD2	2.19	0.43
1:B:109:LYS:HE2	4:B:512:HOH:O	2.19	0.42
1:C:80:HIS:NE2	4:C:303:HOH:O	2.28	0.41

All (5) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:B:393:HOH:O	4:B:564:HOH:O[3_544]	1.85	0.35
4:B:327:HOH:O	4:C:313:HOH:O[3_644]	1.87	0.33
4:B:304:HOH:O	4:B:375:HOH:O[1_655]	1.91	0.29
4:B:448:HOH:O	4:C:512:HOH:O[3_544]	1.98	0.22
1:B:153:LYS:NZ	4:B:307:HOH:O[3_544]	2.18	0.02

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	123/120 (102%)	119 (97%)	4 (3%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	123/120 (102%)	117 (95%)	6 (5%)	0	100	100
1	C	124/120 (103%)	120 (97%)	4 (3%)	0	100	100
All	All	370/360 (103%)	356 (96%)	14 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	109/105 (104%)	107 (98%)	2 (2%)	51	15
1	B	110/105 (105%)	110 (100%)	0	100	100
1	C	107/105 (102%)	105 (98%)	2 (2%)	50	13
All	All	326/315 (104%)	322 (99%)	4 (1%)	76	26

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	109[A]	LYS
1	A	109[B]	LYS
1	C	107[A]	LYS
1	C	107[B]	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

7 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	$\# Z > 2$
3	EDO	B	201	-	3,3,3	0.23	0	2,2,2	0.19	0
2	NO3	B	203	-	1,3,3	0.56	0	0,3,3	-	-
2	NO3	C	201	-	1,3,3	0.17	0	0,3,3	-	-
2	NO3	A	401	-	1,3,3	0.01	0	0,3,3	-	-
2	NO3	B	202	-	1,3,3	0.14	0	0,3,3	-	-
3	EDO	C	202	-	3,3,3	0.32	0	2,2,2	0.24	0
2	NO3	B	204	-	1,3,3	0.10	0	0,3,3	-	-

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
3	EDO	C	202	-	-	1/1/1/1	-
3	EDO	B	201	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	202	EDO	O1-C1-C2-O2

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	203	NO3	1	0
3	C	202	EDO	2	0
2	B	204	NO3	1	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	118/120 (98%)	1.38	23 (19%) 3 4	7, 14, 30, 58	7 (5%)
1	B	120/120 (100%)	1.43	27 (22%) 2 3	8, 14, 31, 35	5 (4%)
1	C	118/120 (98%)	1.62	31 (26%) 1 2	8, 16, 35, 46	8 (6%)
All	All	356/360 (98%)	1.47	81 (22%) 2 3	7, 14, 32, 58	20 (5%)

All (81) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	174	GLY	4.9
1	B	146	LEU	4.2
1	A	128	LYS	4.2
1	C	188	CYS	4.0
1	B	173	ASN	3.9
1	C	108	LEU	3.9
1	C	121	GLU	3.9
1	B	172	ALA	3.8
1	C	132	VAL	3.6
1	B	95	PHE	3.6
1	A	172	ALA	3.5
1	A	189	ASP	3.5
1	A	74	TRP	3.3
1	A	109[A]	LYS	3.2
1	B	80	HIS	3.2
1	A	137[A]	PHE	3.1
1	C	174	GLY	3.1
1	A	72	LYS	3.0
1	C	172	ALA	3.0
1	C	95	PHE	3.0
1	A	168	ALA	3.0
1	B	144	ASP	3.0
1	B	143	GLY	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	133	ILE	2.8
1	B	106	GLU	2.8
1	A	115	TRP	2.8
1	A	108	LEU	2.8
1	A	188	CYS	2.8
1	C	187	LYS	2.7
1	C	104	THR	2.7
1	C	74	TRP	2.7
1	C	173	ASN	2.6
1	C	170	TYR	2.6
1	A	73	ARG	2.6
1	C	92	CYS	2.6
1	C	133	ILE	2.5
1	A	121	GLU	2.5
1	C	165	TYR	2.5
1	C	168	ALA	2.5
1	B	98	PHE	2.5
1	B	124	ARG	2.4
1	C	175[A]	SER	2.4
1	C	126	PRO	2.4
1	C	176	THR	2.4
1	A	136	GLY	2.4
1	B	100	VAL	2.4
1	B	142	ILE	2.4
1	B	126	PRO	2.3
1	B	74	TRP	2.3
1	C	115	TRP	2.3
1	A	92	CYS	2.3
1	C	109	LYS	2.3
1	A	113	TRP	2.3
1	C	122	ASN	2.3
1	A	112	LYS	2.3
1	B	171	PRO	2.3
1	B	108	LEU	2.3
1	B	110	LYS	2.2
1	C	105	GLY	2.2
1	C	124	ARG	2.2
1	C	177	PHE	2.2
1	B	140	ASN	2.2
1	C	128	LYS	2.2
1	B	176	THR	2.2
1	B	147	TYR	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	90	THR	2.1
1	C	97	LEU	2.1
1	B	72[A]	LYS	2.1
1	B	91	VAL	2.1
1	A	131	ASN	2.1
1	C	80	HIS	2.1
1	B	177	PHE	2.1
1	B	127	LEU	2.1
1	B	139	SER	2.1
1	C	179	TRP	2.1
1	A	107	LYS	2.1
1	A	110	LYS	2.1
1	C	146	LEU	2.0
1	B	184	ARG	2.0
1	C	186	ALA	2.0
1	A	76	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 oligosaccharides 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. 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	B-factors(Å ²)	Q<0.9
3	EDO	C	202	4/4	0.61	0.20	35,41,46,50	0
2	NO3	A	401	4/4	0.77	0.13	44,48,56,57	0
2	NO3	B	203	4/4	0.83	0.11	37,38,39,40	0
3	EDO	B	201	4/4	0.87	0.11	19,19,22,23	0
2	NO3	C	201	4/4	0.87	0.11	27,31,33,36	0
2	NO3	B	204	4/4	0.90	0.10	36,39,41,42	0
2	NO3	B	202	4/4	0.94	0.07	19,22,22,25	0

6.5 Other polymers [i](#)

There are no such residues in this entry.