wwPDB X-ray Structure Validation Summary Report (i)

Oct 29, 2024 – 01:38 AM EDT

PDB ID	:	3LIF
Title	:	Crystal Structure of the extracellular domain of the putative histidine kinase rpHK1S-Z16
Authors	:	Zhang, Z.; Hendrickson, W.A.
Deposited on	:	2010-01-24
Resolution	:	2.70 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 (i) 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 (i)) were used in the production of this report:

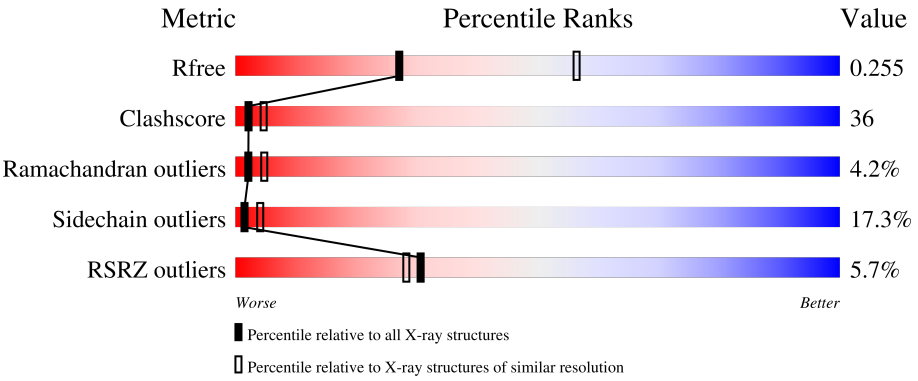
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

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 2.70 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}	164625	3333 (2.70-2.70)
Clashscore	180529	3684 (2.70-2.70)
Ramachandran outliers	177936	3633 (2.70-2.70)
Sidechain outliers	177891	3633 (2.70-2.70)
RSRZ outliers	164620	3333 (2.70-2.70)

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 <=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	254	4% 55% 26% 11%
1	B	254	6% 48% 33% 10% 5%

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
2	MPD	A	1	-	X	X	-
2	MPD	B	2	-	-	X	-
3	CIT	A	294	-	-	X	-

2 Entry composition (i)

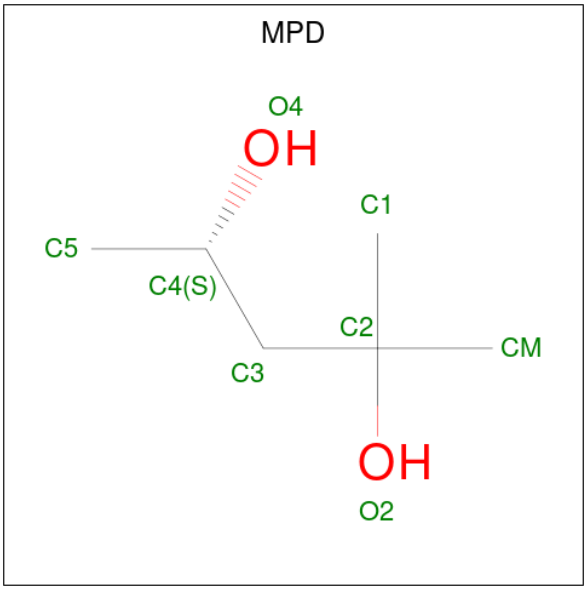
There are 4 unique types of molecules in this entry. The entry contains 3984 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Putative diguanylate cyclase (GGDEF) with PAS/PAC domain.

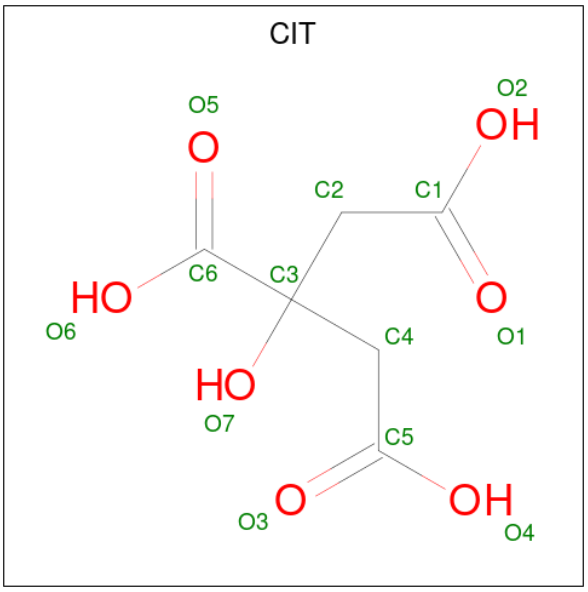
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	243	Total	C	N	O	Se	0	1	0
			1929	1227	336	362	4			
1	B	241	Total	C	N	O	Se	0	1	0
			1920	1224	332	360	4			

- Molecule 2 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula: $C_6H_{14}O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O		0	0
			8	6	2			
2	B	1	Total	C	O		0	0
			8	6	2			

- Molecule 3 is CITRIC ACID (three-letter code: CIT) (formula: $C_6H_8O_7$).



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

- Molecule 4 is water.

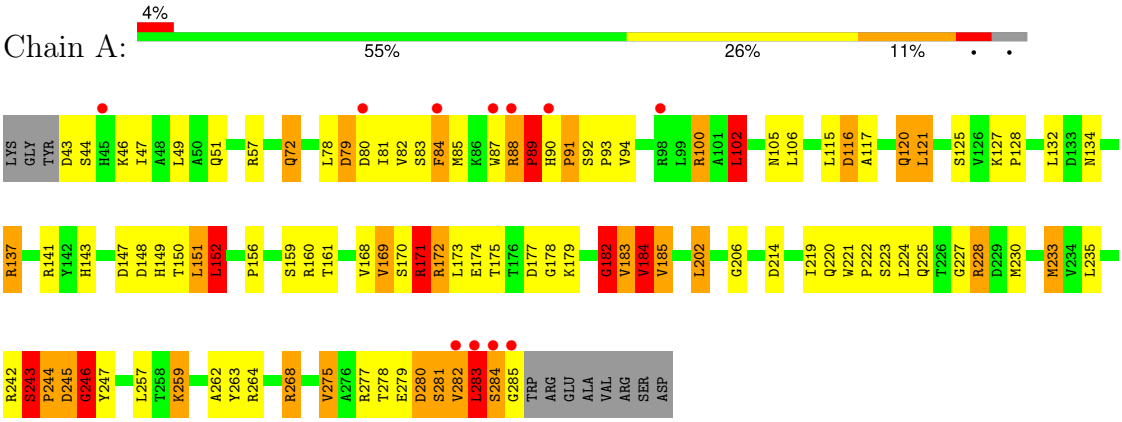
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	57	Total O 57 57	0	0
4	B	36	Total O 36 36	0	0

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.

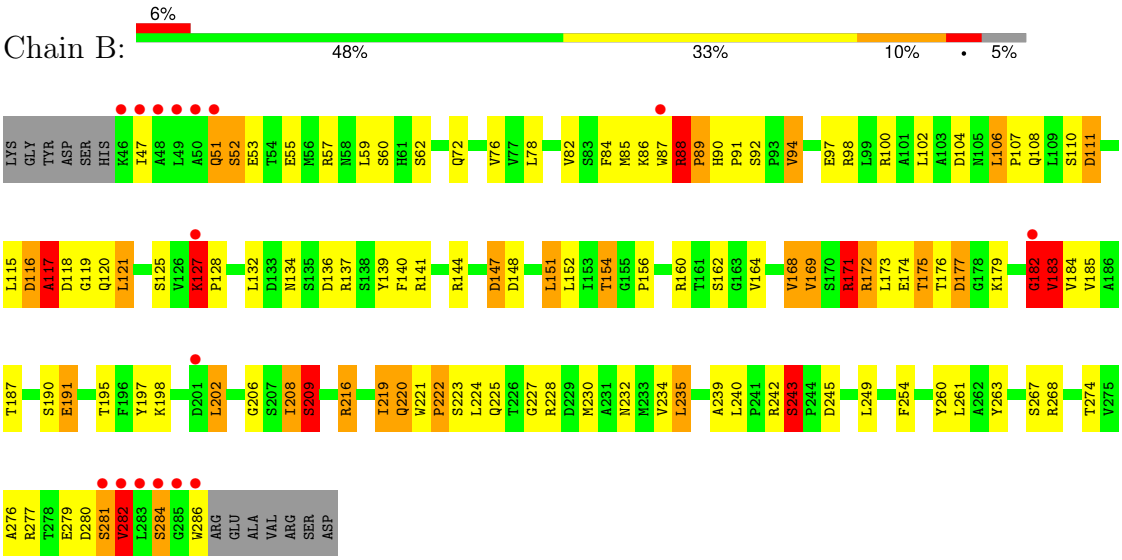
• Molecule 1: Putative diguanylate cyclase (GGDEF) with PAS/PAC domain

Chain A:



• Molecule 1: Putative diguanylate cyclase (GGDEF) with PAS/PAC domain

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 43 3 2	Depositor
Cell constants a, b, c, α, β, γ	184.78Å 184.78Å 184.78Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.39 – 2.70 42.39 – 2.70	Depositor EDS
% Data completeness (in resolution range)	100.0 (42.39-2.70) 99.9 (42.39-2.70)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	9.69 (at 2.69Å)	Xtriage
Refinement program	REFMAC 5.5.0066	Depositor
R, R_{free}	0.205 , 0.256 0.211 , 0.255	Depositor DCC
R_{free} test set	1500 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å^2)	45.8	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 55.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	3984	wwPDB-VP
Average B, all atoms (Å^2)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.88% 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, 0.333 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: MPD, 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	0.99	1/1977 (0.1%)	1.18	16/2685 (0.6%)
1	B	0.94	0/1965	1.09	9/2670 (0.3%)
All	All	0.96	1/3942 (0.0%)	1.13	25/5355 (0.5%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	5
1	B	0	3
All	All	0	8

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	182	GLY	N-CA	5.16	1.53	1.46

The worst 5 of 25 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	184	VAL	CB-CA-C	-9.13	94.06	111.40
1	A	185	VAL	CB-CA-C	-8.38	95.48	111.40
1	A	117	ALA	N-CA-C	-7.29	91.31	111.00
1	A	275	VAL	CB-CA-C	-7.17	97.78	111.40
1	A	184	VAL	CB-CA-C	-7.12	97.86	111.40

There are no chirality outliers.

5 of 8 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	116	ASP	Peptide
1	A	182	GLY	Peptide
1	A	243	SER	Peptide
1	A	244	PRO	Peptide
1	A	88	ARG	Peptide

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	1929	0	1893	133	0
1	B	1920	0	1883	135	0
2	A	8	0	14	18	0
2	B	8	0	14	7	0
3	A	13	0	5	7	0
3	B	13	0	5	4	0
4	A	57	0	0	10	0
4	B	36	0	0	5	0
All	All	3984	0	3814	279	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 36.

The worst 5 of 279 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:B:116:ASP:HB2	1:B:117:ALA:CB	1.38	1.46
1:A:171:ARG:HH11	1:A:171:ARG:CG	1.31	1.36
1:A:230:MSE:HB3	1:A:233:MSE:CE	1.63	1.26
1:B:116:ASP:CB	1:B:117:ALA:HB3	1.64	1.25
1:A:160:ARG:H	2:A:1:MPD:C5	1.52	1.21

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	242/254 (95%)	213 (88%)	18 (7%)	11 (4%)	2 4
1	B	240/254 (94%)	213 (89%)	18 (8%)	9 (4%)	2 6
All	All	482/508 (95%)	426 (88%)	36 (8%)	20 (4%)	2 5

5 of 20 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	280	ASP
1	A	282	VAL
1	B	88	ARG
1	B	127	LYS
1	B	243	SER

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	213/217 (98%)	186 (87%)	27 (13%)	3 9
1	B	211/217 (97%)	165 (78%)	46 (22%)	1 2
All	All	424/434 (98%)	351 (83%)	73 (17%)	1 4

5 of 73 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	183	VAL

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Mol	Chain	Res	Type
1	B	282	VAL
1	B	191	GLU
1	B	220	GLN
1	A	275	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 16 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	158	GLN
1	B	149	HIS
1	A	225	GLN
1	B	143	HIS
1	A	220	GLN

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)

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	$\#\|Z\|>2$
3	CIT	B	294	-	12,12,12	1.46	1 (8%)	17,17,17	2.14	4 (23%)
2	MPD	B	2	-	7,7,7	0.78	0	9,10,10	0.70	0
3	CIT	A	294	-	12,12,12	1.41	1 (8%)	17,17,17	2.65	8 (47%)
2	MPD	A	1	-	7,7,7	1.84	1 (14%)	9,10,10	4.45	6 (66%)

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	CIT	B	294	-	-	1/16/16/16	-
2	MPD	B	2	-	-	0/5/5/5	-
3	CIT	A	294	-	-	6/16/16/16	-
2	MPD	A	1	-	-	5/5/5/5	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1	MPD	O2-C2	4.48	1.55	1.44
3	B	294	CIT	C3-C6	-3.67	1.49	1.53
3	A	294	CIT	C3-C6	2.67	1.56	1.53

The worst 5 of 18 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	1	MPD	CM-C2-C1	7.55	127.54	110.63
2	A	1	MPD	O2-C2-C3	7.47	137.20	109.27
3	A	294	CIT	C4-C3-C2	-5.84	94.32	109.31
2	A	1	MPD	C1-C2-C3	-5.60	86.03	110.20
3	B	294	CIT	C4-C3-C6	-4.81	99.40	110.03

There are no chirality outliers.

5 of 12 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1	MPD	C2-C3-C4-O4
2	A	1	MPD	C2-C3-C4-C5
3	A	294	CIT	O7-C3-C6-O5
3	A	294	CIT	O7-C3-C6-O6

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Mol	Chain	Res	Type	Atoms
3	A	294	CIT	C2-C3-C6-O6

There are no ring outliers.

4 monomers are involved in 36 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	294	CIT	4	0
2	B	2	MPD	7	0
3	A	294	CIT	7	0
2	A	1	MPD	18	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, 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	239/254 (94%)	-0.42	11 (4%)	38	36	18, 33, 57, 90	12 (5%)
1	B	237/254 (93%)	-0.06	16 (6%)	25	23	14, 43, 74, 100	2 (0%)
All	All	476/508 (93%)	-0.24	27 (5%)	30	28	14, 37, 70, 100	14 (2%)

The worst 5 of 27 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	47	ILE	6.3
1	A	87	TRP	5.9
1	B	283	LEU	5.2
1	B	48	ALA	4.8
1	B	51	GLN	4.5

6.2 Non-standard residues in protein, DNA, RNA chains

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates

There are no monosaccharides in this entry.

6.4 Ligands

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, 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	B-factors(Å^2)	Q<0.9
3	CIT	B	294	13/13	0.77	0.18	52,57,59,60	13
2	MPD	A	1	8/8	0.82	0.18	23,30,34,36	0
2	MPD	B	2	8/8	0.88	0.17	39,44,47,48	0
3	CIT	A	294	13/13	0.89	0.10	45,48,54,56	13

6.5 Other polymers (i)

There are no such residues in this entry.