



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 04:48 PM GMT

PDB ID : 1GCK
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 validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

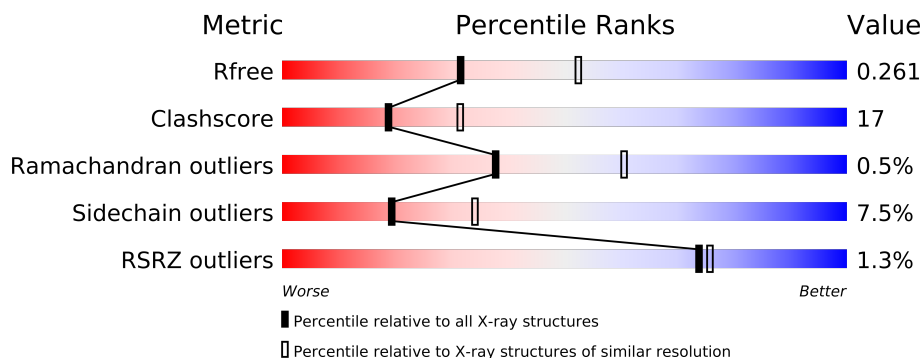
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

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}	66092	2784 (2.50-2.50)
Clashscore	79885	3562 (2.50-2.50)
Ramachandran outliers	78287	3480 (2.50-2.50)
Sidechain outliers	78261	3482 (2.50-2.50)
RSRZ outliers	66119	2785 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	385	
1	B	385	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	ASP	A	414	-	X
3	PLP	A	413	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5960 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

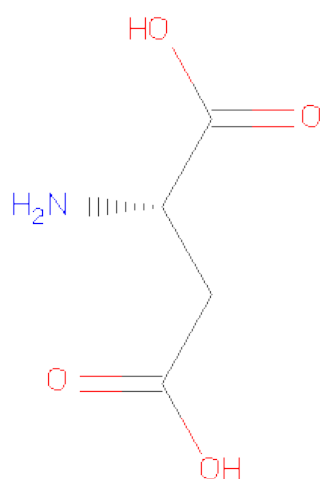
- Molecule 1 is a protein called ASPARTATE AMINOTRANSFERASE.

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:

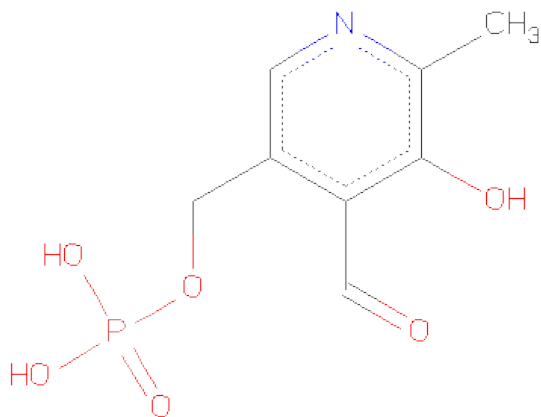
Chain	Residue	Modelled	Actual	Comment	Reference
A	101	SER	LYS	ENGINEERED	UNP Q56232
A	261	ARG	SER	ENGINEERED	UNP Q56232
B	601	SER	LYS	ENGINEERED	UNP Q56232
B	761	ARG	SER	ENGINEERED	UNP Q56232

- Molecule 2 is ASPARTIC ACID (three-letter code: ASP) (formula: C₄H₇NO₄).



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 (three-letter code: PLP) (formula: C₈H₁₀NO₆P).



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		

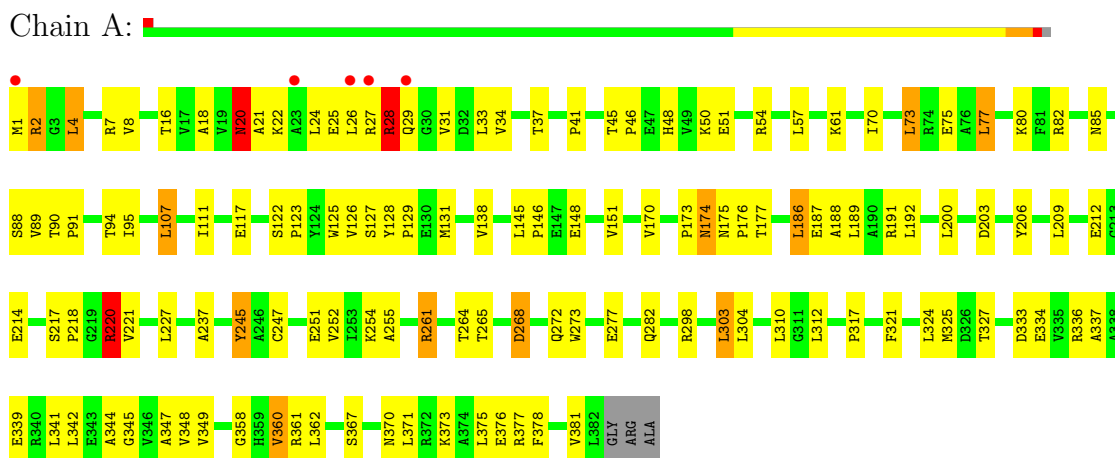
- Molecule 4 is water.

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

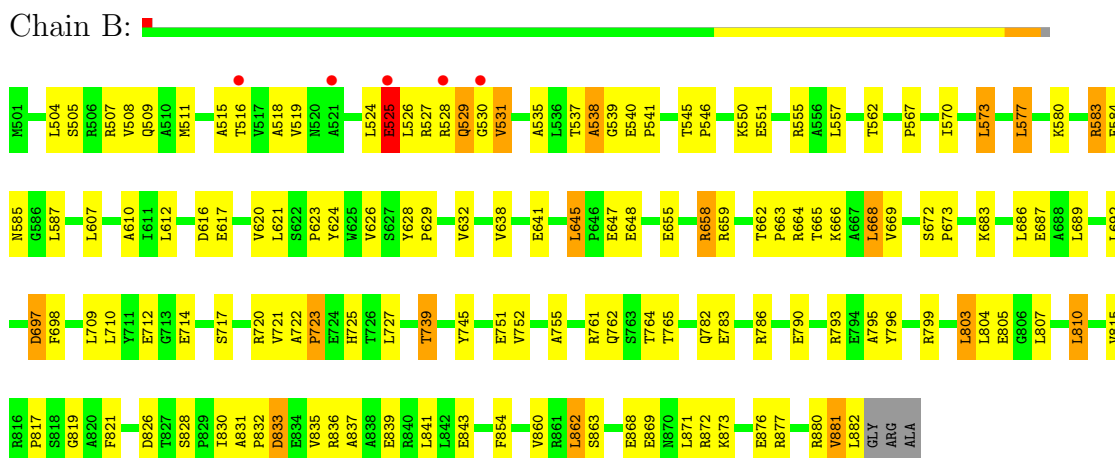
3 Residue-property plots

These plots are drawn for all protein, RNA 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: 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 46.15 – 2.50	Depositor EDS
% Data completeness (in resolution range)	85.3 (8.00-2.50) 89.6 (46.15-2.50)	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	2361 reflections (9.91%)	DCC
Wilson B-factor (Å ²)	34.0	Xtriage
Anisotropy	0.338	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 26.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 24610 reflections	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.

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	0.66	4/3007 (0.1%)	0.89	12/4089 (0.3%)
1	B	0.61	0/3007	0.88	5/4089 (0.1%)
All	All	0.64	4/6014 (0.1%)	0.88	17/8178 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	B	0	1
All	All	0	3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	214	GLU	CG-CD	-9.27	1.38	1.51
1	A	21	ALA	N-CA	8.62	1.63	1.46
1	A	28	ARG	CZ-NH1	6.58	1.41	1.33
1	A	28	ARG	CZ-NH2	5.81	1.40	1.33

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	28	ARG	NE-CZ-NH2	-15.22	112.69	120.30
1	B	529	GLN	N-CA-CB	-8.20	95.85	110.60
1	A	212	GLU	N-CA-CB	7.43	123.97	110.60
1	A	20	ASN	N-CA-CB	6.64	122.56	110.60
1	A	28	ARG	NH1-CZ-NH2	6.59	126.65	119.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	2	ARG	N-CA-CB	6.38	122.08	110.60
1	B	507	ARG	N-CA-CB	-6.36	99.16	110.60
1	A	28	ARG	CG-CD-NE	-6.11	98.97	111.80
1	A	28	ARG	NE-CZ-NH1	-5.79	117.40	120.30
1	A	174	ASN	N-CA-CB	-5.62	100.49	110.60
1	B	511	MET	N-CA-CB	5.50	120.51	110.60
1	A	268	ASP	N-CA-CB	-5.41	100.86	110.60
1	B	712	GLU	N-CA-CB	5.29	120.12	110.60
1	A	220	ARG	N-CA-CB	-5.25	101.14	110.60
1	A	174	ASN	N-CA-C	5.09	124.74	111.00
1	A	20	ASN	CA-C-N	-5.07	106.05	117.20
1	B	833	ASP	N-CA-CB	-5.02	101.56	110.60

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 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2946	0	2964	111	2
1	B	2946	0	2961	103	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	197	2

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 17.

All (197) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(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:NH1	1:B:583:ARG:HG2	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:525:GLU:HG3	1:B:526:LEU:N	2.12	0.63
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:B:764:THR:O	1:B:765:THR:HB	2.03	0.59
1:B:526:LEU:O	1:B:531:VAL:HG22	2.02	0.58

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:227:LEU:HD21	1:A:252:VAL:CG1	2.33	0.57
1:A:128:TYR:HB2	1:A:129:PRO:HD3	1.86	0.57
1:A:7:ARG:NH2	1:B:612:LEU:O	2.38	0.57
1:A:303:LEU:HD23	1:A:375:LEU:HD21	1.87	0.57
1:A:24:LEU:HA	1:A:27:ARG:HB2	1.86	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:77:LEU:CB	1:A:94:THR:HG21	2.32	0.57
1:A:61:LYS:HD3	1:A:273:TRP:CE2	2.40	0.57
1:A:88:SER:HB2	4:A:1019:HOH:O	2.04	0.57
1:B:717:SER:H	1:B:720:ARG:NH1	2.04	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: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:261:ARG:HD2	1:A:261:ARG:C	2.27	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:191:ARG:HH11	1:A:191:ARG:CB	2.21	0.53
1:A:22:LYS:NZ	1:A:339:GLU:OE2	2.40	0.53
1:B:839:GLU:O	1:B:843:GLU:HG3	2.08	0.53
1:A:51:GLU:OE1	1:A:54:ARG:NH1	2.42	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: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:122:SER:HA	1:A:123:PRO:C	2.30	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: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

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
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:860:VAL:HG13	1:B:862:LEU:CD1	2.40	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:HD13	1.93	0.51
1:B:803:LEU:O	1:B:807:LEU:HG	2.11	0.50
1:A:261:ARG:HH22	2:B:914:ASP:CG	2.14	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:B:868:GLU:HA	1:B:871:LEU:HD12	1.94	0.50
1:A:344:ALA:HB2	1:A:381:VAL:CG1	2.42	0.50
1:A:16:THR:O	1:A:20:ASN:OD1	2.29	0.50
1:B:551:GLU:O	1:B:555:ARG:HG3	2.12	0.50
1:B:573:LEU:HD22	1:B:577:LEU:HD22	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:341:LEU:HD12	1:A:348:VAL:HG21	1.94	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: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:A:261:ARG:C	1:A:261:ARG:CD	2.82	0.48
1:B:505:SER:O	1:B:509:GLN:HG3	2.12	0.48
1:B:583:ARG:NH2	1:B:782:GLN:CD	2.67	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:261:ARG:NH2	2:B:914:ASP:CG	2.68	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:655:GLU:O	1:B:659:ARG:HG3	2.14	0.47
1:B:616:ASP:OD1	1:B:666:LYS:NZ	2.47	0.47
1:B:709:LEU:O	1:B:819:GLY:HA2	2.15	0.47
1:B:617:GLU:HG2	1:B:638:VAL:CG2	2.45	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
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:HA	1:A:77:LEU:HD12	1.77	0.45
1:B:658:ARG:HB3	1:B:658:ARG:HH11	1.81	0.45
1:B:515:ALA:HA	1:B:518:ALA:HB3	1.97	0.45
1:A:333:ASP:OD2	1:A:336:ARG:NE	2.49	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:B:672:SER:HA	1:B:673:PRO:C	2.37	0.45
1:A:217:SER:O	1:A:220:ARG:HB3	2.17	0.45
1:A:57:LEU:HD22	1:B:550:LYS:HG2	1.99	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:A:94:THR:HG22	1:A:95:ILE:N	2.32	0.44
1:B:583:ARG:CG	1:B:583:ARG:HH11	2.23	0.44
1:A:2:ARG:NH1	1:B:725:HIS:CE1	2.86	0.44
1:A:8:VAL:HG21	1:B:610:ALA:HA	1.98	0.44
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:N	1:B:862:LEU:HD13	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:HA	1:A:324:LEU:HD23	1.87	0.42
1:B:687:GLU:HG3	1:B:721:VAL:HG21	2.01	0.42
1:A:334:GLU:CG	1:A:358:GLY:H	2.31	0.42
1:B:557:LEU:HA	1:B:557:LEU:HD23	1.80	0.42
1:B:662:THR:HB	1:B:663:PRO:CD	2.49	0.42
1:A:50:LYS:O	1:A:54:ARG:HG3	2.19	0.42
1:A:16:THR:HG21	1:A:349:VAL:HG13	2.02	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:1:MET:O	1:B:698:PHE:HA	2.20	0.42
1:B:862:LEU:N	1:B:862:LEU:CD1	2.82	0.42
1:B:587:LEU:HD11	1:B:717:SER:HB3	2.02	0.42
1:A:70:ILE:HG13	1:A:73:LEU:HB2	2.01	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:HA	1:B:810:LEU:HD12	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:107:LEU:O	1:A:111:ILE:HG12	2.19	0.41
1:A:33:LEU:HD12	1:A:33:LEU:O	2.20	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:B:881:VAL:O	1:B:882:LEU:CB	2.67	0.40
1:B:826:ASP:OD1	1:B:828:SER:HB2	2.21	0.40
1:A:337:ALA:O	1:A:341:LEU:HG	2.21	0.40
1:A:176:PRO:HB3	1:A:361:ARG:HD3	2.02	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	Distance(Å)	Clash(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

5.3.1 Protein backbone ⓘ

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%)	50	73
1	B	380/385 (99%)	351 (92%)	26 (7%)	3 (1%)	27	46
All	All	760/770 (99%)	709 (93%)	47 (6%)	4 (0%)	38	60

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 ⓘ

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%)	285 (93%)	20 (7%)	24	41
1	B	305/306 (100%)	279 (92%)	26 (8%)	15	28
All	All	610/612 (100%)	564 (92%)	46 (8%)	19	34

All (46) 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

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Mol	Chain	Res	Type
1	A	82	ARG
1	A	107	LEU
1	A	186	LEU
1	A	192	LEU
1	A	200	LEU
1	A	209	LEU
1	A	220	ARG
1	A	245	TYR
1	A	254	LYS
1	A	261	ARG
1	A	282	GLN
1	A	303	LEU
1	A	310	LEU
1	A	360	VAL
1	A	376	GLU
1	B	525	GLU
1	B	573	LEU
1	B	577	LEU
1	B	583	ARG
1	B	645	LEU
1	B	647	GLU
1	B	658	ARG
1	B	664	ARG
1	B	668	LEU
1	B	686	LEU
1	B	689	LEU
1	B	692	LEU
1	B	723	PRO
1	B	739	THR
1	B	745	TYR
1	B	751	GLU
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	821	PHE
1	B	833	ASP
1	B	862	LEU
1	B	881	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) 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	585	ASN
1	B	725	HIS

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

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	14,15,16	1.98	6 (42%)	20,22,23	2.12	6 (30%)
2	ASP	A	414	3	8,8,8	1.11	0	10,10,10	2.10	1 (10%)
3	PLP	B	913	2	14,15,16	1.37	3 (21%)	20,22,23	1.54	4 (20%)
2	ASP	B	914	3	8,8,8	0.72	0	10,10,10	2.32	1 (10%)

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	-	0/6/6/8	0/1/1/1
2	ASP	A	414	3	-	0/8/8/8	0/0/0/0
3	PLP	B	913	2	-	0/6/6/8	0/1/1/1
2	ASP	B	914	3	-	0/8/8/8	0/0/0/0

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	413	PLP	P-O3P	-3.59	1.41	1.54
3	A	413	PLP	O3-C3	-3.01	1.29	1.37
3	A	413	PLP	P-O1P	-2.75	1.42	1.51
3	B	913	PLP	C4-C5	2.52	1.43	1.39
3	A	413	PLP	P-O2P	-2.46	1.45	1.54
3	A	413	PLP	C4-C3	-2.32	1.33	1.40
3	B	913	PLP	P-O3P	-2.32	1.46	1.54
3	A	413	PLP	C4-C5	2.17	1.42	1.39
3	B	913	PLP	C3-C2	-2.08	1.39	1.40

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	914	ASP	C-CA-N	6.74	120.53	109.36
2	A	414	ASP	C-CA-N	6.12	119.50	109.36
3	A	413	PLP	O3P-P-O1P	5.15	127.26	110.44
3	A	413	PLP	O2P-P-O1P	-5.12	93.69	110.44
3	B	913	PLP	O2P-P-O4P	-3.53	96.90	106.65
3	B	913	PLP	O3P-P-O1P	3.35	121.41	110.44
3	A	413	PLP	O3P-P-O4P	3.12	115.26	106.65
3	B	913	PLP	C5A-C5-C6	-2.37	114.80	119.28
3	B	913	PLP	C5-C6-N1	-2.12	120.03	123.86
3	A	413	PLP	C3-C4-C5	-2.05	117.41	121.29
3	A	413	PLP	C5-C6-N1	-2.03	120.19	123.86
3	A	413	PLP	P-O4P-C5A	2.02	128.52	121.22

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

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	382/385 (99%)	-0.32	5 (1%) 74 76	6, 19, 47, 76	0
1	B	382/385 (99%)	-0.37	5 (1%) 74 76	5, 17, 46, 70	0
All	All	764/770 (99%)	-0.35	10 (1%) 74 76	5, 18, 48, 76	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	528	ARG	3.9
1	A	1	MET	3.2
1	A	29	GLN	2.9
1	A	23	ALA	2.6
1	B	525	GLU	2.6
1	A	27	ARG	2.6
1	B	516	THR	2.4
1	B	530	GLY	2.3
1	B	521	ALA	2.2
1	A	26	LEU	2.1

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	PLP	A	413	15/16	0.22	3.84	28,34,40,41	0
2	ASP	A	414	9/9	0.26	2.99	47,53,56,58	0
2	ASP	B	914	9/9	0.17	0.32	26,31,37,39	0
3	PLP	B	913	15/16	0.11	-0.27	5,18,20,22	0

6.5 Other polymers ⓘ

There are no such residues in this entry.