



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 09:21 AM UTC

PDB ID : 5LT9 / pdb_00005lt9
Title : Ligand binding domain of Pseudomonas aeruginosa PAO1 amino acid chemoreceptors PctB in complex with L-Arg
Authors : Gavira, J.A.; Rico-Jimenez, M.; Ortega, A.; Conejero-Muriel, M.; Zhulin, I.; Krell, T.
Deposited on : 2016-09-06
Resolution : 3.00 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

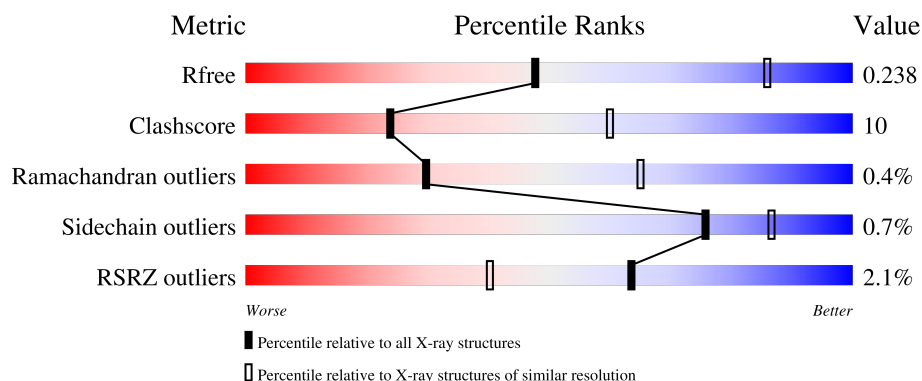
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	290	 67% 14% 19%
1	B	290	 68% 17% 13%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3846 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 Methyl-accepting chemotaxis protein PctB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	234	Total	C	N	O	S	0	3	0
			1830	1158	306	360	6			
1	B	253	Total	C	N	O	S	0	3	0
			1968	1241	332	390	5			

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	9	MET	-	initiating methionine	UNP A0A0F6UK01
A	10	GLY	-	expression tag	UNP A0A0F6UK01
A	11	SER	-	expression tag	UNP A0A0F6UK01
A	12	SER	-	expression tag	UNP A0A0F6UK01
A	13	HIS	-	expression tag	UNP A0A0F6UK01
A	14	HIS	-	expression tag	UNP A0A0F6UK01
A	15	HIS	-	expression tag	UNP A0A0F6UK01
A	16	HIS	-	expression tag	UNP A0A0F6UK01
A	17	HIS	-	expression tag	UNP A0A0F6UK01
A	18	HIS	-	expression tag	UNP A0A0F6UK01
A	19	SER	-	expression tag	UNP A0A0F6UK01
A	20	SER	-	expression tag	UNP A0A0F6UK01
A	21	GLY	-	expression tag	UNP A0A0F6UK01
A	22	LEU	-	expression tag	UNP A0A0F6UK01
A	23	VAL	-	expression tag	UNP A0A0F6UK01
A	24	PRO	-	expression tag	UNP A0A0F6UK01
A	25	ARG	-	expression tag	UNP A0A0F6UK01
A	26	GLY	-	expression tag	UNP A0A0F6UK01
A	27	SER	-	expression tag	UNP A0A0F6UK01
A	28	HIS	-	expression tag	UNP A0A0F6UK01
A	29	MET	-	expression tag	UNP A0A0F6UK01
A	278	ASP	-	expression tag	UNP A0A0F6UK01
A	279	PRO	-	expression tag	UNP A0A0F6UK01
A	280	ASN	-	expression tag	UNP A0A0F6UK01
A	281	SER	-	expression tag	UNP A0A0F6UK01

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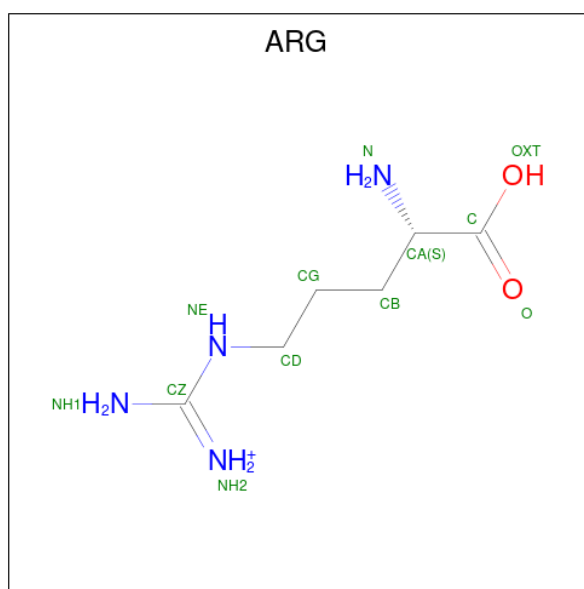
Chain	Residue	Modelled	Actual	Comment	Reference
A	282	SER	-	expression tag	UNP A0A0F6UK01
A	283	SER	-	expression tag	UNP A0A0F6UK01
A	284	VAL	-	expression tag	UNP A0A0F6UK01
A	285	ASP	-	expression tag	UNP A0A0F6UK01
A	286	LYS	-	expression tag	UNP A0A0F6UK01
A	287	LEU	-	expression tag	UNP A0A0F6UK01
A	288	ALA	-	expression tag	UNP A0A0F6UK01
A	289	ALA	-	expression tag	UNP A0A0F6UK01
A	290	ALA	-	expression tag	UNP A0A0F6UK01
A	291	LEU	-	expression tag	UNP A0A0F6UK01
A	292	GLU	-	expression tag	UNP A0A0F6UK01
A	293	HIS	-	expression tag	UNP A0A0F6UK01
A	294	HIS	-	expression tag	UNP A0A0F6UK01
A	295	HIS	-	expression tag	UNP A0A0F6UK01
A	296	HIS	-	expression tag	UNP A0A0F6UK01
A	297	HIS	-	expression tag	UNP A0A0F6UK01
A	298	HIS	-	expression tag	UNP A0A0F6UK01
B	9	MET	-	initiating methionine	UNP A0A0F6UK01
B	10	GLY	-	expression tag	UNP A0A0F6UK01
B	11	SER	-	expression tag	UNP A0A0F6UK01
B	12	SER	-	expression tag	UNP A0A0F6UK01
B	13	HIS	-	expression tag	UNP A0A0F6UK01
B	14	HIS	-	expression tag	UNP A0A0F6UK01
B	15	HIS	-	expression tag	UNP A0A0F6UK01
B	16	HIS	-	expression tag	UNP A0A0F6UK01
B	17	HIS	-	expression tag	UNP A0A0F6UK01
B	18	HIS	-	expression tag	UNP A0A0F6UK01
B	19	SER	-	expression tag	UNP A0A0F6UK01
B	20	SER	-	expression tag	UNP A0A0F6UK01
B	21	GLY	-	expression tag	UNP A0A0F6UK01
B	22	LEU	-	expression tag	UNP A0A0F6UK01
B	23	VAL	-	expression tag	UNP A0A0F6UK01
B	24	PRO	-	expression tag	UNP A0A0F6UK01
B	25	ARG	-	expression tag	UNP A0A0F6UK01
B	26	GLY	-	expression tag	UNP A0A0F6UK01
B	27	SER	-	expression tag	UNP A0A0F6UK01
B	28	HIS	-	expression tag	UNP A0A0F6UK01
B	29	MET	-	expression tag	UNP A0A0F6UK01
B	278	ASP	-	expression tag	UNP A0A0F6UK01
B	279	PRO	-	expression tag	UNP A0A0F6UK01
B	280	ASN	-	expression tag	UNP A0A0F6UK01
B	281	SER	-	expression tag	UNP A0A0F6UK01

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Chain	Residue	Modelled	Actual	Comment	Reference
B	282	SER	-	expression tag	UNP A0A0F6UK01
B	283	SER	-	expression tag	UNP A0A0F6UK01
B	284	VAL	-	expression tag	UNP A0A0F6UK01
B	285	ASP	-	expression tag	UNP A0A0F6UK01
B	286	LYS	-	expression tag	UNP A0A0F6UK01
B	287	LEU	-	expression tag	UNP A0A0F6UK01
B	288	ALA	-	expression tag	UNP A0A0F6UK01
B	289	ALA	-	expression tag	UNP A0A0F6UK01
B	290	ALA	-	expression tag	UNP A0A0F6UK01
B	291	LEU	-	expression tag	UNP A0A0F6UK01
B	292	GLU	-	expression tag	UNP A0A0F6UK01
B	293	HIS	-	expression tag	UNP A0A0F6UK01
B	294	HIS	-	expression tag	UNP A0A0F6UK01
B	295	HIS	-	expression tag	UNP A0A0F6UK01
B	296	HIS	-	expression tag	UNP A0A0F6UK01
B	297	HIS	-	expression tag	UNP A0A0F6UK01
B	298	HIS	-	expression tag	UNP A0A0F6UK01

- Molecule 2 is ARGinine (CCD ID: ARG) (formula: $C_6H_{15}N_4O_2$).

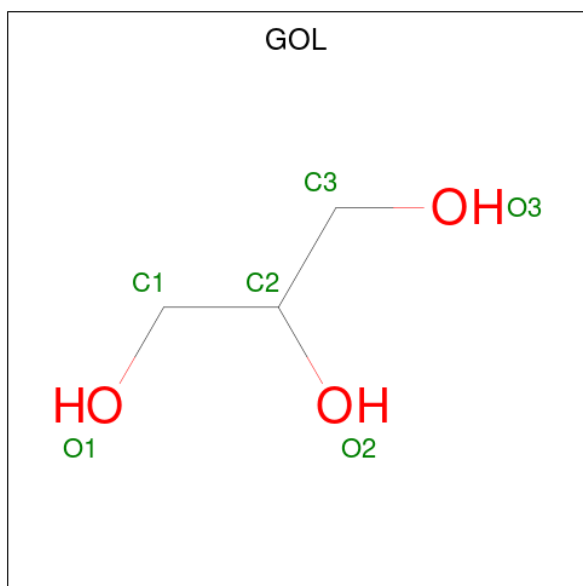


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			12	6	4	2		
2	B	1	Total	C	N	O	0	1
			24	12	8	4		

- Molecule 3 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Na	0	0
			1	1		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	O	0	0
			1	1		
5	B	4	Total	O	0	0
			4	4		

4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	111.73Å 111.73Å 117.55Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	37.35 – 3.00 37.35 – 3.00	Depositor EDS
% Data completeness (in resolution range)	99.9 (37.35-3.00) 94.5 (37.35-3.00)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.07 (at 3.01Å)	Xtriage
Refinement program	PHENIX 1.10_2155	Depositor
R, R_{free}	0.215 , 0.248 0.212 , 0.238	Depositor DCC
R_{free} test set	881 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	79.4	Xtriage
Anisotropy	0.001	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 78.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.038 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3846	wwPDB-VP
Average B, all atoms (Å ²)	86.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.96% 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, NA

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	0.08	0/1871	0.26	0/2546
1	B	0.49	2/2011 (0.1%)	0.57	4/2737 (0.1%)
All	All	0.36	2/3882 (0.1%)	0.45	4/5283 (0.1%)

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	B	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	267	ALA	C-N	17.16	1.58	1.33
1	B	268	TYR	C-N	-8.43	1.22	1.33

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	267	ALA	N-CA-C	-14.64	92.04	110.19
1	B	266	LYS	N-CA-C	-7.29	103.37	111.82
1	B	227	ILE	N-CA-C	-5.41	102.61	109.58
1	B	267	ALA	O-C-N	5.33	129.50	122.36

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	275[A]	ARG	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	1830	0	1789	26	0
1	B	1968	0	1927	51	0
2	A	12	0	12	2	0
2	B	24	0	24	4	0
3	A	1	0	0	0	0
4	B	6	0	8	0	0
5	A	1	0	0	0	0
5	B	4	0	0	0	0
All	All	3846	0	3760	77	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 10.

All (77) 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:264:LYS:O	1:B:267:ALA:HB3	1.76	0.84
1:B:146:GLU:HG2	1:B:149:ILE:HG12	1.61	0.82
1:B:49:MET:HG3	1:B:186:LEU:HD23	1.67	0.77
1:B:115:SER:OG	2:B:301[A]:ARG:NH1	2.19	0.75
1:B:228:GLY:HA2	1:B:229:SER:HB2	1.73	0.70
1:B:228:GLY:CA	1:B:229:SER:HB2	2.24	0.68
1:B:268:TYR:CD2	1:B:271:LEU:HD11	2.29	0.67
1:B:187:ASP:HB3	1:B:188:PHE:C	2.20	0.67
1:B:228:GLY:HA3	1:B:229:SER:C	2.19	0.67
1:B:274:LEU:HD23	1:B:274:LEU:O	1.96	0.66
1:B:268:TYR:O	1:B:271:LEU:HD12	1.96	0.66
1:B:75:LEU:HD13	1:B:102:LEU:HD22	1.79	0.65
1:B:277:SER:O	1:B:279:PRO:HD3	1.96	0.64
1:A:115:SER:OG	2:A:301:ARG:NH1	2.31	0.64
1:A:93:LEU:HD13	1:A:100:THR:HG21	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:65:ILE:HG23	1:A:138:THR:HG21	1.80	0.63
1:A:175:SER:O	1:A:178:THR:OG1	2.16	0.63
1:B:229:SER:N	1:B:230:GLY:HA2	2.16	0.61
1:B:266:LYS:C	1:B:267:ALA:O	2.33	0.61
1:B:263:ASP:O	1:B:267:ALA:HB2	2.00	0.60
1:A:75:LEU:HD13	1:A:102:LEU:HD22	1.82	0.60
1:A:85:GLN:HG3	1:A:112:ARG:HH12	1.68	0.59
1:B:147:PRO:HD2	2:B:301[B]:ARG:N	2.18	0.57
1:B:265:ASP:O	1:B:267:ALA:O	2.22	0.57
1:B:188:PHE:N	1:B:189:GLY:HA2	2.20	0.56
1:B:111:ALA:O	1:B:112:ARG:NH1	2.37	0.56
1:B:233:GLU:HB3	1:B:242:ILE:HG22	1.88	0.56
1:A:41:ASP:N	1:A:41:ASP:OD1	2.36	0.56
1:B:161:ARG:HA	1:B:166:PRO:HA	1.87	0.55
1:B:263:ASP:O	1:B:267:ALA:N	2.40	0.54
1:A:196:LEU:HB3	1:A:205:VAL:HG22	1.89	0.54
1:A:98:LEU:HD11	1:A:175:SER:HB3	1.90	0.53
1:B:77:ARG:NH2	1:B:78[A]:ASP:OD2	2.41	0.53
1:A:179:VAL:O	1:A:183:ILE:HG12	2.07	0.53
1:B:93:LEU:HD13	1:B:100:THR:HG21	1.91	0.53
1:B:146:GLU:HG3	1:B:148:ALA:H	1.74	0.53
1:A:225:PRO:HG2	1:A:243:LEU:HD22	1.91	0.53
1:B:40:GLU:N	1:B:40:GLU:OE2	2.42	0.52
1:B:40:GLU:N	1:B:270:MET:HE2	2.23	0.52
1:A:85:GLN:O	1:A:89:GLU:HG2	2.10	0.52
1:B:136:GLY:O	1:B:161:ARG:NH1	2.42	0.52
1:A:149:ILE:HG13	1:A:151[B]:GLU:HG2	1.94	0.49
1:B:228:GLY:HA3	1:B:229:SER:O	2.12	0.48
1:A:67:LEU:HD13	1:A:93:LEU:HD23	1.95	0.48
1:B:40:GLU:N	1:B:270:MET:CE	2.76	0.48
1:B:76:ALA:HB3	1:B:79:HIS:HD2	1.79	0.47
1:B:109:TYR:OH	2:B:301[A]:ARG:HD2	2.14	0.47
1:B:268:TYR:CG	1:B:271:LEU:HD11	2.50	0.47
1:B:263:ASP:HB3	1:B:266:LYS:HB3	1.97	0.47
1:B:263:ASP:O	1:B:267:ALA:CB	2.63	0.46
1:A:140:LEU:HD22	1:A:156:ILE:HG12	1.98	0.46
1:A:212:VAL:HG12	1:A:213:MET:HG2	1.98	0.45
1:B:149:ILE:HG13	1:B:151:GLU:HG2	1.97	0.45
1:A:99:PHE:CG	2:A:301:ARG:HD2	2.52	0.44
1:B:191:MET:HE3	1:B:191:MET:HB2	1.84	0.44
1:A:109:TYR:HB2	1:A:123:PRO:HG3	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:76:ALA:HB3	1:B:79:HIS:CD2	2.52	0.44
1:A:201:GLY:O	1:A:216:LEU:N	2.52	0.43
1:B:188:PHE:N	1:B:189:GLY:CA	2.82	0.42
1:B:101:TYR:CE2	2:B:301[B]:ARG:HD2	2.54	0.42
1:B:222:ARG:HD2	1:B:223:ASN:H	1.84	0.42
1:B:271:LEU:O	1:B:275[C]:ARG:HG2	2.20	0.42
1:A:43:GLU:HB3	1:A:231:PHE:CE2	2.55	0.42
1:B:268:TYR:CE2	1:B:271:LEU:HD11	2.56	0.41
1:B:210:ASP:OD1	1:B:210:ASP:N	2.53	0.41
1:A:85:GLN:HG3	1:A:112:ARG:NH1	2.33	0.41
1:B:101:TYR:HB3	1:B:111:ALA:CB	2.51	0.41
1:A:87:LEU:O	1:A:90:GLN:NE2	2.45	0.41
1:B:40:GLU:C	1:B:42:LEU:H	2.29	0.40
1:B:78[A]:ASP:N	1:B:78[A]:ASP:OD1	2.53	0.40
1:A:112:ARG:HA	1:A:112:ARG:HD3	1.92	0.40
1:A:57:VAL:HG21	1:A:183:ILE:HD13	2.03	0.40
1:B:122:ASP:HB3	1:B:125:ARG:HG3	2.03	0.40
1:B:251:LEU:HB3	1:B:254:LEU:HB2	2.02	0.40
1:A:75:LEU:HA	1:A:78:ASP:O	2.22	0.40
1:A:229:SER:HA	1:A:230:GLY:HA2	1.68	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	235/290 (81%)	232 (99%)	3 (1%)	0	100	100
1	B	254/290 (88%)	244 (96%)	8 (3%)	2 (1%)	16	50
All	All	489/580 (84%)	476 (97%)	11 (2%)	2 (0%)	30	65

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	279	PRO
1	B	230	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	199/245 (81%)	199 (100%)	0	100	100
1	B	215/245 (88%)	211 (98%)	4 (2%)	50	76
All	All	414/490 (84%)	410 (99%)	4 (1%)	76	84

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	47[A]	HIS
1	B	47[B]	HIS
1	B	270	MET
1	B	271	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	150	HIS
1	B	58	GLN
1	B	239	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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 1 is monoatomic - leaving 4 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
2	ARG	B	301[B]	-	10,11,11	0.76	1 (10%)	9,13,13	1.02	1 (11%)
4	GOL	B	302	-	5,5,5	0.37	0	5,5,5	0.28	0
2	ARG	B	301[A]	-	10,11,11	0.75	1 (10%)	9,13,13	0.89	1 (11%)
2	ARG	A	301	-	10,11,11	0.76	1 (10%)	9,13,13	0.91	1 (11%)

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	ARG	B	301[B]	-	-	4/11/11/11	-
4	GOL	B	302	-	-	2/4/4/4	-
2	ARG	B	301[A]	-	-	4/11/11/11	-
2	ARG	A	301	-	-	2/11/11/11	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	301	ARG	OXT-C	-2.24	1.23	1.30
2	B	301[A]	ARG	OXT-C	-2.22	1.23	1.30
2	B	301[B]	ARG	OXT-C	-2.21	1.23	1.30

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	301[B]	ARG	OXT-C-O	-3.01	117.24	124.08
2	A	301	ARG	OXT-C-O	-2.71	117.93	124.08
2	B	301[A]	ARG	OXT-C-O	-2.48	118.46	124.08

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	301[A]	ARG	N-CA-CB-CG
4	B	302	GOL	O1-C1-C2-C3
2	B	301[A]	ARG	NE-CD-CG-CB
4	B	302	GOL	O1-C1-C2-O2
2	B	301[B]	ARG	O-C-CA-N
2	B	301[A]	ARG	OXT-C-CA-CB
2	B	301[A]	ARG	O-C-CA-CB
2	B	301[B]	ARG	OXT-C-CA-N
2	B	301[B]	ARG	O-C-CA-CB
2	A	301	ARG	O-C-CA-CB
2	A	301	ARG	N-CA-CB-CG
2	B	301[B]	ARG	OXT-C-CA-CB

There are no ring outliers.

3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	301[B]	ARG	2	0
2	B	301[A]	ARG	2	0
2	A	301	ARG	2	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 [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	234/290 (80%)	-0.07	4 (1%)	69 45	42, 77, 124, 140	4 (1%)
1	B	253/290 (87%)	-0.08	6 (2%)	59 36	49, 81, 140, 181	3 (1%)
All	All	487/580 (83%)	-0.08	10 (2%)	63 40	42, 80, 131, 181	7 (1%)

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	229	SER	4.0
1	B	47[A]	HIS	3.7
1	A	191	MET	3.0
1	B	270	MET	2.9
1	B	272	THR	2.7
1	B	40	GLU	2.6
1	B	269	ALA	2.5
1	A	270[A]	MET	2.3
1	A	224	THR	2.2
1	B	278	ASP	2.2

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
4	GOL	B	302	6/6	0.75	0.22	87,98,104,106	0
3	NA	A	302	1/1	0.84	0.24	78,78,78,78	0
2	ARG	B	301[B]	12/12	0.93	0.16	67,76,80,81	12
2	ARG	A	301	12/12	0.93	0.13	64,78,111,115	0
2	ARG	B	301[A]	12/12	0.93	0.16	67,75,80,81	12

6.5 Other polymers [i](#)

There are no such residues in this entry.