



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 07:41 PM UTC

PDB ID : 1WRV / pdb_00001wrv
Title : Crystal Structure of T.th.HB8 Branched-Chain Amino Acid Aminotransferase
Authors : Goto, M.; RIKEN Structural Genomics/Proteomics Initiative (RSGI)
Deposited on : 2004-10-27
Resolution : 1.50 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

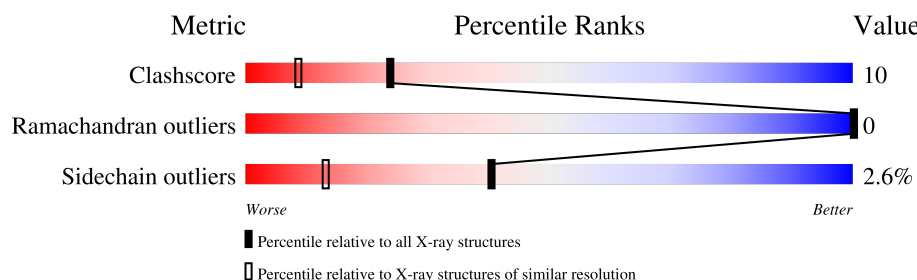
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.50 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))
Clashscore	190562	4235 (1.50-1.50)
Ramachandran outliers	187476	4153 (1.50-1.50)
Sidechain outliers	187428	4150 (1.50-1.50)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	308	
1	B	308	
1	C	308	

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 7387 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.

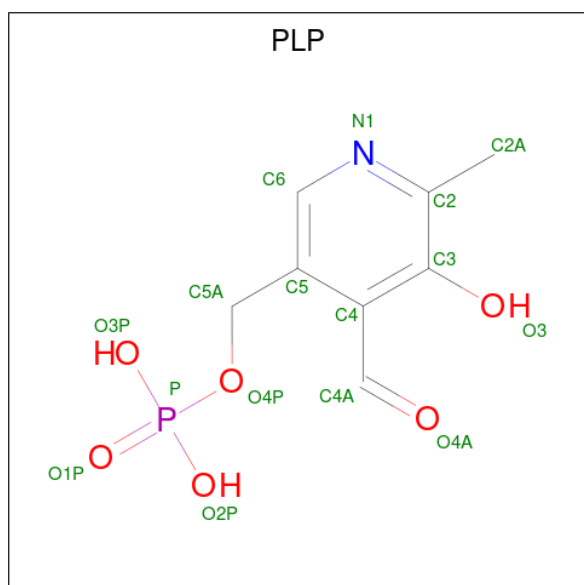
- Molecule 1 is a protein called Branched-Chain Amino Acid Aminotransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	297	Total	C	N	O	S	0	0	0
			2311	1478	401	422	10			
1	B	305	Total	C	N	O	S	0	0	0
			2377	1519	413	435	10			
1	C	295	Total	C	N	O	S	0	0	0
			2299	1470	399	420	10			

- Molecule 2 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

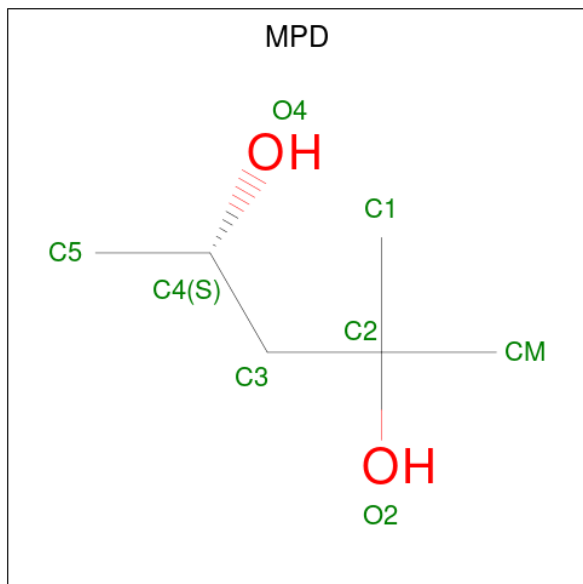
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Cl	0	0
			2	2		
2	B	1	Total	Cl	0	0
			1	1		

- Molecule 3 is PYRIDOXAL-5'-PHOSPHATE (CCD ID: 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		
3	C	1	Total	C	N	O	P	0	0
			15	8	1	5	1		

- Molecule 4 is (4S)-2-METHYL-2,4-PENTANEDIOL (CCD ID: MPD) (formula: C₆H₁₄O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			8	6	2		
4	A	1	Total	C	O	0	0
			8	6	2		
4	B	1	Total	C	O	0	0
			8	6	2		
4	B	1	Total	C	O	0	0
			8	6	2		
4	C	1	Total	C	O	0	0
			8	6	2		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	92	Total	O	0	0
			92	92		
5	B	115	Total	O	0	0
			115	115		

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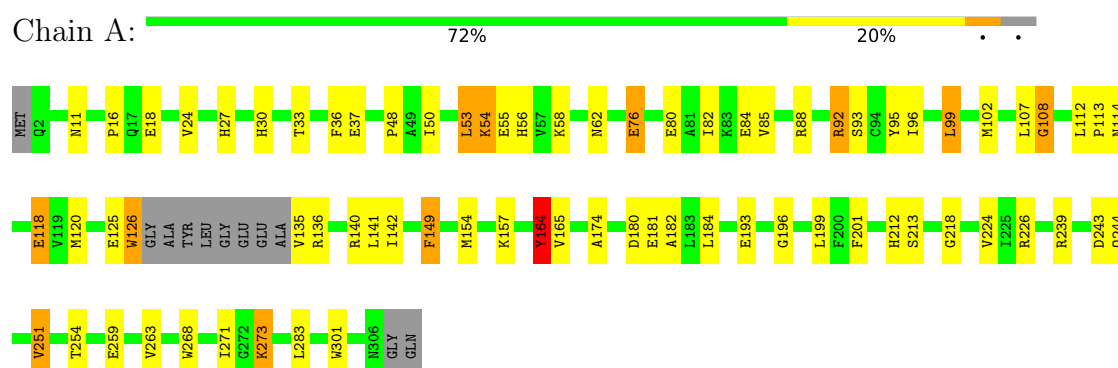
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	105	Total 105	O 105	0	0

3 Residue-property plots

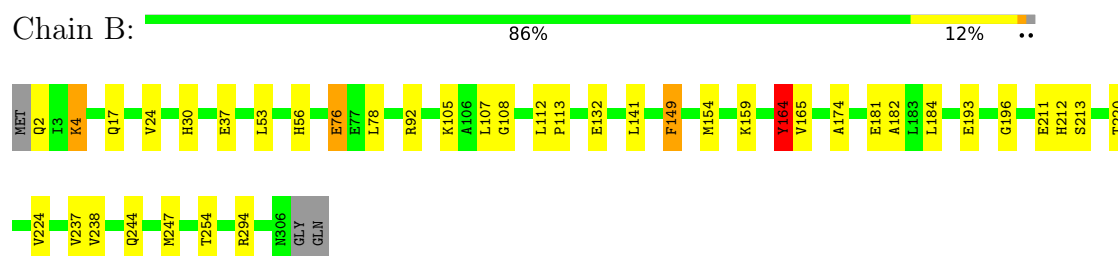
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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.

Note EDS was not executed.

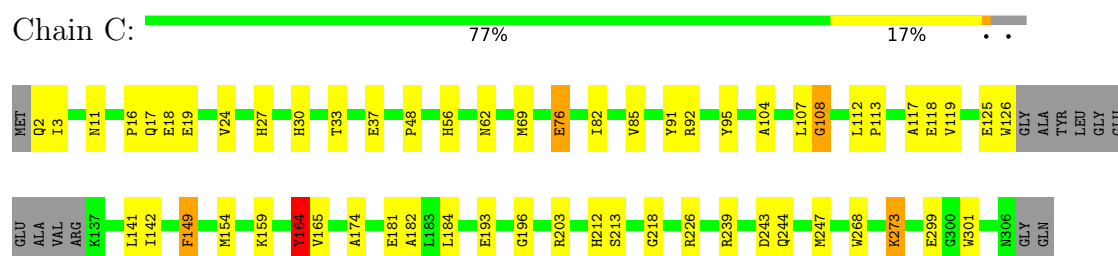
• Molecule 1: Branched-Chain Amino Acid Aminotransferase



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4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	143.41Å 143.41Å 116.22Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	9.99 – 1.50	Depositor
% Data completeness (in resolution range)	88.8 (9.99-1.50)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.205 , 0.218	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7387	wwPDB-VP
Average B, all atoms (Å ²)	13.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PLP, CL, MPD

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	0.37	0/2363	0.93	8/3207 (0.2%)
1	B	0.37	0/2431	0.91	6/3298 (0.2%)
1	C	0.37	0/2351	0.92	9/3190 (0.3%)
All	All	0.37	0/7145	0.92	23/9695 (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	1
1	B	0	1
1	C	0	1
All	All	0	3

There are no bond length outliers.

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	108	GLY	N-CA-C	-9.59	99.78	112.14
1	B	108	GLY	N-CA-C	-9.15	99.56	112.37
1	A	108	GLY	N-CA-C	-9.02	100.51	112.14
1	C	165	VAL	N-CA-C	-7.47	103.18	110.72
1	A	165	VAL	N-CA-C	-7.35	103.30	110.72
1	B	165	VAL	N-CA-C	-7.21	103.44	110.72
1	A	149	PHE	N-CA-C	6.73	118.19	109.64
1	B	149	PHE	N-CA-C	6.66	118.09	109.64
1	C	239	ARG	N-CA-C	-6.50	99.98	110.32
1	C	149	PHE	N-CA-C	6.47	117.86	109.64

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	244	GLN	N-CA-C	-6.03	104.64	112.23
1	C	244	GLN	N-CA-C	-6.02	104.84	111.82
1	A	218	GLY	N-CA-C	5.95	118.31	111.36
1	A	239	ARG	N-CA-C	-5.90	100.93	110.32
1	A	244	GLN	N-CA-C	-5.59	105.33	111.82
1	C	104	ALA	N-CA-C	5.52	118.82	112.97
1	C	164	TYR	N-CA-C	5.35	119.56	113.19
1	C	218	GLY	N-CA-C	5.33	117.60	111.36
1	C	118	GLU	N-CA-C	-5.27	101.94	110.32
1	A	164	TYR	N-CA-C	5.23	119.53	113.20
1	B	164	TYR	N-CA-C	5.19	119.36	113.19
1	B	92	ARG	N-CA-C	-5.15	99.93	108.12
1	A	118	GLU	N-CA-C	-5.08	102.25	110.32

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	164	TYR	Sidechain
1	B	164	TYR	Sidechain
1	C	164	TYR	Sidechain

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	2311	0	2294	67	0
1	B	2377	0	2365	31	0
1	C	2299	0	2283	41	0
2	A	2	0	0	0	0
2	B	1	0	0	0	0
3	A	15	0	7	1	0
3	B	15	0	7	2	0
3	C	15	0	6	2	0
4	A	16	0	28	6	0
4	B	16	0	28	4	0
4	C	8	0	14	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	92	0	0	4	0
5	B	115	0	0	3	0
5	C	105	0	0	2	1
All	All	7387	0	7032	142	1

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 10.

All (142) 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:92:ARG:HE	1:A:92:ARG:H	1.06	0.95
1:C:196:GLY:HA3	4:C:1414:MPD:HM3	1.49	0.92
1:A:196:GLY:HA3	4:A:414:MPD:HM3	1.51	0.91
1:B:220:THR:O	1:B:224:VAL:HG23	1.75	0.87
1:A:92:ARG:H	1:A:92:ARG:NE	1.72	0.87
1:B:196:GLY:HA3	4:B:914:MPD:HM3	1.56	0.86
1:A:251:VAL:HG13	1:A:271:ILE:HG21	1.56	0.86
1:A:251:VAL:HG22	1:A:263:VAL:HB	1.58	0.82
1:B:4:LYS:HD3	1:B:4:LYS:H	1.42	0.82
1:A:84:GLU:HB3	1:A:88:ARG:HH11	1.44	0.81
1:A:92:ARG:HE	1:A:92:ARG:N	1.80	0.79
1:A:84:GLU:HB3	1:A:88:ARG:NH1	2.00	0.76
1:C:142:ILE:HD11	1:C:268:TRP:CZ2	2.22	0.74
1:A:92:ARG:NH2	1:A:125:GLU:HG2	2.03	0.73
1:A:95:TYR:HB2	1:A:126:TRP:CD1	2.23	0.72
1:A:54:LYS:NZ	1:A:55:GLU:HG2	2.05	0.72
1:A:80:GLU:O	1:A:84:GLU:HG3	1.91	0.71
1:B:4:LYS:H	1:B:4:LYS:CD	2.05	0.69
1:C:27:HIS:HD2	1:C:33:THR:OG1	1.76	0.69
1:B:4:LYS:HD3	1:B:4:LYS:N	2.07	0.68
1:B:76:GLU:H	1:B:76:GLU:CD	2.00	0.68
1:C:226:ARG:HD3	1:C:301:TRP:CZ2	2.30	0.67
1:A:58:LYS:HE3	1:C:203:ARG:HH21	1.59	0.66
1:A:27:HIS:HD2	1:A:33:THR:OG1	1.78	0.65
1:C:82:ILE:O	1:C:85:VAL:HG12	1.96	0.65
1:B:224:VAL:HG21	1:B:254:THR:HG21	1.78	0.64
1:A:54:LYS:HZ3	1:A:55:GLU:HG2	1.62	0.64
1:B:212:HIS:HD2	5:B:952:HOH:O	1.81	0.64
1:B:247:MET:HE2	1:C:62:ASN:HB3	1.80	0.63
1:A:82:ILE:O	1:A:85:VAL:HG12	1.99	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:142:ILE:HD11	1:A:268:TRP:CZ2	2.34	0.63
1:B:37:GLU:OE2	1:B:56:HIS:HD2	1.83	0.62
1:C:2:GLN:HE21	1:C:17:GLN:HE21	1.48	0.61
1:C:24:VAL:O	1:C:30:HIS:HE1	1.83	0.61
1:A:24:VAL:O	1:A:30:HIS:HE1	1.85	0.59
1:C:212:HIS:HD2	5:C:1425:HOH:O	1.86	0.59
1:C:16:PRO:HB2	1:C:18:GLU:HG2	1.84	0.59
1:C:48:PRO:HG3	1:C:92:ARG:HA	1.85	0.58
4:A:414:MPD:HM2	4:A:414:MPD:H53	1.86	0.58
1:A:212:HIS:HD2	5:A:1424:HOH:O	1.85	0.57
1:B:2:GLN:HE21	1:B:17:GLN:HE21	1.53	0.57
1:C:76:GLU:H	1:C:76:GLU:CD	2.11	0.57
1:C:37:GLU:OE2	1:C:56:HIS:HD2	1.87	0.56
1:A:27:HIS:HE1	1:A:108:GLY:O	1.88	0.56
1:A:193:GLU:HA	1:A:199:LEU:HD23	1.86	0.56
1:A:37:GLU:OE2	1:A:56:HIS:HD2	1.88	0.55
4:B:914:MPD:H53	4:B:914:MPD:HM2	1.87	0.55
4:B:915:MPD:HM2	4:B:915:MPD:H53	1.88	0.55
1:A:36:PHE:HB3	1:A:99:LEU:HD23	1.88	0.55
1:A:48:PRO:HG3	1:A:92:ARG:HA	1.90	0.54
1:A:58:LYS:HE3	1:C:203:ARG:NH2	2.23	0.54
1:A:50:ILE:HG21	1:A:53:LEU:HG	1.88	0.54
1:A:140:ARG:HG3	1:A:140:ARG:HH11	1.72	0.54
4:C:1414:MPD:HM2	4:C:1414:MPD:H53	1.91	0.53
1:A:58:LYS:CE	1:C:203:ARG:HH21	2.21	0.53
1:B:24:VAL:O	1:B:30:HIS:HE1	1.92	0.53
1:A:92:ARG:HH22	1:A:125:GLU:HG2	1.73	0.53
1:C:91:TYR:CD2	1:C:125:GLU:HG2	2.44	0.52
1:A:141:LEU:CD2	1:A:181:GLU:HB3	2.40	0.52
1:C:27:HIS:HE1	1:C:108:GLY:O	1.93	0.52
1:A:112:LEU:N	1:A:113:PRO:HD2	2.25	0.51
1:C:56:HIS:HE1	5:C:1436:HOH:O	1.93	0.51
1:A:56:HIS:HE1	5:A:1442:HOH:O	1.93	0.51
1:C:159:LYS:NZ	3:C:1413:PLP:O3	2.44	0.51
1:C:141:LEU:CD2	1:C:181:GLU:HB3	2.41	0.51
1:A:62:ASN:HB3	1:C:247:MET:HG2	1.92	0.51
1:A:251:VAL:CG2	1:A:263:VAL:HB	2.37	0.51
1:A:92:ARG:NE	1:A:93:SER:H	2.09	0.50
1:A:135:VAL:HG13	1:A:136:ARG:H	1.77	0.50
1:A:102:MET:HE3	1:A:114:ASN:HB2	1.94	0.50
1:B:56:HIS:HE1	5:B:956:HOH:O	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:82:ILE:HG23	1:A:96:ILE:HG21	1.94	0.50
1:A:135:VAL:HG13	1:A:136:ARG:N	2.27	0.49
1:B:76:GLU:CD	1:B:76:GLU:N	2.69	0.49
1:B:211:GLU:HG2	1:B:237:VAL:HG12	1.95	0.49
4:A:415:MPD:H52	4:A:415:MPD:H11	1.93	0.49
1:C:2:GLN:CG	1:C:3:ILE:H	2.25	0.48
1:A:92:ARG:HH22	1:A:125:GLU:CB	2.27	0.47
1:A:142:ILE:HG23	5:A:1497:HOH:O	2.14	0.47
4:A:415:MPD:HM2	4:A:415:MPD:H53	1.97	0.47
1:B:211:GLU:HG3	1:B:238:VAL:C	2.39	0.47
1:A:140:ARG:NH1	1:A:180:ASP:OD2	2.48	0.47
1:A:76:GLU:H	1:A:76:GLU:CD	2.20	0.47
1:A:273:LYS:HD2	1:A:273:LYS:H	1.81	0.46
1:A:224:VAL:HG13	1:A:283:LEU:HD13	1.98	0.46
1:A:95:TYR:HB2	1:A:126:TRP:NE1	2.31	0.46
1:A:157:LYS:CB	1:C:247:MET:HE1	2.46	0.46
1:A:149:PHE:CG	1:A:154:MET:HB2	2.51	0.45
1:A:201:PHE:HA	1:A:251:VAL:HA	1.98	0.45
1:B:105:LYS:HG3	5:B:1017:HOH:O	2.16	0.45
1:C:2:GLN:HG2	1:C:3:ILE:H	1.80	0.45
1:B:174:ALA:HB2	1:B:182:ALA:HB2	1.98	0.45
1:A:92:ARG:NE	1:A:92:ARG:N	2.49	0.45
1:A:164:TYR:CD1	1:A:164:TYR:N	2.85	0.45
1:B:164:TYR:N	1:B:164:TYR:CD1	2.85	0.45
1:A:226:ARG:HD3	1:A:301:TRP:CZ2	2.52	0.44
1:B:224:VAL:HG21	1:B:254:THR:CG2	2.46	0.44
1:C:18:GLU:HG3	1:C:19:GLU:HG3	1.98	0.44
1:C:164:TYR:CD1	1:C:164:TYR:N	2.85	0.44
1:A:193:GLU:OE1	3:A:413:PLP:N1	2.51	0.44
1:B:212:HIS:O	1:B:213:SER:C	2.61	0.44
1:C:69:MET:HG3	1:C:117:ALA:HB2	1.99	0.44
1:B:2:GLN:NE2	1:B:17:GLN:HE21	2.16	0.43
1:C:226:ARG:HD3	1:C:301:TRP:HZ2	1.81	0.43
1:B:159:LYS:NZ	3:B:913:PLP:O3	2.52	0.43
1:C:141:LEU:HD22	1:C:181:GLU:HB3	1.99	0.43
1:A:92:ARG:HH22	1:A:125:GLU:CG	2.31	0.43
4:B:914:MPD:HM2	4:B:914:MPD:C5	2.48	0.43
1:C:243:ASP:O	1:C:247:MET:HE3	2.18	0.43
1:A:54:LYS:HG2	1:A:55:GLU:N	2.34	0.43
1:B:4:LYS:O	1:B:4:LYS:HG2	2.19	0.43
1:B:193:GLU:OE1	3:B:913:PLP:N1	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:11:ASN:HD21	1:A:118:GLU:HB3	1.83	0.43
1:C:212:HIS:O	1:C:213:SER:C	2.61	0.43
1:C:2:GLN:NE2	1:C:17:GLN:HE21	2.16	0.43
1:B:132:GLU:H	1:B:132:GLU:CD	2.27	0.43
1:C:112:LEU:N	1:C:113:PRO:HD2	2.34	0.42
1:B:107:LEU:HD23	1:B:107:LEU:HA	1.91	0.42
1:C:149:PHE:CG	1:C:154:MET:HB2	2.54	0.42
1:A:54:LYS:N	1:A:54:LYS:HD3	2.35	0.42
1:A:212:HIS:O	1:A:213:SER:C	2.63	0.41
1:B:112:LEU:N	1:B:113:PRO:HD2	2.35	0.41
1:C:193:GLU:OE1	3:C:1413:PLP:N1	2.52	0.41
1:A:92:ARG:NH2	1:A:125:GLU:CG	2.81	0.41
1:C:11:ASN:ND2	1:C:119:VAL:H	2.18	0.41
1:B:141:LEU:CD2	1:B:181:GLU:HB3	2.50	0.41
1:B:294:ARG:CG	1:B:294:ARG:HH11	2.34	0.41
1:C:107:LEU:HD23	1:C:107:LEU:HA	1.90	0.41
1:C:174:ALA:HB2	1:C:182:ALA:HB2	2.03	0.41
4:A:414:MPD:HM2	4:A:414:MPD:C5	2.50	0.41
4:A:414:MPD:H52	5:A:1464:HOH:O	2.20	0.41
1:A:107:LEU:HD23	1:A:107:LEU:HA	1.92	0.41
1:A:27:HIS:CE1	1:A:108:GLY:O	2.71	0.41
1:A:254:THR:HA	1:A:259:GLU:O	2.21	0.41
1:C:95:TYR:HB2	1:C:126:TRP:HD1	1.86	0.41
1:A:174:ALA:HB2	1:A:182:ALA:HB2	2.03	0.41
1:B:149:PHE:CG	1:B:154:MET:HB2	2.56	0.41
1:C:273:LYS:H	1:C:273:LYS:HG3	1.61	0.41
1:A:54:LYS:HZ2	1:A:55:GLU:HG2	1.86	0.40
1:A:99:LEU:HB3	1:A:120:MET:HE3	2.03	0.40
1:A:16:PRO:HB2	1:A:18:GLU:HG2	2.02	0.40
1:A:243:ASP:OD1	1:A:243:ASP:N	2.54	0.40

All (1) 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)
5:C:1485:HOH:O	5:C:1485:HOH:O[8_665]	1.18	1.02

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	293/308 (95%)	286 (98%)	7 (2%)	0	100	100
1	B	303/308 (98%)	295 (97%)	8 (3%)	0	100	100
1	C	291/308 (94%)	285 (98%)	6 (2%)	0	100	100
All	All	887/924 (96%)	866 (98%)	21 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	233/241 (97%)	224 (96%)	9 (4%)	28	5
1	B	239/241 (99%)	234 (98%)	5 (2%)	47	19
1	C	232/241 (96%)	228 (98%)	4 (2%)	53	26
All	All	704/723 (97%)	686 (97%)	18 (3%)	40	13

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	53	LEU
1	A	54	LYS
1	A	76	GLU
1	A	92	ARG
1	A	99	LEU

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Mol	Chain	Res	Type
1	A	126	TRP
1	A	184	LEU
1	A	251	VAL
1	A	273	LYS
1	B	4	LYS
1	B	53	LEU
1	B	76	GLU
1	B	78	LEU
1	B	184	LEU
1	C	76	GLU
1	C	184	LEU
1	C	273	LYS
1	C	299	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	ASN
1	A	27	HIS
1	A	30	HIS
1	A	56	HIS
1	A	62	ASN
1	A	212	HIS
1	B	2	GLN
1	B	30	HIS
1	B	56	HIS
1	B	62	ASN
1	B	166	ASN
1	B	212	HIS
1	C	2	GLN
1	C	11	ASN
1	C	27	HIS
1	C	30	HIS
1	C	56	HIS
1	C	62	ASN
1	C	166	ASN
1	C	212	HIS

5.3.3 RNA ⓘ

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](#)

Of 11 ligands modelled in this entry, 3 are monoatomic - leaving 8 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
3	PLP	A	413	1	15,15,16	1.11	1 (6%)	21,22,23	1.40	3 (14%)
4	MPD	A	414	-	7,7,7	0.73	0	9,10,10	0.78	0
3	PLP	B	913	1	15,15,16	1.52	3 (20%)	21,22,23	1.56	5 (23%)
4	MPD	A	415	-	7,7,7	0.69	0	9,10,10	0.96	0
4	MPD	B	915	-	7,7,7	0.76	0	9,10,10	0.78	0
4	MPD	C	1414	-	7,7,7	0.71	0	9,10,10	0.79	0
4	MPD	B	914	-	7,7,7	0.84	0	9,10,10	0.75	0
3	PLP	C	1413	1	15,15,16	1.42	2 (13%)	21,22,23	1.37	4 (19%)

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	1	-	0/6/6/8	0/1/1/1
4	MPD	A	414	-	-	2/5/5/5	-
3	PLP	B	913	1	-	0/6/6/8	0/1/1/1
4	MPD	A	415	-	-	2/5/5/5	-
4	MPD	B	915	-	-	2/5/5/5	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	MPD	C	1414	-	-	2/5/5/5	-
4	MPD	B	914	-	-	3/5/5/5	-
3	PLP	C	1413	1	-	0/6/6/8	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	1413	PLP	C2A-C2	2.95	1.55	1.50
3	B	913	PLP	C4A-C4	2.94	1.57	1.51
3	B	913	PLP	C2A-C2	2.80	1.54	1.50
3	B	913	PLP	C3-C2	-2.78	1.38	1.41
3	A	413	PLP	C3-C2	-2.35	1.38	1.41
3	C	1413	PLP	C4A-C4	2.27	1.56	1.51

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	913	PLP	O4P-C5A-C5	3.38	115.69	109.36
3	A	413	PLP	O4P-C5A-C5	3.23	115.40	109.36
3	B	913	PLP	O2P-P-O4P	-2.91	99.09	106.67
3	A	413	PLP	O2P-P-O4P	-2.74	99.52	106.67
3	B	913	PLP	C6-C5-C4	2.48	120.13	118.10
3	C	1413	PLP	O4P-C5A-C5	2.44	113.93	109.36
3	C	1413	PLP	O2P-P-O4P	-2.36	100.52	106.67
3	B	913	PLP	C4A-C4-C5	2.28	123.29	120.94
3	C	1413	PLP	C4A-C4-C5	2.26	123.27	120.94
3	C	1413	PLP	C6-C5-C4	2.19	119.89	118.10
3	B	913	PLP	C5-C6-N1	-2.14	120.35	123.83
3	A	413	PLP	C5A-C5-C6	-2.02	116.06	119.36

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	414	MPD	C2-C3-C4-C5
4	A	415	MPD	C2-C3-C4-C5
4	B	914	MPD	C2-C3-C4-C5
4	B	915	MPD	C2-C3-C4-C5
4	C	1414	MPD	C2-C3-C4-C5
4	C	1414	MPD	O2-C2-C3-C4
4	A	414	MPD	O2-C2-C3-C4

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Mol	Chain	Res	Type	Atoms
4	A	415	MPD	O2-C2-C3-C4
4	B	914	MPD	O2-C2-C3-C4
4	B	915	MPD	O2-C2-C3-C4
4	B	914	MPD	C2-C3-C4-O4

There are no ring outliers.

8 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	413	PLP	1	0
4	A	414	MPD	4	0
3	B	913	PLP	2	0
4	A	415	MPD	2	0
4	B	915	MPD	1	0
4	C	1414	MPD	2	0
4	B	914	MPD	3	0
3	C	1413	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

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.