



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 02:22 PM EDT

PDB ID : 3BDL / pdb_00003bdl
Title : Crystal structure of a truncated human Tudor-SN
Authors : Li, C.L.
Deposited on : 2007-11-15
Resolution : 1.90 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

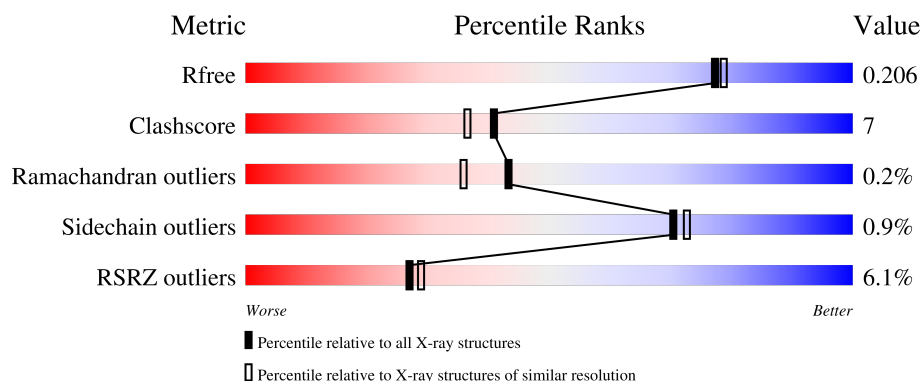
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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 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	570	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4984 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.

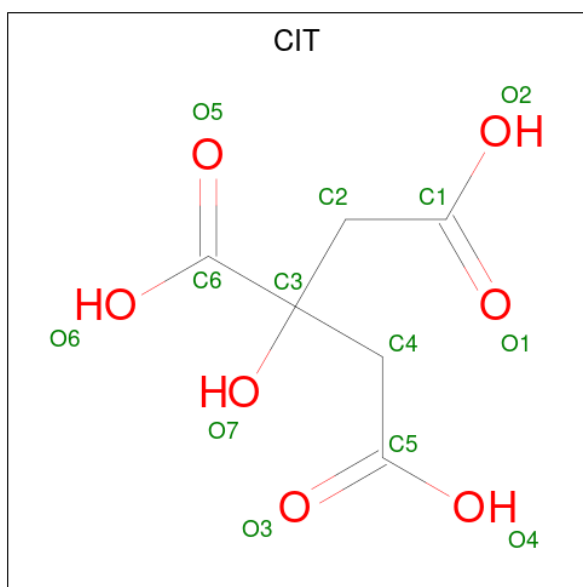
- Molecule 1 is a protein called Staphylococcal nuclease domain-containing protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	537	Total	C	N	O	S	0	0	0
			4242	2689	734	805	14			

There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	294	MET	-	expression tag	UNP Q7KZF4
A	295	GLY	-	expression tag	UNP Q7KZF4
A	296	SER	-	expression tag	UNP Q7KZF4
A	297	SER	-	expression tag	UNP Q7KZF4
A	298	HIS	-	expression tag	UNP Q7KZF4
A	299	HIS	-	expression tag	UNP Q7KZF4
A	300	HIS	-	expression tag	UNP Q7KZF4
A	301	HIS	-	expression tag	UNP Q7KZF4
A	302	HIS	-	expression tag	UNP Q7KZF4
A	303	HIS	-	expression tag	UNP Q7KZF4
A	304	SER	-	expression tag	UNP Q7KZF4
A	305	SER	-	expression tag	UNP Q7KZF4
A	306	GLY	-	expression tag	UNP Q7KZF4
A	307	LEU	-	expression tag	UNP Q7KZF4
A	308	VAL	-	expression tag	UNP Q7KZF4
A	309	PRO	-	expression tag	UNP Q7KZF4
A	310	ARG	-	expression tag	UNP Q7KZF4
A	311	GLY	-	expression tag	UNP Q7KZF4
A	312	SER	-	expression tag	UNP Q7KZF4
A	313	HIS	-	expression tag	UNP Q7KZF4
A	314	MET	-	expression tag	UNP Q7KZF4

- Molecule 2 is CITRIC ACID (CCD ID: CIT) (formula: C₆H₈O₇).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			13	6	7		
2	A	1	Total	C	O	0	0
			13	6	7		
2	A	1	Total	C	O	0	0
			13	6	7		
2	A	1	Total	C	O	0	0
			13	6	7		

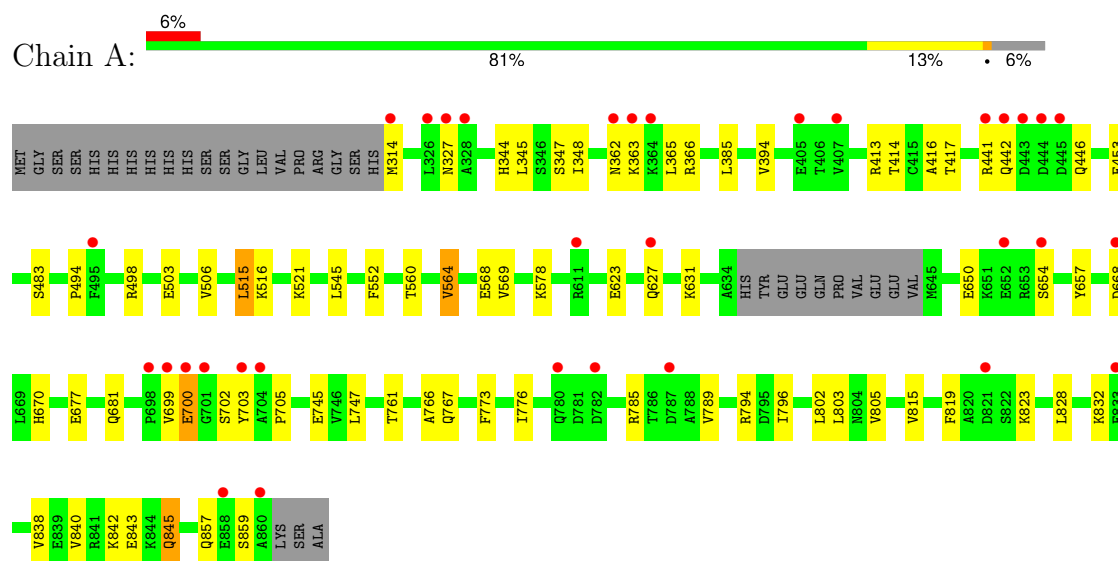
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	690	Total	O	0	0
			690	690		

3 Residue-property plots

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: Staphylococcal nuclease domain-containing protein 1



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	97.41Å 91.93Å 88.04Å 90.00° 91.26° 90.00°	Depositor
Resolution (Å)	23.33 – 1.90 23.33 – 1.90	Depositor EDS
% Data completeness (in resolution range)	93.8 (23.33-1.90) 93.8 (23.33-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.79 (at 1.81Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.175 , 0.209 0.171 , 0.206	Depositor DCC
R_{free} test set	5780 reflections (8.49%)	wwPDB-VP
Wilson B-factor (Å ²)	18.3	Xtriage
Anisotropy	0.247	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 54.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.024 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4984	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.69% 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

5.1 Standard geometry

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

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	0.34	0/4322	0.84	9/5844 (0.2%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	564	VAL	N-CA-C	6.48	120.02	113.47
1	A	385	LEU	N-CA-C	6.15	121.66	114.04
1	A	840	VAL	N-CA-C	5.71	116.95	109.58
1	A	670	HIS	N-CA-C	-5.48	102.08	110.14
1	A	767	GLN	N-CA-C	5.29	119.45	112.89
1	A	796	ILE	N-CA-C	5.17	120.10	109.34
1	A	417	THR	N-CA-C	-5.10	101.00	109.46
1	A	347	SER	N-CA-C	5.06	119.03	112.86
1	A	766	ALA	N-CA-C	-5.05	102.58	110.10

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	4242	0	4264	56	0
2	A	52	0	20	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	690	0	0	11	0
All	All	4984	0	4284	56	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 7.

All (56) 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:503:GLU:H	1:A:521:LYS:HZ2	1.19	0.88
1:A:506:VAL:HG21	1:A:564:VAL:HG22	1.61	0.83
1:A:503:GLU:H	1:A:521:LYS:NZ	1.76	0.83
1:A:314:MET:HE1	1:A:414:THR:HG21	1.58	0.82
1:A:843:GLU:HB3	1:A:845:GLN:HE22	1.50	0.77
1:A:650:GLU:O	1:A:845:GLN:HG3	1.87	0.75
1:A:545:LEU:HD23	3:A:2601:HOH:O	1.90	0.70
1:A:843:GLU:HB3	1:A:845:GLN:NE2	2.11	0.64
1:A:650:GLU:HB2	1:A:845:GLN:HG2	1.79	0.64
1:A:453:GLU:HG3	3:A:2158:HOH:O	1.97	0.63
1:A:506:VAL:HG21	1:A:564:VAL:CG2	2.27	0.63
1:A:761:THR:HG21	3:A:2261:HOH:O	2.00	0.61
1:A:631:LYS:HE2	3:A:2477:HOH:O	2.00	0.60
1:A:700:GLU:H	1:A:700:GLU:CD	2.10	0.59
1:A:838:VAL:H	1:A:857:GLN:NE2	2.01	0.59
1:A:483:SER:OG	1:A:516:LYS:HE3	2.03	0.58
1:A:552:PHE:CE2	1:A:631:LYS:HA	2.40	0.57
1:A:446:GLN:HA	3:A:2198:HOH:O	2.04	0.57
1:A:845:GLN:NE2	1:A:845:GLN:H	2.05	0.55
1:A:745:GLU:CD	1:A:747:LEU:HD21	2.31	0.55
1:A:564:VAL:HG22	1:A:564:VAL:O	2.08	0.53
1:A:506:VAL:CG2	1:A:564:VAL:HG22	2.35	0.53
1:A:552:PHE:CZ	1:A:631:LYS:HA	2.43	0.53
1:A:703:TYR:O	1:A:705:PRO:HD3	2.08	0.53
1:A:650:GLU:H	1:A:845:GLN:NE2	2.06	0.53
1:A:365:LEU:HD23	1:A:366:ARG:N	2.25	0.51
1:A:623:GLU:O	1:A:627:GLN:HG3	2.11	0.51
1:A:365:LEU:HD23	1:A:365:LEU:C	2.37	0.50
1:A:803:LEU:C	1:A:803:LEU:HD23	2.39	0.48
1:A:842:LYS:HB3	1:A:842:LYS:NZ	2.29	0.47
1:A:327:ASN:HB3	3:A:2085:HOH:O	2.14	0.47
1:A:703:TYR:CE2	1:A:705:PRO:HG3	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:515:LEU:HD12	1:A:515:LEU:N	2.30	0.47
1:A:668:ASP:HA	1:A:785:ARG:HD2	1.98	0.46
1:A:564:VAL:CG2	1:A:569:VAL:HG11	2.45	0.46
1:A:773:PHE:HB2	1:A:776:ILE:HD11	1.97	0.46
1:A:859:SER:HB2	3:A:2395:HOH:O	2.17	0.44
1:A:441:ARG:HB3	1:A:442:GLN:NE2	2.33	0.44
1:A:494:PRO:HB2	1:A:498:ARG:HH12	1.83	0.44
1:A:345:LEU:HB2	1:A:348:ILE:HD11	2.00	0.43
1:A:699:VAL:HB	1:A:702:SER:OG	2.18	0.43
1:A:819:PHE:O	1:A:823:LYS:HA	2.19	0.43
1:A:344:HIS:CE1	1:A:413:ARG:HD3	2.53	0.43
1:A:362:ASN:C	1:A:363:LYS:HG2	2.44	0.42
1:A:521:LYS:HD3	3:A:2075:HOH:O	2.18	0.42
1:A:828:LEU:O	1:A:832:LYS:HG2	2.19	0.42
1:A:568:GLU:HG3	3:A:2110:HOH:O	2.18	0.42
1:A:560:THR:O	1:A:564:VAL:HG12	2.19	0.42
1:A:578:LYS:HE2	1:A:702:SER:O	2.20	0.42
1:A:394:VAL:HA	1:A:416:ALA:HB2	2.03	0.41
1:A:681:GLN:HG2	3:A:2346:HOH:O	2.21	0.41
1:A:657:TYR:HB3	1:A:802:LEU:HB3	2.03	0.41
1:A:805:VAL:HA	1:A:815:VAL:HG12	2.01	0.41
1:A:745:GLU:HG2	1:A:747:LEU:HG	2.03	0.40
1:A:794:ARG:NH1	3:A:2471:HOH:O	2.55	0.40
1:A:785:ARG:O	1:A:789:VAL:HG23	2.22	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	533/570 (94%)	520 (98%)	12 (2%)	1 (0%)	43 36

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	654	SER

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	461/490 (94%)	457 (99%)	4 (1%)	70	73

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

Mol	Chain	Res	Type
1	A	515	LEU
1	A	677	GLU
1	A	700	GLU
1	A	845	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	324	GLN
1	A	327	ASN
1	A	344	HIS
1	A	362	ASN
1	A	442	GLN
1	A	450	HIS
1	A	469	HIS
1	A	606	HIS
1	A	627	GLN
1	A	681	GLN
1	A	688	ASN
1	A	696	HIS
1	A	767	GLN
1	A	798	ASN
1	A	818	GLN

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Mol	Chain	Res	Type
1	A	845	GLN
1	A	857	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

4 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	$\# Z > 2$
2	CIT	A	2000	-	12,12,12	0.96	0	17,17,17	1.30	1 (5%)
2	CIT	A	2003	-	12,12,12	0.89	0	17,17,17	1.31	1 (5%)
2	CIT	A	2001	-	12,12,12	1.09	1 (8%)	17,17,17	1.24	1 (5%)
2	CIT	A	2002	-	12,12,12	0.89	0	17,17,17	1.31	1 (5%)

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
2	CIT	A	2000	-	-	0/16/16/16	-
2	CIT	A	2003	-	-	0/16/16/16	-
2	CIT	A	2001	-	-	0/16/16/16	-
2	CIT	A	2002	-	-	0/16/16/16	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	2001	CIT	C2-C3	2.45	1.57	1.54

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	2002	CIT	O6-C6-C3	3.65	120.14	113.14
2	A	2003	CIT	O6-C6-C3	3.59	120.02	113.14
2	A	2000	CIT	O6-C6-C3	3.52	119.89	113.14
2	A	2001	CIT	O6-C6-C3	3.24	119.36	113.14

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	537/570 (94%)	-0.02	33 (6%) 27 29	12, 22, 47, 62	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	699	VAL	4.3
1	A	442	GLN	4.1
1	A	327	ASN	3.8
1	A	821	ASP	3.5
1	A	668	ASP	3.4
1	A	326	LEU	3.0
1	A	701	GLY	3.0
1	A	860	ALA	3.0
1	A	833	GLU	3.0
1	A	314	MET	2.9
1	A	363	LYS	2.9
1	A	703	TYR	2.9
1	A	443	ASP	2.8
1	A	444	ASP	2.8
1	A	704	ALA	2.7
1	A	407	VAL	2.7
1	A	441	ARG	2.6
1	A	654	SER	2.5
1	A	495	PHE	2.4
1	A	364	LYS	2.4
1	A	627	GLN	2.3
1	A	611	ARG	2.3
1	A	328	ALA	2.3
1	A	780	GLN	2.2
1	A	652	GLU	2.2
1	A	858	GLU	2.2
1	A	700	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	445	ASP	2.1
1	A	698	PRO	2.1
1	A	362	ASN	2.1
1	A	782	ASP	2.1
1	A	405	GLU	2.1
1	A	787	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 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($\text{\AA}^2$)	Q<0.9
2	CIT	A	2002	13/13	0.86	0.12	31,38,47,50	0
2	CIT	A	2003	13/13	0.92	0.10	18,31,40,42	0
2	CIT	A	2001	13/13	0.96	0.05	14,20,24,25	0
2	CIT	A	2000	13/13	0.97	0.05	17,20,24,25	0

6.5 Other polymers [i](#)

There are no such residues in this entry.