



Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 03:35 am BST

PDB ID : 1F3T
Title : CRYSTAL STRUCTURE OF TRYPANOSOMA BRUCEI ORNITHINE DE-CARBOXYLASE (ODC) COMPLEXED WITH PUTRESCINE, ODC'S RE-ACTION PRODUCT.
Authors : Jackson, L.K.; Brooks, H.B.; Osterman, A.L.; Goldsmith, E.J.; Phillips, M.A.
Deposited on : 2000-06-06
Resolution : 2.00 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

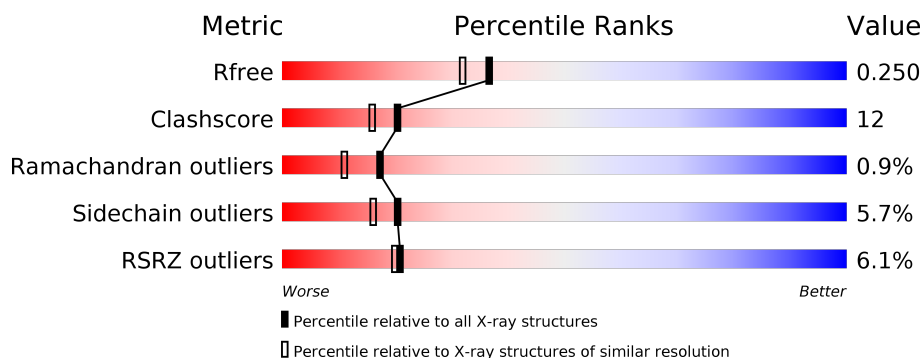
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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	425	<div> <div>7%</div> <div> <div></div> <div>68%</div> <div>17%</div> <div>•</div> <div>11%</div> </div> </div>
1	B	425	<div> <div>5%</div> <div> <div></div> <div>68%</div> <div>19%</div> <div>•</div> <div>10%</div> </div> </div>
1	C	425	<div> <div>6%</div> <div> <div></div> <div>67%</div> <div>17%</div> <div>•</div> <div>13%</div> </div> </div>
1	D	425	<div> <div>4%</div> <div> <div></div> <div>67%</div> <div>17%</div> <div>•</div> <div>14%</div> </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 12201 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 ORNITHINE DECARBOXYLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	378	Total	C	N	O	S	0	0	0
			2965	1905	497	547	16			
1	B	381	Total	C	N	O	S	0	0	0
			2987	1918	500	553	16			
1	C	370	Total	C	N	O	S	0	0	0
			2906	1866	486	538	16			
1	D	367	Total	C	N	O	S	0	0	0
			2879	1852	480	531	16			

There are 8 discrepancies between the modelled and reference sequences:

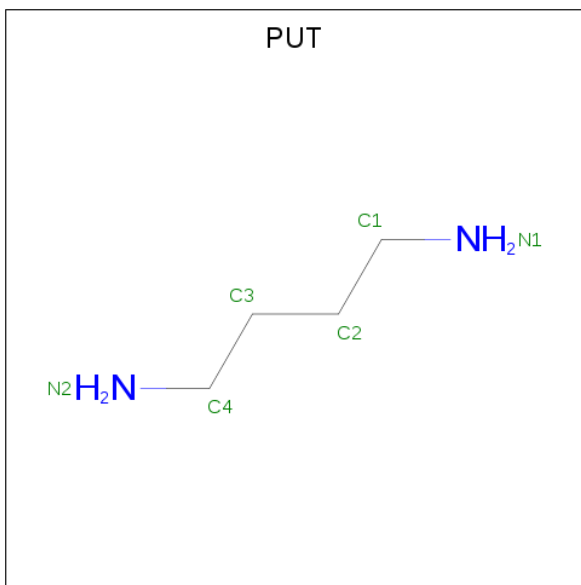
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	CLONING ARTIFACT	UNP P07805
A	2	ALA	-	CLONING ARTIFACT	UNP P07805
B	1	GLY	-	CLONING ARTIFACT	UNP P07805
B	2	ALA	-	CLONING ARTIFACT	UNP P07805
C	1	GLY	-	CLONING ARTIFACT	UNP P07805
C	2	ALA	-	CLONING ARTIFACT	UNP P07805
D	1	GLY	-	CLONING ARTIFACT	UNP P07805
D	2	ALA	-	CLONING ARTIFACT	UNP P07805

- Molecule 2 is PYRIDOXAL-5'-PHOSPHATE (three-letter code: PLP) (formula: C₈H₁₀NO₆P).



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

- Molecule 3 is 1,4-DIAMINOBTANE (three-letter code: PUT) (formula: $C_4H_{12}N_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C N 6 4 2	0	0
3	B	1	Total C N 6 4 2	0	0
3	C	1	Total C N 6 4 2	0	0
3	D	1	Total C N 6 4 2	0	0

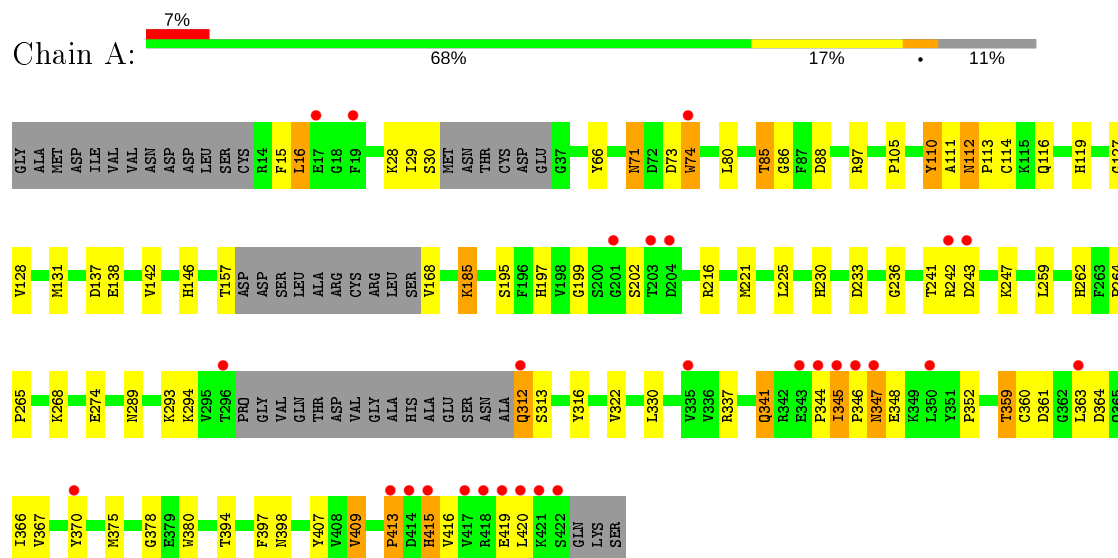
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	81	Total O 81 81	0	0
4	B	110	Total O 110 110	0	0
4	C	77	Total O 77 77	0	0
4	D	112	Total O 112 112	0	0

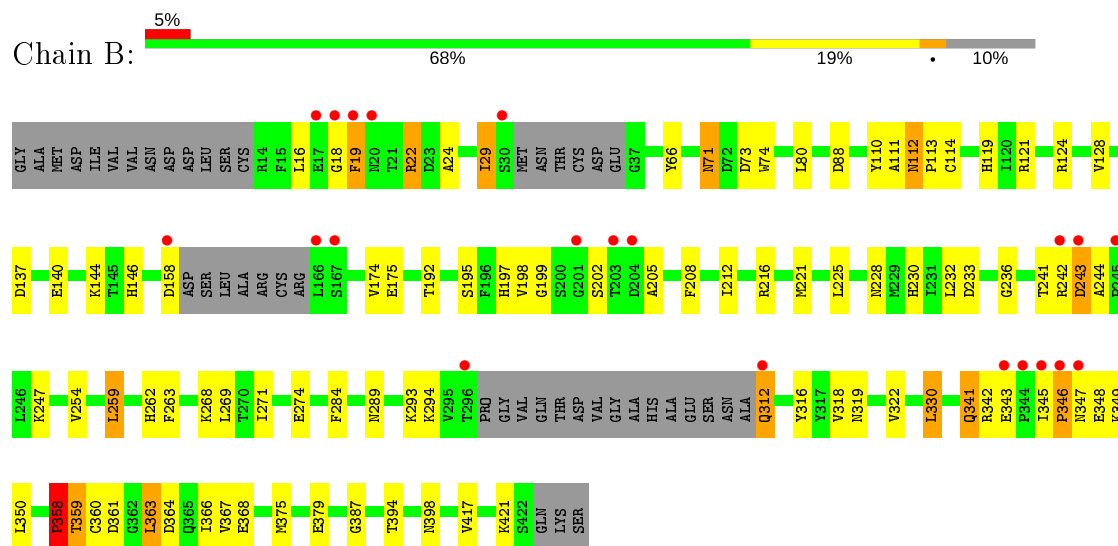
3 Residue-property plots

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: ORNITHINE DECARBOXYLASE

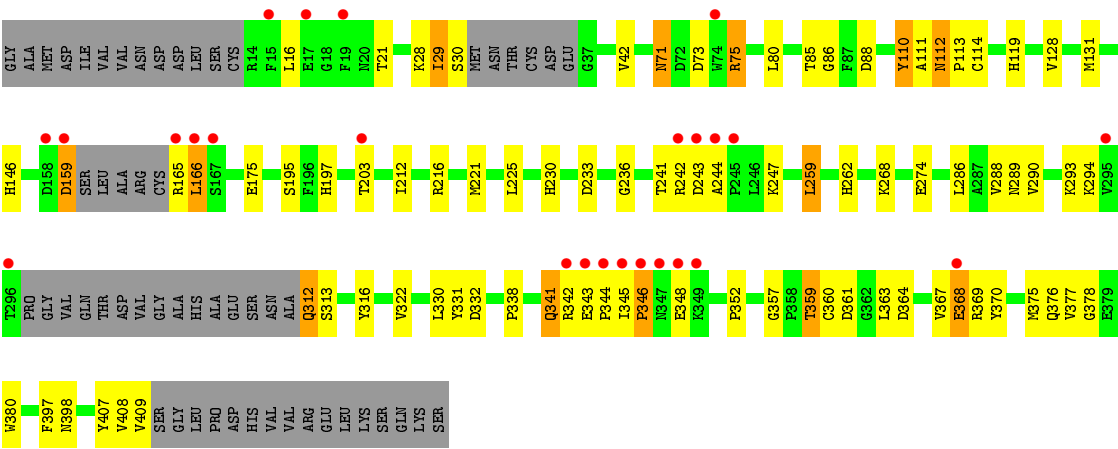


• Molecule 1: ORNITHINE DECARBOXYLASE

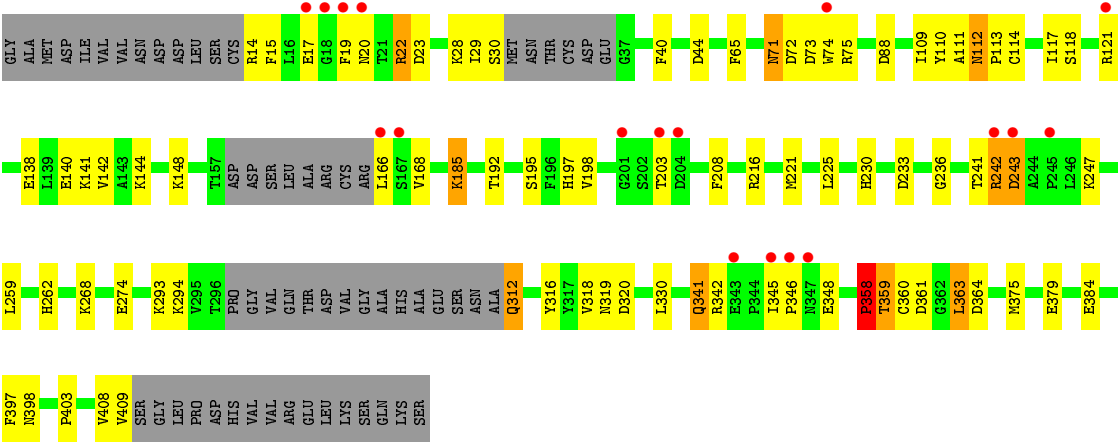


• Molecule 1: ORNITHINE DECARBOXYLASE





● Molecule 1: ORNITHINE DECARBOXYLASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	67.29 Å 151.60 Å 86.41 Å 90.00° 103.18° 90.00°	Depositor
Resolution (Å)	8.00 – 2.00 19.91 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.3 (8.00-2.00) 89.6 (19.91-1.90)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.53 (at 1.90 Å)	Xtriage
Refinement program	CNS 0.9	Depositor
R, R_{free}	0.237 , 0.280 0.215 , 0.250	Depositor DCC
R_{free} test set	2512 reflections (1.98%)	wwPDB-VP
Wilson B-factor (Å ²)	21.1	Xtriage
Anisotropy	0.585	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 54.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	12201	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 46.34 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.1616e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: PUT, 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	0.46	0/3036	0.69	0/4115
1	B	0.50	1/3058 (0.0%)	0.75	2/4145 (0.0%)
1	C	0.45	0/2975	0.67	0/4032
1	D	0.47	0/2948	0.72	1/3996 (0.0%)
All	All	0.47	1/12017 (0.0%)	0.71	3/16288 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	358	PRO	N-CA	5.00	1.55	1.47

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	358	PRO	CA-N-CD	-11.61	95.25	111.50
1	D	358	PRO	CA-N-CD	-8.53	99.56	111.50
1	B	18	GLY	N-CA-C	-7.26	94.94	113.10

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	345	ILE	Mainchain

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2965	0	2931	81	0
1	B	2987	0	2951	70	0
1	C	2906	0	2865	76	0
1	D	2879	0	2844	67	0
2	A	15	0	7	1	0
2	B	15	0	7	1	0
2	C	15	0	7	1	0
2	D	15	0	7	1	0
3	A	6	0	10	1	0
3	B	6	0	10	0	0
3	C	6	0	10	1	0
3	D	6	0	10	1	0
4	A	81	0	0	5	0
4	B	110	0	0	6	0
4	C	77	0	0	2	0
4	D	112	0	0	4	0
All	All	12201	0	11659	292	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 12.

All (292) 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:318:VAL:O	1:B:358:PRO:HD3	1.37	1.18
1:D:318:VAL:O	1:D:358:PRO:HD3	1.55	1.05
1:A:312:GLN:HE21	1:A:312:GLN:HA	1.25	1.01
1:A:409:VAL:HG21	1:A:419:GLU:OE1	1.62	0.98
1:C:359:THR:HG23	1:C:361:ASP:H	1.31	0.96
1:A:359:THR:HG23	1:A:361:ASP:H	1.34	0.93
1:B:16:LEU:CD1	1:B:24:ALA:HB1	1.99	0.91

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:197:HIS:HD2	1:B:236:GLY:H	1.20	0.87
1:A:15:PHE:CE2	1:A:420:LEU:HD22	2.10	0.86
1:D:197:HIS:HD2	1:D:236:GLY:H	1.20	0.86
1:B:22:ARG:HB2	1:B:22:ARG:NH1	1.92	0.84
1:B:312:GLN:HE21	1:B:312:GLN:HA	1.42	0.83
1:A:221:MET:O	1:A:225:LEU:HD23	1.80	0.81
1:D:312:GLN:HE21	1:D:312:GLN:HA	1.45	0.81
1:B:345:ILE:O	1:B:348:GLU:HG3	1.78	0.81
1:A:185:LYS:O	1:A:185:LYS:HE3	1.81	0.80
1:C:221:MET:O	1:C:225:LEU:HD23	1.80	0.80
1:C:197:HIS:HD2	1:C:236:GLY:H	1.29	0.79
1:C:341:GLN:HG2	1:C:380:TRP:HB2	1.64	0.79
1:D:359:THR:HG23	1:D:361:ASP:H	1.47	0.79
1:B:318:VAL:O	1:B:358:PRO:CD	2.28	0.76
1:A:197:HIS:HD2	1:A:236:GLY:H	1.33	0.76
1:B:322:VAL:HB	1:B:330:LEU:HD21	1.67	0.76
1:B:16:LEU:HD11	1:B:24:ALA:HB1	1.68	0.75
1:A:419:GLU:HG3	1:A:420:LEU:HD23	1.69	0.75
1:D:140:GLU:O	1:D:144:LYS:HD3	1.86	0.75
1:D:345:ILE:O	1:D:348:GLU:HG3	1.87	0.75
1:B:216:ARG:HD3	1:B:262:HIS:O	1.86	0.74
1:D:319:ASN:HA	1:D:358:PRO:HG3	1.71	0.73
1:A:345:ILE:O	1:A:348:GLU:HG3	1.89	0.73
1:C:316:TYR:CE2	1:C:375:MET:HB2	2.23	0.72
1:C:21:THR:HG23	1:C:42:VAL:HG11	1.70	0.72
1:B:22:ARG:HH11	1:B:22:ARG:HB2	1.54	0.72
1:C:312:GLN:HA	1:C:312:GLN:HE21	1.55	0.72
1:D:185:LYS:O	1:D:185:LYS:HE3	1.89	0.72
1:A:416:VAL:HA	1:A:419:GLU:OE1	1.90	0.71
1:A:341:GLN:HG2	1:A:380:TRP:HB2	1.72	0.71
1:C:322:VAL:HB	1:C:330:LEU:HD21	1.73	0.70
1:A:359:THR:CG2	1:A:361:ASP:H	2.04	0.70
1:D:216:ARG:HD3	1:D:262:HIS:O	1.91	0.69
1:C:112:ASN:C	1:C:112:ASN:HD22	1.95	0.69
1:B:197:HIS:CD2	1:B:236:GLY:H	2.08	0.68
1:A:195:SER:CB	1:A:233:ASP:HB3	2.24	0.68
1:B:140:GLU:O	1:B:144:LYS:HD3	1.94	0.68
1:D:197:HIS:CD2	1:D:236:GLY:H	2.09	0.68
1:A:416:VAL:O	1:A:420:LEU:HG	1.92	0.68
1:B:359:THR:HG23	1:B:361:ASP:H	1.58	0.67
1:C:80:LEU:HB3	1:C:85:THR:HG21	1.76	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:359:THR:CG2	1:D:361:ASP:H	2.08	0.66
1:B:319:ASN:HA	1:B:358:PRO:HG3	1.78	0.66
1:B:359:THR:HB	1:B:364:ASP:OD2	1.95	0.66
1:B:71:ASN:HD22	1:B:71:ASN:C	1.99	0.66
1:C:290:VAL:HG23	1:C:377:VAL:HA	1.77	0.66
1:B:359:THR:CG2	1:B:361:ASP:H	2.09	0.66
1:A:71:ASN:ND2	1:A:73:ASP:H	1.94	0.65
1:C:359:THR:CG2	1:C:361:ASP:H	2.08	0.65
1:D:195:SER:CB	1:D:233:ASP:HB3	2.25	0.65
1:B:16:LEU:HD13	1:B:24:ALA:HB1	1.79	0.65
1:A:312:GLN:NE2	1:A:312:GLN:HA	2.06	0.65
1:B:366:ILE:HG22	1:B:367:VAL:HG23	1.77	0.65
1:A:71:ASN:C	1:A:71:ASN:HD22	2.00	0.65
1:B:322:VAL:HB	1:B:330:LEU:CD2	2.27	0.65
1:D:44:ASP:HA	1:D:408:VAL:HG13	1.79	0.65
1:A:367:VAL:HG11	1:A:370:TYR:HB2	1.77	0.65
1:A:88:ASP:OD1	1:A:111:ALA:HB3	1.96	0.65
1:D:71:ASN:HD22	1:D:71:ASN:C	1.99	0.65
1:A:85:THR:HG23	1:A:86:GLY:O	1.96	0.64
1:A:71:ASN:HD22	1:A:73:ASP:H	1.45	0.64
1:A:316:TYR:CE2	1:A:375:MET:HB2	2.33	0.64
1:C:21:THR:HG23	1:C:42:VAL:CG1	2.27	0.64
1:A:74:TRP:HD1	1:A:97:ARG:NH2	1.96	0.64
1:C:197:HIS:CD2	1:C:236:GLY:H	2.15	0.63
1:D:185:LYS:HG2	4:D:723:HOH:O	1.98	0.63
1:A:312:GLN:HE21	1:A:312:GLN:CA	2.00	0.62
1:B:341:GLN:O	1:B:341:GLN:NE2	2.32	0.62
1:A:112:ASN:HD22	1:A:112:ASN:C	2.02	0.62
1:C:341:GLN:NE2	1:C:341:GLN:O	2.30	0.62
1:D:88:ASP:OD1	1:D:111:ALA:HB3	1.99	0.62
1:B:230:HIS:HB2	1:B:268:LYS:O	2.00	0.62
1:A:337:ARG:HB3	1:A:337:ARG:HH21	1.63	0.61
1:C:71:ASN:HD22	1:C:73:ASP:H	1.47	0.61
1:A:409:VAL:HG11	1:A:416:VAL:HG13	1.81	0.61
1:C:290:VAL:O	1:C:290:VAL:HG23	1.99	0.61
1:C:359:THR:HB	1:C:364:ASP:OD2	2.01	0.61
1:D:221:MET:O	1:D:225:LEU:HD23	2.01	0.61
1:C:293:LYS:HG2	1:C:294:LYS:N	2.15	0.60
1:B:146:HIS:HD2	4:B:756:HOH:O	1.84	0.60
1:A:322:VAL:HB	1:A:330:LEU:HD21	1.83	0.60
1:C:29:ILE:O	1:C:30:SER:HB3	2.01	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:71:ASN:HD22	1:C:71:ASN:C	2.05	0.60
1:D:44:ASP:HA	1:D:408:VAL:CG1	2.32	0.60
1:D:241:THR:O	1:D:243:ASP:N	2.34	0.59
1:B:312:GLN:HA	1:B:312:GLN:NE2	2.16	0.59
1:C:312:GLN:CA	1:C:312:GLN:HE21	2.15	0.59
1:A:216:ARG:HD3	1:A:262:HIS:O	2.02	0.59
1:D:241:THR:C	1:D:243:ASP:H	2.05	0.59
1:D:112:ASN:C	1:D:112:ASN:HD22	2.05	0.59
1:B:112:ASN:C	1:B:112:ASN:HD22	2.06	0.59
1:A:409:VAL:CG1	1:A:416:VAL:HG13	2.33	0.58
1:C:367:VAL:HG11	1:C:370:TYR:HB2	1.86	0.58
1:B:394:THR:HG21	4:B:799:HOH:O	2.03	0.58
1:B:293:LYS:HG2	1:B:294:LYS:N	2.16	0.58
1:D:312:GLN:NE2	1:D:312:GLN:HA	2.18	0.58
1:D:71:ASN:ND2	1:D:73:ASP:H	2.02	0.57
1:C:71:ASN:ND2	1:C:73:ASP:H	2.02	0.57
1:B:195:SER:CB	1:B:233:ASP:HB3	2.34	0.57
1:A:195:SER:HB3	1:A:233:ASP:HB3	1.86	0.57
1:C:341:GLN:C	1:C:341:GLN:HE21	2.07	0.57
1:D:22:ARG:HA	1:D:22:ARG:HH11	1.69	0.57
1:A:16:LEU:HD21	1:A:28:LYS:HD2	1.87	0.57
1:C:195:SER:CB	1:C:233:ASP:HB3	2.36	0.56
1:C:75:ARG:HH11	1:C:75:ARG:HG2	1.71	0.56
1:D:195:SER:HB3	1:D:233:ASP:HB3	1.86	0.56
1:B:88:ASP:OD1	1:B:111:ALA:HB3	2.06	0.56
1:C:29:ILE:O	1:C:29:ILE:HG22	2.05	0.56
1:C:343:GLU:HA	1:C:343:GLU:OE2	2.05	0.56
3:A:1002:PUT:H11	1:B:360:CYS:SG	2.46	0.56
1:C:146:HIS:HD2	4:C:546:HOH:O	1.88	0.55
1:D:341:GLN:O	1:D:341:GLN:NE2	2.39	0.55
1:A:29:ILE:O	1:A:30:SER:CB	2.54	0.55
1:D:17:GLU:OE2	1:D:17:GLU:HA	2.06	0.55
1:B:241:THR:O	1:B:243:ASP:N	2.38	0.55
1:C:397:PHE:O	1:C:398:ASN:HB2	2.05	0.55
1:B:360:CYS:HB3	4:B:689:HOH:O	2.07	0.55
1:C:322:VAL:HB	1:C:330:LEU:CD2	2.35	0.54
1:D:242:ARG:HG3	4:D:755:HOH:O	2.07	0.54
1:B:241:THR:C	1:B:243:ASP:H	2.12	0.54
1:D:230:HIS:HB2	1:D:268:LYS:O	2.07	0.54
1:C:241:THR:HB	1:C:243:ASP:OD1	2.08	0.53
1:D:29:ILE:O	1:D:30:SER:HB3	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:165:ARG:HB3	1:C:165:ARG:CZ	2.38	0.53
1:D:138:GLU:O	1:D:142:VAL:HG23	2.09	0.53
1:A:289:ASN:HD21	1:A:378:GLY:HA2	1.74	0.53
1:C:216:ARG:HD3	1:C:262:HIS:O	2.08	0.52
1:B:316:TYR:CE2	1:B:375:MET:HB2	2.44	0.52
1:B:71:ASN:ND2	1:B:73:ASP:H	2.07	0.52
1:D:20:ASN:HB2	1:D:23:ASP:HB2	1.92	0.52
1:D:192:THR:HB	1:D:230:HIS:CE1	2.45	0.52
1:A:367:VAL:CG1	1:A:370:TYR:HB2	2.39	0.52
1:C:112:ASN:ND2	1:C:112:ASN:C	2.63	0.51
1:A:157:THR:HG22	1:A:157:THR:O	2.09	0.51
1:C:338:PRO:HD3	1:C:367:VAL:HG21	1.93	0.51
1:B:319:ASN:HA	1:B:358:PRO:CG	2.41	0.51
1:A:360:CYS:HB3	4:A:568:HOH:O	2.10	0.51
1:C:112:ASN:ND2	1:C:114:CYS:H	2.09	0.51
1:C:29:ILE:O	1:C:30:SER:CB	2.59	0.51
1:A:366:ILE:HG22	1:A:367:VAL:HG23	1.94	0.50
1:A:409:VAL:O	1:A:409:VAL:HG22	2.11	0.50
1:A:322:VAL:HB	1:A:330:LEU:CD2	2.41	0.50
1:A:29:ILE:O	1:A:30:SER:HB3	2.12	0.50
1:B:221:MET:O	1:B:225:LEU:HD23	2.10	0.50
1:C:289:ASN:HD21	1:C:378:GLY:HA2	1.76	0.50
1:D:359:THR:HB	1:D:364:ASP:OD2	2.11	0.50
1:D:293:LYS:HG2	1:D:294:LYS:N	2.27	0.50
1:D:71:ASN:HD22	1:D:73:ASP:H	1.59	0.50
1:A:146:HIS:HD2	4:A:505:HOH:O	1.95	0.50
1:C:241:THR:C	1:C:243:ASP:H	2.15	0.50
1:B:66:TYR:CG	1:B:80:LEU:HD13	2.47	0.49
1:A:337:ARG:HB3	1:A:337:ARG:NH2	2.25	0.49
1:B:360:CYS:H	1:B:398:ASN:HD21	1.59	0.49
1:B:128:VAL:O	1:B:146:HIS:HE1	1.95	0.49
1:C:128:VAL:O	1:C:146:HIS:HE1	1.95	0.49
1:A:241:THR:C	1:A:243:ASP:H	2.15	0.49
1:A:359:THR:HB	1:A:364:ASP:OD2	2.12	0.49
1:D:168:VAL:HG12	4:D:859:HOH:O	2.12	0.49
1:A:341:GLN:HG3	4:A:576:HOH:O	2.12	0.49
1:C:322:VAL:CG2	1:C:364:ASP:HA	2.42	0.49
1:C:332:ASP:OD2	3:C:1006:PUT:N2	2.45	0.49
1:D:112:ASN:ND2	1:D:114:CYS:H	2.10	0.49
1:C:159:ASP:N	1:C:159:ASP:OD2	2.46	0.49
1:A:112:ASN:HD22	1:A:113:PRO:N	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:128:VAL:O	1:A:146:HIS:HE1	1.96	0.49
1:C:16:LEU:HD21	1:C:28:LYS:HD2	1.95	0.49
1:D:117:ILE:HG22	1:D:121:ARG:NH2	2.28	0.49
1:C:313:SER:HB3	1:C:352:PRO:HG2	1.95	0.48
1:C:322:VAL:HG22	1:C:364:ASP:HA	1.93	0.48
1:C:312:GLN:HA	1:C:312:GLN:NE2	2.26	0.48
1:B:112:ASN:HD22	1:B:113:PRO:N	2.12	0.48
1:C:119:HIS:HE1	4:D:913:HOH:O	1.96	0.48
1:B:175:GLU:CD	1:B:175:GLU:H	2.15	0.48
1:B:202:SER:HB2	4:B:878:HOH:O	2.12	0.48
1:A:330:LEU:N	1:A:330:LEU:HD22	2.28	0.48
1:A:394:THR:HG21	4:A:650:HOH:O	2.13	0.47
1:B:228:ASN:HB2	4:B:678:HOH:O	2.13	0.47
1:A:112:ASN:ND2	1:A:112:ASN:C	2.67	0.47
1:B:29:ILE:HD12	1:B:341:GLN:OE1	2.15	0.47
1:C:88:ASP:OD1	1:C:111:ALA:HB3	2.15	0.47
1:C:75:ARG:HB2	1:C:407:TYR:CE2	2.49	0.47
1:A:29:ILE:O	1:A:29:ILE:HG22	2.15	0.47
1:B:71:ASN:HD22	1:B:73:ASP:H	1.62	0.47
1:D:360:CYS:H	1:D:398:ASN:HD21	1.62	0.47
1:A:112:ASN:ND2	1:A:114:CYS:H	2.13	0.47
1:C:112:ASN:HD22	1:C:113:PRO:N	2.13	0.47
1:C:344:PRO:C	1:C:345:ILE:HD12	2.36	0.47
1:D:316:TYR:CE2	1:D:375:MET:HB2	2.50	0.46
1:A:195:SER:HB2	1:A:233:ASP:HB3	1.95	0.46
1:C:341:GLN:HG3	4:C:575:HOH:O	2.14	0.46
1:C:85:THR:HG23	1:C:86:GLY:O	2.16	0.46
1:A:241:THR:O	1:A:243:ASP:N	2.41	0.46
1:C:290:VAL:O	1:C:290:VAL:CG2	2.62	0.46
1:D:28:LYS:HD3	1:D:40:PHE:CE1	2.50	0.46
1:A:110:TYR:HB3	1:A:131:MET:HG2	1.98	0.46
1:C:110:TYR:HB3	1:C:131:MET:HG2	1.98	0.46
1:C:290:VAL:HG21	1:C:376:GLN:O	2.16	0.46
1:A:359:THR:HG23	1:A:360:CYS:N	2.30	0.46
1:C:75:ARG:HH11	1:C:75:ARG:CG	2.26	0.45
1:D:185:LYS:HA	1:D:185:LYS:HE3	1.97	0.45
1:D:320:ASP:OD2	1:D:403:PRO:HG2	2.15	0.45
1:D:274:GLU:O	2:D:1007:PLP:H6	2.17	0.45
1:A:293:LYS:HG2	1:A:294:LYS:N	2.30	0.45
1:A:313:SER:HB3	1:A:352:PRO:HG2	1.98	0.45
1:B:174:VAL:HG23	1:B:175:GLU:N	2.30	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:230:HIS:HB2	1:A:268:LYS:O	2.17	0.45
1:A:15:PHE:CD2	1:A:420:LEU:HD22	2.49	0.45
1:C:212:ILE:HD13	1:C:259:LEU:HD13	1.99	0.45
1:B:232:LEU:HD23	1:B:271:ILE:HD12	1.98	0.45
1:C:230:HIS:HB2	1:C:268:LYS:O	2.16	0.45
1:D:195:SER:HB2	1:D:233:ASP:HB3	1.98	0.45
1:A:409:VAL:CG2	1:A:419:GLU:OE1	2.50	0.45
1:D:112:ASN:HD22	1:D:113:PRO:N	2.15	0.45
1:D:14:ARG:HH11	1:D:14:ARG:HG3	1.82	0.45
1:B:198:VAL:HG23	1:B:199:GLY:O	2.17	0.44
1:C:166:LEU:HD13	1:C:166:LEU:N	2.32	0.44
1:C:359:THR:HG23	1:C:360:CYS:N	2.32	0.44
1:D:312:GLN:CA	1:D:312:GLN:HE21	2.15	0.44
1:D:363:LEU:HA	1:D:363:LEU:HD12	1.78	0.44
1:B:345:ILE:H	1:B:348:GLU:CD	2.21	0.44
1:C:331:TYR:CE2	1:D:330:LEU:HD12	2.52	0.44
1:D:198:VAL:HG21	1:D:208:PHE:CE2	2.53	0.44
1:B:198:VAL:HG21	1:B:208:PHE:CE2	2.53	0.44
1:B:192:THR:HB	1:B:230:HIS:CE1	2.53	0.44
1:C:195:SER:HB3	1:C:233:ASP:HB3	1.99	0.44
1:D:15:PHE:HZ	1:D:75:ARG:HE	1.64	0.44
1:D:71:ASN:HD22	1:D:72:ASP:N	2.16	0.44
1:B:349:LYS:HD3	1:B:350:LEU:H	1.83	0.43
1:B:417:VAL:CG1	1:B:421:LYS:HE3	2.47	0.43
1:A:397:PHE:O	1:A:398:ASN:HB2	2.18	0.43
1:C:357:GLY:HA3	1:C:364:ASP:OD1	2.18	0.43
1:A:105:PRO:HB2	1:A:127:GLY:O	2.19	0.43
1:C:225:LEU:HD22	1:C:225:LEU:N	2.34	0.43
1:C:360:CYS:SG	3:D:1008:PUT:H11	2.58	0.43
1:A:197:HIS:CE1	1:A:199:GLY:HA2	2.53	0.43
1:B:19:PHE:CD2	1:B:19:PHE:C	2.90	0.43
1:B:212:ILE:HD13	1:B:259:LEU:HD13	2.00	0.43
1:A:116:GLN:NE2	1:B:319:ASN:HD22	2.17	0.43
1:A:202:SER:HB2	4:A:733:HOH:O	2.17	0.43
1:A:66:TYR:CG	1:A:80:LEU:HD13	2.53	0.43
1:A:312:GLN:NE2	1:A:312:GLN:CA	2.70	0.43
1:A:347:ASN:C	1:A:347:ASN:HD22	2.22	0.43
1:A:409:VAL:HG11	1:A:419:GLU:OE1	2.19	0.43
1:A:168:VAL:O	1:A:168:VAL:HG22	2.18	0.43
1:A:344:PRO:C	1:A:345:ILE:HD12	2.39	0.43
1:B:263:PHE:HB3	1:B:269:LEU:HD21	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:117:ILE:HD11	1:D:141:LYS:HD3	2.01	0.42
1:A:185:LYS:HA	1:A:185:LYS:HD2	1.84	0.42
1:A:413:PRO:O	1:A:415:HIS:N	2.42	0.42
1:B:195:SER:HB3	1:B:233:ASP:HB3	2.01	0.42
1:B:289:ASN:ND2	4:B:673:HOH:O	2.52	0.42
1:A:138:GLU:O	1:A:142:VAL:HG23	2.20	0.42
1:B:121:ARG:HG3	1:B:124:ARG:NH2	2.34	0.42
1:B:205:ALA:HB3	1:B:254:VAL:HG21	2.02	0.42
1:B:342:ARG:HD2	1:B:379:GLU:OE2	2.20	0.42
1:D:342:ARG:HD2	1:D:379:GLU:OE2	2.20	0.42
1:A:195:SER:HA	1:A:233:ASP:O	2.19	0.42
1:A:15:PHE:HA	1:A:407:TYR:O	2.19	0.42
1:D:19:PHE:CE2	1:D:409:VAL:O	2.73	0.42
1:D:19:PHE:HD1	1:D:19:PHE:O	2.02	0.42
1:D:22:ARG:NH2	1:D:384:GLU:OE2	2.52	0.41
1:A:264:PRO:HA	1:A:265:PRO:HD3	1.95	0.41
1:C:75:ARG:NH1	1:C:75:ARG:CG	2.81	0.41
1:B:363:LEU:HA	1:B:363:LEU:HD12	1.81	0.41
1:A:274:GLU:O	2:A:1001:PLP:H6	2.21	0.41
1:D:241:THR:C	1:D:243:ASP:N	2.73	0.41
1:D:397:PHE:O	1:D:398:ASN:HB2	2.20	0.41
1:B:112:ASN:C	1:B:112:ASN:ND2	2.72	0.41
1:B:112:ASN:ND2	1:B:114:CYS:H	2.18	0.41
1:D:166:LEU:HD23	1:D:197:HIS:ND1	2.35	0.41
1:C:368:GLU:OE1	1:C:369:ARG:HG3	2.19	0.41
1:C:75:ARG:HD2	1:C:75:ARG:N	2.36	0.41
1:D:65:PHE:CD2	1:D:109:ILE:HG13	2.56	0.41
1:D:148:LYS:HE2	1:D:148:LYS:HB3	1.86	0.41
1:C:274:GLU:O	2:C:1005:PLP:H6	2.20	0.41
1:C:286:LEU:HG	1:C:288:VAL:HG13	2.03	0.41
1:A:71:ASN:C	1:A:71:ASN:ND2	2.71	0.41
1:B:274:GLU:O	2:B:1003:PLP:H6	2.21	0.41
1:C:342:ARG:NH2	1:C:348:GLU:OE1	2.54	0.41
1:C:367:VAL:CG1	1:C:370:TYR:HB2	2.50	0.41
1:D:29:ILE:O	1:D:30:SER:CB	2.69	0.41
1:B:22:ARG:CZ	1:B:22:ARG:HB2	2.49	0.40
1:D:112:ASN:C	1:D:112:ASN:ND2	2.73	0.40
1:D:118:SER:HA	1:D:121:ARG:NH1	2.36	0.40
1:B:284:PHE:HB2	1:B:387:GLY:CA	2.51	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	370/425 (87%)	351 (95%)	16 (4%)	3 (1%)	19	13
1	B	373/425 (88%)	356 (95%)	13 (4%)	4 (1%)	14	8
1	C	362/425 (85%)	341 (94%)	17 (5%)	4 (1%)	14	8
1	D	359/425 (84%)	345 (96%)	12 (3%)	2 (1%)	25	19
All	All	1464/1700 (86%)	1393 (95%)	58 (4%)	13 (1%)	17	11

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	242	ARG
1	A	346	PRO
1	B	346	PRO
1	C	346	PRO
1	D	242	ARG
1	B	242	ARG
1	C	242	ARG
1	B	29	ILE
1	B	244	ALA
1	C	244	ALA
1	A	413	PRO
1	C	29	ILE
1	D	346	PRO

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	323/362 (89%)	305 (94%)	18 (6%)	21	17
1	B	326/362 (90%)	304 (93%)	22 (7%)	16	11
1	C	316/362 (87%)	298 (94%)	18 (6%)	20	16
1	D	313/362 (86%)	298 (95%)	15 (5%)	25	22
All	All	1278/1448 (88%)	1205 (94%)	73 (6%)	20	16

All (73) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	LEU
1	A	71	ASN
1	A	74	TRP
1	A	85	THR
1	A	110	TYR
1	A	112	ASN
1	A	119	HIS
1	A	137	ASP
1	A	185	LYS
1	A	247	LYS
1	A	259	LEU
1	A	312	GLN
1	A	341	GLN
1	A	347	ASN
1	A	359	THR
1	A	363	LEU
1	A	409	VAL
1	A	415	HIS
1	B	19	PHE
1	B	22	ARG
1	B	71	ASN
1	B	74	TRP
1	B	110	TYR
1	B	112	ASN
1	B	119	HIS
1	B	137	ASP
1	B	158	ASP
1	B	243	ASP
1	B	247	LYS
1	B	259	LEU
1	B	312	GLN
1	B	330	LEU
1	B	341	GLN

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Mol	Chain	Res	Type
1	B	343	GLU
1	B	346	PRO
1	B	347	ASN
1	B	358	PRO
1	B	359	THR
1	B	363	LEU
1	B	368	GLU
1	C	71	ASN
1	C	75	ARG
1	C	110	TYR
1	C	112	ASN
1	C	159	ASP
1	C	166	LEU
1	C	175	GLU
1	C	203	THR
1	C	247	LYS
1	C	259	LEU
1	C	312	GLN
1	C	341	GLN
1	C	346	PRO
1	C	359	THR
1	C	363	LEU
1	C	368	GLU
1	C	408	VAL
1	C	409	VAL
1	D	22	ARG
1	D	71	ASN
1	D	74	TRP
1	D	110	TYR
1	D	112	ASN
1	D	185	LYS
1	D	203	THR
1	D	243	ASP
1	D	247	LYS
1	D	259	LEU
1	D	312	GLN
1	D	341	GLN
1	D	358	PRO
1	D	359	THR
1	D	363	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (44) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	71	ASN
1	A	96	GLN
1	A	112	ASN
1	A	116	GLN
1	A	119	HIS
1	A	146	HIS
1	A	197	HIS
1	A	210	GLN
1	A	230	HIS
1	A	256	ASN
1	A	289	ASN
1	A	312	GLN
1	A	341	GLN
1	B	71	ASN
1	B	112	ASN
1	B	116	GLN
1	B	119	HIS
1	B	146	HIS
1	B	197	HIS
1	B	256	ASN
1	B	289	ASN
1	B	312	GLN
1	B	401	GLN
1	C	71	ASN
1	C	112	ASN
1	C	116	GLN
1	C	119	HIS
1	C	146	HIS
1	C	197	HIS
1	C	210	GLN
1	C	230	HIS
1	C	256	ASN
1	C	262	HIS
1	C	289	ASN
1	C	312	GLN
1	D	71	ASN
1	D	112	ASN
1	D	116	GLN
1	D	146	HIS
1	D	197	HIS
1	D	256	ASN
1	D	289	ASN
1	D	312	GLN

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Mol	Chain	Res	Type
1	D	401	GLN

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 ⓘ

8 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	PUT	C	1006	2	5,5,5	0.35	0	4,4,4	0.65	0
2	PLP	C	1005	3	15,15,16	2.76	4 (26%)	20,22,23	1.32	4 (20%)
2	PLP	A	1001	3	15,15,16	2.36	5 (33%)	20,22,23	1.24	4 (20%)
2	PLP	B	1003	3	15,15,16	2.02	3 (20%)	20,22,23	1.10	3 (15%)
2	PLP	D	1007	3	15,15,16	2.19	4 (26%)	20,22,23	1.58	2 (10%)
3	PUT	D	1008	2	5,5,5	0.33	0	4,4,4	0.55	0
3	PUT	B	1004	2	5,5,5	0.40	0	4,4,4	0.51	0
3	PUT	A	1002	2	5,5,5	0.50	0	4,4,4	0.48	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	PUT	C	1006	2	-	0/3/3/3	-
2	PLP	C	1005	3	-	0/6/6/8	0/1/1/1
2	PLP	A	1001	3	-	0/6/6/8	0/1/1/1
2	PLP	B	1003	3	-	0/6/6/8	0/1/1/1
2	PLP	D	1007	3	-	0/6/6/8	0/1/1/1
3	PUT	D	1008	2	-	0/3/3/3	-
3	PUT	B	1004	2	-	0/3/3/3	-
3	PUT	A	1002	2	-	0/3/3/3	-

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	1005	PLP	C5-C4	7.71	1.49	1.40
2	A	1001	PLP	C5-C4	7.09	1.48	1.40
2	D	1007	PLP	C5-C4	5.70	1.46	1.40
2	B	1003	PLP	C5-C4	5.34	1.46	1.40
2	C	1005	PLP	C3-C2	4.61	1.45	1.40
2	C	1005	PLP	C2A-C2	3.58	1.56	1.50
2	D	1007	PLP	C6-C5	3.52	1.45	1.37
2	A	1001	PLP	C6-C5	3.11	1.44	1.37
2	B	1003	PLP	C6-C5	3.11	1.44	1.37
2	D	1007	PLP	C3-C2	3.10	1.44	1.40
2	D	1007	PLP	C4A-C4	3.02	1.57	1.51
2	C	1005	PLP	C6-C5	2.92	1.43	1.37
2	B	1003	PLP	C3-C2	2.87	1.43	1.40
2	A	1001	PLP	C3-C2	2.76	1.43	1.40
2	A	1001	PLP	C2A-C2	2.62	1.54	1.50
2	A	1001	PLP	C4A-C4	2.34	1.56	1.51

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1007	PLP	C4A-C4-C5	4.55	125.62	120.94
2	D	1007	PLP	C4A-C4-C3	-2.90	115.58	120.50
2	C	1005	PLP	O3-C3-C2	2.40	122.72	117.49
2	B	1003	PLP	C6-N1-C2	2.32	123.47	119.17
2	C	1005	PLP	C5-C6-N1	-2.26	120.05	123.82
2	A	1001	PLP	O4P-P-O1P	2.25	112.80	106.47
2	A	1001	PLP	C6-N1-C2	2.15	123.14	119.17
2	C	1005	PLP	O4P-P-O1P	2.08	112.32	106.47
2	A	1001	PLP	C5-C6-N1	-2.08	120.36	123.82

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1003	PLP	C5-C6-N1	-2.05	120.41	123.82
2	A	1001	PLP	C5A-C5-C6	-2.04	116.01	119.37
2	C	1005	PLP	C6-N1-C2	2.04	122.94	119.17
2	B	1003	PLP	O4P-P-O1P	2.03	112.16	106.47

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

7 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	1006	PUT	1	0
2	C	1005	PLP	1	0
2	A	1001	PLP	1	0
2	B	1003	PLP	1	0
2	D	1007	PLP	1	0
3	D	1008	PUT	1	0
3	A	1002	PUT	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	378/425 (88%)	0.04	28 (7%) 14 13	13, 25, 64, 83	0
1	B	381/425 (89%)	0.02	21 (5%) 25 24	10, 23, 55, 85	0
1	C	370/425 (87%)	0.09	25 (6%) 17 16	13, 26, 63, 88	0
1	D	367/425 (86%)	0.01	18 (4%) 29 28	11, 25, 54, 81	0
All	All	1496/1700 (88%)	0.04	92 (6%) 21 20	10, 25, 60, 88	0

All (92) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	19	PHE	13.6
1	D	19	PHE	9.1
1	B	18	GLY	8.2
1	A	347	ASN	6.8
1	D	166	LEU	6.0
1	D	167	SER	5.6
1	C	74	TRP	5.4
1	B	203	THR	5.3
1	A	345	ILE	5.2
1	B	346	PRO	5.0
1	B	30	SER	5.0
1	A	346	PRO	5.0
1	A	242	ARG	5.0
1	B	347	ASN	5.0
1	B	158	ASP	4.9
1	C	159	ASP	4.6
1	D	74	TRP	4.5
1	A	243	ASP	4.4
1	C	17	GLU	4.4
1	C	343	GLU	4.4
1	C	158	ASP	4.4

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Mol	Chain	Res	Type	RSRZ
1	C	344	PRO	4.3
1	D	242	ARG	4.3
1	A	415	HIS	4.3
1	C	243	ASP	4.3
1	A	421	LYS	4.2
1	C	165	ARG	4.2
1	C	347	ASN	4.1
1	B	345	ILE	3.8
1	D	243	ASP	3.8
1	A	420	LEU	3.7
1	D	18	GLY	3.7
1	A	418	ARG	3.6
1	C	19	PHE	3.6
1	B	201	GLY	3.5
1	C	296	THR	3.5
1	C	203	THR	3.4
1	C	244	ALA	3.4
1	A	344	PRO	3.3
1	B	17	GLU	3.3
1	D	203	THR	3.2
1	A	203	THR	3.2
1	C	345	ILE	3.2
1	D	201	GLY	3.1
1	D	204	ASP	3.1
1	C	346	PRO	3.1
1	A	74	TRP	3.0
1	D	345	ILE	3.0
1	A	417	VAL	2.9
1	B	343	GLU	2.9
1	B	245	PRO	2.9
1	B	242	ARG	2.9
1	A	419	GLU	2.8
1	B	167	SER	2.8
1	B	20	ASN	2.8
1	B	204	ASP	2.8
1	A	413	PRO	2.7
1	C	242	ARG	2.7
1	B	166	LEU	2.7
1	D	20	ASN	2.7
1	B	296	THR	2.6
1	C	368	GLU	2.6
1	A	422	SER	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	204	ASP	2.5
1	A	414	ASP	2.5
1	A	335	VAL	2.5
1	D	346	PRO	2.5
1	A	17	GLU	2.4
1	B	312	GLN	2.4
1	C	349	LYS	2.4
1	C	295	VAL	2.4
1	A	296	THR	2.4
1	B	243	ASP	2.4
1	D	17	GLU	2.3
1	B	344	PRO	2.3
1	D	245	PRO	2.3
1	C	245	PRO	2.2
1	D	343	GLU	2.2
1	D	121	ARG	2.2
1	C	166	LEU	2.2
1	A	350	LEU	2.1
1	D	347	ASN	2.1
1	A	201	GLY	2.1
1	A	19	PHE	2.1
1	C	15	PHE	2.1
1	A	312	GLN	2.1
1	A	363	LEU	2.1
1	C	167	SER	2.1
1	A	343	GLU	2.1
1	C	342	ARG	2.1
1	C	348	GLU	2.1
1	A	370	TYR	2.0

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 [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(\AA^2)	Q<0.9
3	PUT	D	1008	6/6	0.61	0.25	43,50,54,55	0
3	PUT	A	1002	6/6	0.65	0.32	40,46,48,51	0
3	PUT	B	1004	6/6	0.68	0.33	44,51,55,59	0
3	PUT	C	1006	6/6	0.76	0.27	43,49,53,53	0
2	PLP	C	1005	15/16	0.97	0.08	16,22,31,34	0
2	PLP	B	1003	15/16	0.97	0.10	14,22,30,34	0
2	PLP	D	1007	15/16	0.97	0.08	13,23,28,34	0
2	PLP	A	1001	15/16	0.98	0.09	15,23,29,31	0

6.5 Other polymers [i](#)

There are no such residues in this entry.