



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 19, 2026 – 02:24 AM UTC

PDB ID : 6B7Q / pdb_00006b7q
Title : Crystal structure of Legionella effector protein sdeA (lpg2157) aa. 211-910
Authors : Mao, Y.; Akturk, A.; Wasilko, J.
Deposited on : 2017-10-04
Resolution : 2.20 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

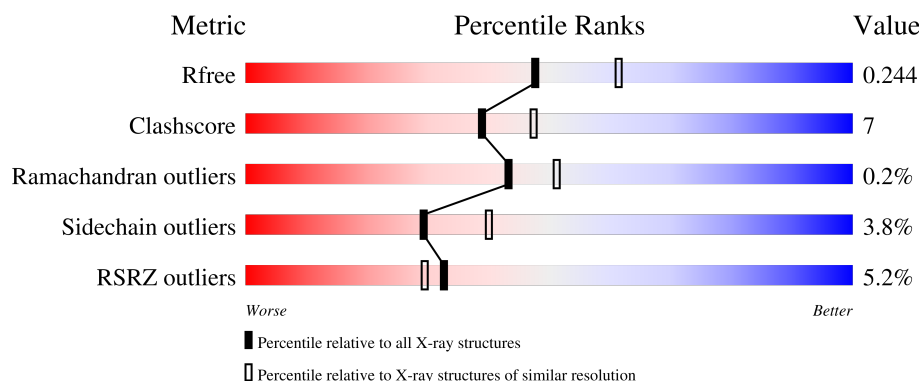
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	705	5% 79% 14% 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	HG	A	1005	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5518 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 SdeA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	667	Total	C	N	O	S	0	0	0
			5338	3362	932	1024	20			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	370	LEU	ILE	conflict	UNP Q6RCR0

- Molecule 2 is MERCURY (II) ION (CCD ID: HG) (formula: Hg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	10	Total	Hg	0	0
			10	10		

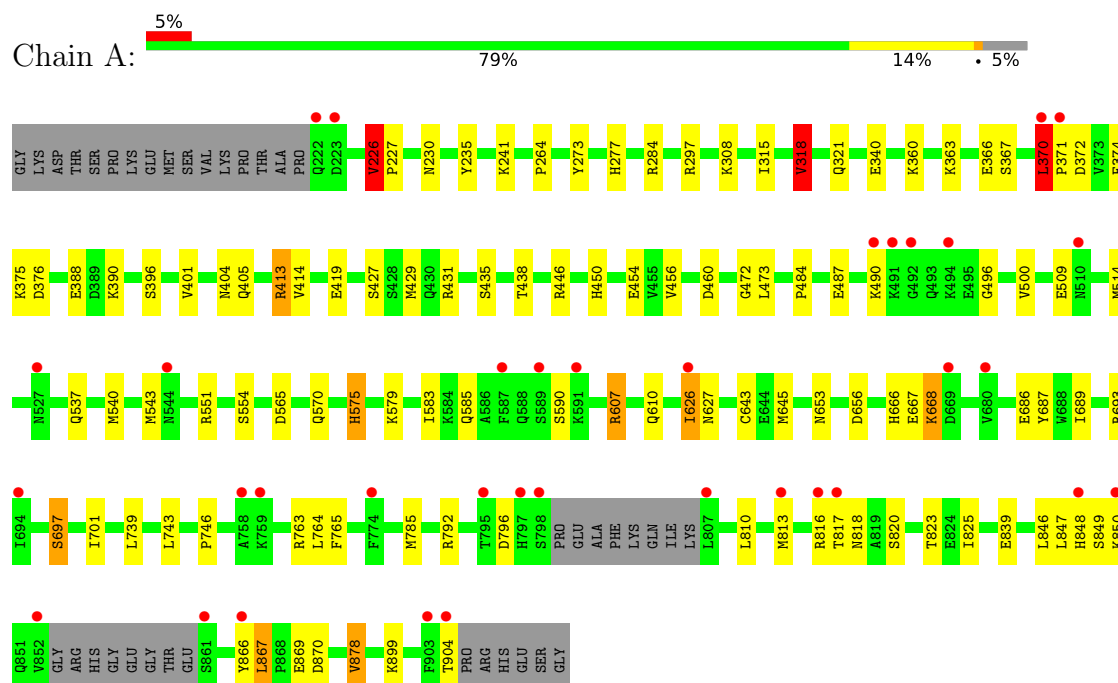
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	170	Total	O	0	0
			170	170		

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: SdeA



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	69.81Å 80.62Å 85.61Å 90.00° 109.87° 90.00°	Depositor
Resolution (Å)	80.51 – 2.20 80.51 – 2.20	Depositor EDS
% Data completeness (in resolution range)	96.5 (80.51-2.20) 96.7 (80.51-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.85 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.195 , 0.241 0.207 , 0.244	Depositor DCC
R_{free} test set	2111 reflections (4.64%)	wwPDB-VP
Wilson B-factor (Å ²)	34.7	Xtriage
Anisotropy	0.583	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 43.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5518	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.26% 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: HG

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	1.08	6/5447 (0.1%)	1.12	12/7358 (0.2%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	374	PHE	C-O	-6.33	1.16	1.24
1	A	226	VAL	N-CA	-5.59	1.39	1.46
1	A	438	THR	N-CA	-5.28	1.39	1.46
1	A	308	LYS	C-O	-5.18	1.17	1.23
1	A	435	SER	N-CA	-5.12	1.40	1.46
1	A	318	VAL	CA-C	5.07	1.58	1.52

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	321	GLN	N-CA-C	-6.88	103.86	111.36
1	A	414	VAL	N-CA-C	6.25	117.37	112.12
1	A	500	VAL	CB-CA-C	-6.05	103.64	110.96
1	A	509	GLU	N-CA-C	-5.45	102.41	110.48
1	A	607	ARG	CB-CA-C	-5.40	102.40	110.88
1	A	490	LYS	N-CA-C	-5.32	103.51	110.53
1	A	372	ASP	N-CA-C	5.27	118.88	112.23
1	A	429	MET	CA-C-N	5.20	127.50	120.38
1	A	429	MET	C-N-CA	5.20	127.50	120.38
1	A	575	HIS	N-CA-C	-5.15	105.58	111.14
1	A	878	VAL	CB-CA-C	-5.14	102.59	110.50
1	A	796	ASP	N-CA-C	5.03	116.94	109.69

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	5338	0	5239	72	0
2	A	10	0	0	2	0
3	A	170	0	0	11	0
All	All	5518	0	5239	72	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 (72) 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:366:GLU:HG2	1:A:370:LEU:CD2	1.47	1.40
1:A:370:LEU:HB3	1:A:371:PRO:CD	1.64	1.27
1:A:366:GLU:CG	1:A:370:LEU:HD23	1.68	1.21
1:A:366:GLU:O	1:A:370:LEU:HB2	1.45	1.13
1:A:370:LEU:HB3	1:A:371:PRO:HD2	1.33	1.07
1:A:366:GLU:CG	1:A:370:LEU:CD2	2.31	0.97
1:A:366:GLU:HG2	1:A:370:LEU:HD21	1.44	0.94
1:A:366:GLU:HG2	1:A:370:LEU:HD23	0.97	0.94
1:A:643:CYS:SG	2:A:1005:HG:HG	1.96	0.82
1:A:645:MET:HE2	1:A:746:PRO:HD3	1.72	0.71
1:A:370:LEU:HB3	1:A:371:PRO:HD3	1.70	0.70
1:A:413:ARG:NH2	3:A:1102:HOH:O	2.23	0.69
1:A:450:HIS:CG	1:A:540:MET:HE2	2.36	0.61
1:A:693:ARG:O	1:A:697:SER:HB3	2.00	0.61
1:A:366:GLU:CB	1:A:370:LEU:HD23	2.31	0.60
1:A:375:LYS:HD3	1:A:376:ASP:OD2	2.01	0.59
1:A:554:SER:HB2	3:A:1142:HOH:O	2.02	0.59
1:A:450:HIS:CE1	1:A:456:VAL:HG21	2.36	0.59
1:A:848:HIS:NE2	3:A:1101:HOH:O	2.22	0.58
1:A:366:GLU:O	1:A:370:LEU:CB	2.37	0.58
1:A:370:LEU:CB	1:A:371:PRO:CD	2.57	0.57
1:A:643:CYS:HG	2:A:1005:HG:HG	1.48	0.56
1:A:846:LEU:HD13	1:A:899:LYS:HB2	1.87	0.56
1:A:241:LYS:CE	1:A:565:ASP:OD1	2.55	0.55

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:816:ARG:O	1:A:818:ASN:OD1	2.25	0.55
1:A:446:ARG:NE	3:A:1107:HOH:O	2.30	0.52
1:A:454:GLU:HB3	3:A:1136:HOH:O	2.10	0.51
1:A:870:ASP:OD2	1:A:899:LYS:NZ	2.42	0.51
1:A:264:PRO:HB3	1:A:273:TYR:CD2	2.45	0.51
1:A:235:TYR:OH	1:A:575:HIS:HD2	1.92	0.51
1:A:570:GLN:HG2	3:A:1108:HOH:O	2.11	0.51
1:A:370:LEU:CB	1:A:371:PRO:HD2	2.24	0.51
1:A:366:GLU:CG	1:A:370:LEU:HD21	2.25	0.50
1:A:668:LYS:CD	1:A:668:LYS:C	2.85	0.49
1:A:450:HIS:ND1	1:A:540:MET:HE2	2.28	0.49
1:A:315:ILE:O	1:A:318:VAL:HG13	2.13	0.48
1:A:869:GLU:HB2	1:A:904:THR:HG22	1.95	0.48
1:A:396:SER:HB2	3:A:1118:HOH:O	2.15	0.47
1:A:764:LEU:HD23	1:A:823:THR:HG22	1.97	0.47
1:A:419:GLU:HG2	1:A:460:ASP:O	2.16	0.46
1:A:763:ARG:NH2	1:A:839:GLU:OE2	2.47	0.46
1:A:765:PHE:CE1	1:A:825:ILE:HD13	2.50	0.46
1:A:401:VAL:O	1:A:405:GLN:HB2	2.16	0.46
1:A:847:LEU:CD2	1:A:867:LEU:CD2	2.94	0.46
1:A:543:MET:HE1	1:A:551:ARG:HD2	1.97	0.45
1:A:666:HIS:ND1	1:A:667:GLU:OE1	2.49	0.45
1:A:226:VAL:HG13	1:A:297:ARG:NH1	2.32	0.44
1:A:820:SER:N	3:A:1113:HOH:O	2.50	0.44
1:A:666:HIS:CE1	1:A:667:GLU:OE1	2.70	0.44
1:A:849:SER:C	1:A:850:LYS:HG2	2.43	0.44
1:A:363:LYS:HD2	1:A:363:LYS:HA	1.78	0.44
1:A:241:LYS:HE3	1:A:565:ASP:OD1	2.18	0.44
1:A:849:SER:HB3	3:A:1185:HOH:O	2.17	0.44
1:A:484:PRO:HB3	1:A:514:MET:CE	2.48	0.43
1:A:626:ILE:HG23	1:A:627:ASN:ND2	2.33	0.43
1:A:230:ASN:C	1:A:230:ASN:OD1	2.59	0.43
1:A:388:GLU:OE2	1:A:390:LYS:NZ	2.49	0.43
1:A:487:GLU:O	1:A:496:GLY:HA3	2.18	0.43
1:A:579:LYS:O	1:A:583:ILE:HD12	2.18	0.43
1:A:686:GLU:O	1:A:687:TYR:C	2.62	0.42
1:A:610:GLN:HG2	1:A:656:ASP:OD1	2.19	0.42
1:A:610:GLN:HG2	1:A:656:ASP:CG	2.45	0.42
1:A:817:THR:HG21	1:A:866:TYR:HD1	1.84	0.42
1:A:472:GLY:O	1:A:473:LEU:C	2.63	0.41
1:A:653:ASN:OD1	1:A:653:ASN:C	2.64	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:739:LEU:O	1:A:743:LEU:HG	2.20	0.41
1:A:277:HIS:HD2	1:A:340:GLU:OE1	2.03	0.41
1:A:847:LEU:HD22	1:A:867:LEU:CD2	2.50	0.41
1:A:792:ARG:HD3	3:A:1151:HOH:O	2.21	0.41
1:A:227:PRO:HD2	3:A:1264:HOH:O	2.21	0.41
1:A:785:MET:HE1	1:A:810:LEU:HD23	2.03	0.40
1:A:366:GLU:HG3	1:A:370:LEU:CD2	2.41	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	661/705 (94%)	639 (97%)	21 (3%)	1 (0%)	43 51

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	370	LEU

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	583/614 (95%)	561 (96%)	22 (4%)	29 40

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	226	VAL
1	A	284	ARG
1	A	318	VAL
1	A	360	LYS
1	A	367	SER
1	A	370	LEU
1	A	404	ASN
1	A	413	ARG
1	A	427	SER
1	A	431	ARG
1	A	537	GLN
1	A	585	GLN
1	A	590	SER
1	A	607	ARG
1	A	626	ILE
1	A	668	LYS
1	A	689	ILE
1	A	697	SER
1	A	701	ILE
1	A	813	MET
1	A	867	LEU
1	A	878	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	245	GLN
1	A	266	ASN
1	A	277	HIS
1	A	348	GLN
1	A	353	GLN
1	A	365	ASN
1	A	377	GLN
1	A	436	GLN
1	A	572	GLN
1	A	575	HIS
1	A	615	ASN
1	A	757	GLN
1	A	818	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 10 ligands modelled in this entry, 10 are 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	667/705 (94%)	0.15	35 (5%) 33 29	22, 45, 85, 113	1 (0%)

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	370	LEU	7.7
1	A	852	VAL	6.1
1	A	807	LEU	4.3
1	A	758	ALA	4.2
1	A	798	SER	3.8
1	A	861	SER	3.8
1	A	491	LYS	3.7
1	A	797	HIS	3.7
1	A	492	GLY	3.3
1	A	817	THR	3.3
1	A	694	ILE	3.2
1	A	816	ARG	3.1
1	A	850	LYS	3.0
1	A	494	LYS	3.0
1	A	371	PRO	2.7
1	A	591	LYS	2.7
1	A	490	LYS	2.7
1	A	527	ASN	2.7
1	A	774	PHE	2.4
1	A	510	ASN	2.4
1	A	795	THR	2.4
1	A	680	VAL	2.4
1	A	222	GLN	2.3
1	A	223	ASP	2.3
1	A	669	ASP	2.3
1	A	626	ILE	2.3
1	A	866	TYR	2.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	544	ASN	2.3
1	A	813	MET	2.2
1	A	904	THR	2.2
1	A	587	PHE	2.2
1	A	759	LYS	2.1
1	A	848	HIS	2.1
1	A	589	SER	2.1
1	A	903	PHE	2.1

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	HG	A	1010	1/1	0.89	0.10	55,55,55,55	1
2	HG	A	1007	1/1	0.91	0.10	71,71,71,71	1
2	HG	A	1001	1/1	0.93	0.07	58,58,58,58	1
2	HG	A	1009	1/1	0.94	0.07	64,64,64,64	1
2	HG	A	1002	1/1	0.94	0.08	60,60,60,60	1
2	HG	A	1006	1/1	0.95	0.06	52,52,52,52	1
2	HG	A	1003	1/1	0.95	0.06	72,72,72,72	1
2	HG	A	1008	1/1	0.95	0.06	43,43,43,43	1
2	HG	A	1004	1/1	0.95	0.07	75,75,75,75	1
2	HG	A	1005	1/1	0.95	0.07	60,60,60,60	0

6.5 Other polymers [i](#)

There are no such residues in this entry.