



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 08:15 PM GMT

PDB ID : 4Y01
Title : Crystal structure of dipeptidyl peptidase 11 (DPP11) from Porphyromonas gingivalis
Authors : Sakamoto, Y.; Suzuki, Y.; Iizuka, I.; Tateoka, C.; Roppongi, S.; Fujimoto, M.; Nonaka, T.; Ogasawara, W.; Tanaka, N.
Deposited on : 2015-02-05
Resolution : 2.46 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-20026982
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-20026982

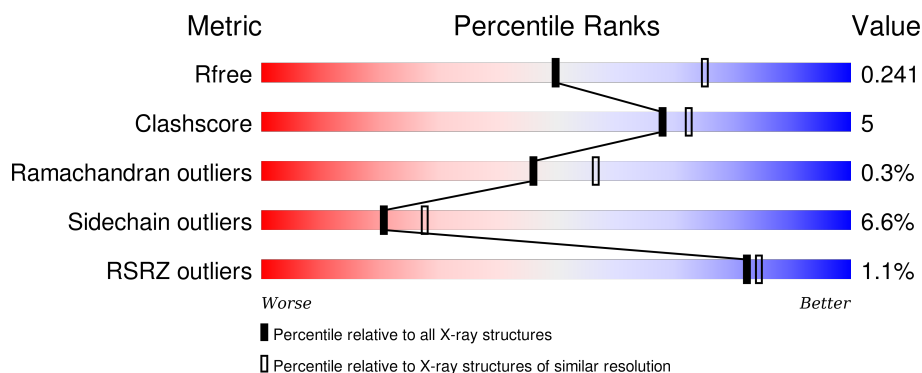
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



The reported resolution of this entry is 2.46 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	4776 (2.50-2.42)
Clashscore	102246	1030 (2.48-2.44)
Ramachandran outliers	100387	1024 (2.48-2.44)
Sidechain outliers	100360	1024 (2.48-2.44)
RSRZ outliers	91569	4787 (2.50-2.42)

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	720	 2% 82% 13% . .
1	B	720	 85% 10% . . .

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	GOL	A	801	-	-	-	X

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 11408 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 Peptidase S46.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	699	Total	C	N	O	S	Se	0	0	0
			5608	3553	970	1058	4	23			
1	B	699	Total	C	N	O	S	Se	0	0	0
			5608	3553	970	1058	4	23			

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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	61	Total	O	0	0
			61	61		

Continued on next page...

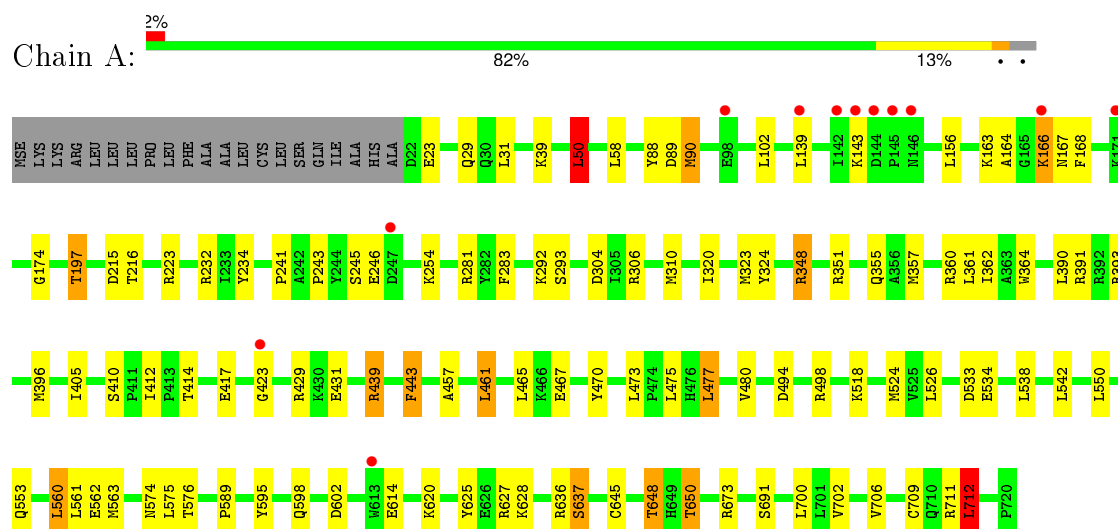
Continued from previous page...

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	125	Total 125	O 125	0	0

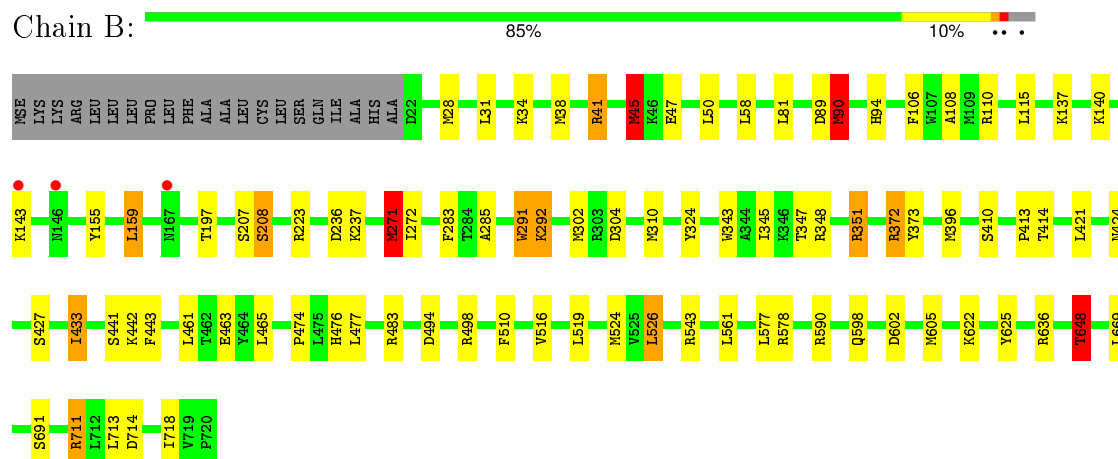
3 Residue-property plots

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: Peptidase S46



• Molecule 1: Peptidase S46



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	99.79Å 115.49Å 147.57Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 2.46 38.91 – 2.46	Depositor EDS
% Data completeness (in resolution range)	99.2 (40.00-2.46) 99.2 (38.91-2.46)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.38 (at 2.45Å)	Xtriage
Refinement program	REFMAC 5.8.0103	Depositor
R, R_{free}	0.184 , 0.233 0.198 , 0.241	Depositor DCC
R_{free} test set	2980 reflections (5.07%)	DCC
Wilson B-factor (Å ²)	36.5	Xtriage
Anisotropy	0.014	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 30.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 61765 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11408	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.77% 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, 0.375 respectively for untwinned datasets, and 0.333, 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	0.72	1/5710 (0.0%)	0.93	11/7676 (0.1%)
1	B	0.75	1/5710 (0.0%)	0.93	13/7676 (0.2%)
All	All	0.73	2/11420 (0.0%)	0.93	24/15352 (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	2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	208	SER	CB-OG	-5.24	1.35	1.42
1	A	293	SER	CB-OG	-5.01	1.35	1.42

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	223	ARG	NE-CZ-NH2	-11.03	114.78	120.30
1	B	223	ARG	NE-CZ-NH1	9.79	125.19	120.30
1	A	223	ARG	NE-CZ-NH2	-8.96	115.82	120.30
1	B	271	MSE	CG-SE-CE	-8.45	80.31	98.90
1	B	41	ARG	NE-CZ-NH1	7.54	124.07	120.30
1	A	50	LEU	CA-CB-CG	7.20	131.86	115.30
1	B	372	ARG	NE-CZ-NH1	6.76	123.68	120.30
1	A	223	ARG	NE-CZ-NH1	6.53	123.56	120.30
1	A	711	ARG	NE-CZ-NH1	6.42	123.51	120.30

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	45	MSE	CG-SE-CE	-6.25	85.15	98.90
1	B	711	ARG	NE-CZ-NH1	6.22	123.41	120.30
1	B	291	TRP	CA-CB-CG	5.87	124.85	113.70
1	A	712	LEU	CA-CB-CG	5.81	128.67	115.30
1	A	306	ARG	NE-CZ-NH1	5.75	123.17	120.30
1	B	351	ARG	NE-CZ-NH1	5.74	123.17	120.30
1	A	391	ARG	NE-CZ-NH1	5.72	123.16	120.30
1	A	223	ARG	CG-CD-NE	-5.51	100.23	111.80
1	B	711	ARG	NE-CZ-NH2	-5.37	117.62	120.30
1	A	391	ARG	NE-CZ-NH2	-5.36	117.62	120.30
1	A	627	ARG	NE-CZ-NH2	-5.33	117.63	120.30
1	A	348	ARG	NE-CZ-NH1	5.32	122.96	120.30
1	B	90	MSE	CG-SE-CE	5.27	110.49	98.90
1	B	236	ASP	CB-CG-OD1	5.14	122.92	118.30
1	B	590	ARG	NE-CZ-NH2	-5.00	117.80	120.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	23	GLU	Peptide
1	A	423	GLY	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	5608	0	5491	50	0
1	B	5608	0	5491	55	0
2	A	6	0	8	0	0
3	A	61	0	0	1	0
3	B	125	0	0	1	0
All	All	11408	0	10990	105	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 5.

All (105) 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:410:SER:CB	1:B:524:MSE:HE3	1.88	1.04
1:A:410:SER:CB	1:A:524:MSE:HE3	1.87	1.04
1:B:302:MSE:HE3	1:B:396:MSE:SE	2.24	0.87
1:B:474:PRO:HG2	1:B:477:LEU:HD12	1.59	0.84
1:B:410:SER:HB2	1:B:524:MSE:CE	2.07	0.84
1:B:310:MSE:HE3	1:B:324:TYR:HD1	1.45	0.82
1:B:31:LEU:HD12	1:B:38:MSE:HE1	1.65	0.79
1:B:410:SER:HB2	1:B:524:MSE:HE3	1.65	0.78
1:B:410:SER:CB	1:B:524:MSE:CE	2.62	0.75
1:A:410:SER:HB2	1:A:524:MSE:HE3	1.68	0.75
1:A:310:MSE:HE3	1:A:324:TYR:HD2	1.50	0.75
1:B:410:SER:HB3	1:B:524:MSE:HE3	1.73	0.65
1:A:410:SER:HB3	1:A:524:MSE:HE3	1.76	0.64
1:B:34:LYS:HB2	1:B:38:MSE:CE	2.28	0.64
1:B:108:ALA:O	1:B:711:ARG:NH1	2.33	0.62
1:A:410:SER:HB2	1:A:524:MSE:CE	2.30	0.61
1:B:302:MSE:HE1	1:B:463:GLU:HG3	1.84	0.58
1:A:443:PHE:CE2	1:A:524:MSE:HE1	2.39	0.58
1:A:174:GLY:HA3	1:A:241:PRO:HG2	1.86	0.57
1:A:89:ASP:HB2	1:A:90:MSE:HE2	1.86	0.56
1:A:414:THR:HG22	1:A:417:GLU:OE1	2.05	0.56
1:B:94:HIS:CG	1:B:115:LEU:HD23	2.40	0.56
1:B:343:TRP:O	1:B:347:THR:HB	2.05	0.56
1:A:310:MSE:HE3	1:A:324:TYR:CD2	2.37	0.56
1:B:421:LEU:HD21	1:B:433:ILE:HD13	1.86	0.56
1:A:412:ILE:O	1:A:439:ARG:NH2	2.39	0.56
1:B:197:THR:HB	3:B:828:HOH:O	2.06	0.55
1:B:302:MSE:HE1	1:B:463:GLU:CG	2.37	0.54
1:B:413:PRO:HD3	1:B:519:LEU:HD21	1.89	0.54
1:B:272:ILE:HD12	1:B:648:THR:HG21	1.89	0.54
1:A:320:ILE:HD13	1:A:323:MSE:CE	2.37	0.54
1:B:413:PRO:HG3	1:B:519:LEU:CD2	2.37	0.54
1:B:292:LYS:HB2	1:B:345:ILE:HD13	1.90	0.54
1:A:429:ARG:NH1	3:A:901:HOH:O	2.36	0.54
1:B:34:LYS:HB2	1:B:38:MSE:HE2	1.89	0.54
1:A:320:ILE:HD13	1:A:323:MSE:HE2	1.89	0.54
1:A:410:SER:OG	1:A:524:MSE:HE3	2.08	0.53
1:B:89:ASP:HB2	1:B:90:MSE:HE3	1.90	0.53
1:B:34:LYS:C	1:B:38:MSE:HE3	2.29	0.53
1:B:577:LEU:O	1:B:578:ARG:HD3	2.09	0.53
1:B:413:PRO:HG3	1:B:519:LEU:HD21	1.91	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:589:PRO:HG3	1:A:595:TYR:CE2	2.45	0.52
1:B:648:THR:HG22	1:B:691:SER:OG	2.08	0.52
1:A:351:ARG:HH12	1:A:355:GLN:HE21	1.56	0.52
1:B:347:THR:CG2	1:B:348:ARG:NH1	2.75	0.50
1:A:281:ARG:HH22	1:A:357:MSE:HE1	1.76	0.50
1:A:31:LEU:HD21	1:A:50:LEU:HD13	1.92	0.50
1:B:155:TYR:CE2	1:B:159:LEU:HD21	2.47	0.50
1:B:285:ALA:HB1	1:B:351:ARG:NH1	2.26	0.49
1:B:28:MSE:HE3	1:B:577:LEU:HG	1.94	0.49
1:B:50:LEU:HD21	1:B:271:MSE:HE1	1.94	0.49
1:A:197:THR:HG22	1:A:243:PRO:HA	1.94	0.49
1:A:405:ILE:HD11	1:A:461:LEU:HD13	1.95	0.48
1:A:457:ALA:O	1:A:461:LEU:HB2	2.13	0.48
1:A:393:ARG:NH2	1:A:534:GLU:OE2	2.46	0.48
1:B:413:PRO:CG	1:B:519:LEU:HD21	2.43	0.47
1:B:89:ASP:HB2	1:B:90:MSE:CE	2.44	0.47
1:A:396:MSE:HE3	1:A:467:GLU:HB3	1.97	0.47
1:A:709:CYS:O	1:A:712:LEU:CD2	2.63	0.47
1:A:88:TYR:CE1	1:A:102:LEU:HD11	2.51	0.46
1:B:413:PRO:CD	1:B:519:LEU:HD21	2.46	0.46
1:B:106:PHE:CZ	1:B:115:LEU:HD13	2.50	0.46
1:A:215:ASP:N	1:A:614:GLU:OE2	2.48	0.46
1:A:414:THR:HG22	1:A:417:GLU:CD	2.36	0.46
1:A:360:ARG:HG2	1:A:563:MSE:HE1	1.98	0.45
1:B:81:LEU:HD13	1:B:669:LEU:HD11	1.99	0.45
1:B:155:TYR:CE2	1:B:159:LEU:CD2	3.00	0.45
1:A:650:THR:HG22	1:A:673:ARG:HG3	1.99	0.45
1:B:648:THR:HG22	1:B:691:SER:CB	2.47	0.44
1:A:164:ALA:HB1	1:A:168:PHE:CD2	2.53	0.44
1:A:232:ARG:HD2	1:A:234:TYR:CE1	2.53	0.44
1:B:34:LYS:CB	1:B:38:MSE:HE2	2.48	0.44
1:B:34:LYS:CB	1:B:38:MSE:CE	2.95	0.44
1:B:648:THR:CG2	1:B:691:SER:OG	2.66	0.43
1:B:372:ARG:NH2	1:B:373:TYR:OH	2.52	0.43
1:A:216:THR:N	1:A:614:GLU:OE2	2.45	0.43
1:A:166:LYS:HA	1:A:167:ASN:HA	1.79	0.43
1:B:410:SER:OG	1:B:524:MSE:HE3	2.17	0.43
1:A:538:LEU:O	1:A:542:LEU:HB2	2.18	0.43
1:A:494:ASP:OD1	1:A:498:ARG:HD3	2.19	0.43
1:A:31:LEU:HD21	1:A:50:LEU:CD1	2.49	0.43
1:A:361:LEU:HD22	1:A:560:LEU:HD13	2.00	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:405:ILE:CD1	1:A:461:LEU:HD13	2.49	0.42
1:A:589:PRO:HG3	1:A:595:TYR:CZ	2.55	0.42
1:B:476:HIS:CD2	1:B:526:LEU:HB3	2.54	0.42
1:B:718:ILE:N	1:B:718:ILE:HD12	2.34	0.42
1:B:272:ILE:CD1	1:B:648:THR:HG21	2.49	0.42
1:B:443:PHE:CE2	1:B:524:MSE:HE1	2.54	0.42
1:B:605:MSE:HG3	1:B:625:TYR:CE2	2.55	0.42
1:A:470:TYR:HA	1:A:473:LEU:HD12	2.01	0.42
1:B:433:ILE:HD12	1:B:433:ILE:HA	1.88	0.42
1:A:477:LEU:O	1:A:480:VAL:HG12	2.20	0.42
1:B:292:LYS:CB	1:B:345:ILE:HD13	2.50	0.42
1:A:702:VAL:O	1:A:706:VAL:HB	2.20	0.41
1:A:648:THR:HG22	1:A:691:SER:HB2	2.02	0.41
1:A:712:LEU:H	1:A:712:LEU:HD22	1.84	0.41
1:A:625:TYR:O	1:A:628:LYS:HD3	2.20	0.41
1:B:510:PHE:CZ	1:B:519:LEU:HD12	2.56	0.41
1:B:272:ILE:HD12	1:B:648:THR:CG2	2.51	0.41
1:A:636:ARG:O	1:A:637:SER:CB	2.69	0.41
1:A:574:ASN:OD1	1:A:576:THR:HB	2.21	0.40
1:A:645:CYS:HA	1:A:691:SER:O	2.21	0.40
1:A:364:TRP:HB2	1:A:563:MSE:HE3	2.04	0.40
1:B:494:ASP:OD1	1:B:498:ARG:HD3	2.22	0.40
1:B:45:MSE:HE1	1:B:58:LEU:HD22	2.03	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	697/720 (97%)	671 (96%)	25 (4%)	1 (0%)	56 71
1	B	697/720 (97%)	671 (96%)	23 (3%)	3 (0%)	39 49

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1394/1440 (97%)	1342 (96%)	48 (3%)	4 (0%)	46 57

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	516	VAL
1	B	648	THR
1	B	143	LYS
1	A	637	SER

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	590/583 (101%)	547 (93%)	43 (7%)	17 23
1	B	590/583 (101%)	555 (94%)	35 (6%)	24 34
All	All	1180/1166 (101%)	1102 (93%)	78 (7%)	21 28

All (78) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	GLN
1	A	39	LYS
1	A	50	LEU
1	A	58	LEU
1	A	90	MSE
1	A	139	LEU
1	A	143	LYS
1	A	156	LEU
1	A	163	LYS
1	A	166	LYS
1	A	197	THR
1	A	245	SER
1	A	246	GLU
1	A	254	LYS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	283	PHE
1	A	292	LYS
1	A	304	ASP
1	A	348	ARG
1	A	362	ILE
1	A	390	LEU
1	A	431	GLU
1	A	439	ARG
1	A	443	PHE
1	A	461	LEU
1	A	465	LEU
1	A	475	LEU
1	A	477	LEU
1	A	518	LYS
1	A	526	LEU
1	A	533	ASP
1	A	550	LEU
1	A	553	GLN
1	A	560	LEU
1	A	561	LEU
1	A	562	GLU
1	A	575	LEU
1	A	598	GLN
1	A	602	ASP
1	A	620	LYS
1	A	648	THR
1	A	650	THR
1	A	700	LEU
1	A	712	LEU
1	B	41	ARG
1	B	45	MSE
1	B	47	GLU
1	B	90	MSE
1	B	110	ARG
1	B	137	LYS
1	B	140	LYS
1	B	159	LEU
1	B	207	SER
1	B	208	SER
1	B	237	LYS
1	B	271	MSE
1	B	283	PHE

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	291	TRP
1	B	292	LYS
1	B	304	ASP
1	B	414	THR
1	B	424	ASN
1	B	427	SER
1	B	433	ILE
1	B	441	SER
1	B	442	LYS
1	B	461	LEU
1	B	465	LEU
1	B	483	ARG
1	B	526	LEU
1	B	543	ARG
1	B	561	LEU
1	B	598	GLN
1	B	602	ASP
1	B	622	LYS
1	B	636	ARG
1	B	648	THR
1	B	713	LEU
1	B	714	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	83	ASN
1	A	117	ASN
1	A	267	ASN
1	A	307	GLN
1	A	355	GLN
1	A	569	GLN
1	A	598	GLN
1	B	83	ASN
1	B	117	ASN
1	B	186	ASN
1	B	307	GLN
1	B	489	GLN
1	B	506	GLN
1	B	553	GLN
1	B	569	GLN
1	B	598	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 ⓘ

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	0.66	0	5,5,5	0.71	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 [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	676/720 (93%)	-0.10	12 (1%) 71 74	20, 42, 71, 115	0
1	B	676/720 (93%)	-0.32	3 (0%) 93 93	16, 34, 61, 99	0
All	All	1352/1440 (93%)	-0.21	15 (1%) 82 84	16, 38, 68, 115	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	171	LYS	4.8
1	A	423	GLY	4.8
1	A	142	ILE	4.2
1	A	143	LYS	4.2
1	A	166	LYS	4.1
1	B	146	ASN	3.3
1	A	146	ASN	2.7
1	A	247	ASP	2.7
1	A	139	LEU	2.5
1	A	145	PRO	2.5
1	B	143	LYS	2.5
1	A	613	TRP	2.4
1	B	167	ASN	2.4
1	A	98	GLU	2.3
1	A	144	ASP	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(\AA^2)	Q<0.9
2	GOL	A	801	6/6	0.81	0.24	3.50	51,56,63,67	0

6.5 Other polymers [i](#)

There are no such residues in this entry.