



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 09:31 AM GMT

PDB ID : 3ISL  
Title : Crystal structure of ureidoglycine-glyoxylate aminotransferase (pucG) from *Bacillus subtilis*  
Authors : Costa, R.; Cendron, L.; Ramazzina, I.; Berni, R.; Peracchi, A.; Percudani, R.; Zanutti, G.  
Deposited on : 2009-08-26  
Resolution : 2.06 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](mailto: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.

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

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
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) : trunk26865

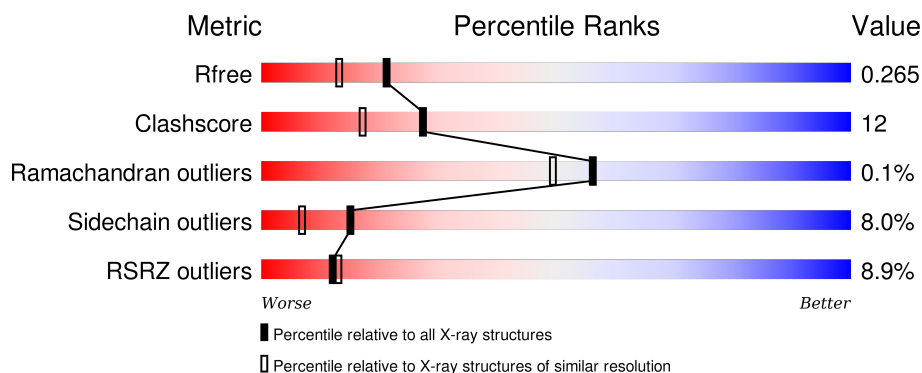
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.06 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	1799 (2.08-2.04)
Clashscore	102246	1910 (2.08-2.04)
Ramachandran outliers	100387	1893 (2.08-2.04)
Sidechain outliers	100360	1893 (2.08-2.04)
RSRZ outliers	91569	1802 (2.08-2.04)

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  $\geq 3$ , 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	416	<div> <div>9%</div> <div>69%</div> <div>20%</div> <div>7%</div> </div>
1	B	416	<div> <div>8%</div> <div>71%</div> <div>19%</div> <div>7%</div> </div>

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	PLP	A	419	-	X	-	-

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 6338 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 Purine catabolism protein pucG.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	387	Total	C	N	O	S	0	0	0
			3001	1899	521	560	21			
1	B	387	Total	C	N	O	S	0	0	0
			3001	1899	521	560	21			

- Molecule 2 is PYRIDOXAL-5'-PHOSPHATE (three-letter code: PLP) (formula:  $C_8H_{10}NO_6P$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			15	8	1	5	1		
2	B	1	Total	C	N	O	P	0	0
			15	8	1	5	1		

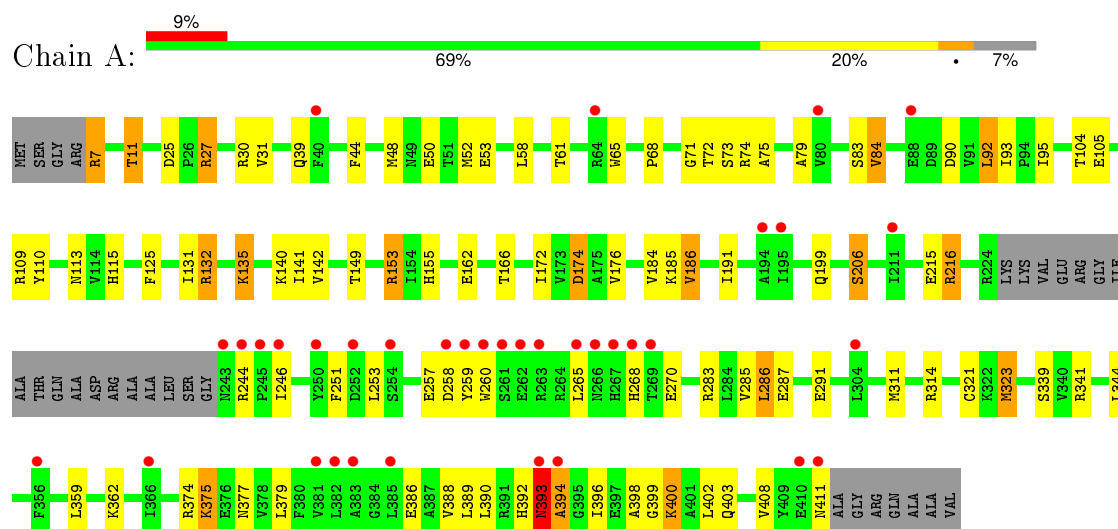
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	153	Total 153	O 153	0	0
3	B	153	Total 153	O 153	0	0

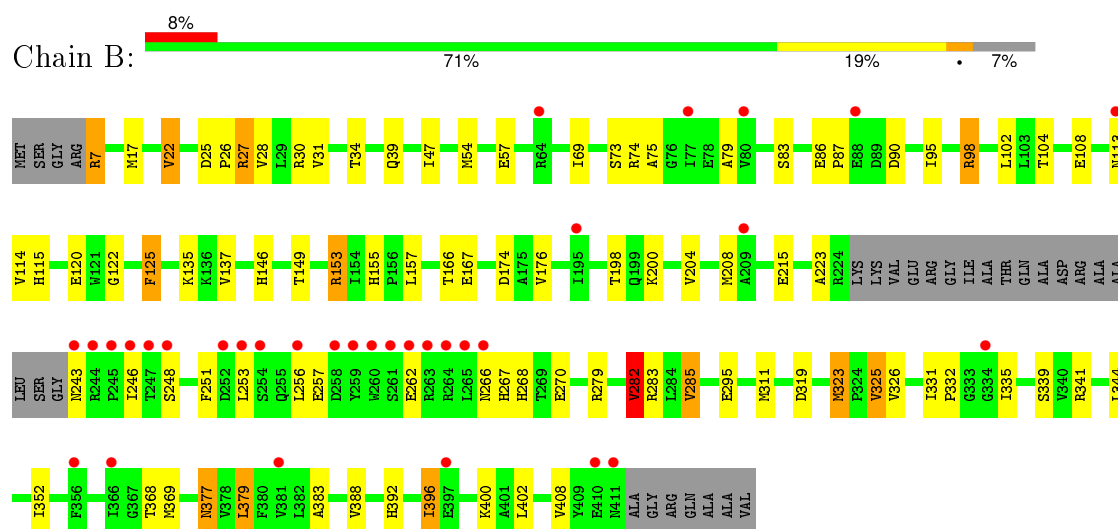
### 3 Residue-property plots [i](#)

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: Purine catabolism protein pucG



- Molecule 1: Purine catabolism protein pucG



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	102.40Å 95.99Å 101.29Å 90.00° 114.23° 90.00°	Depositor
Resolution (Å)	27.69 – 2.06 27.69 – 2.06	Depositor EDS
% Data completeness (in resolution range)	99.5 (27.69-2.06) 99.6 (27.69-2.06)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.53 (at 2.06Å)	Xtriage
Refinement program	REFMAC 5.5.0072	Depositor
R, $R_{free}$	0.193 , 0.247 0.212 , 0.265	Depositor DCC
$R_{free}$ test set	2787 reflections (5.35%)	DCC
Wilson B-factor (Å <sup>2</sup> )	31.0	Xtriage
Anisotropy	0.024	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 37.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 54876 reflections	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	6338	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.13% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: PLP

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	1.12	6/3058 (0.2%)	1.07	18/4143 (0.4%)
1	B	1.08	2/3058 (0.1%)	1.10	18/4143 (0.4%)
All	All	1.10	8/6116 (0.1%)	1.08	36/8286 (0.4%)

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	1
All	All	0	2

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	325	VAL	CB-CG1	-7.11	1.38	1.52
1	B	57	GLU	CB-CG	-6.83	1.39	1.52
1	A	215	GLU	CG-CD	5.62	1.60	1.51
1	A	53	GLU	CG-CD	5.56	1.60	1.51
1	A	135	LYS	CD-CE	5.54	1.65	1.51
1	A	110	TYR	CE2-CZ	5.18	1.45	1.38
1	A	185	LYS	CE-NZ	5.15	1.61	1.49
1	A	53	GLU	CD-OE1	5.05	1.31	1.25

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	27	ARG	NE-CZ-NH2	-14.12	113.24	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	153	ARG	NE-CZ-NH1	12.72	126.66	120.30
1	B	27	ARG	NE-CZ-NH1	11.72	126.16	120.30
1	B	30	ARG	NE-CZ-NH2	-10.34	115.13	120.30
1	A	30	ARG	NE-CZ-NH2	-10.30	115.15	120.30
1	A	27	ARG	NE-CZ-NH2	-10.25	115.18	120.30
1	B	30	ARG	NE-CZ-NH1	10.14	125.37	120.30
1	A	30	ARG	NE-CZ-NH1	9.06	124.83	120.30
1	B	153	ARG	CG-CD-NE	8.65	129.97	111.80
1	B	153	ARG	NE-CZ-NH2	-8.64	115.98	120.30
1	B	153	ARG	CD-NE-CZ	7.27	133.78	123.60
1	A	27	ARG	NE-CZ-NH1	7.21	123.91	120.30
1	A	153	ARG	NE-CZ-NH1	-7.16	116.72	120.30
1	B	285	VAL	CG1-CB-CG2	6.95	122.01	110.90
1	B	282	VAL	CG1-CB-CG2	6.69	121.61	110.90
1	B	174	ASP	CB-CG-OD1	6.65	124.28	118.30
1	A	286	LEU	CB-CG-CD1	6.55	122.13	111.00
1	A	285	VAL	CG1-CB-CG2	6.41	121.16	110.90
1	B	22	VAL	CG1-CB-CG2	6.26	120.92	110.90
1	A	374	ARG	NE-CZ-NH1	-6.07	117.27	120.30
1	A	153	ARG	NE-CZ-NH2	6.05	123.33	120.30
1	A	84	VAL	CG1-CB-CG2	5.95	120.43	110.90
1	B	283	ARG	NE-CZ-NH2	-5.90	117.35	120.30
1	A	135	LYS	CD-CE-NZ	5.86	125.17	111.70
1	A	92	LEU	CB-CG-CD2	5.80	120.87	111.00
1	A	216	ARG	NE-CZ-NH2	-5.62	117.49	120.30
1	A	258	ASP	CB-CG-OD1	5.50	123.25	118.30
1	A	174	ASP	CB-CG-OD1	5.46	123.21	118.30
1	B	22	VAL	N-CA-CB	-5.30	99.83	111.50
1	B	379	LEU	CA-CB-CG	5.29	127.46	115.30
1	A	132	ARG	NE-CZ-NH2	-5.16	117.72	120.30
1	B	379	LEU	CB-CG-CD1	5.15	119.75	111.00
1	B	325	VAL	CG1-CB-CG2	-5.14	102.68	110.90
1	A	142	VAL	CG1-CB-CG2	5.09	119.05	110.90
1	A	71	GLY	N-CA-C	-5.04	100.49	113.10
1	B	285	VAL	CA-CB-CG1	5.04	118.46	110.90

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	393	ASN	Peptide
1	B	125	PHE	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	3001	0	3005	88	0
1	B	3001	0	3005	57	0
2	A	15	0	7	2	0
2	B	15	0	7	2	0
3	A	153	0	0	7	0
3	B	153	0	0	3	0
All	All	6338	0	6024	143	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 12.

All (143) 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:375:LYS:NZ	1:A:375:LYS:HB2	1.55	1.18
1:A:7:ARG:HH11	1:A:7:ARG:CG	1.61	1.14
1:A:375:LYS:HZ3	1:A:375:LYS:HB2	0.90	1.05
1:A:389:LEU:O	1:A:394:ALA:HB2	1.57	1.04
1:A:7:ARG:HG2	1:A:7:ARG:NH1	1.51	1.01
1:B:83:SER:HA	1:B:246:ILE:HD12	1.41	0.99
1:A:375:LYS:HZ3	1:A:375:LYS:CB	1.76	0.94
1:A:339:SER:OG	1:A:392:HIS:HE1	1.50	0.93
1:A:44:PHE:CE2	1:A:48:MET:HE2	2.05	0.90
1:A:393:ASN:HA	1:A:394:ALA:HB3	1.54	0.89
1:A:375:LYS:NZ	1:A:375:LYS:CB	2.23	0.89
1:A:83:SER:HA	1:A:246:ILE:HG12	1.52	0.89
1:A:44:PHE:HD2	1:A:48:MET:HE3	1.39	0.87
1:A:11:THR:HG23	3:B:433:HOH:O	1.74	0.87
1:B:331:ILE:HG23	1:B:335:ILE:HD11	1.57	0.85
1:B:325:VAL:HG11	1:B:369:MET:HG2	1.59	0.84
1:A:44:PHE:CD2	1:A:48:MET:CE	2.60	0.84
1:A:52:MET:HE2	1:A:260:TRP:HB3	1.60	0.82
1:A:44:PHE:CD2	1:A:48:MET:HE3	2.16	0.81
1:B:325:VAL:CG1	1:B:369:MET:HG2	2.11	0.79
1:A:386:GLU:OE1	1:A:398:ALA:O	1.99	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:268:HIS:HE1	1:A:270:GLU:OE2	1.65	0.79
1:A:390:LEU:HA	1:A:394:ALA:HB1	1.66	0.78
1:A:149:THR:HA	1:A:323:MET:HE1	1.66	0.76
1:A:311:MET:CE	1:A:389:LEU:HD12	2.17	0.75
1:A:65:TRP:CD1	3:A:427:HOH:O	2.40	0.74
1:A:79:ALA:HA	1:A:251:PHE:HB3	1.70	0.74
1:A:44:PHE:CE2	1:A:48:MET:CE	2.71	0.74
1:B:7:ARG:CZ	1:B:7:ARG:HB2	2.18	0.74
1:A:52:MET:CE	1:A:260:TRP:HE3	2.02	0.73
1:B:388:VAL:O	1:B:392:HIS:HD2	1.72	0.72
1:B:325:VAL:O	1:B:325:VAL:HG12	1.88	0.72
1:A:153:ARG:HD3	1:A:321:CYS:SG	2.29	0.72
1:A:44:PHE:CD2	1:A:48:MET:HE2	2.25	0.71
1:A:7:ARG:HG2	1:A:7:ARG:HH11	0.69	0.71
1:B:311:MET:O	1:B:396:ILE:HG22	1.89	0.71
1:A:52:MET:HE3	1:A:260:TRP:HE3	1.57	0.70
1:A:390:LEU:HA	1:A:394:ALA:CB	2.22	0.69
1:A:268:HIS:CE1	1:A:270:GLU:OE2	2.46	0.69
1:A:95:ILE:HD11	1:A:104:THR:HG21	1.73	0.69
1:A:65:TRP:HD1	3:A:427:HOH:O	1.75	0.68
1:B:223:ALA:O	1:B:243:ASN:HB3	1.93	0.68
1:B:25:ASP:OD1	1:B:27:ARG:HD3	1.96	0.65
1:B:339:SER:OG	1:B:392:HIS:HE1	1.79	0.65
1:A:74:ARG:HH21	2:A:419:PLP:H5A2	1.62	0.63
1:A:311:MET:HE3	1:A:386:GLU:HA	1.79	0.63
1:A:95:ILE:HD11	1:A:104:THR:CG2	2.30	0.62
1:A:341:ARG:NH1	3:A:483:HOH:O	2.33	0.61
1:B:383:ALA:HB2	1:B:402:LEU:HD11	1.82	0.61
1:A:339:SER:OG	1:A:392:HIS:CE1	2.43	0.60
1:A:393:ASN:HA	1:A:394:ALA:CB	2.30	0.59
1:A:25:ASP:OD1	1:A:27:ARG:HD3	2.03	0.59
1:B:79:ALA:HA	1:B:251:PHE:HB3	1.85	0.58
1:A:393:ASN:CA	1:A:394:ALA:HB3	2.31	0.58
1:A:311:MET:HE2	1:A:389:LEU:HD12	1.83	0.58
1:B:253:LEU:HA	1:B:256:LEU:HB2	1.87	0.56
1:B:341:ARG:HD2	1:B:352:ILE:O	2.05	0.56
1:A:393:ASN:CA	1:A:394:ALA:CB	2.84	0.56
1:B:149:THR:HA	1:B:323:MET:HE1	1.86	0.56
1:B:108:GLU:HG3	1:B:114:VAL:HG21	1.88	0.56
1:A:153:ARG:HH11	1:A:153:ARG:HG2	1.70	0.55
1:A:375:LYS:HZ2	1:A:375:LYS:CB	2.08	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:125:PHE:H	1:B:155:HIS:HE1	1.54	0.54
1:A:311:MET:HE1	1:A:389:LEU:HD12	1.88	0.54
1:B:332:PRO:HD2	1:B:335:ILE:HG12	1.88	0.54
1:B:75:ALA:HB1	1:B:251:PHE:CE1	2.42	0.54
1:A:90:ASP:OD1	1:A:115:HIS:HE1	1.90	0.54
1:A:50:GLU:OE2	1:B:7:ARG:NH2	2.41	0.54
1:A:74:ARG:NH2	2:A:419:PLP:H5A2	2.23	0.53
1:A:61:THR:HB	1:A:186:VAL:HG22	1.91	0.53
1:B:7:ARG:CB	1:B:7:ARG:CZ	2.86	0.53
1:B:325:VAL:CG1	1:B:369:MET:CG	2.83	0.53
1:B:200:LYS:NZ	2:B:419:PLP:O3	2.39	0.53
1:B:146:HIS:CD2	1:B:157:LEU:HD11	2.44	0.53
1:B:95:ILE:HD11	1:B:104:THR:HG21	1.90	0.52
1:B:267:HIS:ND1	1:B:268:HIS:HD2	2.07	0.52
1:A:93:ILE:HD12	1:A:104:THR:HG22	1.90	0.52
1:B:335:ILE:C	1:B:335:ILE:HD12	2.30	0.51
1:B:28:VAL:O	1:B:31:VAL:HG12	2.10	0.51
1:A:79:ALA:CA	1:A:251:PHE:HB3	2.41	0.51
1:A:52:MET:HE2	1:A:260:TRP:CB	2.38	0.51
1:A:259:TYR:HD1	1:A:265:LEU:O	1.94	0.50
1:B:125:PHE:H	1:B:155:HIS:CE1	2.29	0.50
1:A:253:LEU:O	1:A:257:GLU:HG2	2.11	0.50
1:B:54:MET:HB3	1:B:282:VAL:HG22	1.92	0.50
1:A:75:ALA:HB1	1:A:251:PHE:CE1	2.46	0.50
1:A:283:ARG:O	1:A:287:GLU:HG3	2.10	0.49
1:A:216:ARG:HD2	3:A:533:HOH:O	2.11	0.49
1:A:153:ARG:HH11	1:A:153:ARG:CG	2.25	0.49
1:B:388:VAL:O	1:B:392:HIS:CD2	2.61	0.49
1:B:198:THR:HB	1:B:204:VAL:HB	1.94	0.48
1:B:253:LEU:O	1:B:257:GLU:HG2	2.13	0.48
1:A:386:GLU:CD	1:A:398:ALA:O	2.51	0.48
1:B:331:ILE:HG23	1:B:335:ILE:CD1	2.37	0.48
1:B:73:SER:HB3	1:B:176:VAL:HG21	1.95	0.48
1:A:11:THR:HG21	1:B:34:THR:HG21	1.96	0.48
1:B:267:HIS:ND1	1:B:268:HIS:CD2	2.81	0.48
1:A:68:PRO:HD2	1:A:260:TRP:CZ2	2.50	0.47
1:A:141:ILE:HG23	1:A:172:ILE:HD12	1.96	0.47
1:A:153:ARG:NH1	1:A:153:ARG:HG2	2.30	0.46
1:A:411:ASN:C	3:A:506:HOH:O	2.53	0.46
1:B:325:VAL:HG13	1:B:368:THR:O	2.16	0.46
1:B:90:ASP:OD1	1:B:115:HIS:HE1	2.00	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:75:ALA:HB1	1:B:251:PHE:CZ	2.52	0.45
1:A:105:GLU:O	1:A:109:ARG:HG3	2.16	0.45
1:B:74:ARG:HG2	1:B:102:LEU:HD21	1.98	0.45
1:A:140:LYS:NZ	3:A:462:HOH:O	2.38	0.45
1:A:153:ARG:CD	1:A:321:CYS:SG	3.03	0.45
1:A:52:MET:HE2	1:A:260:TRP:HE3	1.77	0.45
1:A:125:PHE:H	1:A:155:HIS:HE1	1.65	0.45
1:B:86:GLU:O	1:B:87:PRO:C	2.54	0.44
1:B:167:GLU:HB2	3:B:443:HOH:O	2.16	0.44
1:B:122:GLY:O	1:B:153:ARG:HD2	2.18	0.44
1:A:7:ARG:CG	1:A:7:ARG:NH1	2.35	0.44
1:A:79:ALA:CB	1:A:251:PHE:HB3	2.47	0.44
1:B:98:ARG:HD3	1:B:98:ARG:C	2.36	0.44
1:A:259:TYR:CD1	1:A:265:LEU:O	2.70	0.44
1:B:325:VAL:CG1	1:B:325:VAL:O	2.60	0.43
1:A:186:VAL:HB	1:A:191:ILE:HB	2.00	0.43
1:A:73:SER:HB3	1:A:176:VAL:CG2	2.48	0.43
1:A:52:MET:HE2	1:A:260:TRP:CE3	2.53	0.43
1:B:73:SER:HB2	2:B:419:PLP:P	2.59	0.43
1:A:359:LEU:HA	1:A:362:LYS:HD3	2.01	0.42
1:B:149:THR:HG22	1:B:326:VAL:HG21	2.02	0.42
1:A:402:LEU:HD23	1:A:402:LEU:HA	1.73	0.42
1:A:399:GLY:O	1:A:403:GLN:HG2	2.18	0.42
1:B:208:MET:CE	1:B:270:GLU:OE1	2.67	0.42
1:A:344:LEU:HA	1:A:344:LEU:HD23	1.90	0.42
1:A:162:GLU:O	1:A:166:THR:HG23	2.19	0.42
1:B:47:ILE:HG23	1:B:279:ARG:HG3	2.00	0.41
1:A:73:SER:HB3	1:A:176:VAL:HG21	2.03	0.41
1:A:131:ILE:HG22	1:A:135:LYS:HE2	2.02	0.41
1:B:325:VAL:HG12	1:B:369:MET:CG	2.50	0.41
1:A:125:PHE:H	1:A:155:HIS:CE1	2.39	0.41
1:B:248:SER:HB3	1:B:251:PHE:HB2	2.02	0.41
1:A:48:MET:HB3	1:A:260:TRP:CZ3	2.54	0.41
1:B:95:ILE:HD11	1:B:104:THR:CG2	2.51	0.41
1:A:388:VAL:O	1:A:392:HIS:HD2	2.03	0.41
1:B:27:ARG:NH2	3:B:458:HOH:O	2.39	0.41
1:A:199:GLN:HB3	1:A:206:SER:HB3	2.02	0.40
1:B:402:LEU:HD23	1:B:402:LEU:HA	1.77	0.40
1:A:400:LYS:HE3	3:A:509:HOH:O	2.21	0.40
1:B:17:MET:SD	1:B:377:ASN:HB3	2.61	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	383/416 (92%)	367 (96%)	15 (4%)	1 (0%)	46	36
1	B	383/416 (92%)	365 (95%)	18 (5%)	0	100	100
All	All	766/832 (92%)	732 (96%)	33 (4%)	1 (0%)	56	49

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	394	ALA

### 5.3.2 Protein sidechains [i](#)

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	318/336 (95%)	292 (92%)	26 (8%)	14	6
1	B	318/336 (95%)	293 (92%)	25 (8%)	15	7
All	All	636/672 (95%)	585 (92%)	51 (8%)	15	7

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	ARG
1	A	11	THR
1	A	31	VAL
1	A	39	GLN
1	A	58	LEU

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Mol	Chain	Res	Type
1	A	72	THR
1	A	84	VAL
1	A	92	LEU
1	A	113	ASN
1	A	132	ARG
1	A	174	ASP
1	A	184	VAL
1	A	186	VAL
1	A	206	SER
1	A	244	ARG
1	A	286	LEU
1	A	291	GLU
1	A	314	ARG
1	A	323	MET
1	A	375	LYS
1	A	377	ASN
1	A	379	LEU
1	A	393	ASN
1	A	396	ILE
1	A	400	LYS
1	A	408	VAL
1	B	7	ARG
1	B	22	VAL
1	B	26	PRO
1	B	39	GLN
1	B	69	ILE
1	B	98	ARG
1	B	113	ASN
1	B	120	GLU
1	B	135	LYS
1	B	137	VAL
1	B	166	THR
1	B	215	GLU
1	B	262	GLU
1	B	266	ASN
1	B	282	VAL
1	B	285	VAL
1	B	295	GLU
1	B	319	ASP
1	B	323	MET
1	B	344	LEU
1	B	377	ASN

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Mol	Chain	Res	Type
1	B	379	LEU
1	B	396	ILE
1	B	400	LYS
1	B	408	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	113	ASN
1	A	115	HIS
1	A	155	HIS
1	A	268	HIS
1	A	347	GLN
1	A	377	ASN
1	A	392	HIS
1	B	39	GLN
1	B	113	ASN
1	B	115	HIS
1	B	155	HIS
1	B	266	ASN
1	B	268	HIS
1	B	377	ASN
1	B	392	HIS

### 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 ⓘ

2 ligands are modelled in this entry.



In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z  > 2$	Counts	RMSZ	# $ Z  > 2$
2	PLP	A	419	1	15,15,16	2.97	8 (53%)	21,22,23	2.57	10 (47%)
2	PLP	B	419	1	15,15,16	2.50	8 (53%)	21,22,23	2.80	8 (38%)

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	PLP	A	419	1	-	0/6/6/8	0/1/1/1
2	PLP	B	419	1	-	0/6/6/8	0/1/1/1

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	419	PLP	C2A-C2	-4.61	1.41	1.50
2	A	419	PLP	C2A-C2	-4.51	1.41	1.50
2	A	419	PLP	O4P-C5A	-2.21	1.35	1.44
2	B	419	PLP	P-O2P	2.28	1.62	1.54
2	B	419	PLP	P-O3P	2.45	1.63	1.54
2	B	419	PLP	C5-C4	2.58	1.43	1.40
2	A	419	PLP	C6-C5	2.68	1.43	1.37
2	B	419	PLP	C6-N1	2.90	1.40	1.34
2	B	419	PLP	C4A-C4	3.20	1.58	1.51
2	A	419	PLP	C4A-C4	3.21	1.58	1.51
2	B	419	PLP	P-O4P	3.26	1.71	1.60
2	A	419	PLP	P-O2P	3.33	1.66	1.54
2	A	419	PLP	P-O3P	3.83	1.68	1.54
2	B	419	PLP	C2-N1	4.03	1.42	1.34
2	A	419	PLP	C6-N1	4.50	1.44	1.34
2	A	419	PLP	C2-N1	6.39	1.47	1.34

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	419	PLP	C2A-C2-C3	-6.49	113.21	121.04
2	B	419	PLP	C4A-C4-C3	-4.98	111.33	120.36
2	A	419	PLP	C5A-C5-C4	-4.21	116.07	121.65
2	B	419	PLP	C5A-C5-C6	-3.80	112.10	119.28
2	B	419	PLP	C5-C6-N1	-2.90	118.82	123.86
2	A	419	PLP	O2P-P-O4P	-2.25	100.09	106.56
2	A	419	PLP	C6-N1-C2	-2.21	114.76	119.28
2	A	419	PLP	C5-C6-N1	-2.06	120.29	123.86
2	A	419	PLP	O4P-P-O1P	2.06	112.38	107.14
2	A	419	PLP	O2P-P-O1P	2.11	117.37	110.58
2	B	419	PLP	O3-C3-C2	2.22	121.51	117.66
2	B	419	PLP	C6-C5-C4	2.36	120.15	118.15
2	A	419	PLP	O3-C3-C2	2.37	121.78	117.66
2	B	419	PLP	O4P-P-O1P	2.51	113.53	107.14
2	A	419	PLP	C2A-C2-N1	3.19	125.02	117.95
2	B	419	PLP	C5A-C5-C4	4.60	127.75	121.65
2	A	419	PLP	C6-C5-C4	5.63	122.92	118.15
2	B	419	PLP	C4A-C4-C5	8.10	129.32	120.88

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	419	PLP	2	0
2	B	419	PLP	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 ⓘ

### 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	387/416 (93%)	0.30	36 (9%) 11 12	14, 26, 42, 54	0
1	B	387/416 (93%)	0.32	33 (8%) 13 15	15, 27, 41, 53	0
All	All	774/832 (93%)	0.31	69 (8%) 12 13	14, 27, 42, 54	0

All (69) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	356	PHE	6.6
1	B	265	LEU	5.8
1	A	263	ARG	5.6
1	B	244	ARG	5.4
1	A	262	GLU	5.3
1	B	246	ILE	5.2
1	B	247	THR	5.0
1	A	266	ASN	4.8
1	B	263	ARG	4.5
1	B	262	GLU	4.5
1	A	356	PHE	4.5
1	B	260	TRP	4.4
1	A	245	PRO	4.4
1	B	245	PRO	4.2
1	A	246	ILE	4.1
1	A	244	ARG	4.1
1	A	259	TYR	4.0
1	B	243	ASN	3.9
1	B	411	ASN	3.8
1	A	261	SER	3.8
1	A	243	ASN	3.7
1	A	411	ASN	3.6
1	A	381	VAL	3.6
1	B	259	TYR	3.6

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Mol	Chain	Res	Type	RSRZ
1	A	88	GLU	3.5
1	B	254	SER	3.5
1	A	254	SER	3.4
1	B	80	VAL	3.3
1	A	265	LEU	3.2
1	A	260	TRP	3.2
1	A	393	ASN	3.1
1	A	385	LEU	3.1
1	B	252	ASP	3.0
1	A	382	LEU	2.9
1	B	334	GLY	2.8
1	B	195	ILE	2.8
1	B	266	ASN	2.7
1	A	366	ILE	2.6
1	B	381	VAL	2.6
1	A	267	HIS	2.6
1	B	248	SER	2.6
1	A	268	HIS	2.6
1	B	64	ARG	2.5
1	B	410	GLU	2.5
1	A	211	ILE	2.5
1	A	64	ARG	2.5
1	B	88	GLU	2.5
1	B	256	LEU	2.4
1	A	195	ILE	2.4
1	A	258	ASP	2.3
1	B	113	ASN	2.3
1	B	397	GLU	2.3
1	A	410	GLU	2.3
1	A	394	ALA	2.2
1	B	264	ARG	2.2
1	B	258	ASP	2.2
1	A	250	TYR	2.2
1	A	383	ALA	2.1
1	A	194	ALA	2.1
1	A	269	THR	2.1
1	A	252	ASP	2.1
1	B	366	ILE	2.1
1	A	304	LEU	2.1
1	B	261	SER	2.1
1	A	40	PHE	2.1
1	B	253	LEU	2.0

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Mol	Chain	Res	Type	RSRZ
1	B	209	ALA	2.0
1	A	80	VAL	2.0
1	B	77	ILE	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, 95<sup>th</sup> 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( $\text{\AA}^2$ )	Q<0.9
2	PLP	A	419	15/16	0.89	0.16	0.59	47,54,57,59	0
2	PLP	B	419	15/16	0.89	0.14	-0.42	45,53,56,56	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.