



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 04:52 PM UTC

PDB ID : 1GD9 / pdb_00001gd9
Title : CRYSTALL STRUCTURE OF PYROCOCCUS PROTEIN-A1
Authors : Ura, H.; Harata, K.; Matsui, I.; Kuramitsu, S.
Deposited on : 2000-09-22
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

<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

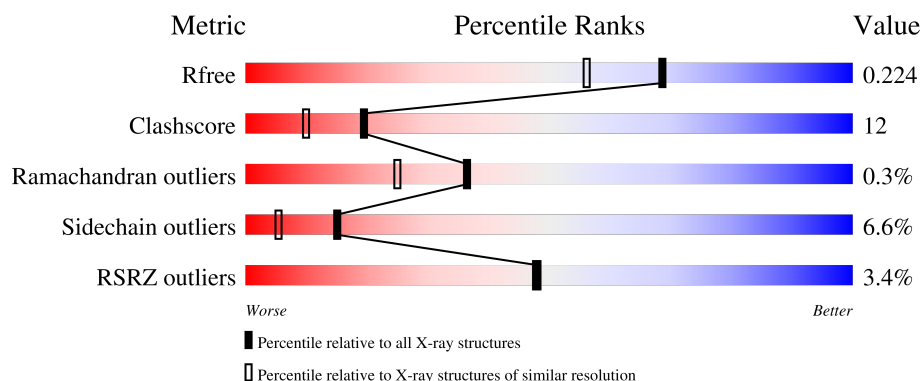
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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}	180053	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (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 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	389	3% 71% 25% .
1	B	389	4% 71% 24% .

2 Entry composition [i](#)

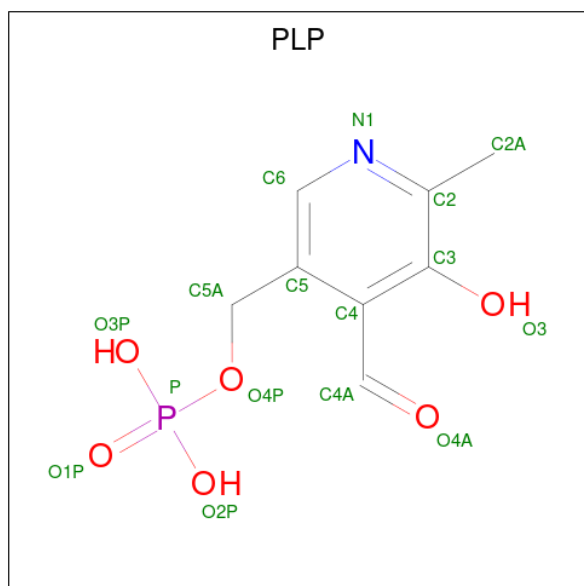
There are 3 unique types of molecules in this entry. The entry contains 6666 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 ASPARTATE AMINOTRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	388	Total	C	N	O	S	0	0	0
			3082	1981	515	572	14			
1	B	388	Total	C	N	O	S	0	0	0
			3082	1981	515	572	14			

- Molecule 2 is PYRIDOXAL-5'-PHOSPHATE (CCD ID: 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		

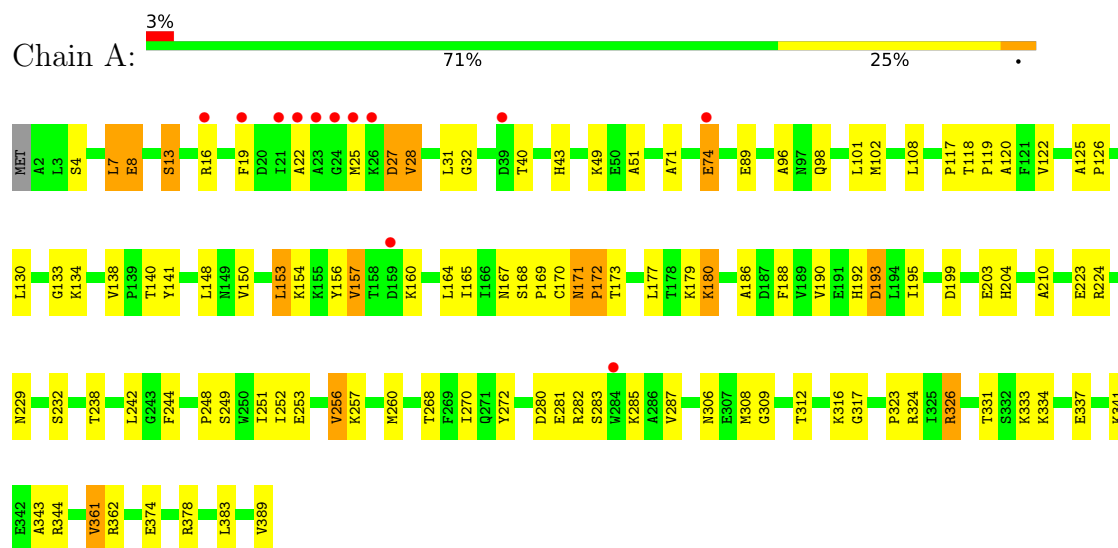
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	246	Total 246	O 246	0	0
3	B	226	Total 226	O 226	0	0

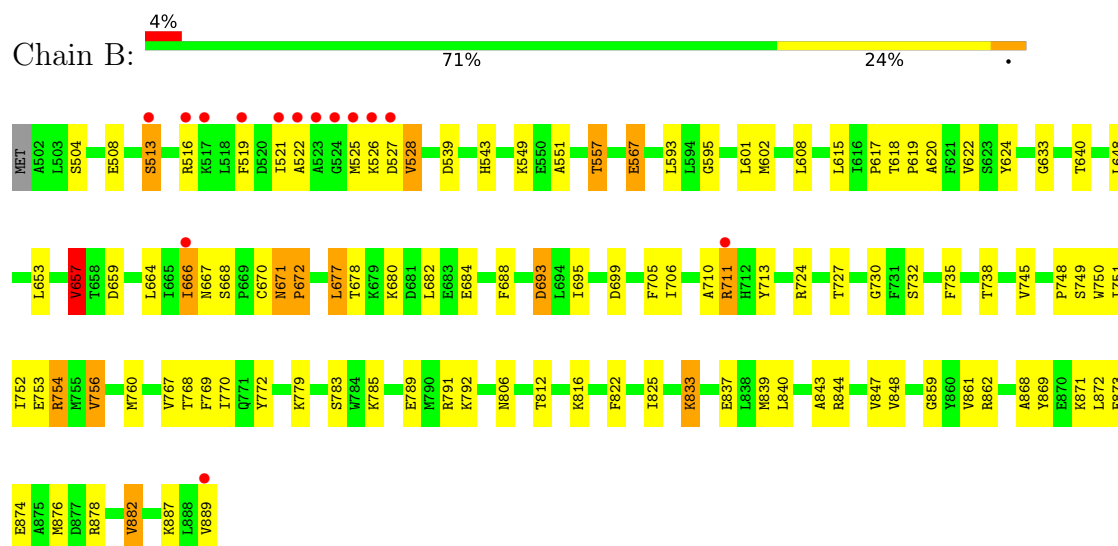
3 Residue-property plots

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: ASPARTATE AMINOTRANSFERASE



• Molecule 1: ASPARTATE AMINOTRANSFERASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	59.67Å 122.28Å 127.17Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 1.80 8.00 – 1.80	Depositor EDS
% Data completeness (in resolution range)	86.4 (8.00-1.80) 94.3 (8.00-1.80)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.54 (at 1.80Å)	Xtriage
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.210 , 0.242 0.196 , 0.224	Depositor DCC
R_{free} test set	8324 reflections (10.03%)	wwPDB-VP
Wilson B-factor (Å ²)	13.0	Xtriage
Anisotropy	0.260	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.58 , 80.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.017 for -h,l,k	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6666	wwPDB-VP
Average B, all atoms (Å ²)	12.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.79% 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: 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	0.50	0/3144	0.97	12/4241 (0.3%)
1	B	0.49	0/3144	1.00	15/4241 (0.4%)
All	All	0.49	0/6288	0.99	27/8482 (0.3%)

There are no bond length outliers.

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	735	PHE	N-CA-C	-9.69	95.41	110.42
1	B	671	ASN	N-CA-C	-8.76	95.97	109.64
1	A	171	ASN	N-CA-C	-8.41	96.52	109.64
1	A	170	CYS	N-CA-C	6.92	121.59	109.96
1	B	670	CYS	N-CA-C	6.68	120.56	109.46
1	B	624	TYR	N-CA-C	6.55	118.11	110.97
1	B	528	VAL	N-CA-C	6.45	117.40	108.89
1	A	195	ILE	N-CA-C	-6.24	99.44	108.36
1	A	122	VAL	N-CA-C	6.20	118.80	111.05
1	B	522	ALA	N-CA-C	-6.11	105.98	113.50
1	A	22	ALA	N-CA-C	-5.91	105.91	113.23
1	B	526	LYS	N-CA-C	5.75	118.12	107.60
1	B	657	VAL	CB-CA-C	-5.70	103.42	111.28
1	A	141	TYR	N-CA-C	5.49	118.42	109.59
1	A	156	TYR	N-CA-C	5.42	120.05	113.38
1	A	268	THR	N-CA-C	5.38	116.83	111.07
1	A	173	THR	N-CA-C	5.37	118.99	112.23
1	B	847	VAL	CA-C-N	-5.34	118.13	123.04
1	B	847	VAL	C-N-CA	-5.34	118.13	123.04
1	B	695	ILE	N-CA-C	-5.29	100.76	108.17
1	B	640	THR	N-CA-C	-5.22	101.37	109.72
1	A	28	VAL	N-CA-C	5.16	115.70	108.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	595	GLY	N-CA-C	-5.07	106.34	112.48
1	B	622	VAL	N-CA-C	5.07	119.19	111.89
1	A	140	THR	N-CA-C	-5.04	101.65	109.72
1	B	768	THR	N-CA-C	5.01	116.43	110.97
1	A	285	LYS	N-CA-C	-5.01	105.90	111.36

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	3082	0	3127	70	0
1	B	3082	0	3127	79	0
2	A	15	0	6	3	0
2	B	15	0	7	1	0
3	A	246	0	0	7	0
3	B	226	0	0	6	0
All	All	6666	0	6267	145	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 (145) 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:257:LYS:HA	1:A:260:MET:HE3	1.43	0.98
1:B:711:ARG:HG2	3:B:1163:HOH:O	1.79	0.82
1:A:193:ASP:HA	1:A:224:ARG:NH2	1.94	0.82
1:B:833:LYS:HE3	1:B:833:LYS:H	1.45	0.80
1:A:25:MET:HE2	1:A:344:ARG:NH1	2.00	0.77
1:B:754:ARG:HH11	1:B:754:ARG:HG3	1.54	0.70
1:A:71:ALA:O	1:A:74:GLU:HG3	1.92	0.69
1:A:331:THR:H	1:A:334:LYS:HD2	1.59	0.68
1:A:4:SER:O	1:A:8:GLU:HG2	1.92	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:306:ASN:HD21	1:A:312:THR:H	1.43	0.66
1:A:252:ILE:O	1:A:256:VAL:HG13	1.96	0.65
1:B:748:PRO:HD2	1:B:751:ILE:HD12	1.77	0.65
1:B:551:ALA:HB2	1:B:772:TYR:CD2	2.32	0.64
1:B:816:LYS:HE3	3:B:1344:HOH:O	1.97	0.64
1:A:51:ALA:HB2	1:A:272:TYR:CD2	2.34	0.62
1:B:754:ARG:HG3	1:B:754:ARG:NH1	2.14	0.62
1:B:504:SER:O	1:B:508:GLU:HG2	2.00	0.61
1:A:43:HIS:HE1	1:A:283:SER:OG	1.82	0.61
1:A:333:LYS:HG2	3:A:1250:HOH:O	2.01	0.61
1:B:543:HIS:HE1	1:B:783:SER:OG	1.83	0.61
1:B:657:VAL:HG13	1:B:688:PHE:HZ	1.66	0.60
1:A:125:ALA:HB3	1:A:126:PRO:HD3	1.83	0.60
1:A:306:ASN:ND2	1:A:312:THR:H	1.99	0.60
1:B:617:PRO:O	1:B:620:ALA:HB2	2.01	0.60
1:B:525:MET:HE2	1:B:844:ARG:NH1	2.17	0.59
1:B:711:ARG:NE	1:B:711:ARG:H	2.01	0.59
1:A:323:PRO:HG2	1:A:361:VAL:HG12	1.85	0.58
1:B:672:PRO:HB3	1:B:862:ARG:HD3	1.84	0.58
1:A:180:LYS:HD3	1:A:180:LYS:H	1.69	0.58
3:A:1010:HOH:O	1:B:557:THR:HG22	2.04	0.58
1:A:232:SER:HB3	1:A:238:THR:HG22	1.85	0.57
1:A:7:LEU:HD13	1:A:130:LEU:HD21	1.86	0.57
1:A:108:LEU:HD21	1:A:133:GLY:HA3	1.86	0.57
1:A:249:SER:O	1:A:253:GLU:HG3	2.05	0.57
1:B:567:GLU:H	1:B:567:GLU:CD	2.13	0.55
1:B:668:SER:HB2	1:B:677:LEU:HD22	1.87	0.55
1:B:680:LYS:O	1:B:684:GLU:HG3	2.06	0.55
1:A:25:MET:HE2	1:A:344:ARG:HH11	1.70	0.55
1:A:309:GLY:HA3	1:A:389:VAL:HG11	1.89	0.54
1:A:157:VAL:HG13	1:A:188:PHE:HZ	1.71	0.54
1:A:167:ASN:ND2	1:A:171:ASN:H	2.05	0.54
1:A:193:ASP:HA	1:A:224:ARG:HH21	1.69	0.54
1:B:667:ASN:ND2	1:B:671:ASN:H	2.05	0.54
1:A:308:MET:O	1:A:389:VAL:HG11	2.07	0.53
1:B:711:ARG:NE	1:B:711:ARG:N	2.55	0.53
1:A:117:PRO:O	1:A:120:ALA:HB2	2.09	0.53
1:A:283:SER:O	1:A:287:VAL:HG23	2.08	0.53
1:B:519:PHE:C	1:B:519:PHE:CD2	2.87	0.53
1:A:193:ASP:HA	1:A:224:ARG:HH22	1.72	0.53
1:B:521:ILE:O	1:B:525:MET:SD	2.68	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:680:LYS:HE3	3:B:1144:HOH:O	2.08	0.52
1:B:648:LEU:HD21	1:B:653:LEU:HD11	1.91	0.52
1:B:711:ARG:N	1:B:711:ARG:HE	2.08	0.52
1:A:19:PHE:C	1:A:19:PHE:CD2	2.88	0.51
1:B:806:ASN:ND2	1:B:812:THR:H	2.08	0.51
1:A:309:GLY:HA3	1:A:389:VAL:CG1	2.41	0.51
1:A:118:THR:HA	1:A:119:PRO:C	2.36	0.51
1:B:859:GLY:HA2	3:B:1450:HOH:O	2.09	0.51
1:B:806:ASN:HD21	1:B:812:THR:H	1.57	0.51
1:B:525:MET:HB3	1:B:844:ARG:NH2	2.26	0.50
1:A:16:ARG:NH1	1:B:760:MET:SD	2.85	0.50
1:B:785:LYS:O	1:B:789:GLU:HG3	2.12	0.49
1:A:148:LEU:HD11	1:A:153:LEU:HD13	1.95	0.49
1:A:280:ASP:OD2	1:A:282:ARG:HD3	2.12	0.49
1:B:868:ALA:HB3	1:B:871:LYS:HG3	1.95	0.49
1:A:343:ALA:O	1:A:378:ARG:HD3	2.13	0.48
1:A:27:ASP:HB2	3:A:1224:HOH:O	2.14	0.48
1:B:749:SER:O	1:B:753:GLU:HG3	2.13	0.48
1:B:667:ASN:HD21	1:B:671:ASN:H	1.61	0.48
1:B:618:THR:HA	1:B:619:PRO:C	2.38	0.48
1:A:117:PRO:HA	1:A:138:VAL:O	2.13	0.48
1:A:169:PRO:HB3	1:A:317:GLY:HA3	1.95	0.48
1:B:657:VAL:HG13	1:B:688:PHE:CZ	2.47	0.48
1:B:528:VAL:HA	1:B:844:ARG:O	2.13	0.47
1:A:40:THR:OG1	1:B:557:THR:HG21	2.14	0.47
1:B:889:VAL:HG23	1:B:889:VAL:OXT	2.14	0.47
1:B:693:ASP:HA	1:B:724:ARG:NH2	2.29	0.47
1:A:167:ASN:HD21	1:A:171:ASN:H	1.62	0.47
1:B:543:HIS:HD2	3:B:1124:HOH:O	1.95	0.47
1:A:43:HIS:HD2	3:A:1189:HOH:O	1.98	0.47
1:A:172:PRO:HB3	1:A:362:ARG:HD3	1.97	0.46
1:A:248:PRO:HD2	1:A:251:ILE:HD12	1.96	0.46
1:B:678:THR:HA	1:B:711:ARG:HH22	1.81	0.46
1:A:134:LYS:HE3	1:A:134:LYS:HB2	1.69	0.46
1:B:711:ARG:H	1:B:711:ARG:HE	1.62	0.46
1:A:171:ASN:HB3	2:A:413:PLP:H2A1	1.98	0.46
1:B:752:ILE:O	1:B:756:VAL:HG13	2.15	0.46
1:A:188:PHE:O	1:A:192:HIS:HD2	1.97	0.46
1:A:344:ARG:NE	3:A:1224:HOH:O	2.48	0.46
1:A:229:ASN:HB3	1:A:244:PHE:CZ	2.51	0.45
1:B:750:TRP:O	1:B:754:ARG:NH1	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:843:ALA:O	1:B:878:ARG:HD3	2.17	0.45
1:B:699:ASP:OD1	2:B:913:PLP:N1	2.50	0.45
1:B:730:GLY:HA3	3:B:1270:HOH:O	2.17	0.44
1:A:117:PRO:HD2	1:A:165:ILE:O	2.17	0.44
1:A:168:SER:HA	1:A:169:PRO:C	2.43	0.44
1:A:13:SER:OG	1:B:760:MET:HE1	2.17	0.44
1:B:789:GLU:HA	1:B:792:LYS:CE	2.48	0.44
1:B:619:PRO:HG2	1:B:822:PHE:CE1	2.53	0.43
1:A:199:ASP:OD1	2:A:413:PLP:N1	2.50	0.43
1:B:672:PRO:CB	1:B:862:ARG:HD3	2.47	0.43
1:B:678:THR:OG1	1:B:680:LYS:HG2	2.18	0.43
1:A:28:VAL:HG22	1:A:344:ARG:HD2	2.01	0.43
1:A:223:GLU:H	1:A:223:GLU:CD	2.26	0.43
1:B:839:MET:HE1	1:B:882:VAL:HG22	2.01	0.43
1:B:682:LEU:HD12	1:B:713:TYR:HB2	2.01	0.43
1:B:666:ILE:HD13	1:B:666:ILE:HA	1.72	0.43
1:A:160:LYS:HA	1:A:160:LYS:HD2	1.91	0.43
1:B:732:SER:HB3	1:B:738:THR:HG22	2.00	0.42
1:A:323:PRO:HD2	1:A:361:VAL:O	2.19	0.42
1:B:608:LEU:HD21	1:B:633:GLY:HA3	2.02	0.42
1:A:324:ARG:HD2	1:A:326:ARG:CZ	2.49	0.42
1:A:150:VAL:O	1:A:154:LYS:HG3	2.18	0.42
1:A:96:ALA:HB3	2:A:413:PLP:H5A2	2.01	0.42
1:B:753:GLU:HB2	1:B:754:ARG:NH1	2.34	0.42
1:B:833:LYS:O	1:B:837:GLU:HG3	2.20	0.42
1:A:32:GLY:HA2	3:A:1342:HOH:O	2.20	0.42
1:B:593:LEU:HD22	1:B:602:MET:HE2	2.02	0.42
1:B:705:PHE:HA	1:B:791:ARG:HG3	2.01	0.42
1:B:843:ALA:HB2	1:B:882:VAL:HG11	2.02	0.42
1:A:186:ALA:O	1:A:190:VAL:HG23	2.20	0.42
1:A:242:LEU:HD22	1:A:270:ILE:HG22	2.01	0.42
1:B:789:GLU:HA	1:B:792:LYS:HE2	2.01	0.42
1:A:210:ALA:HB2	1:A:316:LYS:HB3	2.01	0.41
1:B:727:THR:O	1:B:745:VAL:HA	2.20	0.41
1:B:785:LYS:C	1:B:785:LYS:HD2	2.46	0.41
1:B:615:LEU:HD11	1:B:657:VAL:HG13	2.02	0.41
1:B:513:SER:O	1:B:516:ARG:HB3	2.21	0.41
1:B:525:MET:HE1	1:B:840:LEU:HG	2.02	0.41
1:A:8:GLU:HG2	1:A:8:GLU:H	1.69	0.41
1:A:203:GLU:HG3	1:A:204:HIS:N	2.35	0.41
1:A:333:LYS:O	1:A:337:GLU:HG3	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:49:LYS:HG2	1:B:549:LYS:HG2	2.02	0.41
1:A:98:GLN:HB3	1:A:102:MET:HE2	2.02	0.40
3:A:1010:HOH:O	1:B:557:THR:CG2	2.67	0.40
1:B:825:ILE:C	1:B:825:ILE:HD12	2.46	0.40
1:B:710:ALA:HB2	1:B:816:LYS:HB3	2.03	0.40
1:A:89:GLU:HA	1:A:252:ILE:HG13	2.03	0.40
1:A:179:LYS:HB3	1:A:179:LYS:HE2	1.80	0.40
1:B:521:ILE:O	1:B:521:ILE:HG22	2.20	0.40
1:B:551:ALA:HB3	1:B:769:PHE:HB2	2.03	0.40
1:B:648:LEU:CD2	1:B:666:ILE:HD11	2.52	0.40
1:B:767:VAL:HB	1:B:770:ILE:HB	2.04	0.40
1:B:869:TYR:O	1:B:873:GLU:HG3	2.22	0.40
1:B:872:LEU:O	1:B:876:MET:HG2	2.21	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	386/389 (99%)	378 (98%)	7 (2%)	1 (0%)	36	25
1	B	386/389 (99%)	375 (97%)	10 (3%)	1 (0%)	36	25
All	All	772/778 (99%)	753 (98%)	17 (2%)	2 (0%)	36	25

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	27	ASP
1	B	527	ASP

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	326/327 (100%)	306 (94%)	20 (6%)	17	6
1	B	326/327 (100%)	303 (93%)	23 (7%)	13	4
All	All	652/654 (100%)	609 (93%)	43 (7%)	15	5

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

Mol	Chain	Res	Type
1	A	7	LEU
1	A	8	GLU
1	A	13	SER
1	A	31	LEU
1	A	74	GLU
1	A	101	LEU
1	A	153	LEU
1	A	157	VAL
1	A	164	LEU
1	A	172	PRO
1	A	177	LEU
1	A	180	LYS
1	A	193	ASP
1	A	256	VAL
1	A	281	GLU
1	A	326	ARG
1	A	341	LYS
1	A	361	VAL
1	A	374	GLU
1	A	383	LEU
1	B	513	SER
1	B	539	ASP
1	B	557	THR
1	B	567	GLU
1	B	601	LEU
1	B	657	VAL
1	B	659	ASP

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Mol	Chain	Res	Type
1	B	664	LEU
1	B	666	ILE
1	B	672	PRO
1	B	677	LEU
1	B	693	ASP
1	B	706	ILE
1	B	711	ARG
1	B	754	ARG
1	B	756	VAL
1	B	779	LYS
1	B	833	LYS
1	B	848	VAL
1	B	861	VAL
1	B	874	GLU
1	B	882	VAL
1	B	887	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	43	HIS
1	A	97	ASN
1	A	167	ASN
1	A	192	HIS
1	A	306	ASN
1	B	543	HIS
1	B	597	ASN
1	B	667	ASN
1	B	806	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](#)

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	413	1	15,15,16	1.64	3 (20%)	21,22,23	2.33	6 (28%)
2	PLP	B	913	1	15,15,16	1.63	3 (20%)	21,22,23	2.39	7 (33%)

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	413	1	-	2/6/6/8	0/1/1/1
2	PLP	B	913	1	-	2/6/6/8	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	413	PLP	C3-C2	-3.83	1.37	1.41
2	B	913	PLP	C5-C4	3.31	1.44	1.40
2	B	913	PLP	C3-C2	-3.31	1.37	1.41
2	A	413	PLP	C5-C4	3.02	1.43	1.40
2	B	913	PLP	C2A-C2	2.29	1.54	1.50
2	A	413	PLP	C2A-C2	2.16	1.53	1.50

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	413	PLP	O4P-C5A-C5	7.33	123.10	109.36
2	B	913	PLP	O4P-C5A-C5	7.27	122.98	109.36
2	B	913	PLP	C5A-C5-C6	-3.86	113.06	119.36
2	B	913	PLP	O2P-P-O4P	-3.72	96.97	106.67
2	A	413	PLP	O2P-P-O4P	-3.59	97.31	106.67
2	A	413	PLP	C5A-C5-C6	-3.54	113.59	119.36
2	B	913	PLP	C4A-C4-C5	2.81	123.84	120.94
2	A	413	PLP	C5A-C5-C4	2.62	127.80	122.64
2	B	913	PLP	C5A-C5-C4	2.61	127.80	122.64
2	A	413	PLP	C4A-C4-C5	2.54	123.56	120.94
2	B	913	PLP	C6-N1-C2	2.14	123.08	119.20
2	A	413	PLP	C6-N1-C2	2.09	122.99	119.20
2	B	913	PLP	C5-C6-N1	-2.02	120.54	123.83

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	413	PLP	C4-C5-C5A-O4P
2	A	413	PLP	C6-C5-C5A-O4P
2	B	913	PLP	C4-C5-C5A-O4P
2	B	913	PLP	C6-C5-C5A-O4P

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	413	PLP	3	0
2	B	913	PLP	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

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	388/389 (99%)	-0.41	12 (3%) 51 51	4, 9, 25, 44	0
1	B	388/389 (99%)	-0.38	14 (3%) 46 46	3, 9, 24, 44	0
All	All	776/778 (99%)	-0.39	26 (3%) 48 48	3, 9, 25, 44	0

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	521	ILE	5.6
1	A	26	LYS	5.4
1	B	527	ASP	4.7
1	A	19	PHE	4.4
1	B	523	ALA	4.1
1	B	889	VAL	4.0
1	A	21	ILE	3.8
1	B	524	GLY	3.7
1	B	526	LYS	3.5
1	B	519	PHE	3.5
1	B	711	ARG	3.1
1	A	24	GLY	3.0
1	B	522	ALA	3.0
1	A	23	ALA	2.7
1	A	16	ARG	2.6
1	B	513	SER	2.6
1	A	159	ASP	2.4
1	B	516	ARG	2.4
1	B	666	ILE	2.2
1	B	517	LYS	2.2
1	A	39	ASP	2.2
1	A	284	TRP	2.2
1	B	525	MET	2.1
1	A	74	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	22	ALA	2.1
1	A	25	MET	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	PLP	B	913	15/16	0.94	0.08	8,13,15,15	0
2	PLP	A	413	15/16	0.95	0.07	5,14,16,17	0

6.5 Other polymers [i](#)

There are no such residues in this entry.