



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 10:56 PM UTC

PDB ID : 4TR8 / pdb_00004tr8
Title : Crystal structure of DNA polymerase sliding clamp from *Pseudomonas aeruginosa*
Authors : Olieric, V.; Burnouf, D.; Ennifar, E.; Wolff, P.
Deposited on : 2014-06-15
Resolution : 1.80 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
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

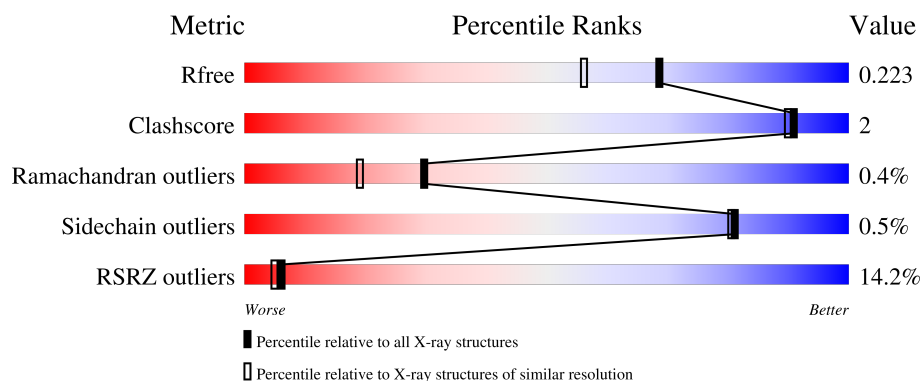
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	383	12% 95% ..
1	B	383	16% 88% 6% 6%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5989 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 DNA polymerase III subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	379	Total	C	N	O	S	0	3	0
			2864	1798	506	550	10			
1	B	361	Total	C	N	O	S	0	1	0
			2692	1692	475	516	9			

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	9985	HIS	-	expression tag	UNP V4MZL6
A	9986	HIS	-	expression tag	UNP V4MZL6
A	9987	HIS	-	expression tag	UNP V4MZL6
A	9988	HIS	-	expression tag	UNP V4MZL6
A	9989	HIS	-	expression tag	UNP V4MZL6
A	9990	HIS	-	expression tag	UNP V4MZL6
A	9991	SER	-	expression tag	UNP V4MZL6
A	9992	SER	-	expression tag	UNP V4MZL6
A	9993	GLY	-	expression tag	UNP V4MZL6
A	9994	LEU	-	expression tag	UNP V4MZL6
A	9995	VAL	-	expression tag	UNP V4MZL6
A	9996	PRO	-	expression tag	UNP V4MZL6
A	9997	ARG	-	expression tag	UNP V4MZL6
A	9998	GLY	-	expression tag	UNP V4MZL6
A	9999	SER	-	expression tag	UNP V4MZL6
A	0	HIS	-	expression tag	UNP V4MZL6
B	-15	HIS	-	expression tag	UNP V4MZL6
B	-14	HIS	-	expression tag	UNP V4MZL6
B	-13	HIS	-	expression tag	UNP V4MZL6
B	-12	HIS	-	expression tag	UNP V4MZL6
B	-11	HIS	-	expression tag	UNP V4MZL6
B	-10	HIS	-	expression tag	UNP V4MZL6
B	-9	SER	-	expression tag	UNP V4MZL6
B	-8	SER	-	expression tag	UNP V4MZL6
B	-7	GLY	-	expression tag	UNP V4MZL6

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-6	LEU	-	expression tag	UNP V4MZL6
B	-5	VAL	-	expression tag	UNP V4MZL6
B	-4	PRO	-	expression tag	UNP V4MZL6
B	-3	ARG	-	expression tag	UNP V4MZL6
B	-2	GLY	-	expression tag	UNP V4MZL6
B	-1	SER	-	expression tag	UNP V4MZL6
B	0	HIS	-	expression tag	UNP V4MZL6

- Molecule 2 is SODIUM ION (CCD ID: NA) (formula: Na).

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

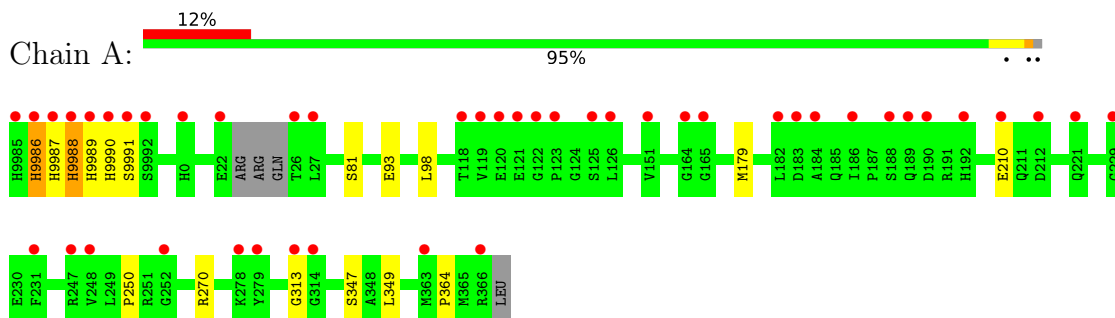
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	259	Total O 259 259	0	0
3	B	173	Total O 173 173	0	0

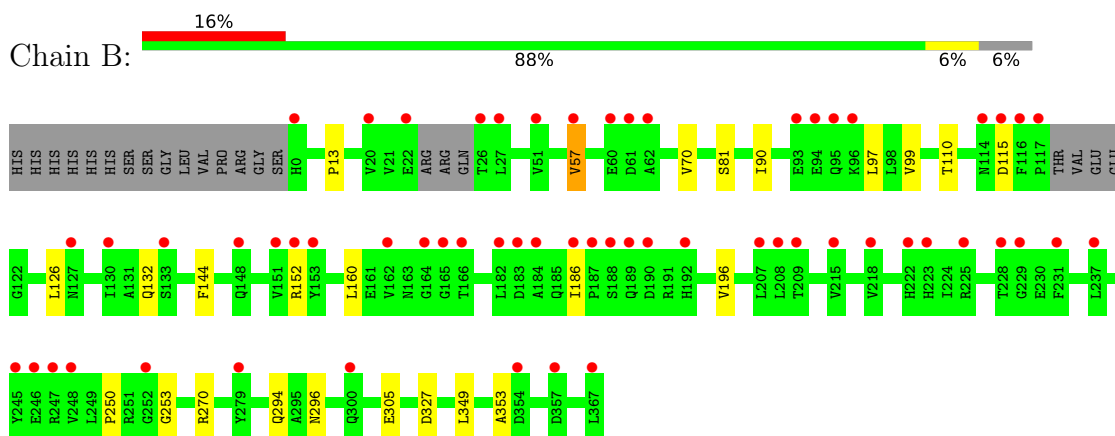
3 Residue-property plots [i](#)

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: DNA polymerase III subunit beta



- Molecule 1: DNA polymerase III subunit beta



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	80.53Å 92.41Å 127.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.90 – 1.80 43.90 – 1.80	Depositor EDS
% Data completeness (in resolution range)	100.0 (43.90-1.80) 83.9 (43.90-1.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.05 (at 1.79Å)	Xtriage
Refinement program	BUSTER 2.10.0	Depositor
R, R_{free}	0.207 , 0.232 (Not available) , 0.223	Depositor DCC
R_{free} test set	4450 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	24.5	Xtriage
Anisotropy	0.486	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 45.4	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	0.95	EDS
Total number of atoms	5989	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.52% 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: NA

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	0.82	1/2913 (0.0%)	1.16	9/3954 (0.2%)
1	B	0.78	0/2729	1.12	6/3705 (0.2%)
All	All	0.80	1/5642 (0.0%)	1.14	15/7659 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	179	MET	SD-CE	-5.11	1.66	1.79

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	313	GLY	CA-C-N	-6.68	115.66	122.27
1	A	313	GLY	C-N-CA	-6.68	115.66	122.27
1	B	353	ALA	N-CA-C	6.00	117.49	111.07
1	A	9990	HIS	CA-C-N	5.74	132.51	121.54
1	A	9990	HIS	C-N-CA	5.74	132.51	121.54
1	B	132	GLN	CA-C-N	5.57	127.75	120.28
1	B	132	GLN	C-N-CA	5.57	127.75	120.28
1	B	115	ASP	CA-CB-CG	5.54	118.14	112.60
1	A	9987	HIS	CA-C-N	5.20	129.65	120.87
1	A	9987	HIS	C-N-CA	5.20	129.65	120.87
1	B	152	ARG	CA-C-N	5.05	127.30	120.38
1	B	152	ARG	C-N-CA	5.05	127.30	120.38
1	A	210	GLU	CA-C-N	5.01	127.25	120.38
1	A	210	GLU	C-N-CA	5.01	127.25	120.38
1	A	9986	HIS	CA-CB-CG	5.01	118.81	113.80

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	2864	0	2841	9	0
1	B	2692	0	2660	13	0
2	A	1	0	0	0	0
3	A	259	0	0	0	0
3	B	173	0	0	0	0
All	All	5989	0	5501	19	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 2.

All (19) 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:9988:HIS:H	1:A:9988:HIS:CD2	1.84	0.94
1:A:9988:HIS:H	1:A:9988:HIS:HD2	1.25	0.82
1:B:294:GLN:HE22	1:B:296:ASN:HD22	1.32	0.74
1:A:81:SER:OG	1:B:270:ARG:HG2	2.04	0.57
1:A:9988:HIS:CD2	1:A:9988:HIS:N	2.62	0.55
1:B:70:VAL:HG12	1:B:110:THR:HG22	1.92	0.51
1:A:270:ARG:HG2	1:B:81:SER:OG	2.11	0.50
1:B:144:PHE:CD1	1:B:327:ASP:HB3	2.49	0.47
1:B:70:VAL:HG11	1:B:97:LEU:HD22	1.97	0.47
1:B:250:PRO:HG2	1:B:349:LEU:HB2	1.99	0.44
1:A:347:SER:HA	1:A:364:PRO:HD3	1.99	0.44
1:A:93:GLU:HG3	1:A:98:LEU:HD22	2.00	0.43
1:A:250:PRO:HG2	1:A:349:LEU:HB2	2.00	0.42
1:B:126:LEU:HD21	1:B:186:ILE:HG21	2.01	0.42
1:B:160:LEU:HD23	1:B:196:VAL:HG23	2.02	0.42
1:B:13:PRO:HG3	1:B:57:VAL:HG22	2.01	0.41
1:B:90:ILE:HG12	1:B:99:VAL:HG22	2.01	0.41
1:A:81:SER:CB	1:B:270:ARG:HG2	2.51	0.41
1:B:270:ARG:HH21	1:B:305:GLU:CD	2.30	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	378/383 (99%)	364 (96%)	12 (3%)	2 (0%)	24	14
1	B	356/383 (93%)	348 (98%)	7 (2%)	1 (0%)	36	25
All	All	734/766 (96%)	712 (97%)	19 (3%)	3 (0%)	30	19

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	9991	SER
1	B	253	GLY
1	A	9989	HIS

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	307/331 (93%)	305 (99%)	2 (1%)	76	73
1	B	282/331 (85%)	281 (100%)	1 (0%)	84	83
All	All	589/662 (89%)	586 (100%)	3 (0%)	81	80

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9986	HIS
1	A	9988	HIS
1	B	57	VAL

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

Mol	Chain	Res	Type
1	A	9986	HIS
1	A	9988	HIS
1	A	6	GLN
1	A	40	GLN
1	A	41	GLN
1	A	222	HIS
1	A	296	ASN
1	B	32	ASN
1	B	40	GLN
1	B	148	GLN
1	B	163	ASN
1	B	285	GLN
1	B	296	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 1 ligands modelled in this entry, 1 is 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	379/383 (98%)	0.58	45 (11%) 9 7	13, 26, 55, 81	3 (0%)
1	B	361/383 (94%)	0.96	60 (16%) 4 3	16, 36, 67, 91	1 (0%)
All	All	740/766 (96%)	0.76	105 (14%) 6 5	13, 31, 63, 91	4 (0%)

All (105) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	9987	HIS	6.3
1	A	252	GLY	6.1
1	A	9988	HIS	5.9
1	B	165	GLY	5.8
1	B	367	LEU	5.5
1	A	212	ASP	5.3
1	A	9991	SER	5.3
1	A	9986	HIS	5.0
1	A	9989	HIS	5.0
1	B	164	GLY	5.0
1	A	229	GLY	4.8
1	B	26	THR	4.5
1	B	0	HIS	4.4
1	A	26	THR	4.4
1	A	164	GLY	4.4
1	B	190	ASP	4.3
1	B	252	GLY	4.3
1	B	187	PRO	4.1
1	B	182	LEU	4.0
1	A	165	GLY	3.9
1	B	151	VAL	3.7
1	A	22	GLU	3.7
1	B	248	VAL	3.7
1	A	313	GLY	3.7

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Mol	Chain	Res	Type	RSRZ
1	B	117	PRO	3.7
1	A	120	GLU	3.6
1	A	121	GLU	3.5
1	B	223	HIS	3.4
1	A	9990	HIS	3.3
1	A	210	GLU	3.3
1	A	182	LEU	3.2
1	A	314	GLY	3.2
1	B	192	HIS	3.2
1	A	123	PRO	3.2
1	B	279	TYR	3.2
1	A	221	GLN	3.2
1	B	60	GLU	3.2
1	B	130	ILE	3.1
1	B	186	ILE	3.1
1	B	246	GLU	3.1
1	B	189	GLN	3.1
1	B	184	ALA	3.0
1	A	126	LEU	3.0
1	A	9992	SER	3.0
1	A	118	THR	3.0
1	A	122	GLY	2.9
1	B	27	LEU	2.9
1	B	166	THR	2.9
1	A	363	MET	2.9
1	B	114	ASN	2.8
1	A	125	SER	2.8
1	B	231	PHE	2.8
1	A	119	VAL	2.8
1	A	188	SER	2.7
1	B	148	GLN	2.7
1	B	93	GLU	2.7
1	B	61	ASP	2.7
1	B	62	ALA	2.6
1	B	133	SER	2.6
1	A	9985	HIS	2.6
1	A	27	LEU	2.6
1	A	247	ARG	2.5
1	A	366	ARG	2.5
1	B	237	LEU	2.5
1	B	127	ASN	2.5
1	B	188	SER	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	183	ASP	2.5
1	B	51	VAL	2.5
1	A	184	ALA	2.4
1	A	278	LYS	2.4
1	B	95	GLN	2.4
1	B	247	ARG	2.3
1	B	245	TYR	2.3
1	B	228	THR	2.3
1	A	186	ILE	2.3
1	B	222	HIS	2.3
1	B	96	LYS	2.3
1	B	218	VAL	2.3
1	B	354	ASP	2.3
1	B	300	GLN	2.3
1	A	183	ASP	2.3
1	B	20	VAL	2.3
1	A	231	PHE	2.3
1	B	152	ARG	2.3
1	A	190	ASP	2.2
1	B	209	THR	2.2
1	B	207	LEU	2.2
1	A	151	VAL	2.2
1	B	57	VAL	2.2
1	B	115	ASP	2.2
1	B	357	ASP	2.2
1	B	229	GLY	2.2
1	B	116	PHE	2.1
1	A	279	TYR	2.1
1	B	153	TYR	2.1
1	B	215	VAL	2.1
1	A	192	HIS	2.1
1	A	189	GLN	2.1
1	B	94	GLU	2.1
1	A	0	HIS	2.1
1	B	162	VAL	2.1
1	B	208	LEU	2.1
1	B	225	ARG	2.1
1	B	22	GLU	2.0
1	A	248	VAL	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	NA	A	1001	1/1	1.00	0.08	11,11,11,11	0

6.5 Other polymers [i](#)

There are no such residues in this entry.