



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 03:14 AM UTC

PDB ID : 1GCK / pdb_00001gck
Title : THERMUS THERMOPHILUS ASPARTATE AMINOTRANSFERASE
DOUBLE MUTANT 1 COMPLEXED WITH ASPARTATE
Authors : Ura, H.; Nakai, T.; Hirotsu, K.; Kuramitsu, S.
Deposited on : 2000-08-04
Resolution : 2.50 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

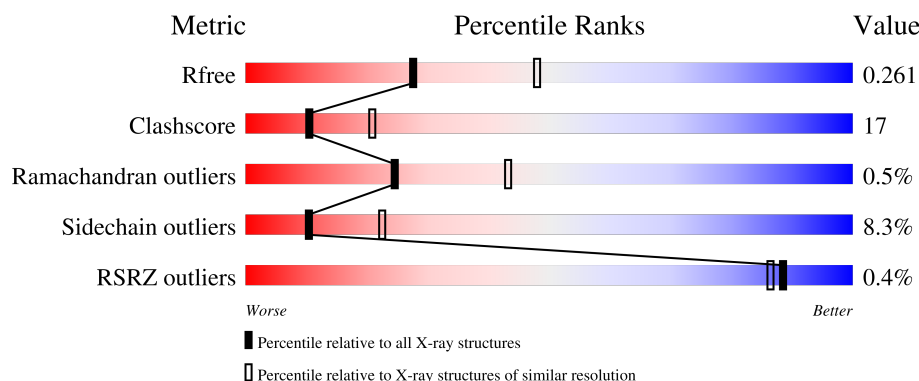
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	385	 60% 32% 6% ..
1	B	385	 57% 34% 6% ..

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	ASP	B	914	-	-	X	-
3	PLP	A	413	-	X	-	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 5960 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.

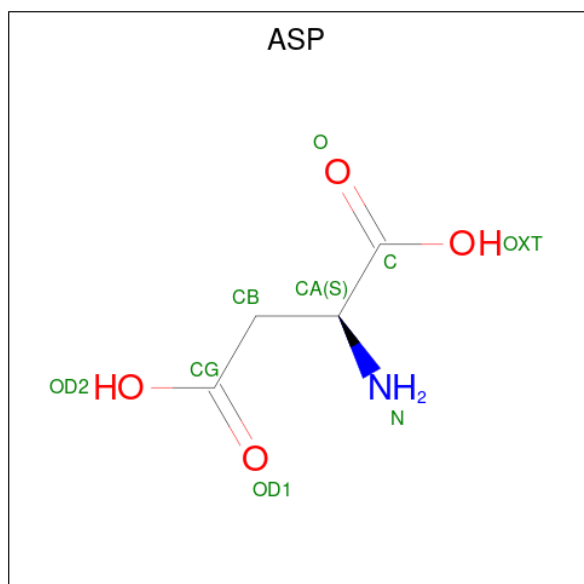
- Molecule 1 is a protein called ASPARTATE AMINOTRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	382	Total	C	N	O	S	0	0	0
			2946	1864	523	551	8			
1	B	382	Total	C	N	O	S	0	0	0
			2946	1864	523	551	8			

There are 4 discrepancies between the modelled and reference sequences:

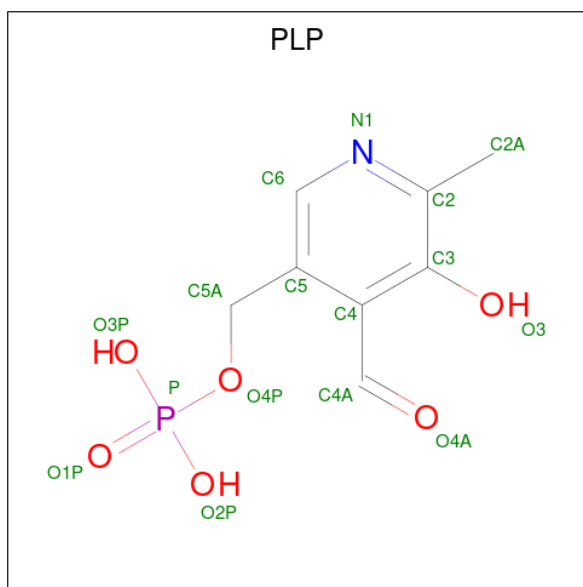
Chain	Residue	Modelled	Actual	Comment	Reference
A	101	SER	LYS	engineered mutation	UNP Q56232
A	261	ARG	SER	engineered mutation	UNP Q56232
B	601	SER	LYS	engineered mutation	UNP Q56232
B	761	ARG	SER	engineered mutation	UNP Q56232

- Molecule 2 is ASPARTIC ACID (CCD ID: ASP) (formula: $C_4H_7NO_4$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			9	4	1	4		
2	B	1	Total	C	N	O	0	0
			9	4	1	4		

- Molecule 3 is PYRIDOXAL-5'-PHOSPHATE (CCD ID: PLP) (formula: $C_8H_{10}NO_6P$).



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

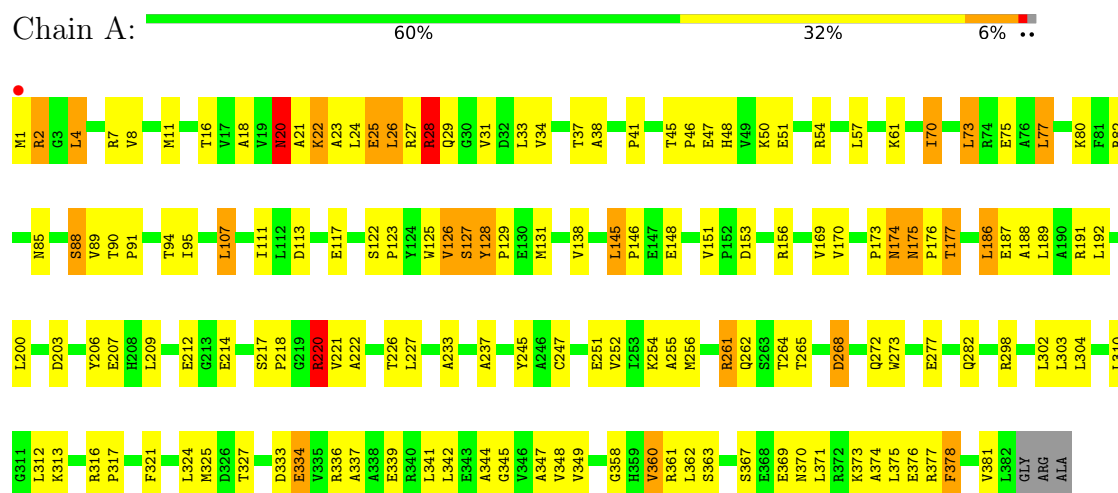
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	9	Total	O	0	0
			9	9		
4	B	11	Total	O	0	0
			11	11		

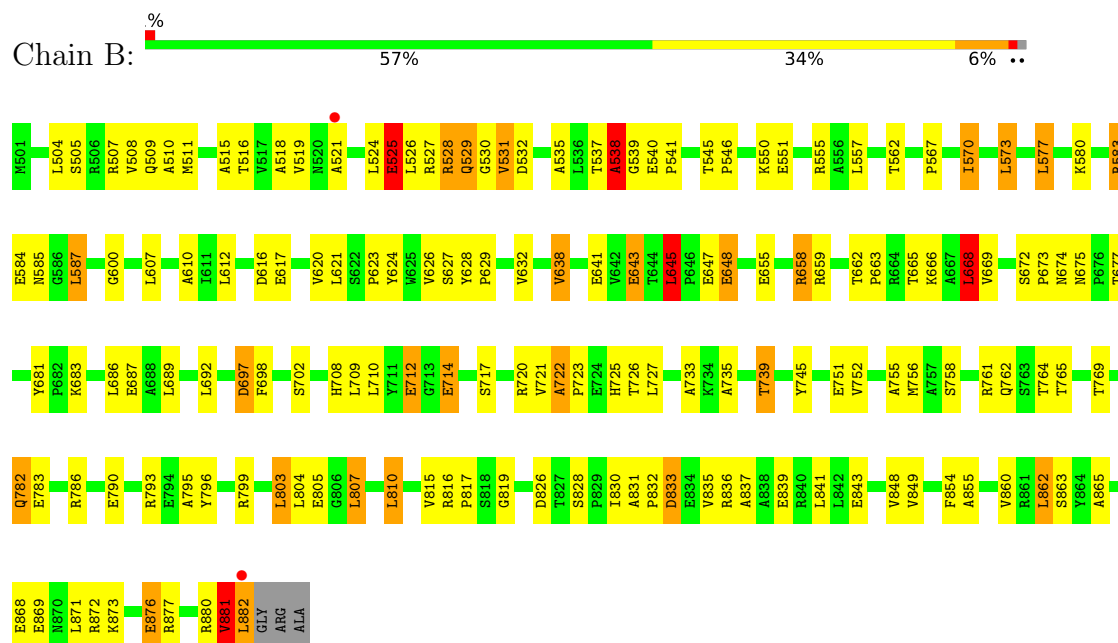
3 Residue-property plots

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	62.75Å 74.89Å 163.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.50 8.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	85.3 (8.00-2.50) 86.7 (8.00-2.50)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.10 (at 2.51Å)	Xtriage
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.205 , 0.284 0.193 , 0.261	Depositor DCC
R_{free} test set	2436 reflections (9.90%)	wwPDB-VP
Wilson B-factor (Å ²)	34.0	Xtriage
Anisotropy	0.338	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 61.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5960	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.45% 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	0.93	14/3007 (0.5%)	1.25	37/4089 (0.9%)
1	B	0.87	6/3007 (0.2%)	1.31	47/4089 (1.1%)
All	All	0.90	20/6014 (0.3%)	1.28	84/8178 (1.0%)

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
1	B	0	1
All	All	0	3

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	21	ALA	N-CA	13.58	1.63	1.46
1	B	529	GLN	N-CA	-7.98	1.36	1.46
1	A	214	GLU	N-CA	-7.64	1.36	1.46
1	A	22	LYS	N-CA	7.57	1.55	1.46
1	A	363	SER	N-CA	-6.81	1.38	1.46
1	A	38	ALA	C-O	-6.32	1.16	1.24
1	A	28	ARG	CZ-NH1	6.32	1.41	1.32
1	A	88	SER	C-O	-6.05	1.16	1.23
1	B	756	MET	SD-CE	-5.93	1.64	1.79
1	A	21	ALA	CA-C	5.91	1.60	1.52
1	A	214	GLU	CG-CD	-5.60	1.38	1.52
1	A	28	ARG	CZ-NH2	5.51	1.40	1.33
1	B	529	GLN	CA-CB	-5.44	1.45	1.53
1	B	848	VAL	N-CA	5.39	1.52	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	233	ALA	N-CA	-5.34	1.40	1.46
1	A	28	ARG	NE-CZ	5.30	1.38	1.33
1	A	256	MET	SD-CE	-5.11	1.66	1.79
1	B	714	GLU	N-CA	-5.07	1.39	1.46
1	A	20	ASN	N-CA	-5.04	1.40	1.46
1	B	538	ALA	N-CA	5.04	1.52	1.46

All (84) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	174	ASN	N-CA-C	9.64	124.74	110.48
1	B	529	GLN	N-CA-CB	-9.42	95.85	110.92
1	B	816	ARG	CA-C-N	9.33	129.28	119.85
1	B	816	ARG	C-N-CA	9.33	129.28	119.85
1	A	38	ALA	N-CA-C	9.29	122.52	109.15
1	B	722	ALA	CA-C-N	8.96	128.70	119.56
1	B	722	ALA	C-N-CA	8.96	128.70	119.56
1	A	212	GLU	N-CA-CB	7.97	123.97	110.49
1	A	20	ASN	N-CA-CB	7.96	122.56	110.70
1	A	316	ARG	CA-C-N	7.76	127.52	119.76
1	A	316	ARG	C-N-CA	7.76	127.52	119.76
1	B	674	ASN	N-CA-C	7.63	122.45	110.17
1	B	532	ASP	CA-C-N	-7.53	112.60	123.00
1	B	532	ASP	C-N-CA	-7.53	112.60	123.00
1	B	782	GLN	N-CA-C	7.51	120.42	111.33
1	B	570	ILE	CA-C-N	7.51	127.88	119.32
1	B	570	ILE	C-N-CA	7.51	127.88	119.32
1	B	521	ALA	CA-C-N	7.44	130.58	120.38
1	B	521	ALA	C-N-CA	7.44	130.58	120.38
1	A	28	ARG	NE-CZ-NH2	-7.24	112.69	119.20
1	B	733	ALA	N-CA-C	7.21	118.78	111.07
1	B	726	THR	N-CA-C	7.04	120.72	109.24
1	B	507	ARG	N-CA-CB	-6.95	99.16	110.14
1	A	113	ASP	N-CA-C	-6.95	100.47	110.40
1	A	226	THR	N-CA-C	6.93	120.23	109.07
1	A	145	LEU	CA-C-N	6.91	126.72	119.05
1	A	145	LEU	C-N-CA	6.91	126.72	119.05
1	A	145	LEU	N-CA-C	6.84	119.75	110.31
1	B	675	ASN	N-CA-C	-6.81	98.51	109.40
1	A	207	GLU	N-CA-C	6.76	119.22	111.11
1	A	25	GLU	N-CA-C	-6.74	105.06	113.28
1	B	510	ALA	CA-C-N	-6.62	113.87	123.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	510	ALA	C-N-CA	-6.62	113.87	123.00
1	A	378	PHE	N-CA-C	-6.60	103.78	110.97
1	B	645	LEU	N-CA-C	6.55	118.90	110.39
1	A	174	ASN	N-CA-CB	-6.45	100.49	110.29
1	B	511	MET	N-CA-CB	6.33	120.51	110.57
1	A	47	GLU	N-CA-C	6.28	118.93	111.71
1	A	2	ARG	N-CA-CB	6.24	122.08	111.66
1	B	677	THR	N-CA-C	6.17	117.68	111.07
1	A	28	ARG	CG-CD-NE	-5.92	98.97	112.00
1	B	881	VAL	N-CA-C	-5.79	105.48	113.00
1	B	865	ALA	N-CA-C	-5.79	103.36	110.65
1	A	25	GLU	N-CA-CB	5.78	119.54	110.46
1	B	627	SER	N-CA-C	5.71	119.43	112.23
1	A	26	LEU	N-CA-C	-5.68	105.46	112.90
1	B	645	LEU	CA-C-N	5.67	125.51	119.28
1	B	645	LEU	C-N-CA	5.67	125.51	119.28
1	A	28	ARG	NH1-CZ-NH2	5.66	126.65	119.30
1	B	643	GLU	N-CA-C	5.64	118.46	110.50
1	A	175	ASN	N-CA-C	-5.64	100.37	109.40
1	B	769	THR	N-CA-C	5.63	118.18	111.71
1	B	849	VAL	N-CA-C	5.62	113.44	107.76
1	A	169	VAL	N-CA-C	5.58	115.63	107.37
1	A	220	ARG	N-CA-CB	-5.57	101.14	110.39
1	A	23	ALA	N-CA-C	5.52	116.98	111.07
1	B	681	TYR	CA-C-N	5.48	125.39	119.85
1	B	681	TYR	C-N-CA	5.48	125.39	119.85
1	B	807	LEU	N-CA-C	-5.44	105.26	111.14
1	A	177	THR	N-CA-C	5.42	118.11	111.82
1	B	674	ASN	N-CA-CB	-5.41	101.82	110.63
1	B	528	ARG	N-CA-C	-5.36	106.72	113.20
1	B	833	ASP	N-CA-CB	-5.33	101.56	110.57
1	B	735	ALA	N-CA-C	5.33	116.77	111.07
1	B	855	ALA	N-CA-C	-5.32	103.89	111.24
1	A	20	ASN	N-CA-C	5.30	118.77	112.72
1	B	712	GLU	N-CA-CB	5.29	120.12	111.08
1	B	600	GLY	N-CA-C	-5.28	107.77	114.16
1	B	876	GLU	N-CA-CB	-5.28	102.42	110.07
1	A	70	ILE	CA-C-N	5.28	125.34	119.32
1	A	70	ILE	C-N-CA	5.28	125.34	119.32
1	A	268	ASP	N-CA-CB	-5.27	100.86	109.56
1	B	668	LEU	CA-C-N	-5.25	116.30	123.12
1	B	668	LEU	C-N-CA	-5.25	116.30	123.12

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	882	LEU	N-CA-CB	-5.25	101.58	110.50
1	B	648	GLU	N-CA-C	-5.18	106.35	113.30
1	B	708	HIS	N-CA-C	5.08	119.46	113.16
1	A	127	SER	N-CA-C	5.07	119.17	112.89
1	A	374	ALA	N-CA-C	-5.03	105.51	111.69
1	A	222	ALA	CA-C-N	5.03	125.30	119.47
1	A	222	ALA	C-N-CA	5.03	125.30	119.47
1	B	702	SER	N-CA-C	5.02	116.81	108.02
1	A	126	VAL	N-CA-C	5.02	119.78	109.34
1	A	128	TYR	N-CA-CB	5.01	117.32	110.11

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	20	ASN	Mainchain
1	A	28	ARG	Sidechain
1	B	525	GLU	Mainchain

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	2946	0	2964	114	2
1	B	2946	0	2961	105	2
2	A	9	0	3	1	0
2	B	9	0	3	5	0
3	A	15	0	6	1	0
3	B	15	0	6	0	0
4	A	9	0	0	1	0
4	B	11	0	0	0	0
All	All	5960	0	5943	201	2

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 17.

All (201) 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:188:ALA:HA	1:A:191:ARG:NH1	1.54	1.19
1:A:188:ALA:HA	1:A:191:ARG:HH12	1.07	0.96
1:B:727:LEU:HD11	1:B:752:VAL:HG11	1.57	0.86
1:B:526:LEU:HG	1:B:531:VAL:HG21	1.57	0.86
1:A:1:MET:HB3	1:B:697:ASP:O	1.77	0.84
1:A:77:LEU:HB3	1:A:94:THR:HG21	1.58	0.84
1:B:830:ILE:HD13	1:B:841:LEU:HD21	1.61	0.83
1:A:173:PRO:HG2	1:A:206:TYR:O	1.80	0.80
1:A:45:THR:HG22	1:B:562:THR:HG21	1.64	0.80
1:A:371:LEU:O	1:A:375:LEU:HG	1.81	0.79
1:A:188:ALA:CA	1:A:191:ARG:HH12	1.92	0.78
1:A:261:ARG:NH1	1:A:265:THR:HG23	2.02	0.74
1:A:227:LEU:HD21	1:A:252:VAL:HG11	1.71	0.73
1:A:2:ARG:HH11	1:B:725:HIS:CD2	2.06	0.73
1:A:37:THR:HG23	1:A:347:ALA:O	1.87	0.73
1:B:583:ARG:HG2	1:B:583:ARG:HH11	1.54	0.72
1:B:583:ARG:HG2	1:B:583:ARG:NH1	2.06	0.71
1:B:539:GLY:HA3	2:B:914:ASP:HA	1.74	0.68
1:B:837:ALA:O	1:B:841:LEU:HG	1.95	0.66
1:A:34:VAL:HG11	1:A:373:LYS:HD2	1.79	0.65
1:B:873:LYS:O	1:B:876:GLU:HB3	1.95	0.65
1:A:2:ARG:NH1	1:B:725:HIS:CD2	2.65	0.65
1:B:717:SER:H	1:B:720:ARG:HH12	1.45	0.64
1:B:868:GLU:O	1:B:872:ARG:HG3	1.98	0.64
1:B:668:LEU:HD12	1:B:669:VAL:N	2.13	0.63
1:A:34:VAL:CG1	1:A:373:LYS:HD2	2.28	0.63
1:B:881:VAL:O	1:B:882:LEU:HB2	1.99	0.62
1:B:835:VAL:O	1:B:839:GLU:HG3	1.99	0.62
1:B:655:GLU:OE1	1:B:658:ARG:NH1	2.33	0.62
1:B:529:GLN:O	1:B:529:GLN:HG2	1.99	0.62
1:A:2:ARG:NH1	1:B:725:HIS:NE2	2.48	0.61
1:B:786:ARG:O	1:B:790:GLU:HG2	1.99	0.61
1:A:373:LYS:O	1:A:377:ARG:HG3	2.01	0.60
1:B:515:ALA:O	1:B:519:VAL:HG23	2.01	0.60
1:A:304:LEU:HD11	1:A:317:PRO:HD2	1.83	0.60
1:A:117:GLU:HG2	1:A:138:VAL:CG2	2.32	0.59
1:A:261:ARG:HD2	1:A:261:ARG:C	2.27	0.59
1:B:526:LEU:O	1:B:531:VAL:HG22	2.02	0.58
1:B:764:THR:O	1:B:765:THR:HB	2.03	0.58
1:A:227:LEU:HD21	1:A:252:VAL:CG1	2.33	0.57
1:A:7:ARG:NH2	1:B:612:LEU:O	2.38	0.57
1:A:128:TYR:HB2	1:A:129:PRO:HD3	1.86	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:24:LEU:HA	1:A:27:ARG:HB2	1.86	0.57
1:A:303:LEU:HD23	1:A:375:LEU:HD21	1.87	0.57
1:B:831:ALA:HB1	1:B:832:PRO:HD2	1.84	0.57
1:A:174:ASN:HD22	1:A:177:THR:H	1.52	0.57
1:A:61:LYS:HD3	1:A:273:TRP:CE2	2.40	0.57
1:A:77:LEU:CB	1:A:94:THR:HG21	2.32	0.57
1:A:88:SER:HB2	4:A:1019:HOH:O	2.04	0.57
1:B:525:GLU:HG3	1:B:526:LEU:N	2.12	0.57
1:B:717:SER:H	1:B:720:ARG:NH1	2.04	0.56
1:A:122:SER:HA	1:A:123:PRO:C	2.30	0.56
1:A:261:ARG:NH2	2:B:914:ASP:OD1	2.39	0.56
1:B:869:GLU:OE1	1:B:872:ARG:NH1	2.39	0.56
1:A:261:ARG:NH1	1:A:265:THR:CG2	2.69	0.56
1:B:683:LYS:O	1:B:687:GLU:HB2	2.06	0.55
1:A:107:LEU:HD21	1:A:227:LEU:HD22	1.88	0.55
1:A:261:ARG:HD2	1:A:261:ARG:O	2.06	0.55
1:A:312:LEU:HD13	1:A:325:MET:SD	2.46	0.55
1:A:261:ARG:HH22	2:B:914:ASP:CG	2.14	0.55
1:B:836:ARG:NH1	1:B:839:GLU:OE1	2.39	0.55
1:B:795:ALA:O	1:B:799:ARG:HG3	2.06	0.55
1:A:186:LEU:HG	1:A:218:PRO:HG3	1.89	0.55
1:A:203:ASP:OD1	3:A:413:PLP:N1	2.40	0.55
1:A:378:PHE:O	1:A:381:VAL:HG22	2.07	0.55
1:B:628:TYR:HB2	1:B:629:PRO:HD3	1.89	0.54
1:A:29:GLN:HB3	1:A:31:VAL:HG23	1.89	0.54
1:A:191:ARG:HH11	1:A:191:ARG:HB3	1.73	0.54
1:A:334:GLU:HG3	1:A:358:GLY:H	1.73	0.53
1:A:45:THR:CG2	1:B:562:THR:HG21	2.36	0.53
1:A:22:LYS:NZ	1:A:339:GLU:OE2	2.40	0.53
1:A:191:ARG:HH11	1:A:191:ARG:CB	2.21	0.53
1:A:51:GLU:OE1	1:A:54:ARG:NH1	2.42	0.53
1:A:261:ARG:C	1:A:261:ARG:CD	2.82	0.53
1:B:839:GLU:O	1:B:843:GLU:HG3	2.08	0.53
1:A:188:ALA:HA	1:A:191:ARG:HH11	1.61	0.53
1:A:367:SER:OG	1:A:370:ASN:ND2	2.41	0.53
1:B:583:ARG:NH2	1:B:782:GLN:CD	2.67	0.53
1:B:524:LEU:HB3	1:B:528:ARG:NH1	2.24	0.53
1:A:77:LEU:HD21	1:A:245:TYR:CB	2.39	0.52
1:A:4:LEU:HD12	1:B:610:ALA:O	2.10	0.52
1:B:525:GLU:O	1:B:529:GLN:HB2	2.09	0.52
1:A:261:ARG:NH2	2:B:914:ASP:CG	2.68	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:877:ARG:O	1:B:881:VAL:HG12	2.09	0.52
1:B:645:LEU:HB2	1:B:648:GLU:HG3	1.91	0.51
1:A:251:GLU:HG2	1:B:504:LEU:HD11	1.92	0.51
1:B:525:GLU:CG	1:B:526:LEU:N	2.72	0.51
1:A:345:GLY:HA3	1:A:377:ARG:HH11	1.75	0.51
1:B:545:THR:HG23	1:B:546:PRO:HD2	1.93	0.51
1:B:860:VAL:HG13	1:B:862:LEU:CD1	2.40	0.51
1:B:860:VAL:HG13	1:B:862:LEU:HD13	1.93	0.51
1:B:803:LEU:O	1:B:807:LEU:HG	2.11	0.50
1:A:170:VAL:HG11	1:A:189:LEU:HD13	1.94	0.50
1:B:877:ARG:HA	1:B:880:ARG:HD2	1.93	0.50
1:A:16:THR:O	1:A:20:ASN:OD1	2.29	0.50
1:A:344:ALA:HB2	1:A:381:VAL:CG1	2.42	0.50
1:B:551:GLU:O	1:B:555:ARG:HG3	2.12	0.50
1:B:868:GLU:HA	1:B:871:LEU:HD12	1.94	0.50
1:B:573:LEU:HD22	1:B:577:LEU:HD22	1.94	0.50
1:A:341:LEU:HD12	1:A:348:VAL:HG21	1.94	0.50
1:B:607:LEU:HG	1:B:727:LEU:HD13	1.94	0.50
1:B:628:TYR:O	1:B:632:VAL:HG23	2.12	0.50
1:A:46:PRO:HB2	1:A:48:HIS:CE1	2.47	0.50
1:A:26:LEU:O	1:A:29:GLN:HB3	2.12	0.49
1:A:41:PRO:HG2	1:A:237:ALA:HA	1.93	0.49
1:A:127:SER:O	1:A:131:MET:HG3	2.13	0.49
1:B:672:SER:HA	1:B:673:PRO:C	2.37	0.49
1:B:545:THR:CG2	1:B:546:PRO:HD2	2.42	0.49
1:A:348:VAL:HG12	1:A:362:LEU:CD2	2.43	0.49
1:A:48:HIS:O	1:A:277:GLU:HG2	2.13	0.49
1:B:527:ARG:HA	1:B:530:GLY:O	2.13	0.48
1:B:505:SER:O	1:B:509:GLN:HG3	2.12	0.48
1:B:580:LYS:HZ1	1:B:585:ASN:HD21	1.61	0.48
1:A:16:THR:CG2	1:A:349:VAL:HG13	2.44	0.48
1:B:541:PRO:HB3	1:B:796:TYR:OH	2.14	0.48
1:A:8:VAL:HG11	1:B:755:ALA:HB1	1.96	0.48
1:A:90:THR:HB	1:A:91:PRO:HD2	1.96	0.47
1:A:125:TRP:NE1	2:A:414:ASP:OXT	2.47	0.47
1:A:2:ARG:HH12	1:B:725:HIS:CE1	2.32	0.47
1:B:537:THR:O	1:B:538:ALA:HB3	2.15	0.47
1:B:621:LEU:O	1:B:624:TYR:HB3	2.15	0.47
1:B:616:ASP:OD1	1:B:666:LYS:NZ	2.47	0.47
1:B:655:GLU:O	1:B:659:ARG:HG3	2.14	0.47
1:B:617:GLU:HG2	1:B:638:VAL:CG2	2.45	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:709:LEU:O	1:B:819:GLY:HA2	2.15	0.47
1:A:191:ARG:NH1	1:A:191:ARG:CB	2.78	0.46
1:B:505:SER:O	1:B:508:VAL:HG12	2.15	0.46
1:B:815:VAL:O	1:B:817:PRO:HD3	2.15	0.46
1:A:321:PHE:HD1	1:A:321:PHE:O	1.98	0.46
1:A:264:THR:O	1:A:265:THR:HB	2.16	0.46
1:A:334:GLU:HG3	1:A:358:GLY:N	2.31	0.46
1:B:540:GLU:HB3	1:B:739:THR:HG21	1.98	0.46
1:B:567:PRO:HA	1:B:765:THR:O	2.16	0.46
1:B:580:LYS:NZ	1:B:585:ASN:HD21	2.14	0.46
1:B:804:LEU:HD23	1:B:807:LEU:HD12	1.98	0.45
1:A:18:ALA:O	1:A:22:LYS:HG3	2.16	0.45
1:A:77:LEU:HD12	1:A:77:LEU:HA	1.77	0.45
1:A:333:ASP:OD2	1:A:336:ARG:NE	2.49	0.45
1:B:515:ALA:HA	1:B:518:ALA:HB3	1.97	0.45
1:B:658:ARG:HH11	1:B:658:ARG:HB3	1.81	0.45
1:B:617:GLU:HG2	1:B:638:VAL:HG22	1.98	0.45
1:A:80:LYS:NZ	1:A:85:ASN:HD21	2.14	0.45
1:B:710:LEU:HD21	1:B:714:GLU:HA	1.99	0.45
1:A:57:LEU:HD22	1:B:550:LYS:HG2	1.99	0.45
1:A:217:SER:O	1:A:220:ARG:HB3	2.17	0.45
1:A:187:GLU:HG2	1:A:221:VAL:HG21	1.98	0.44
1:B:722:ALA:N	1:B:723:PRO:HD3	2.32	0.44
1:B:583:ARG:HH11	1:B:583:ARG:CG	2.23	0.44
1:A:2:ARG:NH1	1:B:725:HIS:CE1	2.86	0.44
1:A:94:THR:HG22	1:A:95:ILE:N	2.32	0.44
1:A:8:VAL:HG21	1:B:610:ALA:HA	1.98	0.44
1:A:11:MET:HE1	1:B:758:SER:HB3	1.99	0.43
1:A:117:GLU:HG2	1:A:138:VAL:HG23	1.99	0.43
1:B:527:ARG:HH12	1:B:535:ALA:HB2	1.84	0.43
1:A:26:LEU:HD12	1:A:342:LEU:HD22	2.01	0.43
1:B:862:LEU:HD13	1:B:862:LEU:N	2.34	0.43
1:B:620:VAL:O	1:B:641:GLU:HA	2.18	0.43
1:A:89:VAL:HG11	1:A:247:CYS:HB2	2.00	0.43
1:B:516:THR:HG23	1:B:626:VAL:HG21	2.00	0.43
1:A:261:ARG:HH11	1:A:265:THR:HG23	1.79	0.43
1:A:324:LEU:HD23	1:A:324:LEU:HA	1.87	0.42
1:A:334:GLU:CG	1:A:358:GLY:H	2.31	0.42
1:B:557:LEU:HD23	1:B:557:LEU:HA	1.80	0.42
1:B:662:THR:HB	1:B:663:PRO:CD	2.49	0.42
1:B:687:GLU:HG3	1:B:721:VAL:HG21	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:16:THR:HG21	1:A:349:VAL:HG13	2.02	0.42
1:A:50:LYS:O	1:A:54:ARG:HG3	2.19	0.42
1:A:1:MET:O	1:B:698:PHE:HA	2.20	0.42
1:B:862:LEU:CD1	1:B:862:LEU:N	2.82	0.42
1:A:70:ILE:HG13	1:A:73:LEU:HB2	2.01	0.42
1:B:587:LEU:HD11	1:B:717:SER:HB3	2.02	0.42
1:B:623:PRO:HA	1:B:854:PHE:O	2.20	0.42
1:A:255:ALA:HB1	1:B:508:VAL:HG21	2.02	0.42
1:B:539:GLY:HA3	2:B:914:ASP:CA	2.44	0.42
1:B:583:ARG:HG3	1:B:584:GLU:N	2.29	0.42
1:B:880:ARG:O	1:B:882:LEU:HG	2.19	0.42
1:A:191:ARG:NH1	1:A:191:ARG:HB2	2.34	0.41
1:A:261:ARG:HA	1:A:265:THR:HA	2.01	0.41
1:B:537:THR:O	1:B:863:SER:HB3	2.21	0.41
1:B:810:LEU:HD12	1:B:810:LEU:HA	1.93	0.41
1:A:261:ARG:O	1:A:261:ARG:CD	2.68	0.41
1:B:570:ILE:HG13	1:B:573:LEU:HB2	2.02	0.41
1:A:1:MET:HA	1:B:665:THR:O	2.20	0.41
1:A:33:LEU:HD12	1:A:33:LEU:O	2.20	0.41
1:A:107:LEU:O	1:A:111:ILE:HG12	2.19	0.41
1:A:95:ILE:O	1:A:245:TYR:HA	2.20	0.41
1:A:327:THR:HG21	1:A:360:VAL:HB	2.02	0.41
1:A:298:ARG:HG2	1:A:298:ARG:HH11	1.84	0.41
1:A:145:LEU:HA	1:A:146:PRO:HD3	1.91	0.41
1:A:77:LEU:HD21	1:A:245:TYR:HB3	2.02	0.41
1:A:148:GLU:O	1:A:151:VAL:HG12	2.21	0.40
1:A:175:ASN:OD1	1:A:361:ARG:NH1	2.51	0.40
1:A:268:ASP:O	1:A:272:GLN:HG3	2.22	0.40
1:A:173:PRO:HG2	1:A:206:TYR:C	2.45	0.40
1:B:826:ASP:OD1	1:B:828:SER:HB2	2.21	0.40
1:B:881:VAL:O	1:B:882:LEU:CB	2.67	0.40
1:A:153:ASP:OD1	1:A:153:ASP:C	2.65	0.40
1:A:176:PRO:HB3	1:A:361:ARG:HD3	2.02	0.40
1:A:337:ALA:O	1:A:341:LEU:HG	2.21	0.40
1:A:367:SER:O	1:A:371:LEU:HG	2.22	0.40
1:B:790:GLU:OE1	1:B:793:ARG:NH1	2.54	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:28:ARG:NH2	1:B:529:GLN:NE2[1_655]	2.02	0.18
1:A:28:ARG:NH1	1:B:529:GLN:NE2[1_655]	2.10	0.10

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	380/385 (99%)	358 (94%)	21 (6%)	1 (0%)	36	55
1	B	380/385 (99%)	351 (92%)	26 (7%)	3 (1%)	16	31
All	All	760/770 (99%)	709 (93%)	47 (6%)	4 (0%)	24	43

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	531	VAL
1	B	697	ASP
1	B	538	ALA
1	A	126	VAL

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	305/306 (100%)	281 (92%)	24 (8%)	11	24
1	B	305/306 (100%)	278 (91%)	27 (9%)	9	20
All	All	610/612 (100%)	559 (92%)	51 (8%)	10	22

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

Mol	Chain	Res	Type
1	A	4	LEU
1	A	25	GLU
1	A	73	LEU
1	A	75	GLU
1	A	77	LEU
1	A	82	ARG
1	A	107	LEU
1	A	156	ARG
1	A	186	LEU
1	A	192	LEU
1	A	200	LEU
1	A	209	LEU
1	A	220	ARG
1	A	254	LYS
1	A	261	ARG
1	A	262	GLN
1	A	282	GLN
1	A	302	LEU
1	A	310	LEU
1	A	313	LYS
1	A	334	GLU
1	A	360	VAL
1	A	369	GLU
1	A	376	GLU
1	B	525	GLU
1	B	573	LEU
1	B	577	LEU
1	B	583	ARG
1	B	587	LEU
1	B	638	VAL
1	B	643	GLU
1	B	645	LEU
1	B	647	GLU
1	B	658	ARG
1	B	668	LEU
1	B	686	LEU
1	B	689	LEU
1	B	692	LEU
1	B	712	GLU
1	B	739	THR
1	B	745	TYR
1	B	751	GLU

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Mol	Chain	Res	Type
1	B	761	ARG
1	B	762	GLN
1	B	783	GLU
1	B	803	LEU
1	B	805	GLU
1	B	810	LEU
1	B	833	ASP
1	B	862	LEU
1	B	881	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	20	ASN
1	A	85	ASN
1	A	174	ASN
1	A	370	ASN
1	B	529	GLN
1	B	585	ASN
1	B	606	ASN

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 [i](#)

4 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
3	PLP	A	413	2	15,15,16	2.20	7 (46%)	21,22,23	4.27	8 (38%)
3	PLP	B	913	2	15,15,16	1.56	4 (26%)	21,22,23	1.74	6 (28%)
2	ASP	B	914	3	7,8,8	0.93	0	6,10,10	1.27	1 (16%)
2	ASP	A	414	3	7,8,8	1.13	0	6,10,10	0.97	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
3	PLP	A	413	2	-	3/6/6/8	0/1/1/1
3	PLP	B	913	2	-	3/6/6/8	0/1/1/1
2	ASP	B	914	3	-	2/8/8/8	-
2	ASP	A	414	3	-	3/8/8/8	-

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	413	PLP	P-O3P	-3.56	1.41	1.54
3	A	413	PLP	C4A-C4	3.38	1.58	1.51
3	A	413	PLP	C3-C4	-3.28	1.33	1.40
3	A	413	PLP	O3-C3	-3.11	1.29	1.36
3	B	913	PLP	C4A-C4	2.76	1.57	1.51
3	A	413	PLP	P-O1P	-2.73	1.42	1.50
3	B	913	PLP	C5-C4	2.66	1.43	1.40
3	A	413	PLP	P-O2P	-2.45	1.45	1.54
3	B	913	PLP	P-O3P	-2.31	1.46	1.54
3	A	413	PLP	C5-C4	2.11	1.42	1.40
3	B	913	PLP	C2-N1	2.03	1.37	1.33

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	413	PLP	C4A-C4-C5	-15.38	105.08	120.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	413	PLP	C4A-C4-C3	8.09	134.01	120.52
3	A	413	PLP	O2P-P-O1P	-4.40	93.69	110.83
3	A	413	PLP	O3P-P-O1P	4.22	127.26	110.83
3	B	913	PLP	O2P-P-O4P	-3.75	96.90	106.67
3	A	413	PLP	O3P-P-O4P	3.29	115.26	106.67
3	B	913	PLP	C6-C5-C4	3.14	120.67	118.10
3	A	413	PLP	C6-C5-C4	2.97	120.53	118.10
3	B	913	PLP	C5A-C5-C6	-2.80	114.80	119.36
3	B	913	PLP	O3P-P-O1P	2.71	121.41	110.83
3	B	913	PLP	C4A-C4-C5	-2.45	118.41	120.94
3	B	913	PLP	C5-C6-N1	-2.33	120.03	123.83
3	A	413	PLP	C5-C6-N1	-2.24	120.19	123.83
2	B	914	ASP	OD1-CG-CB	-2.22	116.04	122.84
3	A	413	PLP	O4P-C5A-C5	2.04	113.18	109.36

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	413	PLP	C5A-O4P-P-O2P
3	A	413	PLP	C5A-O4P-P-O3P
3	B	913	PLP	C5A-O4P-P-O1P
3	B	913	PLP	C5A-O4P-P-O2P
3	B	913	PLP	C5A-O4P-P-O3P
2	B	914	ASP	CA-CB-CG-OD2
3	A	413	PLP	C6-C5-C5A-O4P
2	A	414	ASP	CA-CB-CG-OD1
2	A	414	ASP	CA-CB-CG-OD2
2	B	914	ASP	CA-CB-CG-OD1
2	A	414	ASP	C-CA-CB-CG

There are no ring outliers.

3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	413	PLP	1	0
2	B	914	ASP	5	0
2	A	414	ASP	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	382/385 (99%)	-0.46	1 (0%) 90 87	6, 19, 47, 76	0
1	B	382/385 (99%)	-0.48	2 (0%) 87 85	5, 17, 46, 70	0
All	All	764/770 (99%)	-0.47	3 (0%) 88 86	5, 18, 48, 76	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	521	ALA	2.4
1	B	882	LEU	2.3
1	A	1	MET	2.1

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(Å ²)	Q<0.9
2	ASP	A	414	9/9	0.85	0.14	47,53,56,58	0
3	PLP	A	413	15/16	0.87	0.13	28,34,40,41	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	ASP	B	914	9/9	0.88	0.12	26,31,37,39	0
3	PLP	B	913	15/16	0.98	0.05	5,18,20,22	0

6.5 Other polymers [i](#)

There are no such residues in this entry.