



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 11:34 AM UTC

PDB ID : 5LWF / pdb_00005lwf
Title : Structure of a single domain camelid antibody fragment cAb-G10S in complex with the BlaP beta-lactamase from *Bacillus licheniformis*
Authors : Vettore, N.; Kerff, F.; Pain, C.; Herman, R.; Sauvage, E.; Preumont, S.; Charlier, P.; Dumoulin, M.
Deposited on : 2016-09-16
Resolution : 2.56 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

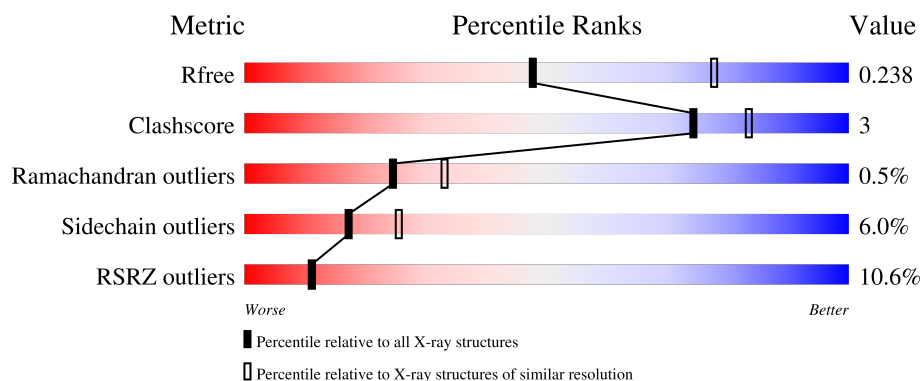
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.56 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	1853 (2.58-2.54)
Clashscore	190562	1897 (2.58-2.54)
Ramachandran outliers	187476	1875 (2.58-2.54)
Sidechain outliers	187428	1875 (2.58-2.54)
RSRZ outliers	180081	1853 (2.58-2.54)

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	273	5% 86% 7% 5%
1	B	273	14% 85% 9% 5%
2	C	131	5% 75% 17% 7%
2	D	131	17% 78% 10% 10%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5927 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 Beta-lactamase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	259	Total	C	N	O	S	0	0	0
			2019	1266	349	401	3			
1	B	258	Total	C	N	O	S	0	0	0
			2014	1263	348	400	3			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	197A	PRO	-	insertion	UNP P00808
A	197B	GLY	-	insertion	UNP P00808
A	296	GLY	-	expression tag	UNP P00808
A	297	PRO	-	expression tag	UNP P00808
A	298	HIS	-	expression tag	UNP P00808
A	299	HIS	-	expression tag	UNP P00808
A	300	HIS	-	expression tag	UNP P00808
A	301	HIS	-	expression tag	UNP P00808
A	302	HIS	-	expression tag	UNP P00808
B	197A	PRO	-	insertion	UNP P00808
B	197B	GLY	-	insertion	UNP P00808
B	296	GLY	-	expression tag	UNP P00808
B	297	PRO	-	expression tag	UNP P00808
B	298	HIS	-	expression tag	UNP P00808
B	299	HIS	-	expression tag	UNP P00808
B	300	HIS	-	expression tag	UNP P00808
B	301	HIS	-	expression tag	UNP P00808
B	302	HIS	-	expression tag	UNP P00808

- Molecule 2 is a protein called Camelid heavy-chain antibody variable fragment cAb-G10S.

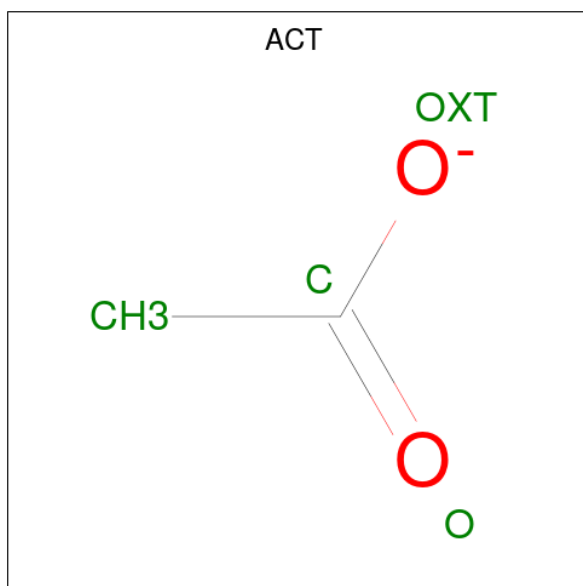
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	122	Total	C	N	O	S	0	0	0
			948	599	163	182	4			

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	118	Total	C	N	O	S	0	0	0
			925	584	159	178	4			

- Molecule 3 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	A	1	Total	C	O	0	0
			4	2	2		
3	C	1	Total	C	O	0	0
			4	2	2		
3	C	1	Total	C	O	0	0
			4	2	2		
3	D	1	Total	C	O	0	0
			4	2	2		

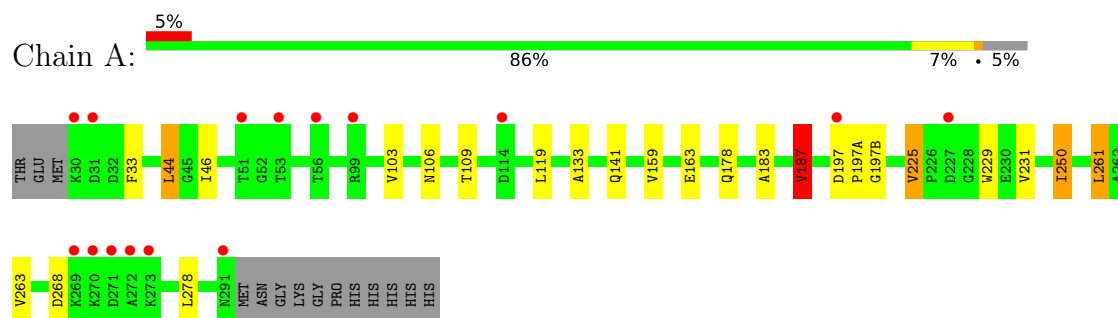
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	1	Total	O	0	0
			1	1		

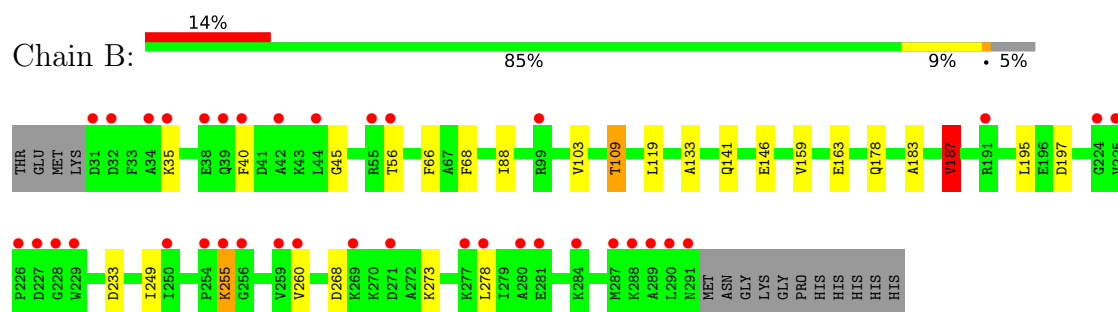
3 Residue-property plots [i](#)

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.

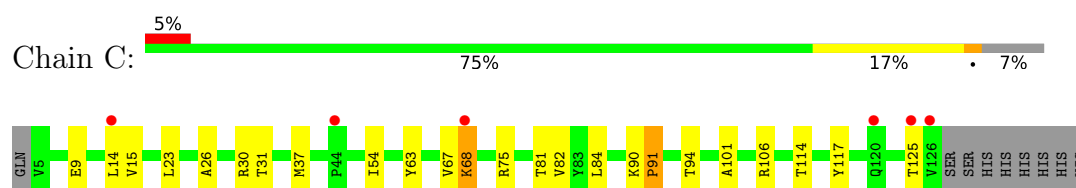
• Molecule 1: Beta-lactamase



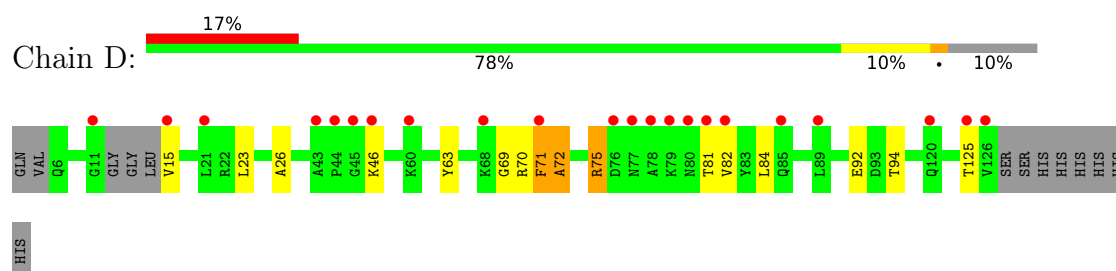
• Molecule 1: Beta-lactamase



• Molecule 2: Camelid heavy-chain antibody variable fragment cAb-G10S



• Molecule 2: Camelid heavy-chain antibody variable fragment cAb-G10S



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	47.28Å 121.69Å 137.03Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.50 – 2.56 40.50 – 2.56	Depositor EDS
% Data completeness (in resolution range)	93.9 (40.50-2.56) 94.3 (40.50-2.56)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.34 (at 2.58Å)	Xtriage
Refinement program	BUSTER 2.10.2	Depositor
R, R_{free}	0.215 , 0.242 (Not available) , 0.238	Depositor DCC
R_{free} test set	1231 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	34.2	Xtriage
Anisotropy	0.535	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 50.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	5927	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.35% 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: ACT

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	0.59	0/2050	1.07	4/2777 (0.1%)
1	B	0.65	0/2045	1.18	5/2770 (0.2%)
2	C	0.64	0/972	1.14	2/1317 (0.2%)
2	D	0.68	0/948	1.12	1/1283 (0.1%)
All	All	0.63	0/6015	1.13	12/8147 (0.1%)

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	197	ASP	CA-CB-CG	-11.04	101.56	112.60
2	D	71	PHE	CA-CB-CG	8.29	122.09	113.80
1	B	197	ASP	CA-CB-CG	-7.58	105.02	112.60
1	B	195	LEU	CA-C-N	6.06	129.22	120.38
1	B	195	LEU	C-N-CA	6.06	129.22	120.38
2	C	91	PRO	CA-C-N	5.89	128.76	120.28
2	C	91	PRO	C-N-CA	5.89	128.76	120.28
1	B	68	PHE	CA-CB-CG	5.84	119.64	113.80
1	B	187	VAL	N-CA-CB	5.62	119.01	110.58
1	A	187	VAL	N-CA-CB	5.62	119.00	110.58
1	A	268	ASP	CA-CB-CG	5.32	117.92	112.60
1	A	197	ASP	CB-CA-C	5.26	117.18	110.13

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	2019	0	2028	12	0
1	B	2014	0	2026	8	0
2	C	948	0	900	9	0
2	D	925	0	873	7	0
3	A	8	0	6	0	0
3	C	8	0	6	0	0
3	D	4	0	3	0	0
4	C	1	0	0	0	0
All	All	5927	0	5842	36	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 (36) 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:109:THR:HG21	1:A:133:ALA:HB3	1.50	0.94
1:A:44:LEU:HD21	1:A:278:LEU:HD11	1.58	0.83
1:B:109:THR:HG21	1:B:133:ALA:HB3	1.70	0.73
2:D:71:PHE:HD1	2:D:72:ALA:H	1.38	0.69
1:A:106:ASN:HB3	1:A:109:THR:CG2	2.27	0.63
1:A:183:ALA:O	1:A:187:VAL:HG13	2.02	0.60
2:D:63:TYR:CE1	2:D:71:PHE:HB2	2.36	0.59
2:C:23:LEU:HD12	2:C:84:LEU:HD23	1.85	0.57
2:D:69:GLY:H	2:D:71:PHE:HD2	1.53	0.56
1:B:183:ALA:O	1:B:187:VAL:HG13	2.05	0.56
2:C:37:MET:HG3	2:C:82:VAL:HG11	1.88	0.55
1:A:106:ASN:HB3	1:A:109:THR:HG23	1.89	0.53
1:B:40:PHE:HB2	1:B:278:LEU:HD13	1.91	0.52
2:C:26:ALA:HA	2:C:81:THR:HG22	1.91	0.51
1:A:33:PHE:CE1	1:A:46:ILE:HD13	2.48	0.48
2:D:23:LEU:HD12	2:D:84:LEU:HD23	1.96	0.47
2:D:26:ALA:HA	2:D:81:THR:HG22	1.97	0.47
1:B:45:GLY:HA3	1:B:66:PHE:HE2	1.80	0.47
1:A:231:VAL:HG22	1:A:250:ILE:HG23	1.97	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:94:THR:HG23	2:D:125:THR:HA	1.99	0.45
1:A:119:LEU:HD21	1:A:141:GLN:HG3	1.98	0.45
2:D:75:ARG:HB3	2:D:82:VAL:HG13	2.00	0.44
2:C:54:ILE:HD13	2:C:75:ARG:HG3	1.99	0.44
1:B:187:VAL:HB	1:B:260:VAL:HG22	2.00	0.44
2:C:94:THR:HG23	2:C:125:THR:HA	1.99	0.44
1:B:45:GLY:HA3	1:B:66:PHE:CE2	2.54	0.43
1:A:261:LEU:HD11	1:A:263:VAL:CG2	2.47	0.43
1:B:119:LEU:HD21	1:B:141:GLN:HG3	2.00	0.43
1:B:255:LYS:HD3	1:B:255:LYS:H	1.85	0.42
2:C:106:ARG:HD3	2:C:114:THR:O	2.19	0.42
2:C:90:LYS:HG2	2:C:91:PRO:HD2	2.02	0.42
1:A:197(A):PRO:CG	1:A:197(A):PRO:O	2.67	0.42
2:C:63:TYR:HB2	2:C:68:LYS:HG2	2.01	0.42
1:A:225:VAL:CG1	1:A:229:TRP:HB2	2.50	0.41
1:A:197(A):PRO:O	1:A:197(A):PRO:HG2	2.20	0.41
2:C:101:ALA:HB3	2:C:117:TYR:HB2	2.02	0.41

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	257/273 (94%)	250 (97%)	5 (2%)	2 (1%)	16	22
1	B	256/273 (94%)	251 (98%)	4 (2%)	1 (0%)	30	39
2	C	120/131 (92%)	115 (96%)	5 (4%)	0	100	100
2	D	114/131 (87%)	106 (93%)	7 (6%)	1 (1%)	14	20
All	All	747/808 (92%)	722 (97%)	21 (3%)	4 (0%)	24	33

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	72	ALA
1	A	197(B)	GLY
1	B	103	VAL
1	A	103	VAL

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	157/230 (68%)	149 (95%)	8 (5%)	21	32
1	B	217/230 (94%)	203 (94%)	14 (6%)	15	22
2	C	96/105 (91%)	89 (93%)	7 (7%)	13	18
2	D	94/105 (90%)	89 (95%)	5 (5%)	20	31
All	All	564/670 (84%)	530 (94%)	34 (6%)	17	25

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	44	LEU
1	A	159	VAL
1	A	163	GLU
1	A	178	GLN
1	A	187	VAL
1	A	225	VAL
1	A	250	ILE
1	A	261	LEU
1	B	35	LYS
1	B	56	THR
1	B	88	ILE
1	B	109	THR
1	B	146	GLU
1	B	159	VAL
1	B	163	GLU
1	B	178	GLN
1	B	187	VAL
1	B	233	ASP

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	249	ILE
1	B	255	LYS
1	B	268	ASP
1	B	273	LYS
2	C	9	GLU
2	C	14	LEU
2	C	15	VAL
2	C	30	ARG
2	C	31	THR
2	C	67	VAL
2	C	68	LYS
2	D	15	VAL
2	D	46	LYS
2	D	70	ARG
2	D	75	ARG
2	D	92	GLU

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	161	ASN
1	A	170	ASN
1	A	173	ASN
1	B	173	ASN
1	B	178	GLN
1	B	245	ASN
2	C	8	GLN
2	C	87	ASN
2	C	120	GLN
2	D	8	GLN
2	D	120	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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

5 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	ACT	C	202	-	3,3,3	1.17	0	3,3,3	0.94	0
3	ACT	C	201	-	3,3,3	1.15	0	3,3,3	0.92	0
3	ACT	A	402	-	3,3,3	1.17	0	3,3,3	0.89	0
3	ACT	A	401	-	3,3,3	1.17	0	3,3,3	0.87	0
3	ACT	D	201	-	3,3,3	1.11	0	3,3,3	0.99	0

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	259/273 (94%)	0.40	15 (5%) 29 29	21, 44, 78, 106	0
1	B	258/273 (94%)	0.77	37 (14%) 6 6	19, 61, 124, 152	0
2	C	122/131 (93%)	0.35	6 (4%) 35 35	16, 44, 70, 81	0
2	D	118/131 (90%)	1.24	22 (18%) 3 3	22, 72, 106, 116	0
All	All	757/808 (93%)	0.65	80 (10%) 11 11	16, 50, 106, 152	0

All (80) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	11	GLY	5.3
2	D	44	PRO	4.6
2	D	78	ALA	4.6
2	C	14	LEU	4.5
1	B	256	GLY	3.9
2	D	77	ASN	3.8
1	A	291	ASN	3.7
1	B	291	ASN	3.6
2	C	120	GLN	3.6
1	B	288	LYS	3.4
1	B	31	ASP	3.4
2	D	126	VAL	3.3
1	B	225	VAL	3.3
2	C	44	PRO	3.2
1	B	278	LEU	3.2
1	B	40	PHE	3.1
2	D	79	LYS	3.1
2	D	43	ALA	3.1
1	A	114	ASP	3.0
2	C	126	VAL	3.0
2	D	81	THR	2.9

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	255	LYS	2.9
1	B	277	LYS	2.9
1	B	44	LEU	2.9
1	B	269	LYS	2.9
2	D	45	GLY	2.8
2	D	89	LEU	2.8
1	B	229	TRP	2.8
1	B	226	PRO	2.8
1	A	272	ALA	2.7
1	B	35	LYS	2.7
1	B	284	LYS	2.7
1	B	280	ALA	2.7
1	B	250	ILE	2.6
2	D	80	ASN	2.6
2	D	68	LYS	2.6
1	B	260	VAL	2.6
2	D	125	THR	2.6
2	D	46	LYS	2.6
1	B	290	LEU	2.5
2	D	15	VAL	2.5
1	B	224	GLY	2.4
2	C	68	LYS	2.4
1	A	227	ASP	2.4
1	A	271	ASP	2.4
2	D	21	LEU	2.4
1	A	99	ARG	2.4
2	D	82	VAL	2.4
2	D	60	LYS	2.4
2	D	120	GLN	2.3
1	B	32	ASP	2.3
1	A	269	LYS	2.3
1	A	31	ASP	2.3
1	B	287	MET	2.3
1	B	271	ASP	2.3
1	B	254	PRO	2.3
1	B	228	GLY	2.2
1	A	197	ASP	2.2
1	B	55	ARG	2.2
1	B	259	VAL	2.2
1	B	39	GLN	2.2
1	A	56	THR	2.2
2	D	85	GLN	2.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	30	LYS	2.2
1	A	270	LYS	2.1
2	C	125	THR	2.1
2	D	76	ASP	2.1
1	B	191	ARG	2.1
1	B	227	ASP	2.1
1	B	289	ALA	2.1
1	B	38	GLU	2.1
1	B	281	GLU	2.1
1	A	273	LYS	2.1
2	D	71	PHE	2.0
1	B	34	ALA	2.0
1	B	99	ARG	2.0
1	A	53	THR	2.0
1	B	42	ALA	2.0
1	A	51	THR	2.0
1	B	56	THR	2.0

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 oligosaccharides 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, 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(Å ²)	Q<0.9
3	ACT	C	202	4/4	0.64	0.23	49,50,50,52	0
3	ACT	A	401	4/4	0.78	0.22	55,56,56,56	0
3	ACT	C	201	4/4	0.83	0.14	33,34,35,35	0
3	ACT	A	402	4/4	0.84	0.12	33,33,34,34	0
3	ACT	D	201	4/4	0.89	0.11	40,40,40,41	0

6.5 Other polymers [i](#)

There are no such residues in this entry.