



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 02:34 AM UTC

PDB ID : 3WOJ / pdb_00003woj
Title : Crystal structure of the DAP BII
Authors : Sakamoto, Y.; Suzuki, Y.; Iizuka, I.; Tateoka, C.; Roppongi, S.; Fujimoto, M.;
Nonaka, T.; Ogasawara, W.; Tanaka, N.
Deposited on : 2013-12-29
Resolution : 2.20 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
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

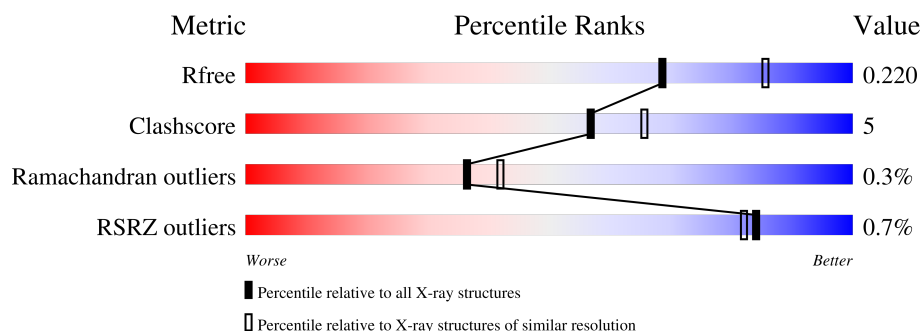
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	698	89% 10%
1	B	698	90% 10%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 11715 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 dipeptidyl aminopeptidase BII.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	698	Total	C	N	O	S	0	0	0
			5381	3404	939	1019	19			
1	B	698	Total	C	N	O	S	0	0	0
			5381	3404	939	1019	19			

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

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

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Zn	0	0
			1	1		
3	B	2	Total	Zn	0	0
			2	2		

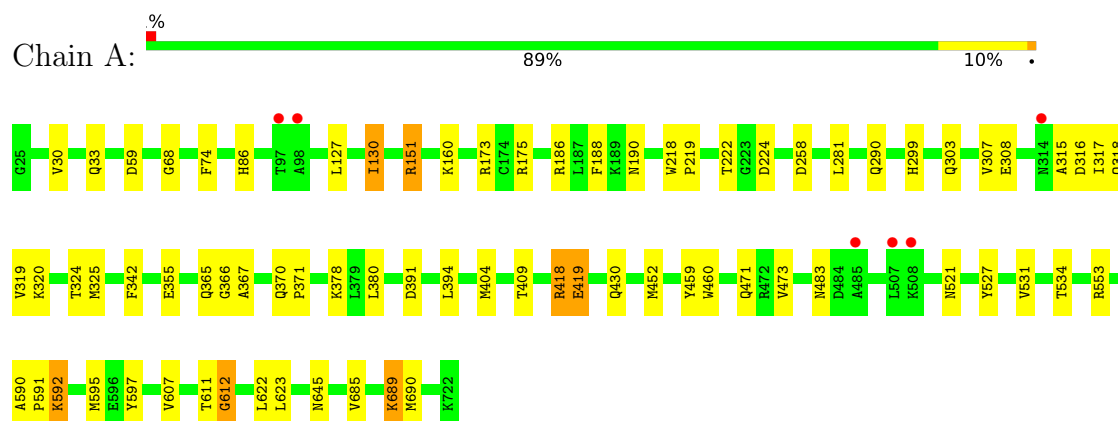
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	462	Total	O	0	0
			462	462		
4	B	446	Total	O	0	0
			446	446		

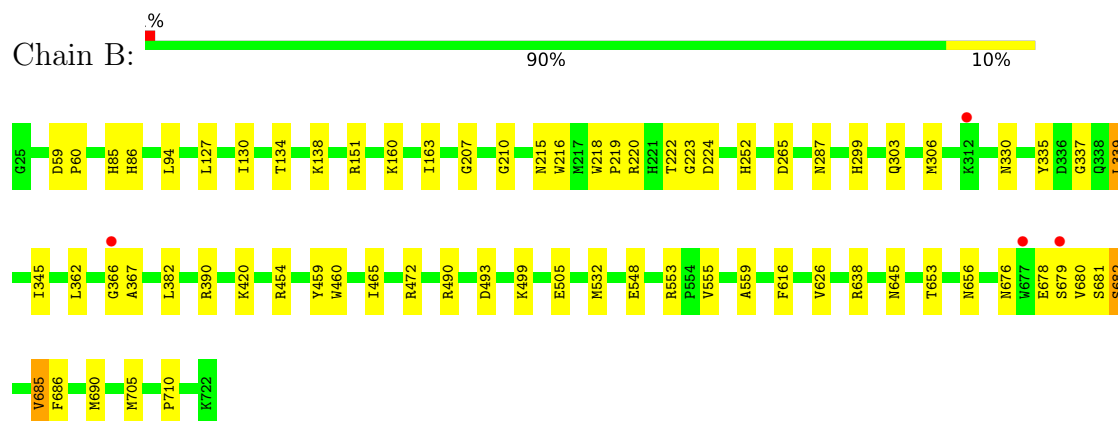
3 Residue-property plots

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: dipeptidyl aminopeptidase BII



• Molecule 1: dipeptidyl aminopeptidase BII



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	76.46Å 130.71Å 170.47Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.59 – 2.20 29.59 – 2.20	Depositor EDS
% Data completeness (in resolution range)	98.5 (29.59-2.20) 98.5 (29.59-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.40 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.165 , 0.220 (Not available) , 0.220	Depositor DCC
R_{free} test set	4318 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	17.7	Xtriage
Anisotropy	0.135	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 33.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11715	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.75% 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, ZN

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	1.22	9/5499 (0.2%)	1.12	13/7451 (0.2%)
1	B	1.22	6/5499 (0.1%)	1.10	12/7451 (0.2%)
All	All	1.22	15/10998 (0.1%)	1.11	25/14902 (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	3
1	B	0	1
All	All	0	4

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	130	ILE	C-O	-5.99	1.18	1.24
1	B	339	LEU	C-O	-5.92	1.17	1.24
1	B	490	ARG	N-CA	5.76	1.53	1.46
1	B	681	SER	C-O	5.69	1.31	1.24
1	A	74	PHE	N-CA	5.52	1.52	1.46
1	A	258	ASP	C-O	-5.48	1.16	1.24
1	B	532	MET	C-O	-5.46	1.19	1.24
1	A	597	TYR	C-O	-5.24	1.17	1.24
1	A	355	GLU	CA-C	5.19	1.59	1.52
1	A	151	ARG	CD-NE	-5.18	1.39	1.46
1	B	626	VAL	C-O	-5.16	1.18	1.24
1	A	342	PHE	N-CA	5.16	1.52	1.46
1	B	682	SER	CA-CB	5.12	1.61	1.53
1	A	59	ASP	CA-C	5.08	1.58	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	689	LYS	C-O	-5.05	1.18	1.24

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	685	VAL	CB-CA-C	-8.69	97.35	110.55
1	A	365	GLN	CA-C-N	-6.55	108.57	121.41
1	A	365	GLN	C-N-CA	-6.55	108.57	121.41
1	B	160	LYS	N-CA-C	6.45	118.31	111.28
1	A	612	GLY	N-CA-C	6.38	128.31	113.18
1	B	555	VAL	N-CA-C	-6.38	104.05	111.00
1	B	330	ASN	N-CA-C	6.28	118.67	111.02
1	A	590	ALA	CA-C-N	-6.25	113.33	119.64
1	A	590	ALA	C-N-CA	-6.25	113.33	119.64
1	B	681	SER	CA-C-N	6.08	128.71	120.38
1	B	681	SER	C-N-CA	6.08	128.71	120.38
1	A	592	LYS	N-CA-C	5.91	118.06	109.07
1	B	163	ILE	N-CA-C	-5.90	104.99	110.53
1	A	324	THR	N-CA-C	-5.75	105.09	111.36
1	B	553	ARG	CA-C-N	-5.49	113.24	119.28
1	B	553	ARG	C-N-CA	-5.49	113.24	119.28
1	A	553	ARG	CA-C-N	-5.46	113.40	119.19
1	A	553	ARG	C-N-CA	-5.46	113.40	119.19
1	A	409	THR	N-CA-C	5.39	116.84	111.07
1	A	418	ARG	CA-C-N	-5.35	116.98	126.45
1	A	418	ARG	C-N-CA	-5.35	116.98	126.45
1	B	337	GLY	N-CA-C	5.06	118.80	112.73
1	A	685	VAL	CB-CA-C	-5.05	104.28	111.40
1	B	680	VAL	N-CA-CB	-5.03	102.04	110.65
1	B	85	HIS	N-CA-C	5.01	117.12	111.11

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	366	GLY	Peptide
1	A	419	GLU	Peptide
1	A	611	THR	Peptide
1	B	366	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	5381	0	5297	56	0
1	B	5381	0	5297	43	0
2	A	24	0	32	1	0
2	B	18	0	23	4	0
3	A	1	0	0	0	0
3	B	2	0	0	0	0
4	A	462	0	0	9	0
4	B	446	0	0	5	0
All	All	11715	0	10649	99	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 (99) 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:591:PRO:HD2	1:A:595:MET:HE1	1.18	1.11
1:A:307:VAL:HG22	1:A:452:MET:HE3	1.42	0.98
1:A:307:VAL:HA	1:A:452:MET:HE1	1.43	0.98
1:A:315:ALA:O	1:A:319:VAL:HG12	1.64	0.96
1:B:224:ASP:OD1	4:B:1323:HOH:O	1.88	0.91
1:B:679:SER:O	1:B:682:SER:HB3	1.73	0.87
1:B:306:MET:HE1	4:B:1161:HOH:O	1.74	0.86
1:A:224:ASP:OD1	4:A:1296:HOH:O	1.94	0.86
1:A:86:HIS:N	4:A:1296:HOH:O	1.88	0.84
1:A:591:PRO:HD2	1:A:595:MET:CE	2.05	0.84
1:A:130:ILE:HD11	1:A:186:ARG:HD3	1.65	0.78
1:A:591:PRO:CD	1:A:595:MET:HE1	2.09	0.74
1:B:94:LEU:HD23	1:B:94:LEU:O	1.88	0.73
1:A:308:GLU:OE1	1:A:325:MET:HE1	1.89	0.72
1:B:86:HIS:N	4:B:1323:HOH:O	1.94	0.72
1:B:685:VAL:O	1:B:685:VAL:HG23	1.85	0.72
1:A:173:ARG:HD2	1:A:175:ARG:HH21	1.54	0.72
1:A:595:MET:C	1:A:595:MET:HE3	2.16	0.71
1:A:307:VAL:HA	1:A:452:MET:CE	2.18	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:689:LYS:O	1:A:690:MET:HE2	1.94	0.68
1:B:682:SER:HA	1:B:685:VAL:O	1.95	0.67
1:A:299:HIS:HD2	1:A:459:TYR:OH	1.78	0.66
1:B:222:THR:H	1:B:645:ASN:HD21	1.41	0.66
1:B:420:LYS:NZ	4:B:1176:HOH:O	2.28	0.66
1:B:685:VAL:O	1:B:685:VAL:CG2	2.40	0.66
1:A:318:GLN:HG2	4:A:1205:HOH:O	1.95	0.66
1:A:418:ARG:O	1:A:419:GLU:HG2	1.95	0.65
1:A:595:MET:HE3	1:A:595:MET:O	1.97	0.65
1:B:299:HIS:HD2	1:B:459:TYR:OH	1.81	0.63
1:A:317:ILE:CG2	1:A:452:MET:HE2	2.29	0.62
1:A:151:ARG:HD3	4:A:1045:HOH:O	2.00	0.62
1:A:592:LYS:H	1:A:595:MET:HE2	1.64	0.62
1:B:287:ASN:OD1	1:B:390:ARG:NH1	2.34	0.60
1:B:94:LEU:HD23	1:B:94:LEU:C	2.27	0.59
1:A:86:HIS:CB	4:A:1296:HOH:O	2.51	0.58
1:B:454:ARG:NH1	1:B:493:ASP:OD1	2.37	0.56
1:B:382:LEU:HD11	1:B:548:GLU:HB3	1.89	0.55
1:A:30:VAL:H	1:A:33:GLN:HE21	1.55	0.55
1:A:130:ILE:HD12	1:A:188:PHE:CZ	2.43	0.54
1:A:418:ARG:O	1:A:419:GLU:CB	2.55	0.53
1:B:130:ILE:HD12	1:B:130:ILE:N	2.24	0.52
1:A:689:LYS:C	1:A:690:MET:HE2	2.35	0.52
1:B:335:TYR:O	1:B:339:LEU:HG	2.09	0.52
1:B:345:ILE:HD12	1:B:685:VAL:HG21	1.92	0.50
1:B:499:LYS:HB2	1:B:505:GLU:HG2	1.93	0.50
1:B:220:ARG:HB3	1:B:645:ASN:HD22	1.77	0.50
1:B:151:ARG:HD3	4:B:923:HOH:O	2.12	0.49
1:A:317:ILE:HG23	1:A:452:MET:HE2	1.94	0.48
1:B:265:ASP:OD2	2:B:802:GOL:H32	2.14	0.48
1:A:418:ARG:O	1:A:419:GLU:CG	2.61	0.48
1:B:676:ASN:O	1:B:679:SER:HB2	2.13	0.48
1:A:222:THR:H	1:A:645:ASN:HD21	1.60	0.48
1:B:678:GLU:HG2	1:B:690:MET:HE3	1.94	0.48
1:A:521:ASN:H	1:A:521:ASN:HD22	1.60	0.47
1:B:345:ILE:CD1	1:B:685:VAL:HG21	2.44	0.47
1:A:218:TRP:CG	1:A:219:PRO:HA	2.49	0.47
1:A:281:LEU:HD23	1:A:380:LEU:CD1	2.45	0.47
1:B:653:THR:H	1:B:656:ASN:HD22	1.62	0.47
1:B:362:LEU:HD21	1:B:559:ALA:HB1	1.98	0.46
1:B:638:ARG:HH12	2:B:802:GOL:H31	1.79	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:281:LEU:HD23	1:A:380:LEU:HD13	1.97	0.46
1:A:303:GLN:NE2	4:A:1310:HOH:O	2.46	0.46
1:B:705:MET:O	1:B:710:PRO:HA	2.16	0.46
1:A:394:LEU:C	1:A:394:LEU:HD23	2.40	0.45
1:A:303:GLN:NE2	1:A:460:TRP:HE1	2.15	0.45
1:B:215:ASN:O	1:B:216:TRP:HB2	2.17	0.45
1:A:404:MET:HA	1:A:404:MET:HE2	1.98	0.45
1:B:686:PHE:CD1	1:B:686:PHE:C	2.95	0.45
1:A:391:ASP:HB2	1:A:471:GLN:HE21	1.82	0.44
1:B:465:ILE:O	1:B:472:ARG:NH2	2.50	0.44
1:A:127:LEU:HA	1:A:190:ASN:ND2	2.33	0.44
1:B:210:GLY:HA2	1:B:616:PHE:CE1	2.53	0.44
1:B:218:TRP:CG	1:B:219:PRO:HA	2.51	0.44
1:B:127:LEU:HD21	1:B:130:ILE:HD11	2.00	0.43
1:A:218:TRP:HA	1:A:219:PRO:C	2.43	0.43
1:A:218:TRP:CD2	1:A:219:PRO:HA	2.54	0.43
1:A:86:HIS:HB3	4:A:1296:HOH:O	2.16	0.43
1:A:483:ASN:O	1:A:483:ASN:ND2	2.52	0.43
1:A:591:PRO:CD	1:A:595:MET:CE	2.84	0.42
1:B:222:THR:H	1:B:645:ASN:ND2	2.14	0.42
1:B:638:ARG:HH12	2:B:802:GOL:C3	2.32	0.42
1:A:160:LYS:HG2	1:A:430:GLN:NE2	2.33	0.42
1:B:303:GLN:NE2	1:B:460:TRP:HE1	2.17	0.42
1:A:370:GLN:N	1:A:371:PRO:CD	2.82	0.42
1:A:378:LYS:HA	1:A:378:LYS:HD3	1.97	0.42
1:A:316:ASP:OD1	1:A:320:LYS:HE2	2.20	0.42
1:A:527:TYR:CE1	1:A:531:VAL:HG11	2.54	0.42
1:A:290:GLN:NE2	2:A:802:GOL:H11	2.35	0.41
1:B:134:THR:HG22	1:B:138:LYS:HD2	2.01	0.41
1:A:607:VAL:HG22	1:A:623:LEU:HD23	2.02	0.41
1:A:690:MET:HE2	1:A:690:MET:HA	2.02	0.41
1:A:318:GLN:CG	4:A:1205:HOH:O	2.63	0.41
1:A:473:VAL:HG13	1:A:534:THR:HG21	2.02	0.41
1:B:218:TRP:HA	1:B:219:PRO:C	2.46	0.41
1:A:622:LEU:HD23	1:A:622:LEU:C	2.46	0.40
1:B:59:ASP:HA	1:B:60:PRO:HA	1.93	0.40
1:B:207:GLY:HA2	1:B:223:GLY:O	2.21	0.40
1:B:252:HIS:CE1	2:B:803:GOL:H32	2.57	0.40
1:A:86:HIS:ND1	4:A:1296:HOH:O	2.37	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	696/698 (100%)	674 (97%)	19 (3%)	3 (0%)	30	34
1	B	696/698 (100%)	677 (97%)	18 (3%)	1 (0%)	48	57
All	All	1392/1396 (100%)	1351 (97%)	37 (3%)	4 (0%)	36	42

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	612	GLY
1	B	367	ALA
1	A	367	ALA
1	A	68	GLY

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 10 ligands modelled in this entry, 3 are monoatomic - leaving 7 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	GOL	A	804	-	5,5,5	0.68	0	5,5,5	1.79	2 (40%)
2	GOL	A	803	-	5,5,5	0.55	0	5,5,5	1.22	1 (20%)
2	GOL	B	802	-	5,5,5	1.33	0	5,5,5	1.47	1 (20%)
2	GOL	A	802	-	5,5,5	0.52	0	5,5,5	0.47	0
2	GOL	B	801	-	5,5,5	0.61	0	5,5,5	0.48	0
2	GOL	B	803	-	5,5,5	0.96	0	5,5,5	1.34	0
2	GOL	A	801	-	5,5,5	0.21	0	5,5,5	1.17	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	804	-	-	2/4/4/4	-
2	GOL	A	803	-	-	0/4/4/4	-
2	GOL	B	802	-	-	2/4/4/4	-
2	GOL	A	802	-	-	4/4/4/4	-
2	GOL	B	801	-	-	0/4/4/4	-
2	GOL	B	803	-	-	3/4/4/4	-
2	GOL	A	801	-	-	0/4/4/4	-

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	804	GOL	C3-C2-C1	-3.05	100.60	111.80
2	A	804	GOL	O2-C2-C1	2.49	119.47	109.18
2	B	802	GOL	C3-C2-C1	2.47	120.87	111.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	803	GOL	C3-C2-C1	-2.04	104.33	111.80

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	802	GOL	O1-C1-C2-C3
2	A	802	GOL	C1-C2-C3-O3
2	B	802	GOL	C1-C2-C3-O3
2	B	803	GOL	C1-C2-C3-O3
2	B	803	GOL	O2-C2-C3-O3
2	A	804	GOL	C1-C2-C3-O3
2	B	802	GOL	O2-C2-C3-O3
2	A	802	GOL	O2-C2-C3-O3
2	A	802	GOL	O1-C1-C2-O2
2	B	803	GOL	O1-C1-C2-O2
2	A	804	GOL	O1-C1-C2-O2

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	802	GOL	3	0
2	A	802	GOL	1	0
2	B	803	GOL	1	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	698/698 (100%)	-0.44	6 (0%) 81 79	6, 18, 40, 60	0
1	B	698/698 (100%)	-0.53	4 (0%) 85 83	6, 17, 35, 70	0
All	All	1396/1396 (100%)	-0.49	10 (0%) 84 82	6, 17, 38, 70	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	97	THR	3.2
1	A	98	ALA	3.1
1	A	507	LEU	3.1
1	A	485	ALA	2.8
1	B	679	SER	2.4
1	A	508	LYS	2.3
1	B	677	TRP	2.2
1	B	312	LYS	2.1
1	B	366	GLY	2.1
1	A	314	ASN	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	GOL	A	804	6/6	0.89	0.14	32,41,44,47	0
2	GOL	B	803	6/6	0.89	0.13	27,32,34,35	0
2	GOL	B	802	6/6	0.90	0.10	18,20,20,22	0
2	GOL	A	803	6/6	0.92	0.09	19,19,23,24	0
2	GOL	A	802	6/6	0.93	0.11	38,40,40,40	0
2	GOL	A	801	6/6	0.96	0.06	12,14,15,18	0
2	GOL	B	801	6/6	0.96	0.06	14,18,20,20	0
3	ZN	A	805	1/1	0.96	0.16	47,47,47,47	0
3	ZN	B	804	1/1	0.98	0.12	36,36,36,36	0
3	ZN	B	805	1/1	1.00	0.03	15,15,15,15	0

6.5 Other polymers [i](#)

There are no such residues in this entry.