



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 20, 2016 – 07:03 AM EDT

PDB ID : 5C9I
Title : Structure of N-acylhomoserine lactone acylase MacQ shortened spacer mutant (delta202-208) in uncleaved form
Authors : Yasutake, Y.; Kusada, H.; Kimura, N.
Deposited on : 2015-06-27
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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20027790
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20027790

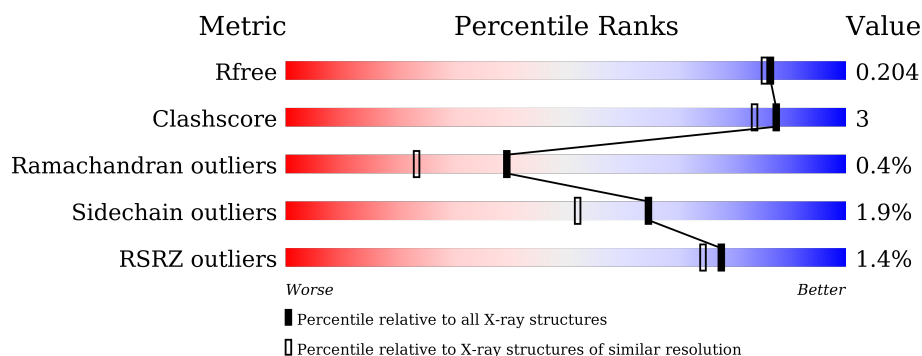
1 Overall quality at a glance

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}	91344	4533 (1.80-1.80)
Clashscore	102246	5383 (1.80-1.80)
Ramachandran outliers	100387	5320 (1.80-1.80)
Sidechain outliers	100360	5319 (1.80-1.80)
RSRZ outliers	91569	4547 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	779	<div> <div>2%</div> <div> <div></div> <div>91%</div> <div>6%</div> <div>..</div> </div> </div>
1	B	779	<div> <div>%</div> <div> <div></div> <div>87%</div> <div>9%</div> <div>..</div> </div> </div>
1	C	779	<div> <div>%</div> <div> <div></div> <div>88%</div> <div>7%</div> <div>..</div> </div> </div>
1	D	779	<div> <div>%</div> <div> <div></div> <div>87%</div> <div>9%</div> <div>..</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 25397 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 Protein related to penicillin acylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	768	Total	C	N	O	S	0	3	0
			5841	3662	1044	1110	25			
1	B	762	Total	C	N	O	S	0	4	0
			5806	3643	1032	1105	26			
1	C	751	Total	C	N	O	S	0	1	0
			5705	3576	1018	1086	25			
1	D	746	Total	C	N	O	S	0	1	0
			5674	3560	1012	1077	25			

There are 88 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	5	SER	-	expression tag	UNP A0A0A1VBK6
A	6	GLY	-	expression tag	UNP A0A0A1VBK6
A	7	GLY	-	expression tag	UNP A0A0A1VBK6
A	8	GLY	-	expression tag	UNP A0A0A1VBK6
A	9	ASP	-	expression tag	UNP A0A0A1VBK6
A	10	GLY	-	expression tag	UNP A0A0A1VBK6
A	11	SER	-	expression tag	UNP A0A0A1VBK6
A	?	-	VAL	deletion	UNP A0A0A1VBK6
A	?	-	GLY	deletion	UNP A0A0A1VBK6
A	?	-	GLY	deletion	UNP A0A0A1VBK6
A	?	-	GLU	deletion	UNP A0A0A1VBK6
A	?	-	LEU	deletion	UNP A0A0A1VBK6
A	?	-	GLY	deletion	UNP A0A0A1VBK6
A	?	-	VAL	deletion	UNP A0A0A1VBK6
A	776	LEU	-	expression tag	UNP A0A0A1VBK6
A	777	GLU	-	expression tag	UNP A0A0A1VBK6
A	778	HIS	-	expression tag	UNP A0A0A1VBK6
A	779	HIS	-	expression tag	UNP A0A0A1VBK6
A	780	HIS	-	expression tag	UNP A0A0A1VBK6
A	781	HIS	-	expression tag	UNP A0A0A1VBK6
A	782	HIS	-	expression tag	UNP A0A0A1VBK6

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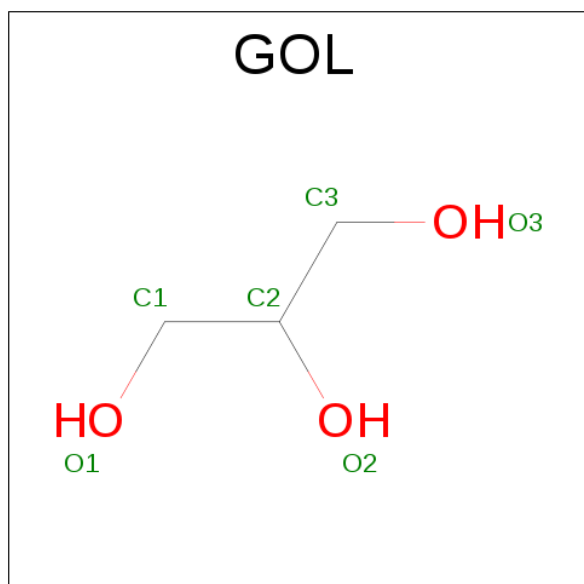
Chain	Residue	Modelled	Actual	Comment	Reference
A	783	HIS	-	expression tag	UNP A0A0A1VBK6
B	5	SER	-	expression tag	UNP A0A0A1VBK6
B	6	GLY	-	expression tag	UNP A0A0A1VBK6
B	7	GLY	-	expression tag	UNP A0A0A1VBK6
B	8	GLY	-	expression tag	UNP A0A0A1VBK6
B	9	ASP	-	expression tag	UNP A0A0A1VBK6
B	10	GLY	-	expression tag	UNP A0A0A1VBK6
B	11	SER	-	expression tag	UNP A0A0A1VBK6
B	?	-	VAL	deletion	UNP A0A0A1VBK6
B	?	-	GLY	deletion	UNP A0A0A1VBK6
B	?	-	GLY	deletion	UNP A0A0A1VBK6
B	?	-	GLU	deletion	UNP A0A0A1VBK6
B	?	-	LEU	deletion	UNP A0A0A1VBK6
B	?	-	GLY	deletion	UNP A0A0A1VBK6
B	?	-	VAL	deletion	UNP A0A0A1VBK6
B	776	LEU	-	expression tag	UNP A0A0A1VBK6
B	777	GLU	-	expression tag	UNP A0A0A1VBK6
B	778	HIS	-	expression tag	UNP A0A0A1VBK6
B	779	HIS	-	expression tag	UNP A0A0A1VBK6
B	780	HIS	-	expression tag	UNP A0A0A1VBK6
B	781	HIS	-	expression tag	UNP A0A0A1VBK6
B	782	HIS	-	expression tag	UNP A0A0A1VBK6
B	783	HIS	-	expression tag	UNP A0A0A1VBK6
C	5	SER	-	expression tag	UNP A0A0A1VBK6
C	6	GLY	-	expression tag	UNP A0A0A1VBK6
C	7	GLY	-	expression tag	UNP A0A0A1VBK6
C	8	GLY	-	expression tag	UNP A0A0A1VBK6
C	9	ASP	-	expression tag	UNP A0A0A1VBK6
C	10	GLY	-	expression tag	UNP A0A0A1VBK6
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C	?	-	GLY	deletion	UNP A0A0A1VBK6
C	?	-	VAL	deletion	UNP A0A0A1VBK6
C	776	LEU	-	expression tag	UNP A0A0A1VBK6
C	777	GLU	-	expression tag	UNP A0A0A1VBK6
C	778	HIS	-	expression tag	UNP A0A0A1VBK6
C	779	HIS	-	expression tag	UNP A0A0A1VBK6
C	780	HIS	-	expression tag	UNP A0A0A1VBK6

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Chain	Residue	Modelled	Actual	Comment	Reference
C	781	HIS	-	expression tag	UNP A0A0A1VBK6
C	782	HIS	-	expression tag	UNP A0A0A1VBK6
C	783	HIS	-	expression tag	UNP A0A0A1VBK6
D	5	SER	-	expression tag	UNP A0A0A1VBK6
D	6	GLY	-	expression tag	UNP A0A0A1VBK6
D	7	GLY	-	expression tag	UNP A0A0A1VBK6
D	8	GLY	-	expression tag	UNP A0A0A1VBK6
D	9	ASP	-	expression tag	UNP A0A0A1VBK6
D	10	GLY	-	expression tag	UNP A0A0A1VBK6
D	11	SER	-	expression tag	UNP A0A0A1VBK6
D	?	-	VAL	deletion	UNP A0A0A1VBK6
D	?	-	GLY	deletion	UNP A0A0A1VBK6
D	?	-	GLY	deletion	UNP A0A0A1VBK6
D	?	-	GLU	deletion	UNP A0A0A1VBK6
D	?	-	LEU	deletion	UNP A0A0A1VBK6
D	?	-	GLY	deletion	UNP A0A0A1VBK6
D	?	-	VAL	deletion	UNP A0A0A1VBK6
D	776	LEU	-	expression tag	UNP A0A0A1VBK6
D	777	GLU	-	expression tag	UNP A0A0A1VBK6
D	778	HIS	-	expression tag	UNP A0A0A1VBK6
D	779	HIS	-	expression tag	UNP A0A0A1VBK6
D	780	HIS	-	expression tag	UNP A0A0A1VBK6
D	781	HIS	-	expression tag	UNP A0A0A1VBK6
D	782	HIS	-	expression tag	UNP A0A0A1VBK6
D	783	HIS	-	expression tag	UNP A0A0A1VBK6

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		

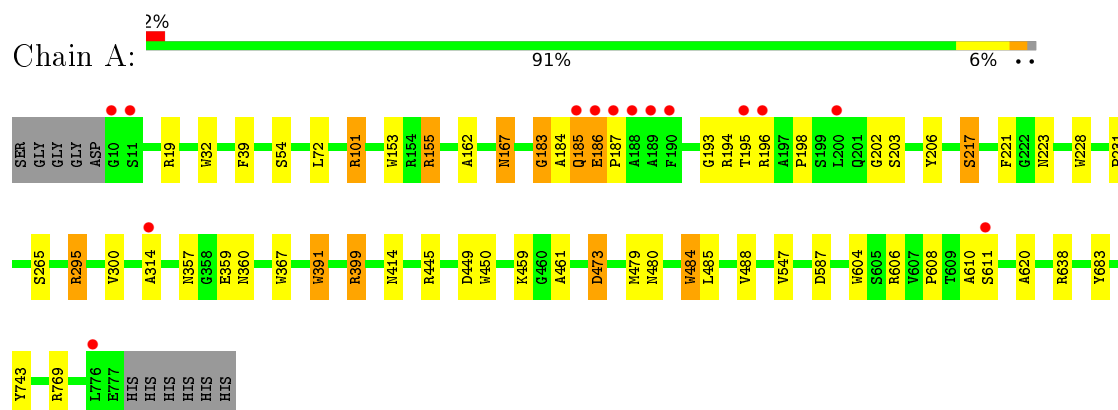
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	658	Total	O	0	0
			658	658		
3	B	636	Total	O	0	0
			636	636		
3	C	518	Total	O	0	0
			518	518		
3	D	553	Total	O	0	0
			553	553		

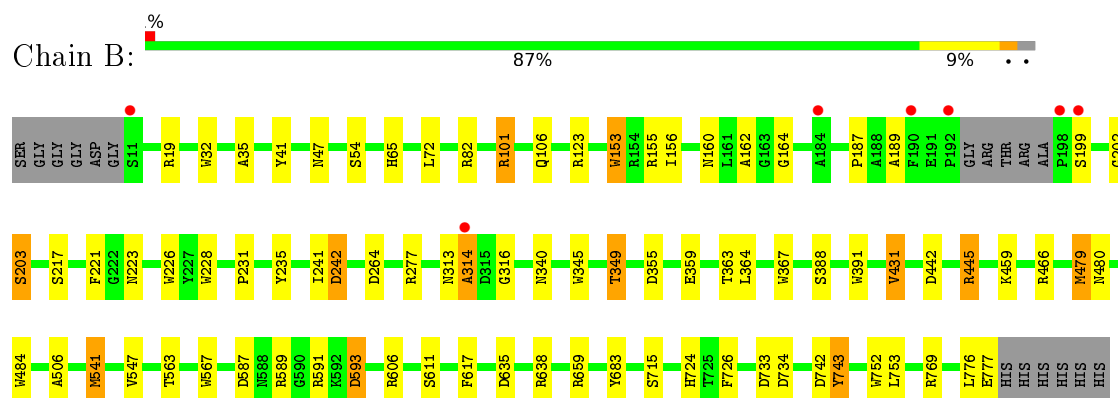
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

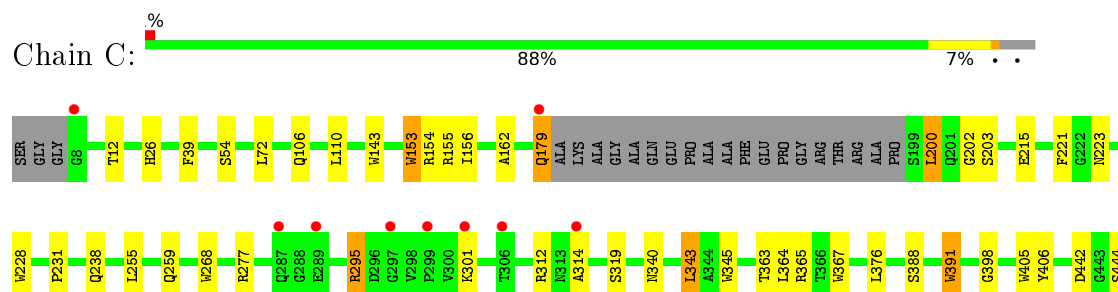
- Molecule 1: Protein related to penicillin acylase



- Molecule 1: Protein related to penicillin acylase

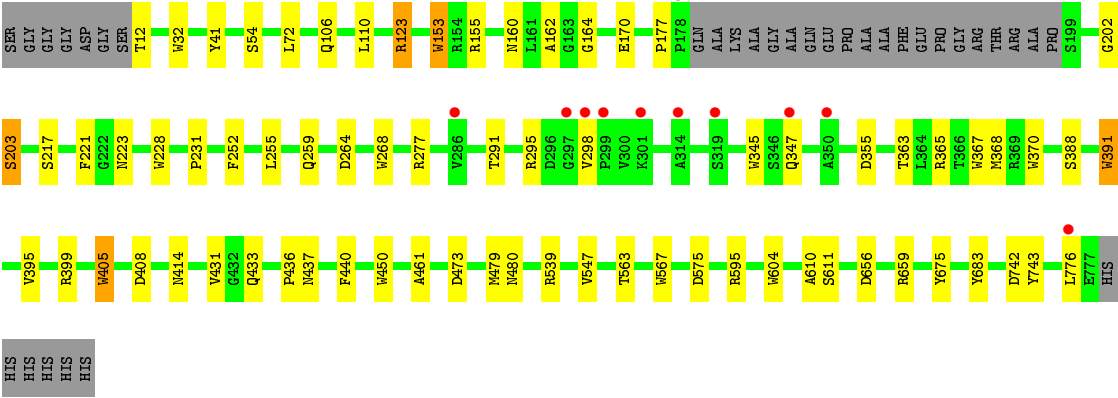
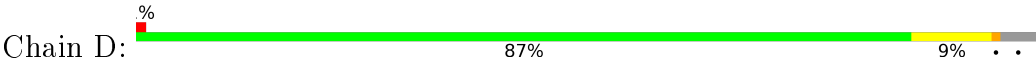


- Molecule 1: Protein related to penicillin acylase





● Molecule 1: Protein related to penicillin acylase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	102.72Å 137.77Å 121.52Å 90.00° 111.45° 90.00°	Depositor
Resolution (Å)	50.00 – 1.80 48.05 – 1.80	Depositor EDS
% Data completeness (in resolution range)	99.8 (50.00-1.80) 99.9 (48.05-1.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.52 (at 1.79Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.166 , 0.206 0.164 , 0.204	Depositor DCC
R_{free} test set	14548 reflections (5.28%)	DCC
Wilson B-factor (Å ²)	18.4	Xtriage
Anisotropy	0.264	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 43.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	25397	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 14.74% 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

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	1.09	9/5997 (0.2%)	1.00	15/8168 (0.2%)
1	B	1.11	12/5964 (0.2%)	1.03	21/8121 (0.3%)
1	C	1.02	11/5853 (0.2%)	0.95	14/7970 (0.2%)
1	D	1.08	14/5822 (0.2%)	0.98	19/7929 (0.2%)
All	All	1.07	46/23636 (0.2%)	0.99	69/32188 (0.2%)

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	1
1	B	0	3
All	All	0	4

All (46) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	752	TRP	CD2-CE2	8.40	1.51	1.41
1	B	593	ASP	CB-CG	8.14	1.68	1.51
1	D	450	TRP	CD2-CE2	7.20	1.50	1.41
1	B	228	TRP	CD2-CE2	7.12	1.49	1.41
1	D	567	TRP	CD2-CE2	7.06	1.49	1.41
1	B	484	TRP	CD2-CE2	6.92	1.49	1.41
1	D	347	GLN	C-O	6.74	1.36	1.23
1	B	217	SER	CB-OG	6.71	1.50	1.42
1	D	391	TRP	CD2-CE2	6.64	1.49	1.41
1	C	391	TRP	CD2-CE2	6.38	1.49	1.41
1	B	153	TRP	CD2-CE2	6.37	1.49	1.41
1	A	228	TRP	CD2-CE2	6.35	1.49	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	32	TRP	CD2-CE2	6.33	1.49	1.41
1	D	405	TRP	CD2-CE2	6.31	1.49	1.41
1	C	367	TRP	CD2-CE2	6.23	1.48	1.41
1	C	345	TRP	CD2-CE2	6.23	1.48	1.41
1	D	268	TRP	CD2-CE2	6.09	1.48	1.41
1	A	391	TRP	CD2-CE2	6.08	1.48	1.41
1	B	226	TRP	CD2-CE2	6.06	1.48	1.41
1	D	367	TRP	CD2-CE2	5.83	1.48	1.41
1	C	604	TRP	CD2-CE2	5.80	1.48	1.41
1	A	367	TRP	CD2-CE2	5.73	1.48	1.41
1	B	367	TRP	CD2-CE2	5.72	1.48	1.41
1	B	567	TRP	CD2-CE2	5.71	1.48	1.41
1	B	235	TYR	CE1-CZ	5.69	1.46	1.38
1	A	604	TRP	CD2-CE2	5.66	1.48	1.41
1	A	32	TRP	CD2-CE2	5.59	1.48	1.41
1	C	153	TRP	CD2-CE2	5.58	1.48	1.41
1	D	370	TRP	CG-CD1	5.55	1.44	1.36
1	C	268	TRP	CD2-CE2	5.51	1.48	1.41
1	A	484	TRP	CD2-CE2	5.46	1.48	1.41
1	C	228	TRP	CD2-CE2	5.45	1.47	1.41
1	B	345	TRP	CD2-CE2	5.42	1.47	1.41
1	C	405	TRP	CD2-CE2	5.39	1.47	1.41
1	A	153	TRP	CD2-CE2	5.38	1.47	1.41
1	C	143	TRP	CD2-CE2	5.35	1.47	1.41
1	D	228	TRP	CD2-CE2	5.33	1.47	1.41
1	C	450	TRP	CD2-CE2	5.32	1.47	1.41
1	D	675	TYR	CG-CD2	5.31	1.46	1.39
1	C	473	ASP	CB-CG	-5.28	1.40	1.51
1	D	153	TRP	CD2-CE2	5.25	1.47	1.41
1	D	345	TRP	CD2-CE2	5.23	1.47	1.41
1	A	206	TYR	CE1-CZ	5.18	1.45	1.38
1	A	450	TRP	CD2-CE2	5.12	1.47	1.41
1	D	32	TRP	CD2-CE2	5.09	1.47	1.41
1	D	604	TRP	CD2-CE2	5.07	1.47	1.41

All (69) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	541	MET	CG-SD-CE	-14.40	77.16	100.20
1	A	295	ARG	NE-CZ-NH2	-10.08	115.26	120.30
1	D	264	ASP	CB-CG-OD1	8.58	126.02	118.30
1	B	466	ARG	NE-CZ-NH2	-8.36	116.12	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	277	ARG	NE-CZ-NH2	-8.22	116.19	120.30
1	C	595	ARG	NE-CZ-NH1	8.01	124.31	120.30
1	D	659	ARG	NE-CZ-NH2	7.98	124.29	120.30
1	D	473	ASP	CB-CG-OD1	-7.93	111.17	118.30
1	A	19	ARG	NE-CZ-NH1	-7.80	116.40	120.30
1	B	277	ARG	NE-CZ-NH2	-7.72	116.44	120.30
1	B	19	ARG	NE-CZ-NH1	-7.58	116.51	120.30
1	C	473	ASP	CB-CG-OD1	-7.22	111.81	118.30
1	D	659	ARG	NE-CZ-NH1	-7.16	116.72	120.30
1	B	635	ASP	CB-CG-OD1	7.13	124.72	118.30
1	C	466	ARG	NE-CZ-NH1	-7.05	116.77	120.30
1	B	264	ASP	CB-CG-OD1	6.99	124.59	118.30
1	B	638	ARG	NE-CZ-NH1	-6.94	116.83	120.30
1	D	277	ARG	NE-CZ-NH2	-6.82	116.89	120.30
1	D	277	ARG	NE-CZ-NH1	6.79	123.69	120.30
1	D	595	ARG	NE-CZ-NH2	-6.68	116.96	120.30
1	A	587	ASP	CB-CG-OD1	6.65	124.29	118.30
1	A	19	ARG	NE-CZ-NH2	6.65	123.62	120.30
1	B	466	ARG	NE-CZ-NH1	6.52	123.56	120.30
1	C	595	ARG	NE-CZ-NH2	-6.46	117.07	120.30
1	D	595	ARG	NE-CZ-NH1	6.45	123.53	120.30
1	D	295	ARG	NE-CZ-NH2	-6.45	117.08	120.30
1	A	473	ASP	CB-CA-C	-6.21	97.98	110.40
1	C	466	ARG	NE-CZ-NH2	6.18	123.39	120.30
1	B	101	ARG	NE-CZ-NH1	-6.16	117.22	120.30
1	B	355	ASP	CB-CG-OD2	-6.16	112.76	118.30
1	D	399	ARG	NE-CZ-NH2	-6.09	117.26	120.30
1	B	734	ASP	CB-CG-OD2	-6.05	112.86	118.30
1	B	742	ASP	CB-CG-OD1	5.92	123.63	118.30
1	B	242	ASP	CB-CG-OD1	5.92	123.63	118.30
1	A	449	ASP	CB-CG-OD1	5.89	123.60	118.30
1	C	449	ASP	CB-CG-OD1	5.83	123.55	118.30
1	C	110	LEU	CB-CG-CD2	-5.83	101.10	111.00
1	B	606	ARG	NE-CZ-NH2	5.80	123.20	120.30
1	C	277	ARG	NE-CZ-NH1	5.80	123.20	120.30
1	D	539	ARG	NE-CZ-NH1	-5.79	117.41	120.30
1	A	155	ARG	NE-CZ-NH2	-5.77	117.41	120.30
1	D	264	ASP	CB-CG-OD2	-5.62	113.24	118.30
1	D	355	ASP	CB-CG-OD1	5.57	123.31	118.30
1	D	252	PHE	CB-CG-CD2	-5.57	116.90	120.80
1	B	606	ARG	NE-CZ-NH1	-5.55	117.53	120.30
1	B	659	ARG	NE-CZ-NH1	-5.51	117.54	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	606	ARG	NE-CZ-NH1	-5.49	117.56	120.30
1	D	742	ASP	CB-CG-OD1	5.47	123.23	118.30
1	B	593	ASP	CB-CG-OD1	5.46	123.22	118.30
1	C	200	LEU	CA-CB-CG	5.39	127.69	115.30
1	A	186	GLU	C-N-CD	-5.35	108.82	120.60
1	A	101	ARG	NE-CZ-NH1	-5.30	117.65	120.30
1	A	399	ARG	NE-CZ-NH2	5.28	122.94	120.30
1	A	638	ARG	NE-CZ-NH1	-5.27	117.66	120.30
1	A	295	ARG	NE-CZ-NH1	5.24	122.92	120.30
1	C	575	ASP	CB-CG-OD1	5.21	122.99	118.30
1	D	123	ARG	NE-CZ-NH2	-5.21	117.70	120.30
1	B	431	VAL	CG1-CB-CG2	5.20	119.23	110.90
1	D	575	ASP	CB-CG-OD1	5.18	122.97	118.30
1	C	480[A]	ASN	CB-CA-C	-5.18	100.05	110.40
1	C	480[B]	ASN	CB-CA-C	-5.18	100.05	110.40
1	B	445	ARG	NE-CZ-NH1	-5.17	117.72	120.30
1	A	445	ARG	NE-CZ-NH1	-5.11	117.75	120.30
1	D	656	ASP	CB-CG-OD1	5.09	122.88	118.30
1	D	408	ASP	CB-CG-OD1	5.08	122.87	118.30
1	B	355	ASP	CB-CG-OD1	5.07	122.86	118.30
1	B	587	ASP	CB-CG-OD1	5.06	122.85	118.30
1	A	769	ARG	NE-CZ-NH1	-5.05	117.78	120.30
1	C	154	ARG	NE-CZ-NH1	5.04	122.82	120.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	183	GLY	Peptide
1	B	199	SER	Peptide
1	B	313	ASN	Peptide
1	B	314	ALA	Peptide

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	5841	0	5582	30	0
1	B	5806	0	5544	35	0
1	C	5705	0	5442	30	0
1	D	5674	0	5419	29	0
2	A	6	0	8	0	0
3	A	658	0	0	3	0
3	B	636	0	0	7	0
3	C	518	0	0	2	0
3	D	553	0	0	4	0
All	All	25397	0	21995	119	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 (119) 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:185:GLN:O	1:A:186:GLU:HG2	1.63	0.97
1:C:179:GLN:HA	1:C:179:GLN:OE1	1.63	0.95
1:A:185:GLN:HE21	1:A:185:GLN:HA	1.35	0.91
1:A:202:GLY:O	1:A:480[A]:ASN:ND2	2.10	0.85
1:B:563:THR:HG23	3:B:1262:HOH:O	1.79	0.82
1:B:202:GLY:O	1:B:480[A]:ASN:ND2	2.12	0.81
1:C:365:ARG:HH11	1:C:365:ARG:HB3	1.46	0.80
1:C:215:GLU:HB3	3:C:1196:HOH:O	1.83	0.79
1:C:365:ARG:HH11	1:C:365:ARG:CB	1.95	0.78
1:A:185:GLN:O	1:A:186:GLU:CG	2.30	0.78
1:B:202:GLY:C	1:B:480[A]:ASN:OD1	2.23	0.76
1:D:202:GLY:C	1:D:480[A]:ASN:OD1	2.25	0.75
1:C:12:THR:HG22	1:C:776:LEU:HD23	1.71	0.73
1:C:179:GLN:CA	1:C:179:GLN:OE1	2.40	0.70
1:C:221:PHE:CZ	1:C:223:ASN:HB2	2.27	0.70
1:A:202:GLY:C	1:A:480[A]:ASN:OD1	2.30	0.69
1:D:202:GLY:CA	1:D:480[A]:ASN:OD1	2.40	0.69
1:C:365:ARG:HB3	1:C:365:ARG:NH1	2.06	0.69
1:A:194:ARG:HA	1:B:189:ALA:HB3	1.76	0.68
1:A:185:GLN:HA	1:A:185:GLN:NE2	2.10	0.66
1:D:170:GLU:HG3	3:D:805:HOH:O	1.95	0.66
1:D:202:GLY:O	1:D:480[A]:ASN:ND2	2.30	0.65
1:D:255:LEU:HD12	1:D:259:GLN:OE1	1.98	0.64
1:B:101:ARG:NH1	1:B:359:GLU:OE2	2.30	0.64
1:B:442:ASP:OD2	1:B:445:ARG:HD3	1.98	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:202:GLY:CA	1:B:480[A]:ASN:OD1	2.47	0.62
1:C:106:GLN:HG3	1:C:153:TRP:CH2	2.34	0.62
1:D:12:THR:HG22	1:D:776:LEU:HA	1.82	0.61
1:B:479:MET:O	1:B:480[A]:ASN:HB2	2.01	0.61
1:D:170:GLU:CG	3:D:805:HOH:O	2.49	0.58
1:D:221:PHE:CZ	1:D:223:ASN:HB2	2.39	0.58
1:B:777:GLU:HG2	3:B:1155:HOH:O	2.04	0.58
1:C:202:GLY:O	1:C:480[A]:ASN:ND2	2.36	0.57
1:B:589:ARG:NH1	1:B:591:ARG:NH1	2.53	0.57
1:B:65:HIS:HE1	3:B:977:HOH:O	1.86	0.57
1:D:110:LEU:HG	1:D:368:MET:SD	2.45	0.57
1:B:541:MET:HE3	3:B:1011:HOH:O	2.06	0.56
1:C:340:ASN:HB3	1:C:343:LEU:HD22	1.86	0.55
1:A:295:ARG:HG3	1:A:300:VAL:HG21	1.87	0.55
1:D:610:ALA:O	1:D:611:SER:HB2	2.06	0.55
1:A:101:ARG:HD2	3:A:1429:HOH:O	2.07	0.55
1:D:479:MET:O	1:D:480[A]:ASN:HB2	2.07	0.55
1:D:431:VAL:HG21	1:D:440:PHE:CD2	2.42	0.55
1:A:221:PHE:CZ	1:A:223:ASN:HB2	2.42	0.55
1:B:541:MET:CE	3:B:1011:HOH:O	2.55	0.54
1:A:202:GLY:CA	1:A:480[A]:ASN:OD1	2.55	0.54
1:C:202:GLY:CA	1:C:480[A]:ASN:OD1	2.56	0.54
1:B:349:THR:HB	3:B:1327:HOH:O	2.07	0.54
1:C:445:ARG:HB3	1:C:447:GLU:CD	2.28	0.53
1:A:185:GLN:HE21	1:A:185:GLN:CA	2.08	0.53
1:D:170:GLU:CD	3:D:805:HOH:O	2.46	0.53
1:B:363:THR:HG22	1:B:388:SER:O	2.09	0.53
1:C:156:ILE:HD13	1:C:364:LEU:HD22	1.91	0.53
1:B:82:ARG:HD3	1:B:189:ALA:HA	1.90	0.52
1:A:265:SER:OG	1:A:399:ARG:NH1	2.37	0.52
1:D:563:THR:HG23	3:D:1102:HOH:O	2.11	0.51
1:A:610:ALA:O	1:A:611:SER:HB2	2.10	0.50
1:C:459:LYS:HD3	1:C:459:LYS:N	2.26	0.50
1:A:54:SER:HB3	1:A:231:PRO:HB3	1.93	0.50
1:B:202:GLY:O	1:B:203:SER:HB2	2.11	0.50
1:D:436:PRO:O	1:D:437:ASN:HB2	2.11	0.50
1:D:162:ALA:HB3	1:D:391:TRP:CE2	2.47	0.50
1:B:221:PHE:CZ	1:B:223:ASN:HB2	2.47	0.49
1:D:363:THR:HG22	1:D:388:SER:O	2.12	0.49
1:B:106:GLN:HG3	1:B:153:TRP:CH2	2.48	0.49
1:C:442:ASP:OD1	1:C:444:SER:HB2	2.12	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:202:GLY:C	1:C:480[A]:ASN:OD1	2.52	0.48
1:D:160:ASN:HD21	1:D:363:THR:CG2	2.27	0.48
1:D:202:GLY:O	1:D:203:SER:HB2	2.14	0.48
1:B:479:MET:O	1:B:480[A]:ASN:CB	2.60	0.47
1:A:399:ARG:HD3	3:A:1387:HOH:O	2.15	0.47
1:A:485:LEU:HD22	1:A:488:VAL:HA	1.96	0.47
1:B:547:VAL:HG22	1:B:683:TYR:CZ	2.50	0.47
1:D:41:TYR:CD1	1:D:123:ARG:HD3	2.50	0.46
1:A:162:ALA:HB3	1:A:391:TRP:CE2	2.50	0.46
1:B:776:LEU:O	1:B:777:GLU:HB2	2.16	0.46
1:A:167:ASN:OD1	1:A:198:PRO:HA	2.16	0.46
1:C:295:ARG:HG2	3:C:1174:HOH:O	2.15	0.46
1:C:54:SER:HB3	1:C:231:PRO:HB3	1.97	0.45
1:B:41:TYR:CD1	1:B:123:ARG:HD3	2.52	0.45
1:B:506:ALA:HA	1:B:617:PHE:CD1	2.52	0.45
1:A:217:SER:HB2	3:A:1192:HOH:O	2.17	0.45
1:A:620:ALA:HB1	1:C:497:ALA:HB1	1.98	0.45
1:A:101:ARG:HH12	1:A:359:GLU:CD	2.20	0.45
1:C:376:LEU:HB2	1:C:398:GLY:HA3	1.97	0.45
1:A:608:PRO:HB3	1:B:316:GLY:HA3	2.00	0.44
1:D:395:VAL:HA	1:D:405:TRP:O	2.17	0.44
1:B:726:PHE:HA	1:B:743:TYR:CE1	2.52	0.44
1:C:255:LEU:HD12	1:C:259:GLN:OE1	2.17	0.44
1:B:35:ALA:HB2	1:B:241:ILE:HG12	2.00	0.43
1:D:255:LEU:HD12	1:D:259:GLN:CD	2.37	0.43
1:A:414:ASN:HB2	1:A:461:ALA:C	2.39	0.43
1:C:312:ARG:HG2	1:C:312:ARG:HH11	1.84	0.43
1:A:183:GLY:H	1:B:340:ASN:HD21	1.66	0.43
1:C:445:ARG:HB3	1:C:447:GLU:OE1	2.19	0.43
1:D:479:MET:O	1:D:480[A]:ASN:CB	2.67	0.43
1:D:547:VAL:HG22	1:D:683:TYR:CZ	2.53	0.43
1:A:196:ARG:HH11	1:A:196:ARG:HG2	1.84	0.43
1:B:47:ASN:HA	1:B:733:ASP:OD1	2.19	0.43
1:C:363:THR:HG22	1:C:388:SER:O	2.18	0.43
1:B:54:SER:HB3	1:B:231:PRO:HB3	2.01	0.43
1:C:459:LYS:CD	1:C:459:LYS:N	2.82	0.42
1:C:26:HIS:HA	1:C:238:GLN:HB3	2.02	0.42
1:D:177:PRO:HG3	1:D:291:THR:CG2	2.49	0.42
1:A:484:TRP:CE3	1:A:485:LEU:HG	2.54	0.42
1:B:715:SER:HB2	3:B:1115:HOH:O	2.19	0.41
1:B:162:ALA:HB3	1:B:391:TRP:CE2	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:506:ALA:HA	1:C:617:PHE:CD1	2.55	0.41
1:D:414:ASN:HB2	1:D:461:ALA:C	2.40	0.41
1:A:357:ASN:HA	1:A:360:ASN:HB3	2.01	0.41
1:A:193:GLY:HA3	1:B:187:PRO:O	2.21	0.41
1:D:54:SER:HB3	1:D:231:PRO:HB3	2.01	0.41
1:B:156:ILE:HD13	1:B:364:LEU:HD22	2.01	0.41
1:D:106:GLN:HG3	1:D:153:TRP:CH2	2.55	0.41
1:A:547:VAL:HG22	1:A:683:TYR:CZ	2.55	0.41
1:D:160:ASN:HD21	1:D:363:THR:HG21	1.85	0.40
1:B:724:HIS:HA	1:B:753:LEU:O	2.21	0.40
1:C:162:ALA:HB3	1:C:391:TRP:CE2	2.57	0.40
1:C:406:TYR:C	1:C:406:TYR:CD2	2.94	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	769/779 (99%)	739 (96%)	26 (3%)	4 (0%)	34	17
1	B	762/779 (98%)	730 (96%)	28 (4%)	4 (0%)	34	17
1	C	748/779 (96%)	716 (96%)	30 (4%)	2 (0%)	46	29
1	D	743/779 (95%)	715 (96%)	26 (4%)	2 (0%)	46	29
All	All	3022/3116 (97%)	2900 (96%)	110 (4%)	12 (0%)	39	23

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	184	ALA
1	A	187	PRO
1	B	314	ALA

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Mol	Chain	Res	Type
1	A	203	SER
1	A	314	ALA
1	B	203	SER
1	C	203	SER
1	D	203	SER
1	B	242	ASP
1	C	314	ALA
1	D	164	GLY
1	B	164	GLY

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	590/595 (99%)	579 (98%)	11 (2%)	65	52
1	B	588/595 (99%)	577 (98%)	11 (2%)	65	52
1	C	578/595 (97%)	564 (98%)	14 (2%)	57	41
1	D	575/595 (97%)	568 (99%)	7 (1%)	78	71
All	All	2331/2380 (98%)	2288 (98%)	43 (2%)	65	54

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	39	PHE
1	A	72	LEU
1	A	155	ARG
1	A	167	ASN
1	A	185	GLN
1	A	195	THR
1	A	217	SER
1	A	459	LYS
1	A	473	ASP
1	A	479	MET
1	A	743	TYR
1	B	72	LEU

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Mol	Chain	Res	Type
1	B	155	ARG
1	B	160	ASN
1	B	349	THR
1	B	431	VAL
1	B	459	LYS
1	B	479	MET
1	B	593	ASP
1	B	611	SER
1	B	743	TYR
1	B	769	ARG
1	C	39	PHE
1	C	72	LEU
1	C	155	ARG
1	C	179	GLN
1	C	200	LEU
1	C	295	ARG
1	C	301	LYS
1	C	319	SER
1	C	343	LEU
1	C	459	LYS
1	C	541	MET
1	C	555	ASP
1	C	743	TYR
1	C	774	LYS
1	D	72	LEU
1	D	155	ARG
1	D	217	SER
1	D	298	VAL
1	D	365	ARG
1	D	433	GLN
1	D	743	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	185	GLN
1	B	65	HIS
1	B	176	GLN
1	D	160	ASN

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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

1 ligand is 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	$\# Z > 2$
2	GOL	A	801	-	5,5,5	0.19	0	5,5,5	1.06	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
2	GOL	A	801	-	-	0/4/4/4	0/0/0/0

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

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 ⓘ

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	768/779 (98%)	-0.62	14 (1%) 71 67	11, 17, 34, 66	0
1	B	762/779 (97%)	-0.65	7 (0%) 85 83	11, 17, 31, 88	0
1	C	751/779 (96%)	-0.45	9 (1%) 81 78	12, 20, 46, 74	0
1	D	746/779 (95%)	-0.50	11 (1%) 76 72	11, 19, 42, 64	0
All	All	3027/3116 (97%)	-0.56	41 (1%) 78 74	11, 18, 40, 88	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	198	PRO	6.5
1	B	192	PRO	5.7
1	A	11	SER	4.5
1	B	314	ALA	4.5
1	C	179	GLN	4.3
1	B	199	SER	4.0
1	C	287	GLN	4.0
1	A	200	LEU	3.8
1	A	314	ALA	3.7
1	D	776	LEU	3.7
1	C	301	LYS	3.7
1	B	190	PHE	3.5
1	D	319	SER	3.2
1	A	10	GLY	3.0
1	A	611	SER	2.9
1	C	8	GLY	2.9
1	A	186	GLU	2.8
1	A	190	PHE	2.8
1	A	196	ARG	2.8
1	A	187	PRO	2.8
1	D	298	VAL	2.7

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Mol	Chain	Res	Type	RSRZ
1	D	299	PRO	2.7
1	C	314	ALA	2.7
1	A	188	ALA	2.6
1	D	178	PRO	2.6
1	D	314	ALA	2.6
1	D	297	GLY	2.5
1	A	185	GLN	2.5
1	A	776	LEU	2.4
1	A	195	THR	2.4
1	C	297	GLY	2.4
1	D	286	VAL	2.2
1	C	299	PRO	2.2
1	D	347	GLN	2.2
1	D	301	LYS	2.2
1	B	184	ALA	2.1
1	B	11	SER	2.0
1	C	306	THR	2.0
1	A	189	ALA	2.0
1	D	350	ALA	2.0
1	C	289	GLU	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
2	GOL	A	801	6/6	0.96	0.08	-	26,29,32,34	0

6.5 Other polymers

There are no such residues in this entry.