



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 12:54 PM UTC

PDB ID : 7AMS / pdb_00007ams
Title : Crystal structure of the complex of the KFN mutant of HuJovi-1 Fab with human TRBC2
Authors : Ferrari, M.; Bulek, A.; Bughda, R.; Jha, R.; Welin, M.; Svensson, A.; Logan, D.T.; Sewell, A.; Onuoha, S.; Pule, M.
Deposited on : 2020-10-09
Resolution : 2.42 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

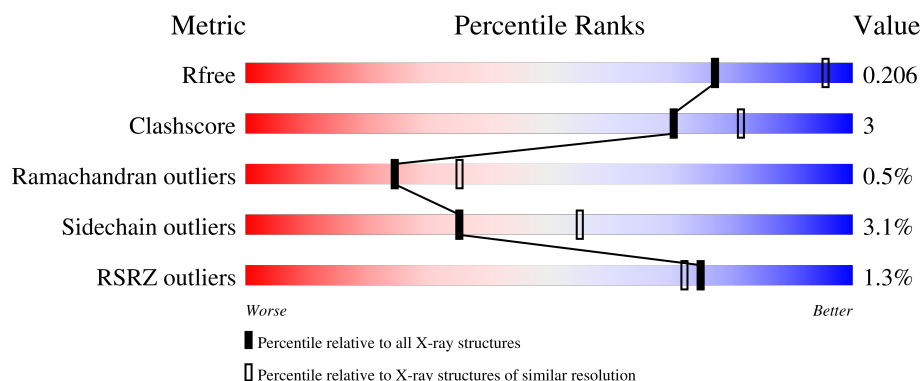
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.42 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	6062 (2.44-2.40)
Clashscore	190562	6562 (2.44-2.40)
Ramachandran outliers	187476	6481 (2.44-2.40)
Sidechain outliers	187428	6482 (2.44-2.40)
RSRZ outliers	180081	6066 (2.44-2.40)

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	204	2% 87% 11% ..
2	B	246	2% 89% 9% .
3	H	225	83% 12% ..
4	L	219	90% 8% .

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
6	CL	L	504	-	-	X	-

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 7492 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 Human A6 T-cell receptor alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	201	Total	C	N	O	S	0	0	0
			1551	961	254	328	8			

- Molecule 2 is a protein called Human A6 T-cell receptor beta chain TRBC2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	240	Total	C	N	O	S	0	1	0
			1910	1200	333	367	10			

- Molecule 3 is a protein called Human Jovi-1 Fab, KFN mutant, heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	H	218	Total	C	N	O	S	0	2	0
			1683	1065	287	320	11			

- Molecule 4 is a protein called Human Jovi-1 Fab, KFN mutant, light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	L	217	Total	C	N	O	S	0	1	0
			1693	1061	292	335	5			

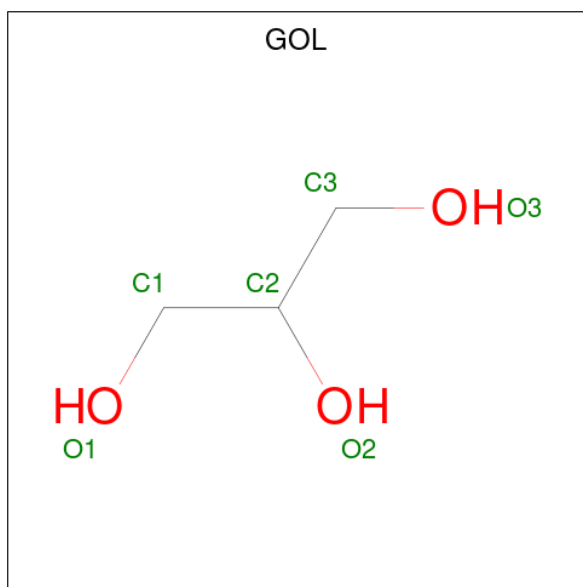
- Molecule 5 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	H	1	Total	Zn	0	0
			1	1		
5	L	2	Total	Zn	0	0
			2	2		

- Molecule 6 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	H	1	Total Cl 1 1	0	0
6	L	1	Total Cl 1 1	0	0

- Molecule 7 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	L	1	Total C O 6 3 3	0	0

- Molecule 8 is water.

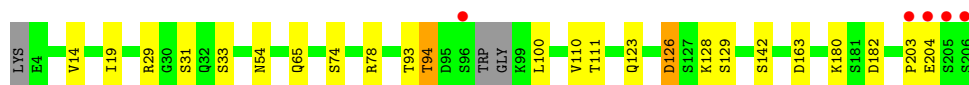
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	124	Total O 124 124	0	0
8	B	160	Total O 160 160	0	0
8	H	182	Total O 182 182	0	0
8	L	178	Total O 178 178	0	0

3 Residue-property plots [i](#)

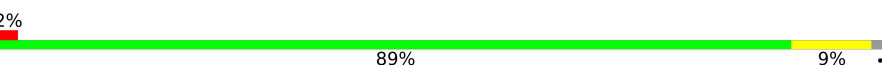
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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.

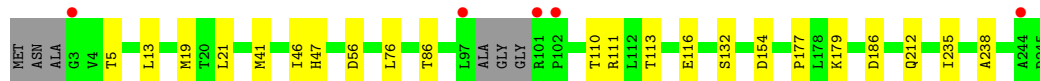
- Molecule 1: Human A6 T-cell receptor alpha chain

Chain A: 



- Molecule 2: Human A6 T-cell receptor beta chain TRBC2

Chain B: 



- Molecule 3: Human Jovi-1 Fab, KFN mutant, heavy chain

Chain H: 



- Molecule 4: Human Jovi-1 Fab, KFN mutant, light chain

Chain L: 



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	114.84Å 91.16Å 123.32Å 90.00° 94.09° 90.00°	Depositor
Resolution (Å)	47.64 – 2.42 47.64 – 2.42	Depositor EDS
% Data completeness (in resolution range)	99.8 (47.64-2.42) 99.8 (47.64-2.42)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.75 (at 2.42Å)	Xtriage
Refinement program	BUSTER 2.11.7	Depositor
R, R_{free}	0.180 , 0.225 (Not available) , 0.206	Depositor DCC
R_{free} test set	2360 reflections (4.87%)	wwPDB-VP
Wilson B-factor (Å ²)	34.8	Xtriage
Anisotropy	1.080	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 65.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7492	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.38% 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: GOL, ZN, 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	0.75	1/1582 (0.1%)	1.02	2/2146 (0.1%)
2	B	0.73	0/1962	1.01	0/2671
3	H	0.70	0/1725	1.04	5/2341 (0.2%)
4	L	0.73	0/1733	1.01	0/2355
All	All	0.73	1/7002 (0.0%)	1.02	7/9513 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	94	THR	CA-C	6.06	1.57	1.52

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	163	ASP	CA-CB-CG	6.43	119.03	112.60
1	A	126	ASP	CA-CB-CG	6.06	118.66	112.60
3	H	50	PHE	CA-CB-CG	5.82	119.62	113.80
3	H	89	ASP	CA-CB-CG	5.21	117.81	112.60
3	H	197	SER	CA-C-N	5.19	129.87	122.08
3	H	197	SER	C-N-CA	5.19	129.87	122.08
3	H	136	ALA	N-CA-C	5.18	116.22	109.64

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	1551	0	1451	11	0
2	B	1910	0	1811	10	0
3	H	1683	0	1635	17	0
4	L	1693	0	1651	14	0
5	H	1	0	0	0	0
5	L	2	0	0	0	0
6	H	1	0	0	0	0
6	L	1	0	0	3	0
7	L	6	0	8	0	0
8	A	124	0	0	0	0
8	B	160	0	0	0	0
8	H	182	0	0	0	0
8	L	178	0	0	0	0
All	All	7492	0	6556	45	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 3.

All (45) 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)
3:H:11:VAL:HB	3:H:158:PRO:HG3	1.63	0.80
3:H:175:HIS:CE1	4:L:143:ASN:HD21	2.03	0.75
3:H:175:HIS:HE1	4:L:143:ASN:HD21	1.42	0.68
4:L:66:ARG:HD2	4:L:82:ARG:O	2.02	0.59
1:A:19:ILE:HG12	1:A:78:ARG:HA	1.86	0.58
2:B:13:LEU:HD21	2:B:19[B]:MET:HG3	1.86	0.57
2:B:21:LEU:HD12	2:B:76:LEU:HD23	1.85	0.57
3:H:53:PRO:HA	3:H:72:ARG:HD2	1.87	0.56
2:B:21:LEU:HD22	2:B:110:THR:HG21	1.89	0.54
4:L:215:ASN:O	4:L:216:ARG:HB2	2.08	0.54
4:L:29:LEU:HD22	4:L:95:GLN:HG3	1.92	0.52
3:H:67:ARG:NH1	3:H:90:ASP:OD2	2.39	0.52
3:H:175:HIS:CE1	4:L:143:ASN:ND2	2.76	0.52
1:A:128:LYS:HZ2	2:B:238:ALA:HA	1.76	0.51
3:H:36:TRP:CE2	3:H:81:MET:HB2	2.47	0.50
3:H:130:PRO:HB3	3:H:156:TYR:HB3	1.93	0.50
2:B:154:ASP:OD1	2:B:177:PRO:HG2	2.12	0.49
4:L:216:ARG:CB	6:L:504:CL:CL	2.98	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:46:ILE:HG22	2:B:47:HIS:CD2	2.49	0.47
1:A:126:ASP:HB3	1:A:129:SER:O	2.15	0.46
1:A:31:SER:HA	1:A:94:THR:HG22	1.97	0.46
3:H:163:VAL:HG22	3:H:209:VAL:HG22	1.97	0.46
4:L:216:ARG:HB3	6:L:504:CL:CL	2.53	0.45
4:L:216:ARG:HB2	6:L:504:CL:CL	2.53	0.45
2:B:186:ASP:HB2	3:H:28:LYS:NZ	2.32	0.45
4:L:118:PRO:HB3	4:L:144:PHE:HB3	1.99	0.45
2:B:86:THR:HG23	2:B:113:THR:HA	2.00	0.44
3:H:103:ASN:O	3:H:104:PHE:HB2	2.17	0.44
2:B:186:ASP:HB2	3:H:28:LYS:HZ3	1.82	0.44
1:A:33:SER:HB2	1:A:93:THR:CG2	2.48	0.43
4:L:24:ARG:HD3	4:L:75:ASP:OD1	2.19	0.43
3:H:62:GLU:HA	3:H:65:ARG:HG3	1.99	0.43
1:A:123:GLN:OE1	1:A:203:PRO:HB3	2.19	0.43
1:A:14:VAL:O	1:A:110:VAL:HA	2.19	0.43
3:H:53:PRO:HB3	3:H:72:ARG:HG3	2.01	0.42
1:A:93:THR:HB	1:A:100:LEU:HD12	2.02	0.41
1:A:65:GLN:HB2	1:A:74:SER:HB2	2.03	0.41
1:A:111:THR:HB	1:A:142:SER:HB3	2.01	0.41
3:H:50:PHE:CE2	3:H:59:GLN:HB2	2.56	0.41
3:H:189:LEU:C	3:H:189:LEU:HD12	2.45	0.41
1:A:180:LYS:HE2	1:A:180:LYS:HB3	1.89	0.41
3:H:175:HIS:HE1	4:L:143:ASN:ND2	2.14	0.41
4:L:125:PRO:HD3	4:L:137:VAL:HG22	2.03	0.41
4:L:202:THR:HG22	4:L:209:PRO:HG3	2.03	0.40
2:B:212:GLN:HG3	2:B:235:ILE:CG2	2.52	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	197/204 (97%)	189 (96%)	6 (3%)	2 (1%)	12	18
2	B	237/246 (96%)	230 (97%)	7 (3%)	0	100	100
3	H	216/225 (96%)	209 (97%)	6 (3%)	1 (0%)	24	35
4	L	216/219 (99%)	213 (99%)	2 (1%)	1 (0%)	24	35
All	All	866/894 (97%)	841 (97%)	21 (2%)	4 (0%)	24	35

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	L	216	ARG
1	A	54	ASN
3	H	199	SER
1	A	204	GLU

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	180/182 (99%)	178 (99%)	2 (1%)	65	81
2	B	209/210 (100%)	202 (97%)	7 (3%)	33	53
3	H	186/190 (98%)	176 (95%)	10 (5%)	20	33
4	L	195/196 (100%)	189 (97%)	6 (3%)	35	55
All	All	770/778 (99%)	745 (97%)	25 (3%)	35	54

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	ARG
1	A	182	ASP
2	B	5	THR
2	B	41	MET
2	B	56	ASP
2	B	111	ARG
2	B	116	GLU

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Mol	Chain	Res	Type
2	B	132	SER
2	B	179	LYS
3	H	1	GLN
3	H	56	ASP
3	H	60	SER
3	H	63	ARG
3	H	72	ARG
3	H	88	SER
3	H	111	PHE
3	H	159	GLU
3	H	189	LEU
3	H	197	SER
4	L	22	SER
4	L	39	HIS
4	L	128[A]	GLU
4	L	128[B]	GLU
4	L	168	VAL
4	L	216	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	39	GLN
1	A	65	GLN
1	A	101	GLN
2	B	37	GLN
2	B	138	HIS
2	B	176	GLN
3	H	100	ASN
3	H	175	HIS
3	H	210	ASN
4	L	98	HIS
4	L	143	ASN

5.3.3 RNA ⓘ

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 6 ligands modelled in this entry, 5 are 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	$\# Z > 2$
7	GOL	L	501	-	5,5,5	0.18	0	5,5,5	0.32	0

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
7	GOL	L	501	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	L	501	GOL	C1-C2-C3-O3
7	L	501	GOL	O2-C2-C3-O3

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	201/204 (98%)	0.11	5 (2%) 58 55	36, 51, 75, 90	0
2	B	240/246 (97%)	-0.10	5 (2%) 63 60	26, 48, 73, 88	1 (0%)
3	H	218/225 (96%)	-0.23	1 (0%) 87 86	21, 44, 61, 75	2 (0%)
4	L	217/219 (99%)	-0.36	0 100 100	26, 42, 54, 63	1 (0%)
All	All	876/894 (97%)	-0.15	11 (1%) 75 72	21, 45, 70, 90	4 (0%)

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	97	LEU	3.9
1	A	96	SER	3.2
2	B	101	ARG	3.0
1	A	205	SER	2.7
1	A	206	SER	2.5
1	A	203	PRO	2.5
3	H	225	LYS	2.4
1	A	204	GLU	2.3
2	B	244	ALA	2.2
2	B	3	GLY	2.2
2	B	102	PRO	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
6	CL	L	504	1/1	0.91	0.14	87,87,87,87	0
6	CL	H	302	1/1	0.93	0.09	75,75,75,75	0
7	GOL	L	501	6/6	0.95	0.08	42,43,44,44	0
5	ZN	L	503	1/1	0.99	0.15	45,45,45,45	0
5	ZN	L	502	1/1	0.99	0.04	41,41,41,41	0
5	ZN	H	301	1/1	1.00	0.03	44,44,44,44	0

6.5 Other polymers [i](#)

There are no such residues in this entry.