



Full wwPDB X-ray Structure Validation Report ⓘ

May 26, 2020 – 05:19 pm BST

PDB ID : 6DND
Title : Crystal structure of wild-type (WT) human Glutamate oxaloacetate transaminase 1 (GOT1)
Authors : Assar, Z.; Holt, M.C.; Stein, A.J.; Lairson, L.; Lyssiotis, C.A.
Deposited on : 2018-06-06
Resolution : 2.10 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

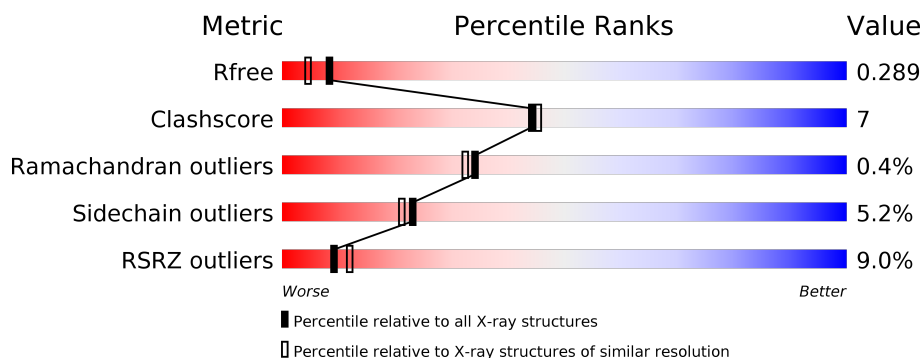
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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}	130704	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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 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	411	<div> <div>13%</div> <div> <div></div> <div>88%</div> <div>10%</div> </div> </div>
1	B	411	<div> <div>5%</div> <div> <div></div> <div>85%</div> <div>12%</div> </div> </div>

2 Entry composition [i](#)

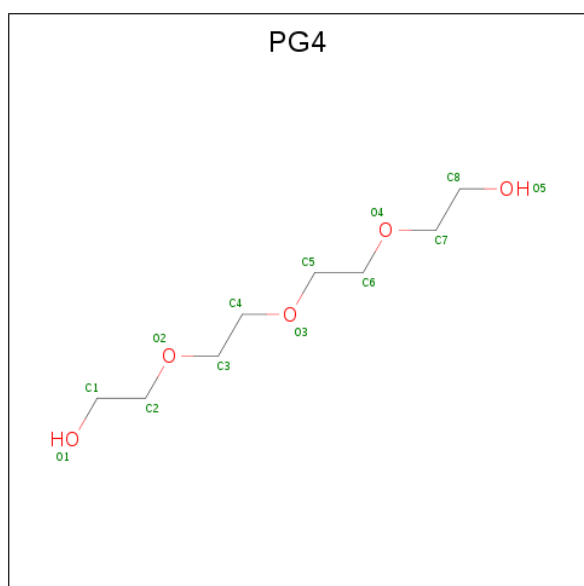
There are 4 unique types of molecules in this entry. The entry contains 6628 atoms, of which 0 are hydrogens and 0 are deuteriums.

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, cytoplasmic.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	409	Total	C	N	O	S	0	4	0
			3237	2070	562	596	9			
1	B	410	Total	C	N	O	S	0	1	0
			3244	2070	564	601	9			

- Molecule 2 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula: C₈H₁₈O₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			7	4	3		
2	B	1	Total	C	O	0	0
			7	4	3		

- 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
			16	8	1	6	1		
3	B	1	Total	C	N	O	P	0	0
			16	8	1	6	1		

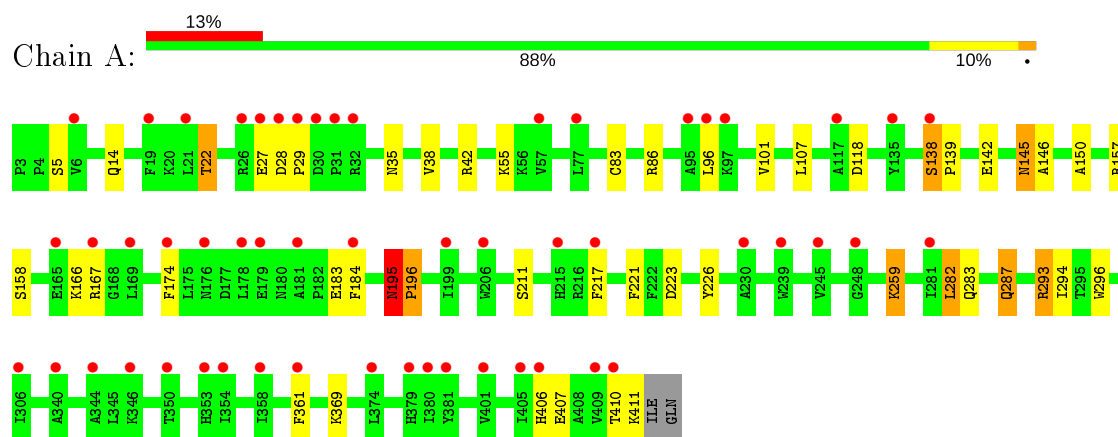
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	35	Total	O	0	0
			35	35		
4	B	66	Total	O	0	0
			66	66		

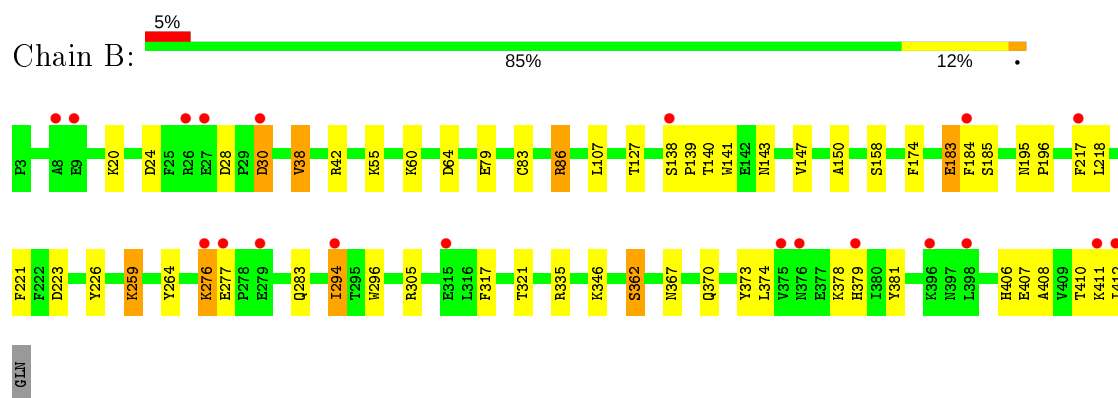
3 Residue-property plots [i](#)

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 ($\text{RSRZ} > 2$). 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, cytoplasmic



- Molecule 1: Aspartate aminotransferase, cytoplasmic



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	91.04Å 96.86Å 97.87Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	68.85 – 2.10 48.43 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.9 (68.85-2.10) 99.9 (48.43-2.10)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.22 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, R_{free}	0.208 , 0.269 0.231 , 0.289	Depositor DCC
R_{free} test set	2570 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	50.6	Xtriage
Anisotropy	0.049	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 52.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.016 for -h,l,k	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6628	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.71% 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: PG4, 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	0.68	0/3332	0.79	6/4533 (0.1%)
1	B	0.75	1/3328 (0.0%)	0.79	4/4525 (0.1%)
All	All	0.72	1/6660 (0.0%)	0.79	10/9058 (0.1%)

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
All	All	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	79	GLU	CD-OE1	6.00	1.32	1.25

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	195	ASN	C-N-CD	-6.66	105.96	120.60
1	A	287[A]	GLN	CA-CB-CG	-5.96	100.28	113.40
1	A	287[B]	GLN	CA-CB-CG	-5.96	100.28	113.40
1	B	86	ARG	NE-CZ-NH2	5.45	123.02	120.30
1	A	118	ASP	CB-CG-OD1	5.41	123.17	118.30
1	B	42	ARG	NE-CZ-NH1	5.40	123.00	120.30
1	A	42	ARG	NE-CZ-NH1	5.33	122.97	120.30
1	B	305	ARG	NE-CZ-NH1	5.33	122.96	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	157	ARG	NE-CZ-NH1	5.06	122.83	120.30
1	B	64	ASP	CB-CG-OD2	-5.01	113.79	118.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	195	ASN	Peptide
1	B	195	ASN	Peptide

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	3237	0	3147	43	0
1	B	3244	0	3179	47	0
2	A	7	0	8	0	0
2	B	7	0	9	0	0
3	A	16	0	7	3	0
3	B	16	0	7	4	0
4	A	35	0	0	2	0
4	B	66	0	0	2	0
All	All	6628	0	6357	85	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 7.

All (85) 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:30:ASP:CG	1:B:379:HIS:CD2	1.97	1.37
1:A:293:ARG:HD3	4:A:631:HOH:O	1.24	1.29
1:B:30:ASP:OD2	1:B:379:HIS:CD2	1.94	1.19
1:A:22:THR:HG22	1:A:35:ASN:HD21	1.04	1.15
1:B:259:LYS:HE2	3:B:502:PLP:O4A	1.46	1.13

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:138:SER:HB3	1:B:139:PRO:HD3	1.25	1.09
1:B:138:SER:HB3	1:B:139:PRO:CD	1.87	1.05
1:A:138:SER:HB3	1:A:139:PRO:HD3	1.39	1.04
1:A:138:SER:HB3	1:A:139:PRO:CD	1.88	1.03
1:B:30:ASP:CG	1:B:379:HIS:HD2	1.43	1.00
1:B:30:ASP:OD2	1:B:379:HIS:CG	2.17	0.97
1:B:28:ASP:OD2	1:B:381:TYR:OH	1.81	0.95
1:A:22:THR:HG22	1:A:35:ASN:ND2	1.85	0.91
1:A:293:ARG:HG3	1:A:293:ARG:HH21	1.34	0.90
1:B:259:LYS:CE	3:B:502:PLP:O4A	2.27	0.82
1:B:30:ASP:OD2	1:B:379:HIS:CB	2.28	0.81
1:B:30:ASP:CB	1:B:379:HIS:CD2	2.65	0.80
1:B:138:SER:CB	1:B:139:PRO:HD3	2.11	0.79
1:B:408:ALA:O	1:B:411:LYS:O	2.04	0.76
1:A:22:THR:CG2	1:A:35:ASN:HD21	1.94	0.74
1:B:30:ASP:OD1	1:B:379:HIS:HD2	1.70	0.74
1:A:28:ASP:OD2	1:A:29:PRO:HD2	1.90	0.72
1:B:30:ASP:OD2	1:B:379:HIS:HB2	1.89	0.72
1:A:406:HIS:O	1:A:410:THR:HG23	1.90	0.70
1:A:195:ASN:HD22	1:A:196:PRO:CA	2.04	0.70
1:B:184:PHE:HA	1:B:217:PHE:O	1.94	0.65
1:B:259:LYS:HD2	1:B:264:TYR:HE1	1.61	0.64
1:A:138:SER:CB	1:A:139:PRO:HD3	2.22	0.64
1:A:195:ASN:HD22	1:A:196:PRO:HA	1.64	0.62
1:B:259:LYS:HD2	1:B:264:TYR:CE1	2.36	0.61
1:B:183:GLU:HG2	1:B:217:PHE:HD2	1.65	0.60
1:A:293:ARG:NH2	1:A:293:ARG:HG3	2.08	0.60
1:A:166:LYS:HG2	1:A:166:LYS:O	2.03	0.58
1:A:294:ILE:HD13	1:B:150:ALA:CB	2.38	0.54
1:A:195:ASN:C	1:A:195:ASN:HD22	2.11	0.54
1:B:407:GLU:O	1:B:411:LYS:HG3	2.08	0.54
1:B:373:TYR:HB2	1:B:412:ILE:CD1	2.38	0.54
1:B:30:ASP:OD1	1:B:379:HIS:CD2	2.52	0.52
1:A:407:GLU:O	1:A:411:LYS:HG3	2.10	0.52
1:B:317:PHE:O	1:B:321:THR:HG23	2.10	0.51
1:A:145[A]:ASN:C	1:A:145[A]:ASN:ND2	2.63	0.51
1:A:259:LYS:NZ	3:A:502:PLP:C4A	2.74	0.51
1:A:294:ILE:HD13	1:B:150:ALA:HB1	1.93	0.51
1:B:226:TYR:OH	3:B:502:PLP:O3	2.27	0.51
1:A:226:TYR:CZ	1:A:361[B]:PHE:HE2	2.29	0.50
1:B:406:HIS:O	1:B:410:THR:HG23	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:150:ALA:HB2	1:B:294:ILE:HD12	1.94	0.50
1:A:138:SER:CB	1:A:139:PRO:CD	2.65	0.50
1:B:38:VAL:HG13	4:B:662:HOH:O	2.13	0.49
1:A:107:LEU:HD11	1:B:107:LEU:HD11	1.95	0.48
1:B:378:LYS:HE2	1:B:407:GLU:OE1	2.13	0.48
1:B:60:LYS:HD2	4:B:633:HOH:O	2.14	0.48
1:A:145[A]:ASN:HD22	1:A:145[A]:ASN:C	2.17	0.48
1:B:30:ASP:OD2	1:B:379:HIS:HD2	1.56	0.47
1:A:145[A]:ASN:HD22	1:A:146:ALA:N	2.13	0.47
1:B:335:ARG:HH12	1:B:362:SER:HB2	1.80	0.47
1:B:83:CYS:SG	1:B:86:ARG:NH2	2.85	0.47
1:B:30:ASP:HB2	1:B:379:HIS:CD2	2.50	0.47
1:A:138:SER:HB3	1:A:139:PRO:HD2	1.85	0.46
1:A:259:LYS:HZ1	3:A:502:PLP:C4A	2.28	0.46
1:B:185:SER:HB2	1:B:218:LEU:HD22	1.97	0.46
1:B:373:TYR:HB2	1:B:412:ILE:HD12	1.97	0.46
1:A:293:ARG:CD	4:A:631:HOH:O	2.08	0.45
1:A:83:CYS:SG	1:A:86:ARG:NH1	2.88	0.45
1:A:28:ASP:HA	1:A:29:PRO:HD3	1.75	0.44
1:B:276:LYS:HB3	1:B:276:LYS:HE3	1.36	0.44
1:B:138:SER:CB	1:B:139:PRO:CD	2.62	0.44
1:A:107:LEU:HD23	1:A:296:TRP:CE2	2.53	0.43
1:A:184:PHE:HA	1:A:217:PHE:O	2.18	0.43
1:B:140:THR:HG23	1:B:141:TRP:O	2.18	0.43
1:B:223:ASP:OD2	3:B:502:PLP:N1	2.51	0.43
1:A:223:ASP:OD2	3:A:502:PLP:N1	2.52	0.43
1:A:282:LEU:HD12	1:A:282:LEU:HA	1.76	0.43
1:A:96:LEU:HD23	1:A:101:VAL:HG21	2.01	0.42
1:B:107:LEU:HD23	1:B:296:TRP:CE2	2.53	0.42
1:A:195:ASN:ND2	1:A:195:ASN:C	2.72	0.41
1:A:226:TYR:CZ	1:A:361[B]:PHE:CE2	3.08	0.41
1:A:293:ARG:NH2	1:A:293:ARG:CG	2.73	0.41
1:B:143:ASN:O	1:B:147:VAL:HG23	2.21	0.41
1:A:226:TYR:OH	1:A:361[B]:PHE:HE2	2.04	0.41
1:A:195:ASN:HD22	1:A:196:PRO:N	2.19	0.40
1:A:5:SER:HB2	1:B:276:LYS:HB2	2.03	0.40
1:B:367:ASN:H	1:B:370:GLN:NE2	2.18	0.40
1:B:20:LYS:HE3	1:B:24:ASP:OD1	2.20	0.40
1:A:217:PHE:N	1:A:217:PHE:CD1	2.89	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	411/411 (100%)	393 (96%)	16 (4%)	2 (0%)	29	26
1	B	409/411 (100%)	395 (97%)	13 (3%)	1 (0%)	47	49
All	All	820/822 (100%)	788 (96%)	29 (4%)	3 (0%)	34	32

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	196	PRO
1	B	196	PRO
1	A	138	SER

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	343/350 (98%)	321 (94%)	22 (6%)	17	14
1	B	348/350 (99%)	332 (95%)	16 (5%)	27	26
All	All	691/700 (99%)	653 (94%)	38 (6%)	23	19

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	GLN
1	A	22	THR
1	A	27	GLU

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Mol	Chain	Res	Type
1	A	38	VAL
1	A	55	LYS
1	A	142	GLU
1	A	145[A]	ASN
1	A	145[B]	ASN
1	A	158	SER
1	A	167	ARG
1	A	174	PHE
1	A	183	GLU
1	A	195	ASN
1	A	211	SER
1	A	221	PHE
1	A	259	LYS
1	A	282	LEU
1	A	283	GLN
1	A	287[A]	GLN
1	A	287[B]	GLN
1	A	293	ARG
1	A	369	LYS
1	B	30	ASP
1	B	38	VAL
1	B	55	LYS
1	B	127	THR
1	B	158	SER
1	B	174	PHE
1	B	183	GLU
1	B	221	PHE
1	B	259	LYS
1	B	276	LYS
1	B	277	GLU
1	B	283	GLN
1	B	294	ILE
1	B	346	LYS
1	B	362	SER
1	B	374	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	12	GLN
1	A	14	GLN
1	A	128	ASN

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Mol	Chain	Res	Type
1	A	195	ASN
1	B	128	ASN
1	B	352	ASN
1	B	370	GLN
1	B	376	ASN
1	B	379	HIS
1	B	406	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 ⓘ

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	B	502	-	16,16,16	2.89	6 (37%)	20,23,23	1.97	6 (30%)
3	PLP	A	502	-	16,16,16	2.97	3 (18%)	20,23,23	2.11	7 (35%)
2	PG4	B	501	-	6,6,12	0.66	0	5,5,11	0.88	0
2	PG4	A	501	-	6,6,12	0.61	0	5,5,11	0.68	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	B	502	-	-	4/8/8/8	0/1/1/1
3	PLP	A	502	-	-	3/8/8/8	0/1/1/1
2	PG4	B	501	-	-	3/4/4/10	-
2	PG4	A	501	-	-	2/4/4/10	-

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	502	PLP	C4-C5	7.55	1.51	1.42
3	B	502	PLP	C3-C2	7.23	1.48	1.40
3	A	502	PLP	C3-C2	6.16	1.47	1.40
3	A	502	PLP	C4-C3	5.95	1.49	1.40
3	B	502	PLP	C4-C5	5.55	1.49	1.42
3	B	502	PLP	C4-C3	5.10	1.48	1.40
3	B	502	PLP	P-O3P	-2.71	1.44	1.54
3	B	502	PLP	P-O2P	-2.57	1.44	1.54
3	B	502	PLP	P-O1P	-2.21	1.43	1.50

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	502	PLP	O4P-C5A-C5	5.32	119.48	109.35
3	A	502	PLP	O4P-C5A-C5	4.62	118.16	109.35
3	A	502	PLP	C4-C3-C2	-4.58	117.36	120.19
3	A	502	PLP	C6-N1-C2	3.49	125.63	119.17
3	B	502	PLP	C3-C4-C5	-3.13	115.86	118.26
3	B	502	PLP	C4-C3-C2	-3.03	118.31	120.19
3	B	502	PLP	O3P-P-O2P	2.66	117.80	107.64
3	B	502	PLP	O3P-P-O4P	-2.37	100.44	106.73
3	A	502	PLP	C2A-C2-N1	2.30	122.17	117.67
3	A	502	PLP	C5A-C5-C6	-2.30	115.60	119.37
3	A	502	PLP	C2A-C2-C3	-2.18	118.19	120.89
3	B	502	PLP	C3-C4-C4A	2.18	122.99	119.90
3	A	502	PLP	O2P-P-O1P	2.15	119.10	110.68

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	502	PLP	C4-C5-C5A-O4P
3	B	502	PLP	C6-C5-C5A-O4P
3	A	502	PLP	C4-C5-C5A-O4P
3	A	502	PLP	C6-C5-C5A-O4P
3	B	502	PLP	C3-C4-C4A-O4A
2	B	501	PG4	O1-C1-C2-O2
2	A	501	PG4	O3-C5-C6-O4
2	B	501	PG4	C4-C3-O2-C2
2	B	501	PG4	C1-C2-O2-C3
3	B	502	PLP	C5-C4-C4A-O4A
3	A	502	PLP	C5A-O4P-P-O2P
2	A	501	PG4	C3-C4-O3-C5

There are no ring outliers.

2 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	502	PLP	4	0
3	A	502	PLP	3	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	409/411 (99%)	0.86	54 (13%) 3 4	40, 60, 82, 99	0
1	B	410/411 (99%)	0.46	20 (4%) 29 35	36, 52, 74, 91	0
All	All	819/822 (99%)	0.66	74 (9%) 9 12	36, 56, 81, 99	0

All (74) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	179	GLU	5.7
1	A	217	PHE	5.5
1	A	27	GLU	4.8
1	B	138	SER	4.6
1	B	412	ILE	4.3
1	B	379	HIS	4.3
1	A	26	ARG	4.3
1	A	30	ASP	4.2
1	A	138	SER	3.8
1	A	28	ASP	3.8
1	A	181	ALA	3.5
1	B	30	ASP	3.5
1	B	9	GLU	3.4
1	B	277	GLU	3.4
1	A	346	LYS	3.4
1	A	29	PRO	3.4
1	B	375	VAL	3.4
1	A	281	ILE	3.3
1	B	27	GLU	3.2
1	A	31	PRO	3.1
1	A	410	THR	3.0
1	A	32	ARG	3.0
1	A	215	HIS	2.9
1	A	409	VAL	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	26	ARG	2.9
1	B	315	GLU	2.9
1	B	376	ASN	2.8
1	A	405	ILE	2.8
1	B	8	ALA	2.8
1	B	184	PHE	2.8
1	A	361[A]	PHE	2.7
1	A	380	ILE	2.7
1	B	279	GLU	2.7
1	A	19	PHE	2.6
1	A	96	LEU	2.6
1	A	199	ILE	2.6
1	A	239	TRP	2.6
1	A	178	LEU	2.6
1	A	401	VAL	2.6
1	A	350	THR	2.6
1	A	57	VAL	2.5
1	A	353	HIS	2.5
1	A	176	ASN	2.5
1	A	117	ALA	2.5
1	A	358	ILE	2.4
1	B	398	LEU	2.4
1	A	340	ALA	2.4
1	A	248	GLY	2.4
1	B	396	LYS	2.4
1	A	230	ALA	2.4
1	A	135	TYR	2.4
1	B	294	ILE	2.3
1	A	374	LEU	2.3
1	A	97	LYS	2.3
1	A	344	ALA	2.3
1	A	6	VAL	2.3
1	B	276	LYS	2.2
1	A	406	HIS	2.2
1	A	306	ILE	2.2
1	A	354	ILE	2.2
1	A	95	ALA	2.2
1	A	167	ARG	2.2
1	A	206	TRP	2.2
1	A	245	VAL	2.1
1	A	77	LEU	2.1
1	B	217	PHE	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	411	LYS	2.1
1	A	21	LEU	2.1
1	A	381	TYR	2.1
1	A	379	HIS	2.1
1	A	169	LEU	2.1
1	A	174	PHE	2.1
1	A	165	GLU	2.0
1	A	184	PHE	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. 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	PG4	A	501	7/13	0.70	0.16	67,70,74,76	0
2	PG4	B	501	7/13	0.80	0.21	51,59,74,83	0
3	PLP	B	502	16/16	0.97	0.16	37,62,71,71	10
3	PLP	A	502	16/16	0.97	0.22	39,73,97,105	10

6.5 Other polymers [i](#)

There are no such residues in this entry.