



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 01:39 am GMT

PDB ID : 2COI
Title : Crystal structure of oxidized human cytosolic branched-chain aminotransferase complexed with gabapentin
Authors : Goto, M.
Deposited on : 2005-05-17
Resolution : 1.90 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

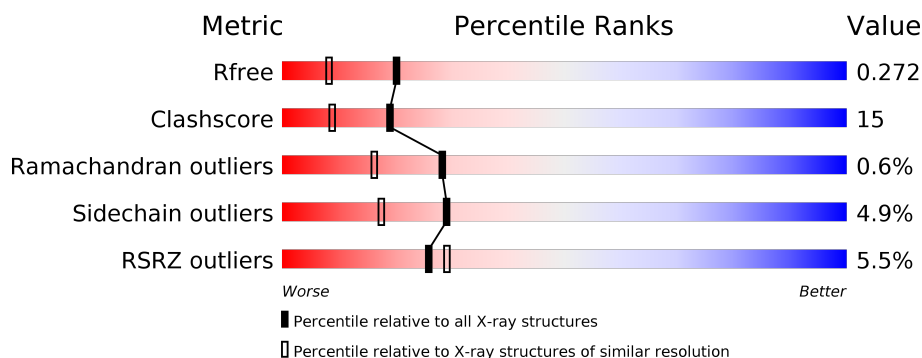
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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}	100719	5047 (1.90-1.90)
Clashscore	112137	5731 (1.90-1.90)
Ramachandran outliers	110173	5669 (1.90-1.90)
Sidechain outliers	110143	5670 (1.90-1.90)
RSRZ outliers	101464	5100 (1.90-1.90)

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. 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	386	<div> <div>7%</div> <div> <div></div> <div>64%</div> <div>26%</div> <div>• 7%</div> </div> </div>
1	B	386	<div> <div>3%</div> <div> <div></div> <div>67%</div> <div>23%</div> <div>• 8%</div> </div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6258 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 aminotransferase 1, cytosolic.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	358	Total	C	N	O	S	0	0	0
			2809	1811	462	521	15			
1	B	356	Total	C	N	O	S	0	0	0
			2814	1813	466	520	15			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	379	ARG	SER	ENGINEERED	UNP P54687
B	379	ARG	SER	ENGINEERED	UNP P54687

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

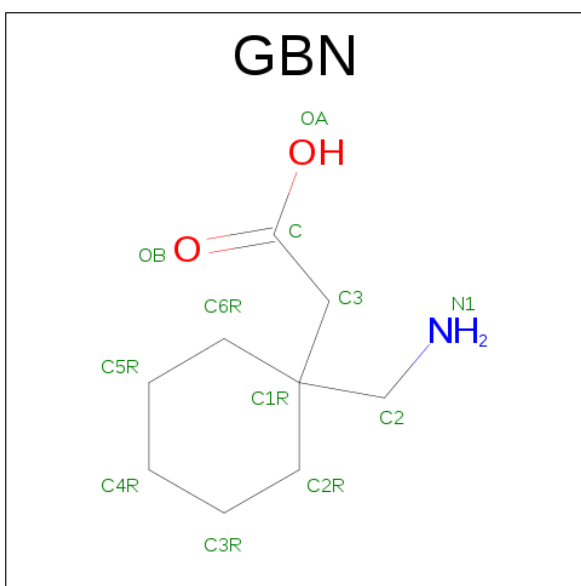
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Mg	0	0
			1	1		

- 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 [1-(AMINOMETHYL)CYCLOHEXYL]ACETIC ACID (three-letter code: GBN) (formula: C₉H₁₇NO₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			12	9	1	2		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total	C	N	O	0	0
			12	9	1	2		

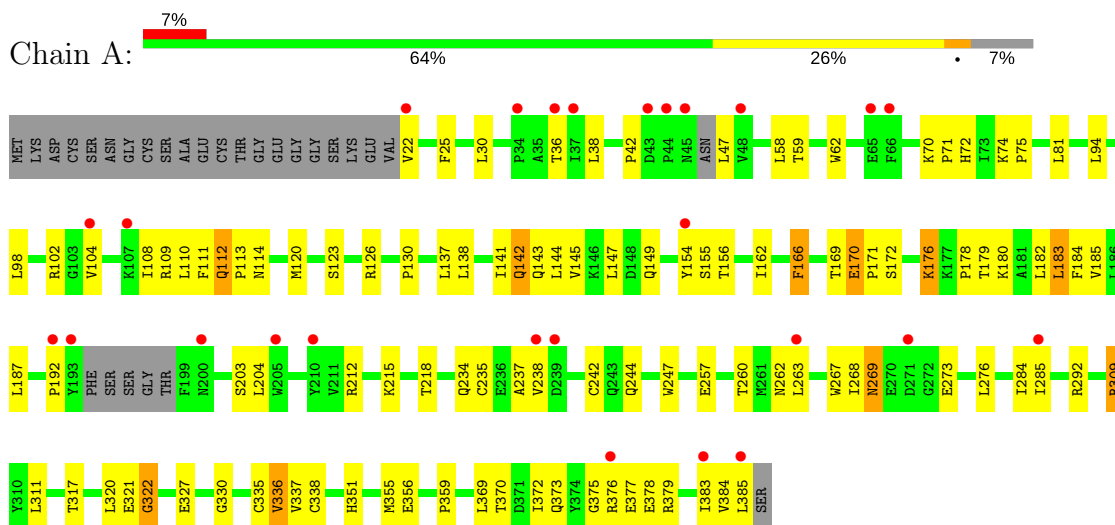
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	217	Total	O	0	0
			217	217		
5	B	363	Total	O	0	0
			363	363		

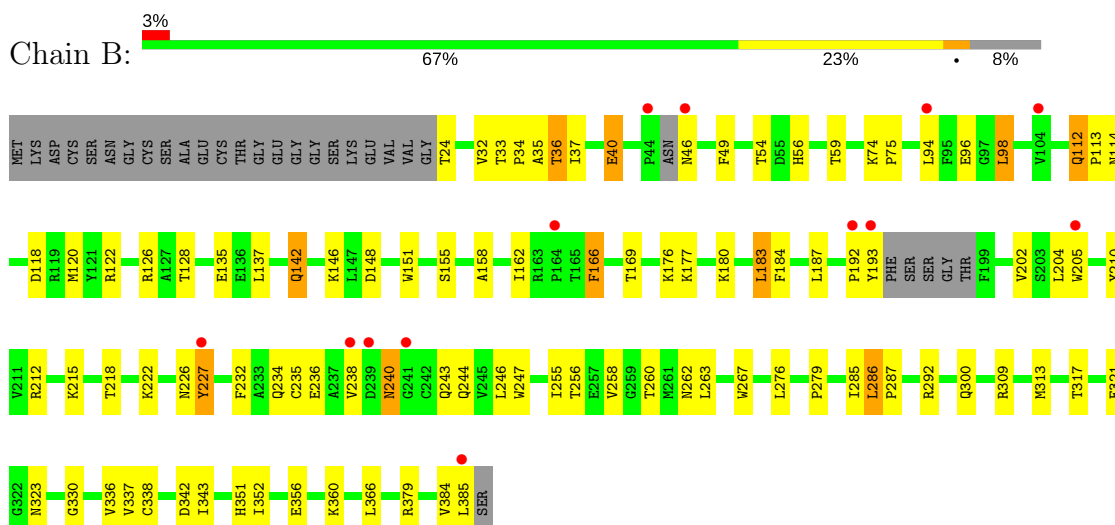
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: branched chain aminotransferase 1, cytosolic



- Molecule 1: branched chain aminotransferase 1, cytosolic



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	66.72Å 107.69Å 110.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	33.37 – 1.90 33.36 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.6 (33.37-1.90) 99.7 (33.36-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.64 (at 1.89Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.224 , 0.272 0.224 , 0.272	Depositor DCC
R_{free} test set	6386 reflections (10.12%)	DCC
Wilson B-factor (Å ²)	23.1	Xtriage
Anisotropy	0.890	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 58.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.019 for -h,l,k	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6258	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.78% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GBN, MG, 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	0.32	0/2880	0.61	0/3920
1	B	0.37	0/2885	0.64	0/3923
All	All	0.34	0/5765	0.63	0/7843

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	B	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	227	TYR	Sidechain

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	2809	0	2760	88	0
1	B	2814	0	2783	76	0
2	B	1	0	0	0	0
3	A	15	0	7	1	0
3	B	15	0	7	1	0
4	A	12	0	16	1	0
4	B	12	0	16	1	0
5	A	217	0	0	7	0
5	B	363	0	0	11	0
All	All	6258	0	5589	164	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 15.

All (164) 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:B:98:LEU:HD13	1:B:162:ILE:HB	1.46	0.97
1:B:35:ALA:H	1:B:56:HIS:HE1	1.07	0.96
1:A:284:ILE:HG13	1:A:285:ILE:HD12	1.50	0.92
1:A:144:LEU:HD13	1:A:187:LEU:HD11	1.64	0.77
1:B:246:LEU:HD11	1:B:255:ILE:HG23	1.67	0.76
1:A:114:ASN:HB2	5:A:627:HOH:O	1.87	0.74
1:B:40:GLU:H	1:B:40:GLU:CD	1.92	0.73
1:A:170:GLU:HG2	1:A:179:THR:H	1.54	0.73
1:B:351:HIS:HE1	1:B:356:GLU:OE1	1.71	0.73
1:A:269:ASN:HD21	1:A:273:GLU:H	1.38	0.71
1:B:74:LYS:HB2	1:B:75:PRO:HD2	1.72	0.71
1:B:126:ARG:HG3	5:B:1588:HOH:O	1.91	0.71
1:A:269:ASN:ND2	1:A:273:GLU:H	1.88	0.70
1:B:35:ALA:H	1:B:56:HIS:CE1	2.00	0.70
1:B:36:THR:HG23	1:B:37:ILE:HG13	1.75	0.69
1:A:142:GLN:NE2	1:A:385:LEU:HD13	2.08	0.69
1:A:369:LEU:O	1:A:373:GLN:HG3	1.95	0.67
1:B:317:THR:O	1:B:321:GLU:HG3	1.97	0.65
1:A:110:LEU:HB3	1:A:113:PRO:HG3	1.78	0.65
1:A:112:GLN:N	1:A:113:PRO:HD3	2.11	0.65
1:A:126:ARG:NH1	1:A:218:THR:HG22	2.13	0.64
1:A:170:GLU:HG3	1:A:178:PRO:HA	1.81	0.62
1:B:183:LEU:HD12	1:B:184:PHE:N	2.16	0.61
1:A:144:LEU:CD1	1:A:187:LEU:HD11	2.32	0.59
1:A:370:THR:HG22	1:A:376:ARG:HH22	1.66	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:170:GLU:OE1	1:A:171:PRO:HD2	2.02	0.59
1:B:24:THR:N	5:B:1614:HOH:O	2.36	0.59
1:B:343:ILE:HD12	1:B:352:ILE:HD11	1.83	0.59
1:A:155:SER:OG	1:A:156:THR:N	2.33	0.58
1:B:114:ASN:HB2	5:B:1627:HOH:O	2.02	0.58
1:B:118:ASP:HB3	5:B:1559:HOH:O	2.03	0.58
1:A:309:ARG:HH11	1:A:311:LEU:HD23	1.69	0.58
1:A:110:LEU:HB3	1:A:113:PRO:CG	2.34	0.58
1:A:138:LEU:O	1:A:142:GLN:HG2	2.03	0.57
1:A:384:VAL:HG12	1:A:385:LEU:N	2.19	0.57
1:B:343:ILE:CD1	1:B:352:ILE:HD11	2.34	0.57
1:A:120:MET:HE1	1:A:183:LEU:HD21	1.86	0.57
1:A:284:ILE:HG13	1:A:285:ILE:CD1	2.31	0.57
1:B:98:LEU:CD1	1:B:162:ILE:HB	2.27	0.57
1:A:58:LEU:O	1:A:184:PHE:HA	2.06	0.56
1:B:35:ALA:N	1:B:56:HIS:HE1	1.90	0.56
1:B:40:GLU:N	1:B:40:GLU:CD	2.59	0.56
1:A:317:THR:O	1:A:321:GLU:HG2	2.07	0.55
1:B:342:ASP:OD2	1:B:351:HIS:HD2	1.89	0.55
1:B:336:VAL:HG13	1:B:337:VAL:N	2.21	0.55
1:A:108:ILE:CD1	1:A:149:GLN:HB2	2.37	0.55
1:A:120:MET:CE	1:A:183:LEU:HD21	2.36	0.55
1:B:169:THR:OG1	1:B:180:LYS:HG2	2.06	0.55
1:A:234:GLN:O	1:A:237:ALA:HB3	2.07	0.55
1:A:384:VAL:HG12	1:A:385:LEU:H	1.72	0.54
1:B:210:TYR:HH	1:B:232:PHE:HE2	1.53	0.54
1:B:204:LEU:CD2	1:B:244:GLN:HB2	2.37	0.54
1:B:202:VAL:HA	1:B:243:GLN:OE1	2.07	0.54
1:B:212:ARG:HB2	1:B:247:TRP:CD2	2.43	0.54
1:A:182:LEU:C	1:A:182:LEU:HD13	2.29	0.54
1:A:58:LEU:HB3	1:A:185:VAL:HB	1.90	0.54
1:A:141:ILE:O	1:A:145:VAL:HG23	2.07	0.53
1:B:384:VAL:HG12	1:B:385:LEU:HG	1.89	0.53
1:B:258:VAL:CG2	1:B:263:LEU:HD23	2.38	0.53
1:B:112:GLN:HA	1:B:112:GLN:HE21	1.74	0.53
1:A:112:GLN:HA	1:A:112:GLN:HE21	1.72	0.53
1:A:30:LEU:CD1	1:A:143:GLN:HG3	2.40	0.52
1:A:351:HIS:HE1	1:A:356:GLU:OE1	1.93	0.52
1:B:336:VAL:HG13	1:B:337:VAL:H	1.75	0.52
1:A:170:GLU:CG	1:A:179:THR:H	2.23	0.52
1:A:263:LEU:HD23	1:A:263:LEU:C	2.30	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:36:THR:HG23	5:A:613:HOH:O	2.09	0.52
1:A:372:ILE:HD13	1:A:379:ARG:HD3	1.92	0.51
1:B:300:GLN:NE2	1:B:379:ARG:HH22	2.08	0.51
1:B:205:TRP:CH2	1:B:236:GLU:OE2	2.64	0.51
1:A:74:LYS:HB2	1:A:75:PRO:CD	2.41	0.51
1:B:146:LYS:HE2	5:B:1700:HOH:O	2.11	0.50
1:B:49:PHE:HB2	5:B:1659:HOH:O	2.11	0.50
1:A:317:THR:HA	1:A:320:LEU:HD12	1.94	0.50
1:B:33:THR:O	1:B:75:PRO:HD3	2.10	0.50
1:A:335:CYS:O	1:A:336:VAL:HB	2.12	0.50
1:A:94:LEU:HG	1:A:166:PHE:HB3	1.93	0.50
1:B:236:GLU:O	1:B:240:ASN:HB2	2.11	0.50
1:A:111:PHE:C	1:A:113:PRO:HD3	2.33	0.50
1:B:193:TYR:HA	5:B:1659:HOH:O	2.11	0.50
1:B:205:TRP:CZ3	1:B:236:GLU:OE2	2.65	0.50
1:A:327:GLU:OE1	1:A:359:PRO:HB2	2.11	0.49
1:B:98:LEU:N	1:B:98:LEU:HD12	2.26	0.49
1:B:118:ASP:O	1:B:122:ARG:HG3	2.12	0.49
1:B:215:LYS:HE3	5:B:1774:HOH:O	2.12	0.49
1:B:56:HIS:HD2	1:B:148:ASP:OD2	1.94	0.49
1:A:269:ASN:C	1:A:269:ASN:HD22	2.14	0.49
1:A:38:LEU:HD22	1:A:38:LEU:N	2.27	0.49
1:B:351:HIS:CE1	1:B:356:GLU:OE1	2.59	0.48
1:A:235:CYS:O	1:A:238:VAL:HG22	2.13	0.48
1:B:135:GLU:HG2	5:B:1574:HOH:O	2.13	0.48
1:A:336:VAL:HG12	1:A:337:VAL:HG23	1.94	0.48
1:A:59:THR:OG1	1:A:74:LYS:HE2	2.13	0.48
1:B:263:LEU:HD13	1:B:263:LEU:C	2.34	0.47
1:B:255:ILE:HD11	1:B:313:MET:SD	2.54	0.47
1:A:180:LYS:HE2	1:A:180:LYS:HB3	1.76	0.47
1:A:180:LYS:HG3	5:A:615:HOH:O	2.13	0.47
1:B:218:THR:HG23	1:B:226:ASN:OD1	2.15	0.47
1:B:155:SER:HB2	1:B:158:ALA:HB3	1.97	0.47
1:B:222:LYS:NZ	3:B:410:PLP:O3	2.46	0.47
1:A:25:PHE:HB2	1:A:71:PRO:HD3	1.97	0.47
1:A:94:LEU:HD12	1:A:94:LEU:C	2.35	0.47
1:B:128:THR:HB	1:B:177:LYS:HG3	1.97	0.46
1:B:279:PRO:HA	1:B:309:ARG:O	2.15	0.46
1:A:102:ARG:HD2	1:A:154:TYR:HA	1.98	0.46
1:B:180:LYS:HE2	5:B:1499:HOH:O	2.15	0.46
1:A:112:GLN:N	1:A:113:PRO:CD	2.79	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:144:LEU:HD11	1:A:187:LEU:HG	1.98	0.45
1:B:227:TYR:CD1	1:B:227:TYR:N	2.85	0.45
1:B:142:GLN:HE21	1:B:142:GLN:HB3	1.55	0.45
1:B:360:LYS:NZ	1:B:360:LYS:HB3	2.31	0.45
1:B:122:ARG:HH11	1:B:122:ARG:HG2	1.82	0.45
1:A:98:LEU:HG	1:A:162:ILE:HB	1.99	0.45
1:A:212:ARG:HB2	1:A:247:TRP:CD2	2.52	0.45
1:A:260:THR:OG1	4:A:420:GBN:N1	2.50	0.45
1:B:56:HIS:CD2	1:B:148:ASP:OD2	2.70	0.45
1:A:30:LEU:HD11	1:A:143:GLN:HG3	1.98	0.45
1:A:172:SER:OG	1:A:176:LYS:HE2	2.18	0.44
1:B:218:THR:HG22	5:B:1775:HOH:O	2.17	0.44
1:B:96:GLU:HG3	1:B:120:MET:HB2	2.00	0.44
1:A:130:PRO:HG2	1:A:166:PHE:HZ	1.83	0.44
1:B:59:THR:OG1	1:B:74:LYS:HE2	2.18	0.43
1:B:205:TRP:CZ3	1:B:236:GLU:HG2	2.52	0.43
1:A:98:LEU:N	1:A:98:LEU:HD23	2.32	0.43
1:A:257:GLU:OE1	3:A:410:PLP:N1	2.52	0.43
1:B:336:VAL:HG11	1:B:366:LEU:CD1	2.48	0.43
1:A:62:TRP:CD2	1:A:130:PRO:HG3	2.54	0.42
1:A:47:LEU:N	5:A:499:HOH:O	2.51	0.42
1:A:215:LYS:HD2	1:A:284:ILE:HG23	2.00	0.42
1:A:378:GLU:OE1	1:A:378:GLU:HA	2.20	0.42
1:B:234:GLN:HE21	1:B:238:VAL:HG23	1.84	0.42
1:B:258:VAL:HG22	1:B:263:LEU:HD23	2.00	0.42
1:A:292:ARG:NH1	5:A:452:HOH:O	2.53	0.42
1:A:262:ASN:O	1:A:330:GLY:HA2	2.20	0.42
1:A:335:CYS:O	1:A:336:VAL:CB	2.67	0.42
1:B:94:LEU:HG	1:B:166:PHE:HB3	2.02	0.42
1:B:32:VAL:HG12	1:B:34:PRO:HD3	2.01	0.42
1:B:255:ILE:HG22	1:B:263:LEU:HG	2.02	0.42
1:A:204:LEU:CD2	1:A:244:GLN:HB2	2.49	0.42
1:B:292:ARG:O	1:B:292:ARG:HD3	2.19	0.42
1:B:54:THR:HG22	1:B:151:TRP:CD2	2.55	0.42
1:A:70:LYS:O	1:A:72:HIS:HD2	2.02	0.41
1:A:74:LYS:HB2	1:A:75:PRO:HD2	2.01	0.41
1:B:112:GLN:N	1:B:113:PRO:CD	2.83	0.41
1:A:182:LEU:HD12	1:A:184:PHE:HD2	1.84	0.41
1:B:260:THR:OG1	4:B:420:GBN:N1	2.53	0.41
1:B:256:THR:HB	1:B:285:ILE:HD11	2.03	0.41
1:B:54:THR:OG1	1:B:187:LEU:HB2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:203:SER:HB3	1:A:242:CYS:HA	2.02	0.41
1:B:204:LEU:HD23	1:B:244:GLN:HB2	2.01	0.41
1:A:22:VAL:HG13	1:A:22:VAL:O	2.20	0.41
1:A:268:ILE:O	1:A:268:ILE:HG22	2.21	0.41
1:A:320:LEU:C	1:A:322:GLY:H	2.24	0.41
1:A:355:MET:HG2	5:A:475:HOH:O	2.20	0.41
1:A:81:LEU:HD22	1:A:169:THR:HG22	2.03	0.41
1:A:104:VAL:HG21	1:A:376:ARG:HG2	2.03	0.41
1:B:262:ASN:O	1:B:330:GLY:HA2	2.20	0.41
1:A:42:PRO:HB2	1:A:47:LEU:HD21	2.03	0.41
1:B:286:LEU:HA	1:B:287:PRO:HD3	1.91	0.40
1:A:182:LEU:HD13	1:A:183:LEU:N	2.35	0.40
1:A:109:ARG:HA	1:A:383:ILE:O	2.21	0.40
1:A:94:LEU:HD22	1:A:123:SER:CB	2.52	0.40
1:A:375:GLY:C	1:A:377:GLU:H	2.22	0.40
1:A:130:PRO:HD3	5:A:449:HOH:O	2.20	0.40

There are no symmetry-related clashes.

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	352/386 (91%)	320 (91%)	29 (8%)	3 (1%)	20	8
1	B	350/386 (91%)	332 (95%)	17 (5%)	1 (0%)	44	34
All	All	702/772 (91%)	652 (93%)	46 (7%)	4 (1%)	28	16

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	192	PRO
1	B	192	PRO

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Mol	Chain	Res	Type
1	A	322	GLY
1	A	336	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	305/334 (91%)	292 (96%)	13 (4%)	33	22
1	B	308/334 (92%)	291 (94%)	17 (6%)	25	14
All	All	613/668 (92%)	583 (95%)	30 (5%)	29	17

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	112	GLN
1	A	137	LEU
1	A	142	GLN
1	A	147	LEU
1	A	166	PHE
1	A	170	GLU
1	A	176	LYS
1	A	183	LEU
1	A	267	TRP
1	A	269	ASN
1	A	276	LEU
1	A	309	ARG
1	A	338	CYS
1	B	36	THR
1	B	40	GLU
1	B	46	ASN
1	B	98	LEU
1	B	112	GLN
1	B	137	LEU
1	B	142	GLN
1	B	166	PHE

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Mol	Chain	Res	Type
1	B	176	LYS
1	B	183	LEU
1	B	235	CYS
1	B	240	ASN
1	B	267	TRP
1	B	276	LEU
1	B	286	LEU
1	B	323	ASN
1	B	338	CYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	72	HIS
1	A	112	GLN
1	A	116	ASN
1	A	142	GLN
1	A	254	GLN
1	A	269	ASN
1	A	300	GLN
1	A	351	HIS
1	B	56	HIS
1	B	112	GLN
1	B	143	GLN
1	B	200	ASN
1	B	243	GLN
1	B	254	GLN
1	B	262	ASN
1	B	300	GLN
1	B	351	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 1 is monoatomic - leaving 4 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	410	1	15,15,16	1.28	2 (13%)	20,22,23	1.60	4 (20%)
4	GBN	A	420	-	8,12,12	1.99	2 (25%)	12,16,16	7.33	5 (41%)
3	PLP	B	410	1	15,15,16	1.11	0	20,22,23	1.37	3 (15%)
4	GBN	B	420	-	8,12,12	1.82	2 (25%)	12,16,16	7.43	5 (41%)

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	410	1	-	0/6/6/8	0/1/1/1
4	GBN	A	420	-	-	0/6/18/18	0/1/1/1
3	PLP	B	410	1	-	0/6/6/8	0/1/1/1
4	GBN	B	420	-	-	0/6/18/18	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	410	PLP	C3-C2	-2.93	1.38	1.40
4	B	420	GBN	C2R-C3R	2.32	1.58	1.52
3	A	410	PLP	C2A-C2	2.42	1.54	1.50
4	A	420	GBN	C2R-C3R	2.48	1.59	1.52
4	B	420	GBN	C3-C1R	2.91	1.58	1.55
4	A	420	GBN	C3-C1R	3.38	1.59	1.55

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	420	GBN	C2-C1R-C3	-21.99	87.55	110.44
4	A	420	GBN	C2-C1R-C3	-21.67	87.88	110.44
4	A	420	GBN	C3R-C2R-C1R	-4.11	107.74	113.60
4	B	420	GBN	C3R-C2R-C1R	-4.00	107.89	113.60
3	A	410	PLP	O2P-P-O4P	-2.18	100.94	106.73
4	A	420	GBN	C6R-C1R-C2	2.03	112.68	109.07
3	B	410	PLP	O3P-P-O1P	2.06	118.56	110.50
3	A	410	PLP	C4A-C4-C5	2.18	123.05	120.86
3	B	410	PLP	C4A-C4-C5	2.38	123.26	120.86
4	B	420	GBN	C6R-C1R-C2	2.40	113.35	109.07
3	A	410	PLP	O3P-P-O1P	2.50	120.28	110.50
3	B	410	PLP	O4P-C5A-C5	2.52	114.38	109.32
3	A	410	PLP	O4P-C5A-C5	4.06	117.48	109.32
4	A	420	GBN	C1R-C3-C	8.10	128.72	116.73
4	B	420	GBN	C1R-C3-C	8.29	128.99	116.73
4	A	420	GBN	C2R-C1R-C2	9.05	125.19	109.07
4	B	420	GBN	C2R-C1R-C2	9.06	125.20	109.07

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	410	PLP	1	0
4	A	420	GBN	1	0
3	B	410	PLP	1	0
4	B	420	GBN	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	358/386 (92%)	0.69	26 (7%) 16 18	17, 34, 55, 82	0
1	B	356/386 (92%)	0.24	13 (3%) 42 46	14, 24, 40, 61	0
All	All	714/772 (92%)	0.47	39 (5%) 26 29	14, 28, 50, 82	0

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	193	TYR	8.4
1	A	22	VAL	7.6
1	A	192	PRO	5.5
1	A	36	THR	5.1
1	B	385	LEU	4.6
1	B	193	TYR	4.2
1	B	192	PRO	3.6
1	A	37	ILE	3.2
1	B	46	ASN	3.2
1	A	66	PHE	3.0
1	A	205	TRP	3.0
1	B	227	TYR	2.8
1	A	376	ARG	2.8
1	A	44	PRO	2.7
1	B	104	VAL	2.6
1	B	239	ASP	2.6
1	A	107	LYS	2.5
1	A	385	LEU	2.5
1	A	104	VAL	2.4
1	B	44	PRO	2.4
1	B	205	TRP	2.3
1	A	43	ASP	2.3
1	A	45	ASN	2.3
1	A	383	ILE	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	34	PRO	2.3
1	A	285	ILE	2.3
1	A	65	GLU	2.3
1	A	200	ASN	2.2
1	A	210	TYR	2.2
1	A	238	VAL	2.2
1	A	239	ASP	2.2
1	A	48	VAL	2.2
1	B	238	VAL	2.1
1	B	241	GLY	2.1
1	A	263	LEU	2.1
1	A	154	TYR	2.1
1	A	271	ASP	2.1
1	B	94	LEU	2.0
1	B	164	PRO	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 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	LLDF	B-factors(\AA^2)	Q < 0.9
4	GBN	B	420	12/12	0.90	0.22	1.24	15,26,29,32	0
3	PLP	A	410	15/16	0.97	0.17	0.10	20,25,30,32	0
4	GBN	A	420	12/12	0.90	0.17	-0.06	24,33,35,36	0
3	PLP	B	410	15/16	0.98	0.15	-0.12	12,19,21,22	0
2	MG	B	1430	1/1	0.91	0.14	-	31,31,31,31	0

6.5 Other polymers ⓘ

There are no such residues in this entry.