



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 11:40 am BST

PDB ID : 4OBU
Title : Ruminococcus gnavus tryptophan decarboxylase RUMGNA_01526 (apo)
Authors : Van Benschoten, A.H.; Fraser, J.S.
Deposited on : 2014-01-07
Resolution : 2.80 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

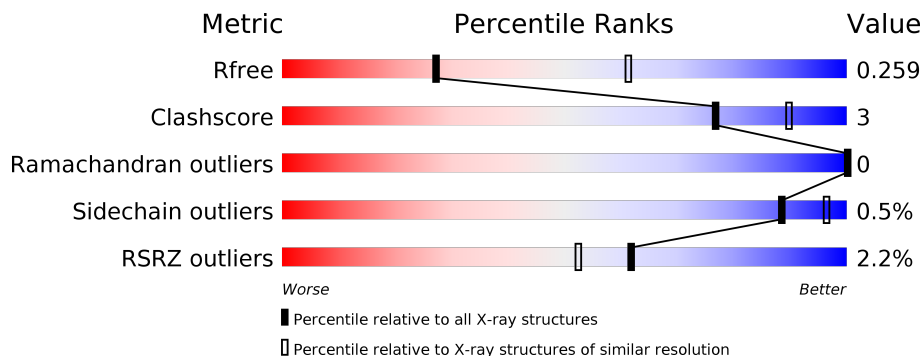
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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	490	 2% 87% 7% 6%
1	B	490	 2% 86% 8% 6%
1	C	490	 2% 84% 11% 6%
1	E	490	 2% 87% 7% 6%
1	F	490	 2% 86% 9% 6%
1	G	490	 2% 87% 8% 6%

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Mol	Chain	Length	Quality of chain
1	H	490	 <p>3% 86% 9% 6%</p>
1	U	490	 <p>2% 88% 6% 6%</p>

2 Entry composition

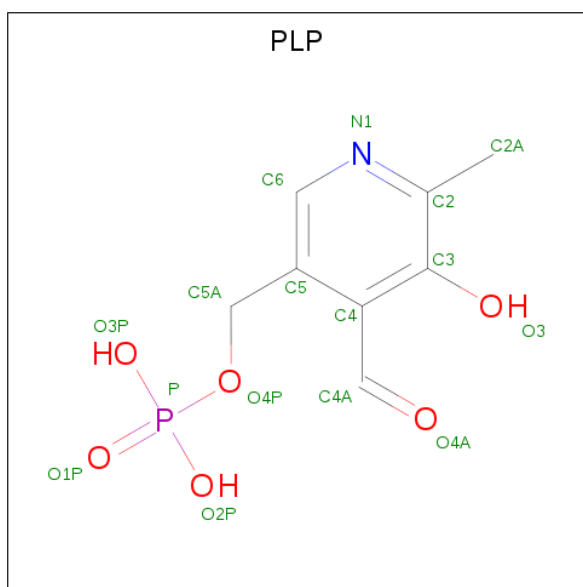
There are 3 unique types of molecules in this entry. The entry contains 29598 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Pyridoxal-dependent decarboxylase domain protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	E	462	3633	2310	618	687	18	0	0	0
1	A	462	3633	2310	618	687	18	0	0	0
1	U	462	3633	2310	618	687	18	0	0	0
1	G	462	3633	2310	618	687	18	0	0	0
1	H	462	3633	2310	618	687	18	0	0	0
1	F	462	3633	2310	618	687	18	0	0	0
1	B	462	3633	2310	618	687	18	0	0	0
1	C	462	3633	2310	618	687	18	0	0	0

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



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	E	1	Total 15	8	1	5	1	0	0
2	A	1	Total 15	8	1	5	1	0	0
2	U	1	Total 15	8	1	5	1	0	0
2	G	1	Total 15	8	1	5	1	0	0
2	H	1	Total 15	8	1	5	1	0	0
2	F	1	Total 15	8	1	5	1	0	0
2	B	1	Total 15	8	1	5	1	0	0
2	C	1	Total 15	8	1	5	1	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	E	56	Total 56	O 56	0	0
3	A	49	Total 49	O 49	0	0
3	U	57	Total 57	O 57	0	0
3	G	46	Total 46	O 46	0	0

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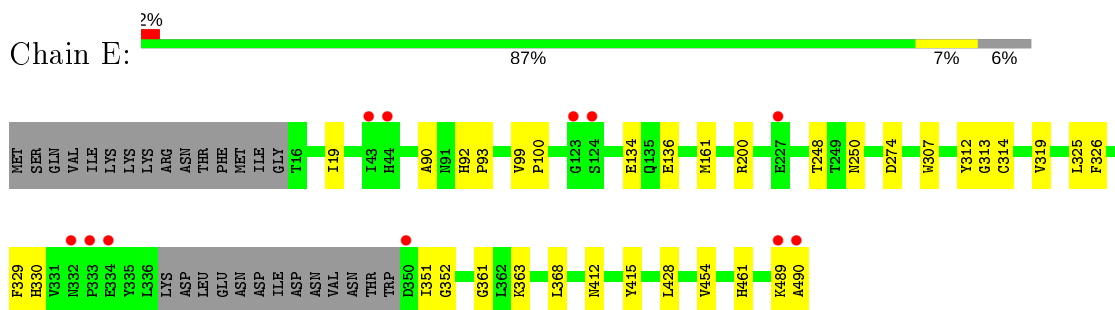
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	H	52	Total 52	O 52	0	0
3	F	73	Total 73	O 73	0	0
3	B	45	Total 45	O 45	0	0
3	C	36	Total 36	O 36	0	0

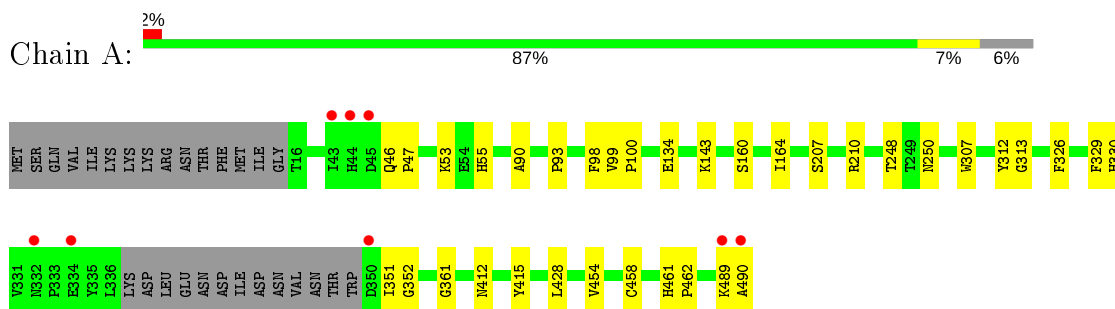
3 Residue-property plots [i](#)

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 ($RSRZ > 2$). 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.

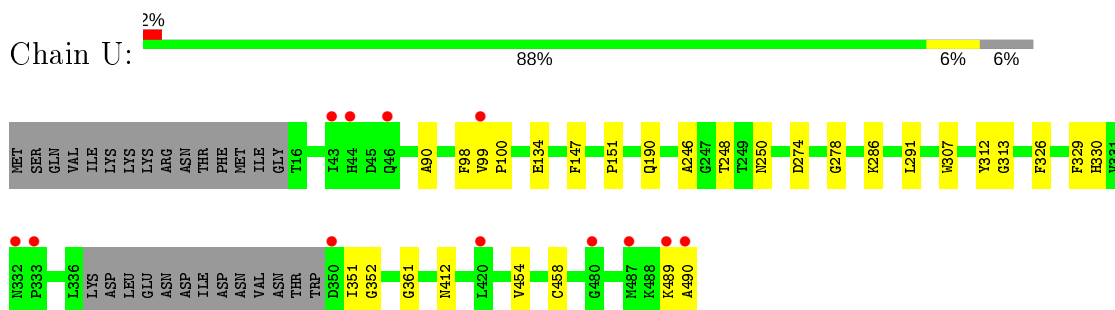
- Molecule 1: Pyridoxal-dependent decarboxylase domain protein



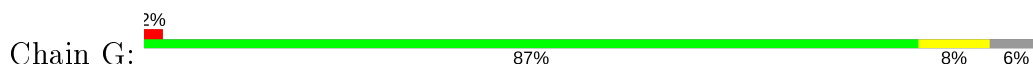
- Molecule 1: Pyridoxal-dependent decarboxylase domain protein

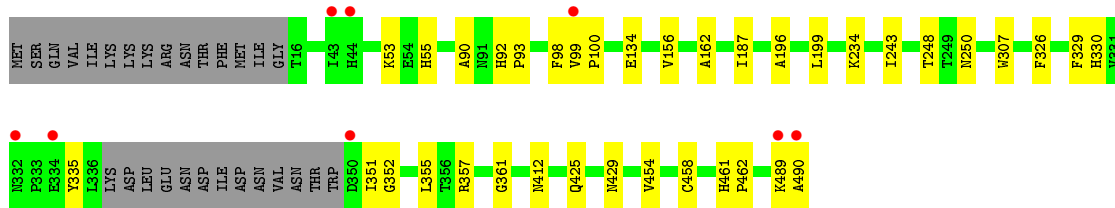


- Molecule 1: Pyridoxal-dependent decarboxylase domain protein

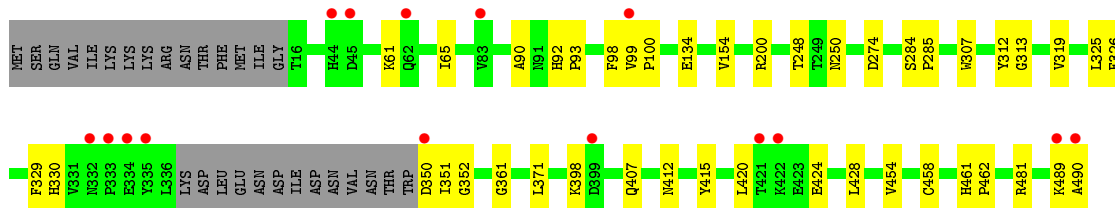
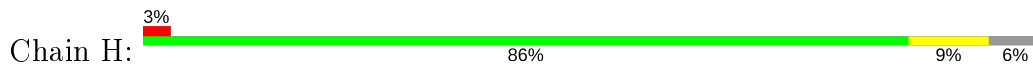


- Molecule 1: Pyridoxal-dependent decarboxylase domain protein

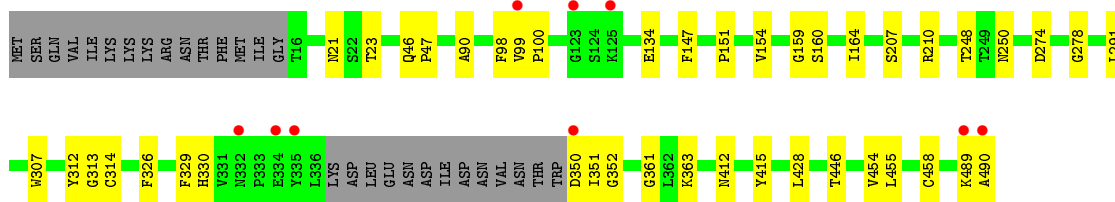
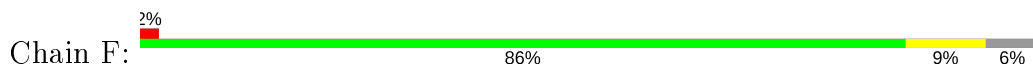




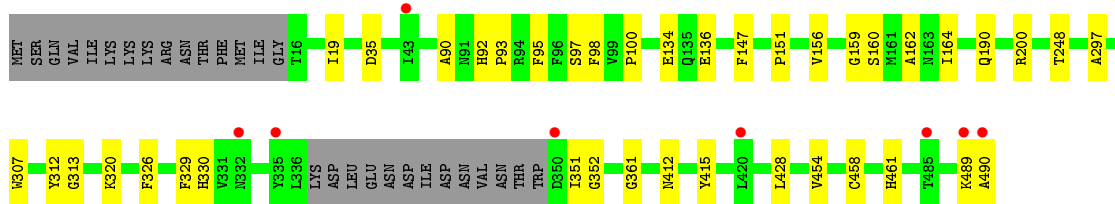
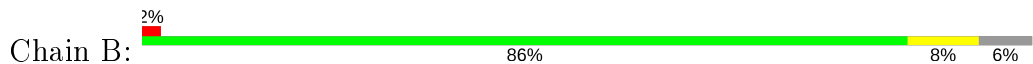
● Molecule 1: Pyridoxal-dependent decarboxylase domain protein



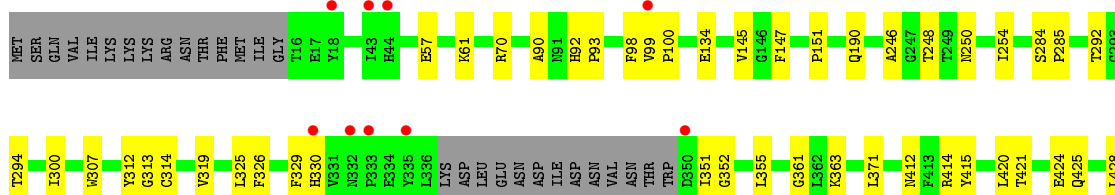
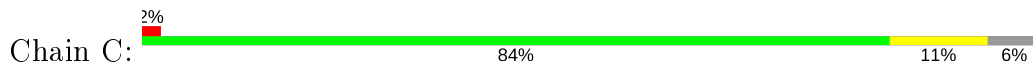
● Molecule 1: Pyridoxal-dependent decarboxylase domain protein



● Molecule 1: Pyridoxal-dependent decarboxylase domain protein



● Molecule 1: Pyridoxal-dependent decarboxylase domain protein





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	58.63Å 145.77Å 165.07Å 72.85° 88.84° 88.30°	Depositor
Resolution (Å)	79.37 – 2.80 79.37 – 2.80	Depositor EDS
% Data completeness (in resolution range)	94.7 (79.37-2.80) 80.1 (79.37-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.88 (at 2.82Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1168)	Depositor
R, R_{free}	0.231 , 0.257 0.232 , 0.259	Depositor DCC
R_{free} test set	6088 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	35.7	Xtrriage
Anisotropy	0.146	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , -10.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtrriage
Estimated twinning fraction	0.105 for h,-k,-l	Xtrriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	29598	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 24.28 % 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 3.9546e-03.*

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: 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	0.21	0/3713	0.39	0/5031
1	B	0.21	0/3713	0.40	0/5031
1	C	0.22	0/3713	0.40	0/5031
1	E	0.21	0/3713	0.40	0/5031
1	F	0.21	0/3713	0.40	0/5031
1	G	0.21	0/3713	0.40	0/5031
1	H	0.21	0/3713	0.40	0/5031
1	U	0.21	0/3713	0.40	0/5031
All	All	0.21	0/29704	0.40	0/40248

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	3633	0	3604	19	0
1	B	3633	0	3604	21	1
1	C	3633	0	3604	34	0
1	E	3633	0	3604	21	0
1	F	3633	0	3604	22	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	G	3633	0	3604	21	1
1	H	3633	0	3604	28	0
1	U	3633	0	3604	17	0
2	A	15	0	7	0	0
2	B	15	0	7	1	0
2	C	15	0	7	0	0
2	E	15	0	7	1	0
2	F	15	0	7	2	0
2	G	15	0	7	0	0
2	H	15	0	7	1	0
2	U	15	0	7	1	0
3	A	49	0	0	1	0
3	B	45	0	0	1	0
3	C	36	0	0	7	0
3	E	56	0	0	1	0
3	F	73	0	0	1	1
3	G	46	0	0	1	0
3	H	52	0	0	3	0
3	U	57	0	0	0	0
All	All	29598	0	28888	176	2

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 3.

All (176) 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:C:421:THR:O	3:C:629:HOH:O	1.89	0.89
1:B:326:PHE:O	1:B:330:HIS:ND1	2.20	0.74
1:H:326:PHE:O	1:H:330:HIS:ND1	2.20	0.74
1:F:326:PHE:O	1:F:330:HIS:ND1	2.23	0.71
1:G:326:PHE:O	1:G:330:HIS:ND1	2.23	0.70
1:C:425:GLN:N	3:C:629:HOH:O	2.23	0.70
1:E:161:MET:SD	3:E:650:HOH:O	2.49	0.70
1:A:326:PHE:O	1:A:330:HIS:ND1	2.26	0.68
1:H:481:ARG:NH2	3:H:611:HOH:O	2.27	0.67
1:U:326:PHE:O	1:U:330:HIS:ND1	2.27	0.67
1:E:326:PHE:O	1:E:330:HIS:ND1	2.27	0.67
1:C:420:LEU:HB3	3:C:629:HOH:O	1.96	0.65
1:C:326:PHE:O	1:C:330:HIS:ND1	2.30	0.65
1:G:357:ARG:NH1	3:G:602:HOH:O	2.30	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:134:GLU:HG2	1:H:361:GLY:HA3	1.86	0.57
1:E:200:ARG:NH2	1:H:330:HIS:O	2.37	0.57
1:C:425:GLN:HG3	3:C:629:HOH:O	2.04	0.56
1:B:90:ALA:HB2	1:B:100:PRO:HB2	1.87	0.55
1:A:90:ALA:HB2	1:A:100:PRO:HB2	1.89	0.55
1:U:329:PHE:CG	1:U:352:GLY:HA3	2.42	0.55
1:E:412:ASN:HB3	1:E:454:VAL:HG11	1.89	0.54
1:E:134:GLU:HG2	1:E:361:GLY:HA3	1.89	0.53
1:F:90:ALA:HB2	1:F:100:PRO:HB2	1.89	0.53
1:C:145:VAL:O	1:C:294:THR:OG1	2.25	0.52
1:C:424:GLU:HG3	3:C:629:HOH:O	2.08	0.52
1:C:70:ARG:NH2	3:C:608:HOH:O	2.41	0.52
1:G:90:ALA:HB2	1:G:100:PRO:HB2	1.91	0.52
1:B:412:ASN:HB3	1:B:454:VAL:HG11	1.92	0.51
1:F:415:TYR:HD2	1:F:428:LEU:HD21	1.76	0.51
1:E:330:HIS:O	1:H:200:ARG:NH2	2.44	0.51
1:E:90:ALA:HB2	1:E:100:PRO:HB2	1.93	0.50
1:C:329:PHE:CG	1:C:352:GLY:HA3	2.46	0.50
1:U:412:ASN:HB3	1:U:454:VAL:HG11	1.93	0.50
1:F:21:ASN:ND2	3:F:621:HOH:O	2.35	0.49
1:C:414:ARG:NH2	1:C:425:GLN:OE1	2.46	0.48
1:G:98:PHE:HB3	1:G:458:CYS:SG	2.54	0.48
1:F:412:ASN:HB3	1:F:454:VAL:HG11	1.94	0.48
1:U:274:ASP:OD2	2:U:501:PLP:N1	2.47	0.48
1:B:134:GLU:HG2	1:B:361:GLY:HA3	1.96	0.48
1:A:412:ASN:HB3	1:A:454:VAL:HG11	1.96	0.48
1:F:207:SER:O	1:F:210:ARG:NH1	2.47	0.47
1:G:134:GLU:HG2	1:G:361:GLY:HA3	1.96	0.47
1:B:415:TYR:HD2	1:B:428:LEU:HD21	1.79	0.47
1:C:190:GLN:HB3	1:C:248:THR:HG21	1.97	0.47
1:H:248:THR:HG22	1:H:250:ASN:H	1.79	0.47
1:H:312:TYR:HA	1:H:313:GLY:HA2	1.58	0.47
1:H:329:PHE:CG	1:H:352:GLY:HA3	2.50	0.47
1:E:274:ASP:OD2	2:E:501:PLP:N1	2.48	0.47
1:F:274:ASP:OD2	2:F:501:PLP:N1	2.48	0.47
1:F:159:GLY:HA3	2:F:501:PLP:H5A2	1.96	0.47
1:H:489:LYS:HA	1:H:490:ALA:HA	1.59	0.46
1:C:415:TYR:HD2	1:C:428:LEU:HD21	1.80	0.46
1:H:93:PRO:HA	1:H:461:HIS:CG	2.51	0.46
1:C:489:LYS:HA	1:C:490:ALA:HA	1.58	0.46
1:E:329:PHE:CG	1:E:352:GLY:HA3	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:196:ALA:HA	1:G:199:LEU:HD12	1.97	0.46
1:A:329:PHE:CG	1:A:352:GLY:HA3	2.50	0.46
1:F:278:GLY:O	1:F:291:LEU:HD11	2.15	0.46
1:A:312:TYR:HA	1:A:313:GLY:HA2	1.57	0.46
1:G:93:PRO:HA	1:G:461:HIS:CG	2.50	0.46
1:B:19:ILE:HG23	1:C:371:LEU:HD13	1.97	0.46
1:E:415:TYR:HD2	1:E:428:LEU:HD21	1.80	0.46
1:B:312:TYR:HA	1:B:313:GLY:HA2	1.57	0.46
1:A:160:SER:O	1:A:164:ILE:HG13	2.16	0.46
1:B:329:PHE:CG	1:B:352:GLY:HA3	2.51	0.46
1:H:461:HIS:HA	1:H:462:PRO:HD3	1.83	0.46
1:H:90:ALA:HB2	1:H:100:PRO:HB2	1.98	0.46
1:A:415:TYR:HD2	1:A:428:LEU:HD21	1.80	0.46
1:A:93:PRO:HA	1:A:461:HIS:CG	2.51	0.46
1:E:93:PRO:HA	1:E:461:HIS:CG	2.51	0.46
1:A:489:LYS:HA	1:A:490:ALA:HA	1.60	0.45
1:H:154:VAL:HB	1:H:350:ASP:HB3	1.98	0.45
1:E:312:TYR:HA	1:E:313:GLY:HA2	1.58	0.45
1:G:489:LYS:HA	1:G:490:ALA:HA	1.59	0.45
1:C:98:PHE:HB3	1:C:458:CYS:SG	2.56	0.45
1:C:147:PHE:HB3	1:C:151:PRO:HB3	1.99	0.45
1:F:489:LYS:HA	1:F:490:ALA:HA	1.59	0.45
1:G:329:PHE:CG	1:G:352:GLY:HA3	2.52	0.45
1:A:98:PHE:HB3	1:A:458:CYS:SG	2.56	0.45
1:U:99:VAL:HB	1:U:100:PRO:HD3	1.98	0.45
1:B:93:PRO:HA	1:B:461:HIS:CG	2.51	0.45
1:C:412:ASN:HB3	1:C:454:VAL:HG11	1.99	0.45
1:G:187:ILE:HG22	1:G:243:ILE:HB	1.98	0.45
1:C:314:CYS:HB2	1:C:363:LYS:HE3	2.00	0.44
1:B:200:ARG:NH2	1:C:330:HIS:O	2.50	0.44
1:H:407:GLN:OE1	3:H:651:HOH:O	2.21	0.44
1:U:90:ALA:HB2	1:U:100:PRO:HB2	1.98	0.44
1:U:190:GLN:HB3	1:U:248:THR:HG21	2.00	0.44
1:U:246:ALA:HB1	1:U:291:LEU:HD22	1.98	0.44
1:A:53:LYS:HD2	1:A:55:HIS:HE1	1.82	0.44
1:F:99:VAL:HB	1:F:100:PRO:HD3	2.00	0.44
1:C:90:ALA:HB2	1:C:100:PRO:HB2	1.98	0.44
1:C:145:VAL:HG11	1:C:300:ILE:HG21	2.00	0.44
1:U:312:TYR:HA	1:U:313:GLY:HA2	1.58	0.44
1:C:246:ALA:HA	1:C:254:ILE:HG23	2.00	0.43
1:F:154:VAL:HB	1:F:350:ASP:HB3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:412:ASN:HB3	1:G:454:VAL:HG11	2.00	0.43
1:U:134:GLU:HG2	1:U:361:GLY:HA3	1.99	0.43
1:C:312:TYR:HA	1:C:313:GLY:HA2	1.60	0.43
1:E:319:VAL:HG21	1:E:325:LEU:HG	2.00	0.43
1:F:248:THR:HG22	1:F:250:ASN:H	1.83	0.43
1:F:329:PHE:CG	1:F:352:GLY:HA3	2.52	0.43
1:G:234:LYS:HB3	1:G:234:LYS:HE2	1.91	0.43
1:G:330:HIS:H	1:G:330:HIS:HD1	1.67	0.43
1:A:248:THR:HG22	1:A:250:ASN:H	1.82	0.43
1:B:147:PHE:HB3	1:B:151:PRO:HB3	2.00	0.43
1:B:489:LYS:HA	1:B:490:ALA:HA	1.60	0.43
1:E:314:CYS:HB2	1:E:363:LYS:HE3	2.01	0.43
1:C:134:GLU:HG2	1:C:361:GLY:HA3	2.00	0.43
1:E:92:HIS:HA	1:E:93:PRO:HD3	1.91	0.43
1:B:297:ALA:O	1:B:320:LYS:HE3	2.19	0.43
1:G:92:HIS:HA	1:G:93:PRO:HD3	1.90	0.43
1:C:99:VAL:HB	1:C:100:PRO:HD3	2.00	0.43
1:G:355:LEU:HA	1:G:355:LEU:HD12	1.91	0.43
1:E:99:VAL:HB	1:E:100:PRO:HD3	2.01	0.42
1:G:53:LYS:HD2	1:G:55:HIS:HE1	1.84	0.42
1:G:99:VAL:HB	1:G:100:PRO:HD3	2.01	0.42
1:E:19:ILE:HG23	1:H:371:LEU:HD13	2.02	0.42
1:U:286:LYS:HB2	1:U:286:LYS:HE3	1.85	0.42
1:B:159:GLY:HA3	2:B:501:PLP:H5A2	2.02	0.42
1:F:312:TYR:HA	1:F:313:GLY:HA2	1.58	0.42
1:F:314:CYS:HB2	1:F:363:LYS:HE3	2.00	0.42
1:H:319:VAL:HG21	1:H:325:LEU:HG	2.01	0.42
1:E:368:LEU:HB3	1:H:65:ILE:HG23	2.02	0.42
1:U:98:PHE:HB3	1:U:458:CYS:SG	2.58	0.42
1:U:489:LYS:HA	1:U:490:ALA:HA	1.59	0.42
1:H:412:ASN:HB3	1:H:454:VAL:HG11	2.01	0.42
1:H:420:LEU:HB3	1:H:424:GLU:HG3	2.02	0.42
1:B:156:VAL:HG21	1:B:162:ALA:HB2	2.02	0.42
1:E:248:THR:HG22	1:E:250:ASN:H	1.85	0.42
1:H:284:SER:HA	1:H:285:PRO:HD2	1.94	0.42
1:H:92:HIS:HA	1:H:93:PRO:HD3	1.91	0.42
1:A:134:GLU:HG2	1:A:361:GLY:HA3	2.00	0.42
1:B:98:PHE:HB3	1:B:458:CYS:SG	2.60	0.41
1:B:136:GLU:OE2	3:B:622:HOH:O	2.21	0.41
1:B:160:SER:O	1:B:164:ILE:HG13	2.20	0.41
1:A:461:HIS:HA	1:A:462:PRO:HD3	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:319:VAL:HG21	1:C:325:LEU:HG	2.02	0.41
1:F:147:PHE:HB3	1:F:151:PRO:HB3	2.02	0.41
1:G:248:THR:HG22	1:G:250:ASN:H	1.85	0.41
1:H:98:PHE:HB3	1:H:458:CYS:SG	2.60	0.41
1:C:355:LEU:HA	1:C:355:LEU:HD12	1.93	0.41
1:F:98:PHE:HB3	1:F:458:CYS:SG	2.60	0.41
1:C:92:HIS:HA	1:C:93:PRO:HD3	1.91	0.41
1:C:284:SER:HA	1:C:285:PRO:HD2	1.93	0.41
1:H:274:ASP:OD2	2:H:501:PLP:N1	2.54	0.41
1:U:278:GLY:O	1:U:291:LEU:HD11	2.20	0.41
1:H:415:TYR:HD2	1:H:428:LEU:HD21	1.86	0.41
1:A:330:HIS:HD1	1:A:330:HIS:H	1.69	0.41
1:C:292:THR:N	3:C:601:HOH:O	2.53	0.41
1:F:446:THR:HG22	1:F:455:LEU:HD23	2.01	0.41
1:A:46:GLN:HA	1:A:47:PRO:HD3	1.96	0.41
1:F:134:GLU:HG2	1:F:361:GLY:HA3	2.02	0.41
1:G:156:VAL:HG21	1:G:162:ALA:HB2	2.02	0.41
1:U:330:HIS:HD1	1:U:330:HIS:H	1.68	0.41
1:A:99:VAL:HB	1:A:100:PRO:HD3	2.02	0.41
1:H:99:VAL:HB	1:H:100:PRO:HD3	2.01	0.41
1:U:147:PHE:HB3	1:U:151:PRO:HB3	2.02	0.41
1:B:95:PHE:CE2	1:B:97:SER:HB3	2.56	0.41
1:G:461:HIS:HA	1:G:462:PRO:HD3	1.84	0.41
1:B:190:GLN:HB3	1:B:248:THR:HG21	2.03	0.40
1:C:428:LEU:O	1:C:432:ILE:HG13	2.20	0.40
1:E:489:LYS:HA	1:E:490:ALA:HA	1.59	0.40
1:A:143:LYS:NZ	3:A:609:HOH:O	2.42	0.40
1:E:136:GLU:OE2	1:H:61:LYS:HG2	2.21	0.40
1:C:248:THR:HG22	1:C:250:ASN:H	1.87	0.40
1:C:57:GLU:O	1:C:61:LYS:HG3	2.22	0.40
1:H:398:LYS:O	3:H:611:HOH:O	2.22	0.40
1:U:248:THR:HG22	1:U:250:ASN:H	1.85	0.40
1:B:92:HIS:HA	1:B:93:PRO:HD3	1.91	0.40
1:C:93:PRO:HA	1:C:461:HIS:CG	2.56	0.40
1:F:160:SER:O	1:F:164:ILE:HG13	2.21	0.40
1:G:425:GLN:O	1:G:429:ASN:ND2	2.52	0.40
1:A:207:SER:O	1:A:210:ARG:NH1	2.54	0.40
1:F:46:GLN:HA	1:F:47:PRO:HD3	1.98	0.40
1:H:330:HIS:H	1:H:330:HIS:HD1	1.68	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:23:THR:OG1	1:B:35:ASP:OD2[1_655]	2.09	0.11
1:G:335:TYR:OH	3:F:611:HOH:O[1_455]	2.12	0.08

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	458/490 (94%)	441 (96%)	17 (4%)	0	100	100
1	B	458/490 (94%)	441 (96%)	17 (4%)	0	100	100
1	C	458/490 (94%)	441 (96%)	17 (4%)	0	100	100
1	E	458/490 (94%)	441 (96%)	17 (4%)	0	100	100
1	F	458/490 (94%)	441 (96%)	17 (4%)	0	100	100
1	G	458/490 (94%)	441 (96%)	17 (4%)	0	100	100
1	H	458/490 (94%)	442 (96%)	16 (4%)	0	100	100
1	U	458/490 (94%)	442 (96%)	16 (4%)	0	100	100
All	All	3664/3920 (94%)	3530 (96%)	134 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	395/422 (94%)	393 (100%)	2 (0%)	88	96

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	395/422 (94%)	393 (100%)	2 (0%)	88	96
1	C	395/422 (94%)	393 (100%)	2 (0%)	88	96
1	E	395/422 (94%)	393 (100%)	2 (0%)	88	96
1	F	395/422 (94%)	393 (100%)	2 (0%)	88	96
1	G	395/422 (94%)	393 (100%)	2 (0%)	88	96
1	H	395/422 (94%)	393 (100%)	2 (0%)	88	96
1	U	395/422 (94%)	393 (100%)	2 (0%)	88	96
All	All	3160/3376 (94%)	3144 (100%)	16 (0%)	88	96

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	307	TRP
1	E	351	ILE
1	A	307	TRP
1	A	351	ILE
1	U	307	TRP
1	U	351	ILE
1	G	307	TRP
1	G	351	ILE
1	H	307	TRP
1	H	351	ILE
1	F	307	TRP
1	F	351	ILE
1	B	307	TRP
1	B	351	ILE
1	C	307	TRP
1	C	351	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	55	HIS
1	G	55	HIS
1	H	407	GLN

5.3.3 RNA [i](#)

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

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
2	PLP	C	501	-	15,15,16	1.01	1 (6%)	20,22,23	1.34	2 (10%)
2	PLP	E	501	-	15,15,16	1.02	1 (6%)	20,22,23	1.45	2 (10%)
2	PLP	F	501	-	15,15,16	1.00	1 (6%)	20,22,23	1.43	2 (10%)
2	PLP	G	501	-	15,15,16	0.98	1 (6%)	20,22,23	1.35	2 (10%)
2	PLP	U	501	-	15,15,16	1.02	1 (6%)	20,22,23	1.27	2 (10%)
2	PLP	H	501	-	15,15,16	1.01	1 (6%)	20,22,23	1.35	2 (10%)
2	PLP	A	501	-	15,15,16	1.02	1 (6%)	20,22,23	1.38	2 (10%)
2	PLP	B	501	-	15,15,16	1.01	1 (6%)	20,22,23	1.41	2 (10%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	PLP	C	501	-	-	3/6/6/8	0/1/1/1
2	PLP	E	501	-	-	3/6/6/8	0/1/1/1
2	PLP	F	501	-	-	2/6/6/8	0/1/1/1
2	PLP	G	501	-	-	3/6/6/8	0/1/1/1
2	PLP	U	501	-	-	2/6/6/8	0/1/1/1
2	PLP	H	501	-	-	3/6/6/8	0/1/1/1
2	PLP	A	501	-	-	3/6/6/8	0/1/1/1
2	PLP	B	501	-	-	2/6/6/8	0/1/1/1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	501	PLP	C2-N1	2.33	1.38	1.33
2	E	501	PLP	C2-N1	2.32	1.38	1.33
2	A	501	PLP	C2-N1	2.32	1.38	1.33
2	U	501	PLP	C2-N1	2.32	1.38	1.33
2	C	501	PLP	C2-N1	2.28	1.38	1.33
2	B	501	PLP	C2-N1	2.28	1.38	1.33
2	F	501	PLP	C2-N1	2.28	1.38	1.33
2	G	501	PLP	C2-N1	2.20	1.38	1.33

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	501	PLP	C4A-C4-C3	-3.83	114.01	120.50
2	E	501	PLP	C4A-C4-C3	-3.65	114.31	120.50
2	H	501	PLP	C4A-C4-C3	-3.56	114.46	120.50
2	B	501	PLP	O4P-C5A-C5	3.56	116.13	109.35
2	C	501	PLP	C4A-C4-C3	-3.46	114.63	120.50
2	A	501	PLP	C4A-C4-C3	-3.43	114.68	120.50
2	G	501	PLP	C4A-C4-C3	-3.38	114.76	120.50
2	B	501	PLP	C4A-C4-C3	-3.38	114.77	120.50
2	E	501	PLP	O4P-C5A-C5	3.26	115.56	109.35
2	A	501	PLP	O4P-C5A-C5	3.15	115.36	109.35
2	U	501	PLP	C4A-C4-C3	-3.12	115.22	120.50
2	F	501	PLP	O4P-C5A-C5	2.98	115.03	109.35
2	G	501	PLP	O4P-C5A-C5	2.94	114.95	109.35
2	C	501	PLP	O4P-C5A-C5	2.89	114.85	109.35
2	U	501	PLP	O4P-C5A-C5	2.71	114.51	109.35
2	H	501	PLP	O4P-C5A-C5	2.69	114.47	109.35

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	501	PLP	C4-C5-C5A-O4P
2	C	501	PLP	C6-C5-C5A-O4P
2	E	501	PLP	C4-C5-C5A-O4P
2	E	501	PLP	C6-C5-C5A-O4P
2	F	501	PLP	C4-C5-C5A-O4P
2	F	501	PLP	C6-C5-C5A-O4P
2	G	501	PLP	C4-C5-C5A-O4P
2	G	501	PLP	C6-C5-C5A-O4P
2	H	501	PLP	C4-C5-C5A-O4P
2	A	501	PLP	C4-C5-C5A-O4P
2	A	501	PLP	C6-C5-C5A-O4P
2	B	501	PLP	C4-C5-C5A-O4P
2	B	501	PLP	C6-C5-C5A-O4P
2	U	501	PLP	C5A-O4P-P-O1P
2	H	501	PLP	C6-C5-C5A-O4P
2	C	501	PLP	C5A-O4P-P-O1P
2	E	501	PLP	C5A-O4P-P-O1P
2	G	501	PLP	C5A-O4P-P-O1P
2	H	501	PLP	C5A-O4P-P-O1P
2	A	501	PLP	C5A-O4P-P-O1P
2	U	501	PLP	C5A-O4P-P-O3P

There are no ring outliers.

5 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	501	PLP	1	0
2	F	501	PLP	2	0
2	U	501	PLP	1	0
2	H	501	PLP	1	0
2	B	501	PLP	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	462/490 (94%)	-0.03	8 (1%) 70 63	1, 14, 63, 107	0
1	B	462/490 (94%)	0.01	8 (1%) 70 63	1, 14, 63, 110	0
1	C	462/490 (94%)	-0.01	11 (2%) 59 49	1, 15, 60, 109	0
1	E	462/490 (94%)	-0.08	11 (2%) 59 49	1, 15, 62, 113	0
1	F	462/490 (94%)	-0.01	9 (1%) 66 59	1, 13, 62, 112	0
1	G	462/490 (94%)	-0.01	8 (1%) 70 63	1, 14, 60, 111	0
1	H	462/490 (94%)	0.04	15 (3%) 47 37	1, 16, 64, 105	0
1	U	462/490 (94%)	-0.01	12 (2%) 56 46	1, 15, 64, 110	0
All	All	3696/3920 (94%)	-0.01	82 (2%) 62 52	1, 15, 63, 113	0

All (82) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	490	ALA	11.1
1	C	490	ALA	10.9
1	B	490	ALA	9.2
1	H	490	ALA	8.6
1	A	332	ASN	7.3
1	G	490	ALA	6.7
1	F	332	ASN	6.6
1	G	332	ASN	6.1
1	E	332	ASN	6.1
1	H	99	VAL	6.0
1	F	489	LYS	6.0
1	B	420	LEU	5.5
1	H	489	LYS	5.3
1	G	43	ILE	4.6
1	C	489	LYS	4.6
1	A	489	LYS	4.5

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Mol	Chain	Res	Type	RSRZ
1	E	44	HIS	4.5
1	G	489	LYS	4.5
1	U	350	ASP	4.5
1	B	489	LYS	4.5
1	B	350	ASP	4.4
1	H	332	ASN	4.4
1	C	350	ASP	4.1
1	G	350	ASP	3.9
1	U	490	ALA	3.9
1	C	18	TYR	3.9
1	B	332	ASN	3.8
1	A	350	ASP	3.7
1	H	335	TYR	3.7
1	E	43	ILE	3.6
1	U	43	ILE	3.5
1	C	43	ILE	3.4
1	F	350	ASP	3.4
1	G	99	VAL	3.4
1	G	44	HIS	3.4
1	U	489	LYS	3.3
1	F	335	TYR	3.3
1	A	490	ALA	3.2
1	F	123	GLY	3.1
1	E	334	GLU	3.0
1	E	350	ASP	3.0
1	C	335	TYR	2.9
1	C	332	ASN	2.9
1	H	62	GLN	2.9
1	U	46	GLN	2.8
1	H	421	THR	2.8
1	U	420	LEU	2.8
1	C	99	VAL	2.8
1	U	44	HIS	2.7
1	A	43	ILE	2.7
1	F	334	GLU	2.7
1	U	333	PRO	2.7
1	U	487	MET	2.7
1	A	45	ASP	2.7
1	A	334	GLU	2.7
1	H	350	ASP	2.7
1	H	334	GLU	2.6
1	G	334	GLU	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	485	THR	2.5
1	H	44	HIS	2.5
1	C	333	PRO	2.5
1	B	335	TYR	2.5
1	H	399	ASP	2.5
1	A	44	HIS	2.4
1	U	332	ASN	2.4
1	U	480	GLY	2.4
1	E	490	ALA	2.4
1	E	489	LYS	2.3
1	F	99	VAL	2.3
1	E	123	GLY	2.3
1	H	45	ASP	2.3
1	H	333	PRO	2.2
1	E	124	SER	2.2
1	E	333	PRO	2.2
1	C	330	HIS	2.2
1	H	83	VAL	2.2
1	F	125	LYS	2.2
1	H	422	LYS	2.1
1	B	43	ILE	2.1
1	C	44	HIS	2.1
1	E	227	GLU	2.1
1	U	99	VAL	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(\AA^2)	Q<0.9
2	PLP	E	501	15/16	0.88	0.27	24,44,79,80	0
2	PLP	A	501	15/16	0.89	0.31	2,49,70,78	0
2	PLP	F	501	15/16	0.90	0.18	6,31,65,66	0
2	PLP	U	501	15/16	0.92	0.21	4,38,74,80	0
2	PLP	B	501	15/16	0.94	0.26	8,43,64,67	0
2	PLP	G	501	15/16	0.95	0.24	6,32,73,76	0
2	PLP	H	501	15/16	0.95	0.24	7,28,73,74	0
2	PLP	C	501	15/16	0.96	0.19	10,35,66,66	0

6.5 Other polymers [i](#)

There are no such residues in this entry.