



Full wwPDB X-ray Structure Validation Report ⓘ

May 19, 2020 – 06:57 am BST

PDB ID : 5ZL6
Title : Histidine Racemase from *Leuconostoc mesenteroides* subsp. sake NBRC 102480
Authors : Adachi, M.; Shimizu, R.; Oikawa, T.
Deposited on : 2018-03-27
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

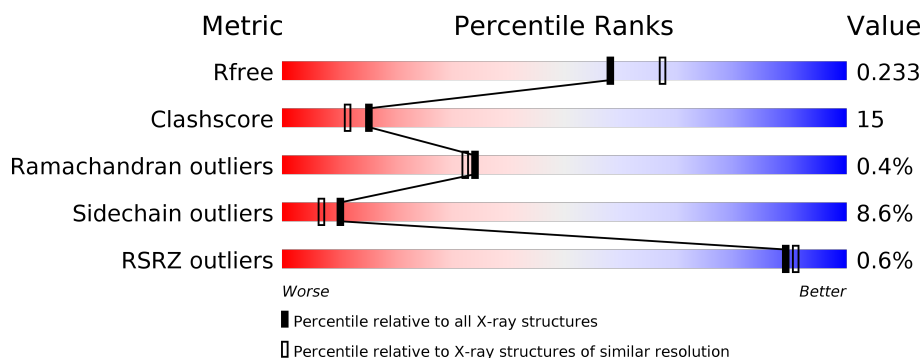
1 Overall quality at a glance ⓘ

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	381	<div> <div>%</div> <div> <div></div> <div>76%</div> <div>18%</div> <div>• •</div> </div> </div>
1	B	381	<div> <div>%</div> <div> <div></div> <div>72%</div> <div>20%</div> <div>• • •</div> </div> </div>
1	C	381	<div> <div>%</div> <div> <div></div> <div>74%</div> <div>17%</div> <div>5% • •</div> </div> </div>
1	D	381	<div> <div></div> <div> <div>75%</div> <div>18%</div> <div>5% •</div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	PLP	B	402	-	-	X	-

2 Entry composition [i](#)

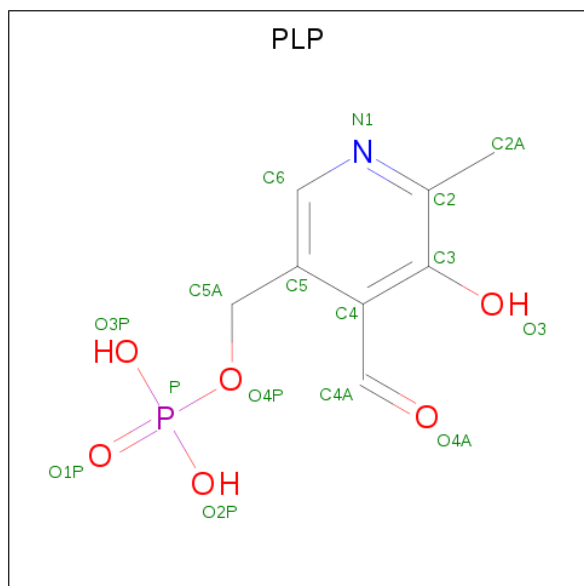
There are 3 unique types of molecules in this entry. The entry contains 13605 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Histidine racemase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	372	Total	C	N	O	S	0	0	0
			2831	1798	487	534	12			
1	B	372	Total	C	N	O	S	0	0	0
			2831	1798	487	534	12			
1	C	372	Total	C	N	O	S	0	0	0
			2831	1798	487	534	12			
1	D	372	Total	C	N	O	S	0	0	0
			2831	1798	487	534	12			

- 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		

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

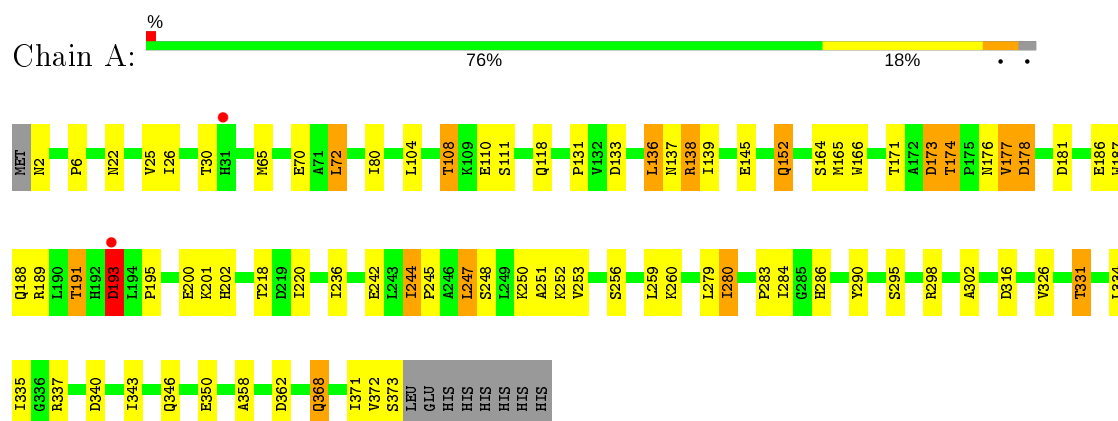
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	494	Total	O	0	0
			494	494		
3	B	553	Total	O	0	0
			553	553		
3	C	568	Total	O	0	0
			568	568		
3	D	576	Total	O	0	0
			576	576		

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.

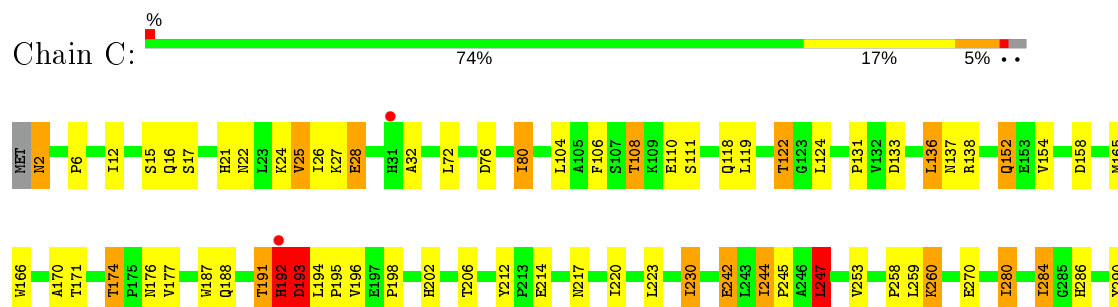
• Molecule 1: Histidine racemase



• Molecule 1: Histidine racemase

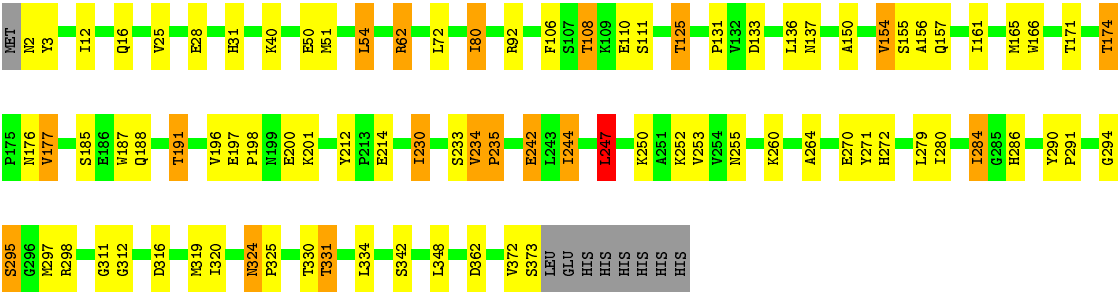


• Molecule 1: Histidine racemase





● Molecule 1: Histidine racemase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	68.05Å 142.47Å 78.70Å 90.00° 99.30° 90.00°	Depositor
Resolution (Å)	33.60 – 2.10 34.10 – 2.10	Depositor EDS
% Data completeness (in resolution range)	97.8 (33.60-2.10) 94.6 (34.10-2.10)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.90 (at 2.10Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.174 , 0.235 0.175 , 0.233	Depositor DCC
R_{free} test set	4208 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	9.4	Xtriage
Anisotropy	1.450	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 50.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	13605	wwPDB-VP
Average B, all atoms (Å ²)	13.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.46% 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: 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	0.27	0/2888	0.50	1/3928 (0.0%)
1	B	0.28	0/2888	0.51	2/3928 (0.1%)
1	C	0.28	0/2888	0.48	1/3928 (0.0%)
1	D	0.28	0/2888	0.51	1/3928 (0.0%)
All	All	0.28	0/11552	0.50	5/15712 (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	2
1	B	0	2
1	D	0	1
All	All	0	5

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	247	LEU	CA-CB-CG	8.46	134.77	115.30
1	B	247	LEU	CA-CB-CG	8.31	134.41	115.30
1	C	247	LEU	CA-CB-CG	7.31	132.12	115.30
1	D	247	LEU	CA-CB-CG	7.01	131.42	115.30
1	B	247	LEU	CB-CG-CD1	5.05	119.59	111.00

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	173	ASP	Peptide
1	A	178	ASP	Peptide
1	B	173	ASP	Peptide
1	B	178	ASP	Peptide
1	D	234	VAL	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	2831	0	2844	66	2
1	B	2831	0	2842	90	2
1	C	2831	0	2844	88	1
1	D	2831	0	2842	90	0
2	A	15	0	7	0	0
2	B	30	0	14	9	1
2	C	15	0	7	1	0
2	D	30	0	14	3	0
3	A	494	0	0	22	2
3	B	553	0	0	50	3
3	C	568	0	0	41	4
3	D	576	0	0	40	4
All	All	13605	0	11414	339	10

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 (339) 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:D:31:HIS:ND1	3:D:502:HOH:O	1.62	1.32
1:D:2:ASN:N	3:D:501:HOH:O	1.58	1.27
1:B:2:ASN:ND2	3:B:502:HOH:O	1.74	1.21
1:B:180:VAL:HG23	3:B:501:HOH:O	1.48	1.14
1:B:72:LEU:HG	3:B:785:HOH:O	1.48	1.12
2:B:402:PLP:P	3:B:503:HOH:O	2.09	1.07
1:B:31:HIS:HE1	3:B:656:HOH:O	1.37	1.07

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:235:PRO:HD2	3:D:531:HOH:O	1.56	1.05
1:B:171:THR:OG1	3:B:501:HOH:O	1.73	1.05
1:D:234:VAL:O	1:D:235:PRO:O	1.73	1.04
1:C:230:ILE:HB	3:C:938:HOH:O	1.57	1.03
1:C:298:ARG:NH2	3:C:501:HOH:O	1.88	1.03
1:B:178:ASP:HB2	3:B:719:HOH:O	1.58	1.03
1:C:286:HIS:NE2	3:C:502:HOH:O	1.93	1.01
1:D:212:TYR:HA	3:D:581:HOH:O	1.61	1.00
1:C:286:HIS:CE1	3:C:502:HOH:O	2.14	1.00
1:B:219:ASP:OD2	3:B:504:HOH:O	1.82	0.98
2:B:402:PLP:O4P	3:B:503:HOH:O	1.79	0.98
1:C:187:TRP:O	1:C:191:THR:HB	1.64	0.96
1:D:187:TRP:O	1:D:191:THR:HB	1.64	0.95
1:D:212:TYR:CD1	3:D:581:HOH:O	2.19	0.95
1:B:138:ARG:HD2	3:B:885:HOH:O	1.67	0.94
1:D:295:SER:HB2	3:D:559:HOH:O	1.68	0.92
1:B:31:HIS:CE1	3:B:656:HOH:O	2.13	0.91
1:A:2:ASN:N	3:A:502:HOH:O	2.01	0.90
1:D:372:VAL:O	1:D:373:SER:HB2	1.69	0.89
2:B:402:PLP:C5A	3:B:503:HOH:O	2.20	0.88
1:C:171:THR:O	1:C:174:THR:HG22	1.73	0.87
1:C:2:ASN:HA	3:C:533:HOH:O	1.74	0.86
1:C:174:THR:HG23	1:C:176:ASN:H	1.38	0.86
1:A:108:THR:HG22	1:A:111:SER:H	1.40	0.86
1:A:110:GLU:HG2	3:A:596:HOH:O	1.76	0.86
1:A:133:ASP:OD2	1:A:137:ASN:HA	1.75	0.85
1:B:284:ILE:HD11	1:B:290:TYR:HB2	1.58	0.85
1:A:186:GLU:OE1	3:A:501:HOH:O	1.95	0.85
1:B:33:LYS:HE3	3:B:521:HOH:O	1.75	0.84
1:C:191:THR:HG22	3:C:939:HOH:O	1.78	0.84
1:D:250:LYS:HE2	3:D:507:HOH:O	1.79	0.82
1:C:260:LYS:HE2	3:C:655:HOH:O	1.79	0.81
1:C:315:MET:HG3	3:C:502:HOH:O	1.79	0.81
1:D:373:SER:C	3:D:616:HOH:O	2.18	0.81
1:A:250:LYS:HE2	3:A:638:HOH:O	1.79	0.80
1:A:178:ASP:HA	1:A:181:ASP:HB2	1.64	0.80
1:B:133:ASP:OD2	1:B:137:ASN:HA	1.82	0.80
1:C:24:LYS:HE2	3:C:777:HOH:O	1.81	0.79
1:D:62:ARG:HD3	3:D:587:HOH:O	1.83	0.79
2:B:402:PLP:O1P	3:B:503:HOH:O	1.98	0.79
1:A:136:LEU:HD13	1:A:138:ARG:HG2	1.63	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:402:PLP:C6	3:B:503:HOH:O	2.30	0.78
1:A:80:ILE:HD12	3:C:923:HOH:O	1.83	0.78
1:B:178:ASP:OD1	1:B:178:ASP:C	2.23	0.77
1:C:196:VAL:HG22	3:C:870:HOH:O	1.84	0.77
1:C:76:ASP:OD2	3:C:503:HOH:O	2.01	0.76
1:C:108:THR:HG22	1:C:111:SER:H	1.50	0.76
1:D:212:TYR:HD1	3:D:581:HOH:O	1.63	0.76
1:D:234:VAL:O	1:D:235:PRO:C	2.23	0.76
1:D:242:GLU:HG3	3:D:533:HOH:O	1.86	0.76
1:B:13:GLU:OE1	3:B:505:HOH:O	2.04	0.76
1:D:252:LYS:HZ3	1:D:331:THR:HG22	1.52	0.74
2:B:402:PLP:H2A2	3:B:989:HOH:O	1.87	0.74
1:C:340:ASP:HB2	3:C:506:HOH:O	1.87	0.74
1:C:28:GLU:OE2	3:C:505:HOH:O	2.06	0.73
1:B:152:GLN:HG2	3:B:944:HOH:O	1.87	0.73
1:C:293:ASN:OD1	3:C:504:HOH:O	2.05	0.73
1:D:133:ASP:OD2	1:D:137:ASN:HA	1.89	0.72
1:D:284:ILE:HD11	1:D:290:TYR:HB2	1.71	0.72
1:A:302:ALA:O	3:A:503:HOH:O	2.07	0.72
2:B:402:PLP:C5	3:B:503:HOH:O	2.38	0.72
1:C:298:ARG:NH2	3:C:510:HOH:O	2.23	0.72
1:D:264:ALA:HA	1:D:271:TYR:O	1.88	0.71
1:B:174:THR:HG23	1:B:176:ASN:H	1.55	0.71
1:D:31:HIS:CE1	3:D:502:HOH:O	2.20	0.71
1:A:2:ASN:HB3	3:A:502:HOH:O	1.91	0.70
1:D:244:ILE:O	1:D:244:ILE:HG13	1.92	0.70
1:B:312:GLY:HA3	3:B:538:HOH:O	1.91	0.70
1:B:138:ARG:HG3	1:B:139:ILE:H	1.57	0.70
1:D:108:THR:HG22	1:D:111:SER:H	1.55	0.70
1:C:27:LYS:HE2	3:C:507:HOH:O	1.92	0.70
1:A:189:ARG:HB2	3:A:595:HOH:O	1.92	0.69
1:C:340:ASP:O	3:C:506:HOH:O	2.11	0.68
1:B:100:ARG:NH1	3:B:515:HOH:O	2.26	0.68
1:B:312:GLY:CA	3:B:538:HOH:O	2.42	0.68
1:A:80:ILE:CD1	3:C:923:HOH:O	2.41	0.68
1:B:108:THR:HG22	1:B:111:SER:H	1.58	0.68
1:B:173:ASP:OD1	3:B:506:HOH:O	2.11	0.68
1:B:136:LEU:HD13	1:B:138:ARG:HG2	1.76	0.67
1:A:2:ASN:CB	3:A:502:HOH:O	2.44	0.66
1:C:196:VAL:CG2	3:C:870:HOH:O	2.39	0.66
1:B:187:TRP:O	1:B:191:THR:HB	1.96	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:31:HIS:ND1	3:B:521:HOH:O	2.28	0.66
2:B:402:PLP:O3	3:B:507:HOH:O	2.15	0.65
1:C:32:ALA:O	3:C:507:HOH:O	2.14	0.65
1:B:178:ASP:HA	1:B:181:ASP:HB2	1.78	0.65
1:D:295:SER:OG	1:D:311:GLY:O	2.08	0.65
1:B:196:VAL:HG12	3:B:815:HOH:O	1.96	0.64
1:D:234:VAL:C	3:D:531:HOH:O	2.35	0.64
1:C:316:ASP:HB2	3:C:515:HOH:O	1.98	0.64
1:D:212:TYR:CA	3:D:581:HOH:O	2.33	0.64
1:D:362:ASP:OD1	3:D:504:HOH:O	2.14	0.64
1:B:28:GLU:HG2	3:B:842:HOH:O	1.98	0.63
1:C:217:ASN:HB2	3:C:891:HOH:O	1.98	0.63
1:A:372:VAL:O	1:A:373:SER:C	2.37	0.63
1:A:187:TRP:O	1:A:191:THR:HB	1.98	0.62
1:D:294:GLY:O	1:D:297:MET:HB2	1.99	0.62
1:D:177:VAL:HG13	1:D:212:TYR:HE1	1.63	0.62
1:A:248:SER:HB2	1:A:335:ILE:O	1.99	0.62
1:A:152:GLN:HE22	1:A:195:PRO:HG3	1.64	0.62
1:C:2:ASN:CA	3:C:533:HOH:O	2.41	0.62
1:D:214:GLU:HA	1:D:214:GLU:OE1	1.99	0.62
1:D:108:THR:HG23	1:D:110:GLU:OE2	2.00	0.61
1:A:174:THR:HG22	1:A:176:ASN:H	1.65	0.61
1:D:157:GLN:NE2	3:D:509:HOH:O	2.34	0.61
1:B:180:VAL:N	3:B:501:HOH:O	2.33	0.61
1:B:80:ILE:HD12	3:B:849:HOH:O	2.01	0.60
1:B:292:PHE:CE1	3:B:578:HOH:O	2.51	0.60
1:C:119:LEU:O	1:C:122:THR:HG22	2.02	0.60
1:C:316:ASP:OD2	1:D:40:LYS:HD2	2.02	0.60
1:D:177:VAL:HG13	1:D:212:TYR:CE1	2.37	0.59
1:D:242:GLU:CG	3:D:533:HOH:O	2.47	0.59
1:C:284:ILE:HD11	1:C:290:TYR:HB2	1.84	0.59
1:A:193:ASP:HA	3:A:598:HOH:O	2.02	0.59
1:A:6:PRO:HA	1:A:368:GLN:HG3	1.84	0.59
1:B:372:VAL:O	1:B:373:SER:C	2.40	0.59
1:D:212:TYR:CB	3:D:581:HOH:O	2.51	0.59
1:B:31:HIS:CE1	3:B:521:HOH:O	2.55	0.58
1:C:21:HIS:CE1	1:C:242:GLU:OE1	2.56	0.58
1:C:196:VAL:HG12	3:C:905:HOH:O	2.03	0.58
1:D:150:ALA:O	1:D:154:VAL:HG12	2.02	0.58
1:B:176:ASN:C	1:B:177:VAL:O	2.40	0.58
1:B:28:GLU:OE1	3:B:508:HOH:O	2.17	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:337:ARG:CD	3:C:671:HOH:O	2.51	0.57
1:A:259:LEU:HD22	1:A:279:LEU:HB2	1.86	0.57
1:A:295:SER:HB2	3:A:659:HOH:O	2.04	0.57
1:C:258:PRO:O	3:C:508:HOH:O	2.17	0.57
1:C:192:HIS:O	1:C:193:ASP:C	2.42	0.57
1:B:371:ILE:HB	1:D:80:ILE:HD11	1.87	0.57
1:D:242:GLU:CD	3:D:533:HOH:O	2.42	0.57
1:D:242:GLU:HG3	3:D:1011:HOH:O	2.04	0.57
1:A:30:THR:OG1	1:A:218:THR:CB	2.53	0.56
1:A:171:THR:C	1:A:173:ASP:H	2.09	0.56
1:A:191:THR:O	1:A:191:THR:CG2	2.52	0.56
1:B:293:ASN:O	3:B:510:HOH:O	2.17	0.56
1:B:80:ILE:CD1	3:B:849:HOH:O	2.53	0.56
1:D:230:ILE:HG13	1:D:230:ILE:O	2.06	0.56
1:A:30:THR:OG1	1:A:218:THR:HB	2.06	0.56
1:C:133:ASP:OD2	1:C:137:ASN:HA	2.06	0.55
1:C:2:ASN:N	3:C:521:HOH:O	2.38	0.55
1:D:50:GLU:O	1:D:54:LEU:HD12	2.06	0.55
1:B:214:GLU:HB2	3:B:742:HOH:O	2.06	0.55
1:C:298:ARG:CZ	3:C:501:HOH:O	2.41	0.55
1:A:177:VAL:O	1:A:178:ASP:CG	2.45	0.55
2:D:402:PLP:H5A2	3:D:541:HOH:O	2.07	0.55
1:D:156:ALA:C	3:D:518:HOH:O	2.45	0.55
1:D:196:VAL:HG12	3:D:795:HOH:O	2.06	0.55
1:A:118:GLN:O	1:D:92:ARG:NH1	2.40	0.55
1:A:80:ILE:O	1:A:80:ILE:HG13	2.06	0.55
1:C:21:HIS:NE2	1:C:242:GLU:OE1	2.40	0.54
1:C:206:THR:OG1	1:C:223:LEU:HD12	2.07	0.54
1:B:193:ASP:HA	3:B:675:HOH:O	2.06	0.54
1:C:230:ILE:HG13	1:C:230:ILE:O	2.06	0.54
1:B:66:ALA:O	3:B:509:HOH:O	2.17	0.54
1:B:24:LYS:HB2	3:B:732:HOH:O	2.08	0.54
1:D:51:MET:HA	1:D:54:LEU:CD1	2.37	0.53
1:D:51:MET:HA	1:D:54:LEU:HD12	1.91	0.53
1:C:259:LEU:HD21	1:C:321:PHE:CE1	2.43	0.53
1:D:331:THR:HG23	3:D:822:HOH:O	2.08	0.53
1:B:174:THR:CG2	1:B:176:ASN:H	2.20	0.52
1:D:174:THR:HG23	1:D:176:ASN:H	1.75	0.52
1:B:131:PRO:HB3	1:B:166:TRP:CZ2	2.45	0.52
1:B:138:ARG:HG3	1:B:139:ILE:N	2.23	0.52
1:A:110:GLU:HB2	3:A:717:HOH:O	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:248:SER:HB2	1:B:335:ILE:O	2.10	0.51
1:A:138:ARG:HG3	1:A:139:ILE:H	1.76	0.51
1:B:264:ALA:HA	1:B:271:TYR:O	2.10	0.51
1:C:12:ILE:HG23	1:C:247:LEU:HD22	1.91	0.51
1:B:261:LYS:HD2	3:B:673:HOH:O	2.10	0.51
1:C:25:VAL:HG21	1:C:242:GLU:OE1	2.11	0.51
1:D:331:THR:CG2	3:D:822:HOH:O	2.59	0.51
1:C:15:SER:OG	1:C:373:SER:HB2	2.11	0.51
1:D:312:GLY:HA3	3:D:752:HOH:O	2.11	0.51
2:D:402:PLP:C5A	3:D:541:HOH:O	2.59	0.51
1:D:196:VAL:HG11	3:D:878:HOH:O	2.10	0.50
1:D:372:VAL:O	1:D:373:SER:CB	2.48	0.50
1:B:138:ARG:CD	3:B:885:HOH:O	2.42	0.50
1:D:253:VAL:HG23	1:D:330:THR:O	2.11	0.50
1:D:324:ASN:HB2	1:D:325:PRO:HD2	1.93	0.50
1:C:136:LEU:HD13	1:C:138:ARG:HB2	1.93	0.50
1:A:251:ALA:HB1	1:A:283:PRO:HD2	1.93	0.50
1:A:252:LYS:NZ	1:A:331:THR:HG22	2.27	0.50
1:A:362:ASP:OD1	3:A:506:HOH:O	2.20	0.50
1:B:295:SER:HB2	3:B:671:HOH:O	2.11	0.50
1:D:174:THR:CG2	1:D:176:ASN:H	2.25	0.50
1:A:174:THR:CG2	1:A:176:ASN:H	2.24	0.49
1:A:178:ASP:OD1	1:A:178:ASP:C	2.50	0.49
1:C:230:ILE:CG2	3:C:938:HOH:O	2.60	0.49
1:C:108:THR:HG23	1:C:110:GLU:HG2	1.95	0.49
1:B:244:ILE:HG13	1:B:244:ILE:O	2.12	0.49
1:C:214:GLU:HB2	3:C:799:HOH:O	2.11	0.49
1:C:206:THR:OG1	2:C:401:PLP:O3P	2.24	0.49
1:C:286:HIS:HA	1:C:290:TYR:O	2.12	0.49
1:A:253:VAL:HG11	1:A:280:ILE:HD12	1.94	0.49
1:B:101:ASN:HB2	3:B:934:HOH:O	2.11	0.49
1:B:112:ILE:HG21	1:B:154:VAL:HG12	1.94	0.49
1:B:312:GLY:HA3	3:B:776:HOH:O	2.13	0.49
1:C:242:GLU:CD	1:C:242:GLU:H	2.14	0.49
1:D:131:PRO:HB3	1:D:166:TRP:CZ2	2.46	0.49
1:C:165:MET:HE2	1:C:198:PRO:HG3	1.95	0.49
1:C:352:THR:OG1	1:C:354:SER:HB2	2.13	0.49
1:D:31:HIS:HA	3:D:730:HOH:O	2.12	0.49
1:B:253:VAL:HG22	1:B:280:ILE:HG23	1.95	0.49
1:D:234:VAL:HA	3:D:590:HOH:O	2.13	0.49
1:A:298:ARG:NE	3:A:525:HOH:O	2.46	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:252:LYS:HZ1	1:B:331:THR:HG22	1.77	0.48
1:C:16:GLN:HB3	1:C:373:SER:C	2.33	0.48
1:D:312:GLY:HA3	3:D:819:HOH:O	2.12	0.48
1:B:150:ALA:O	1:B:154:VAL:HG13	2.14	0.48
1:B:178:ASP:HA	1:B:181:ASP:CB	2.42	0.48
1:B:312:GLY:N	3:B:538:HOH:O	2.46	0.48
1:C:24:LYS:CE	3:C:777:HOH:O	2.50	0.48
1:C:2:ASN:N	3:C:531:HOH:O	2.45	0.48
1:B:12:ILE:HG23	1:B:247:LEU:HD22	1.94	0.48
1:A:371:ILE:HB	1:C:80:ILE:HD11	1.93	0.48
1:D:125:THR:HG23	1:D:161:ILE:HD13	1.95	0.48
1:C:362:ASP:OD1	3:C:509:HOH:O	2.20	0.48
1:B:252:LYS:NZ	1:B:331:THR:HG22	2.28	0.48
1:A:177:VAL:O	1:A:178:ASP:OD1	2.31	0.48
1:C:2:ASN:N	3:C:533:HOH:O	2.46	0.48
1:B:337:ARG:NH1	3:B:539:HOH:O	2.46	0.48
1:B:284:ILE:HD12	1:B:288:ASP:HB2	1.96	0.47
1:B:259:LEU:HD21	1:B:321:PHE:CE1	2.49	0.47
1:D:12:ILE:HG23	1:D:247:LEU:HD22	1.96	0.47
1:A:108:THR:HG22	1:A:111:SER:N	2.20	0.47
1:C:131:PRO:HB3	1:C:166:TRP:CZ2	2.50	0.47
1:B:92:ARG:HG2	1:C:118:GLN:OE1	2.15	0.47
1:D:252:LYS:NZ	1:D:331:THR:HG22	2.27	0.47
1:A:244:ILE:HD12	1:A:245:PRO:HD2	1.96	0.46
1:A:337:ARG:NH1	1:A:340:ASP:HA	2.30	0.46
1:D:212:TYR:CG	3:D:581:HOH:O	2.58	0.46
1:A:236:ILE:HG13	3:A:673:HOH:O	2.15	0.46
2:B:402:PLP:H6	3:B:503:HOH:O	2.04	0.46
1:C:171:THR:HB	1:C:174:THR:HG21	1.96	0.46
1:D:106:PHE:O	1:D:131:PRO:HD2	2.15	0.46
1:A:252:LYS:HZ1	1:A:331:THR:HG22	1.78	0.46
1:A:22:ASN:O	1:A:26:ILE:HG13	2.15	0.46
1:B:260:LYS:HE2	1:B:261:LYS:O	2.15	0.46
1:D:165:MET:CE	1:D:201:LYS:HE3	2.45	0.46
1:A:2:ASN:HA	3:A:667:HOH:O	2.15	0.46
1:B:202:HIS:ND1	1:B:222:ARG:HD3	2.31	0.46
1:A:65:MET:HE2	1:A:70:GLU:HB3	1.98	0.45
1:C:122:THR:CG2	1:C:124:LEU:H	2.29	0.45
1:B:253:VAL:HG11	1:B:280:ILE:HD12	1.97	0.45
1:A:286:HIS:HA	1:A:290:TYR:O	2.16	0.45
1:A:295:SER:CB	3:A:659:HOH:O	2.62	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:152:GLN:HG3	3:C:581:HOH:O	2.17	0.45
1:D:16:GLN:HA	1:D:54:LEU:HD23	1.99	0.45
1:D:298:ARG:HD2	3:D:508:HOH:O	2.16	0.45
1:A:72:LEU:HG	3:A:887:HOH:O	2.17	0.45
1:B:144:ARG:O	1:B:148:VAL:HG23	2.17	0.45
1:A:131:PRO:HB3	1:A:166:TRP:CZ2	2.52	0.45
1:B:178:ASP:OD1	1:B:179:TYR:N	2.49	0.45
1:D:280:ILE:HG13	1:D:320:ILE:HB	1.98	0.45
1:D:2:ASN:CA	3:D:501:HOH:O	2.38	0.45
1:B:138:ARG:CG	3:B:885:HOH:O	2.64	0.44
1:D:191:THR:O	1:D:191:THR:HG23	2.16	0.44
1:A:256:SER:OG	1:A:326:VAL:O	2.26	0.44
1:D:187:TRP:NE1	1:D:191:THR:HG21	2.33	0.44
1:D:214:GLU:HB2	3:D:770:HOH:O	2.17	0.44
1:C:119:LEU:O	1:C:122:THR:CG2	2.65	0.44
1:C:106:PHE:O	1:C:131:PRO:HD2	2.18	0.44
1:C:253:VAL:HG11	1:C:280:ILE:HD12	1.99	0.44
1:D:286:HIS:HA	1:D:290:TYR:O	2.17	0.44
1:A:202:HIS:HB3	1:A:220:ILE:HB	1.99	0.44
1:B:113:LYS:O	1:B:117:GLU:HG3	2.18	0.43
3:A:864:HOH:O	1:B:254:VAL:HG12	2.17	0.43
1:D:234:VAL:HG12	1:D:235:PRO:HD3	2.00	0.43
1:A:165:MET:O	1:A:201:LYS:HA	2.18	0.43
1:B:169:PHE:HA	1:B:183:GLN:HE22	1.83	0.43
1:C:230:ILE:CB	3:C:938:HOH:O	2.35	0.43
1:C:259:LEU:HD21	1:C:321:PHE:HE1	1.81	0.43
1:D:165:MET:HE2	1:D:198:PRO:HG3	2.00	0.43
1:D:242:GLU:CD	1:D:242:GLU:H	2.21	0.43
1:D:62:ARG:CD	3:D:587:HOH:O	2.51	0.43
1:D:291:PRO:HG2	1:D:348:LEU:HG	2.01	0.43
1:D:279:LEU:HD11	1:D:319:MET:HB3	2.01	0.43
1:B:280:ILE:HG22	1:B:280:ILE:O	2.18	0.43
1:D:191:THR:O	1:D:191:THR:CG2	2.66	0.42
1:C:22:ASN:O	1:C:26:ILE:HD13	2.19	0.42
1:C:314:ALA:HB3	1:C:317:GLN:O	2.19	0.42
1:A:358:ALA:HB2	1:B:292:PHE:CD1	2.55	0.42
1:C:192:HIS:NE2	3:C:511:HOH:O	2.26	0.42
1:C:253:VAL:HG22	1:C:280:ILE:HG23	2.00	0.42
1:B:177:VAL:HG13	1:B:212:TYR:CE1	2.54	0.42
1:A:191:THR:O	1:A:191:THR:HG23	2.20	0.42
1:A:331:THR:HG23	3:A:725:HOH:O	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:260:LYS:HG2	3:D:961:HOH:O	2.19	0.42
1:C:298:ARG:HD3	3:C:790:HOH:O	2.20	0.42
1:D:197:GLU:OE2	1:D:201:LYS:NZ	2.48	0.42
1:B:8:ARG:HD2	1:B:283:PRO:HB2	2.01	0.42
1:B:278:TYR:CE2	1:B:325:PRO:HD3	2.55	0.42
1:D:50:GLU:O	1:D:54:LEU:CD1	2.66	0.42
1:A:298:ARG:HB2	1:A:343:ILE:HD11	2.02	0.41
1:C:17:SER:H	1:C:373:SER:CB	2.33	0.41
1:D:250:LYS:CE	3:D:507:HOH:O	2.54	0.41
1:B:322:VAL:HA	3:B:755:HOH:O	2.18	0.41
1:D:171:THR:HB	1:D:174:THR:HG21	2.02	0.41
1:C:244:ILE:HA	1:C:245:PRO:HD3	1.88	0.41
1:B:142:LYS:HD2	1:B:142:LYS:HA	1.89	0.41
1:C:136:LEU:HD23	1:C:170:ALA:HB2	2.02	0.41
1:C:17:SER:H	1:C:373:SER:HB3	1.84	0.41
1:C:347:ASP:HB3	3:C:915:HOH:O	2.20	0.41
1:A:256:SER:OG	1:A:280:ILE:HD11	2.20	0.41
1:A:346:GLN:HE21	1:A:350:GLU:HG3	1.84	0.41
1:B:331:THR:HG23	3:B:805:HOH:O	2.20	0.41
1:C:194:LEU:HB3	1:C:195:PRO:HD2	2.02	0.41
1:C:202:HIS:HB3	1:C:220:ILE:HB	2.02	0.41
1:C:22:ASN:O	1:C:26:ILE:CD1	2.68	0.41
1:D:233:SER:C	1:D:234:VAL:O	2.52	0.41
1:C:24:LYS:HE3	1:C:24:LYS:HB2	1.78	0.41
1:D:272:HIS:HD2	2:D:402:PLP:O2P	2.04	0.41
1:A:164:SER:HB3	1:A:200:GLU:HG3	2.02	0.41
1:B:2:ASN:HA	3:B:817:HOH:O	2.21	0.41
1:D:62:ARG:NH2	1:D:200:GLU:OE2	2.48	0.41
1:A:316:ASP:HB2	3:A:534:HOH:O	2.21	0.40
1:B:113:LYS:HD3	1:B:153:GLU:OE2	2.21	0.40
1:B:171:THR:HB	1:B:174:THR:HG21	2.03	0.40
1:B:244:ILE:HA	1:B:245:PRO:HD3	1.92	0.40
1:B:25:VAL:HG13	1:B:241:PHE:CG	2.57	0.40
1:C:122:THR:HG23	1:C:124:LEU:H	1.86	0.40
1:C:177:VAL:HG22	1:C:212:TYR:OH	2.21	0.40
1:A:108:THR:HG21	3:A:881:HOH:O	2.21	0.40
1:B:191:THR:HG23	1:B:191:THR:O	2.20	0.40
1:C:331:THR:HG23	3:C:800:HOH:O	2.22	0.40
1:C:6:PRO:HA	1:C:368:GLN:HG3	2.03	0.40
1:D:235:PRO:CD	3:D:531:HOH:O	2.33	0.40
1:B:206:THR:OG1	1:B:223:LEU:HD12	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:3:TYR:HB2	3:D:605:HOH:O	2.19	0.40
1:A:145:GLU:HG2	3:A:777:HOH:O	2.21	0.40
1:B:259:LEU:HD21	1:B:321:PHE:HE1	1.87	0.40
1:C:372:VAL:HG23	1:C:373:SER:N	2.36	0.40

All (10) 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)
3:B:891:HOH:O	3:C:980:HOH:O[1_554]	1.63	0.57
3:B:1042:HOH:O	3:D:1054:HOH:O[2_546]	1.91	0.29
1:A:242:GLU:OE2	1:B:31:HIS:NE2[1_655]	1.99	0.21
3:A:674:HOH:O	3:C:847:HOH:O[2_656]	2.01	0.19
1:C:298:ARG:NE	2:B:402:PLP:O1P[1_556]	2.13	0.07
1:A:242:GLU:OE1	1:B:31:HIS:NE2[1_655]	2.13	0.07
3:D:751:HOH:O	3:D:945:HOH:O[2_546]	2.14	0.06
3:C:567:HOH:O	3:D:737:HOH:O[1_655]	2.15	0.05
3:B:957:HOH:O	3:C:935:HOH:O[2_546]	2.16	0.04
3:A:564:HOH:O	3:D:518:HOH:O[2_545]	2.16	0.04

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	370/381 (97%)	354 (96%)	15 (4%)	1 (0%)	41	41
1	B	370/381 (97%)	354 (96%)	14 (4%)	2 (0%)	29	26
1	C	370/381 (97%)	358 (97%)	10 (3%)	2 (0%)	29	26
1	D	370/381 (97%)	356 (96%)	13 (4%)	1 (0%)	41	41
All	All	1480/1524 (97%)	1422 (96%)	52 (4%)	6 (0%)	34	32

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	193	ASP
1	B	193	ASP
1	C	192	HIS
1	D	235	PRO
1	B	177	VAL
1	C	193	ASP

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	304/313 (97%)	284 (93%)	20 (7%)	16	14
1	B	304/313 (97%)	279 (92%)	25 (8%)	11	8
1	C	304/313 (97%)	274 (90%)	30 (10%)	8	5
1	D	304/313 (97%)	275 (90%)	29 (10%)	8	5
All	All	1216/1252 (97%)	1112 (91%)	104 (9%)	10	7

All (104) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	VAL
1	A	72	LEU
1	A	104	LEU
1	A	108	THR
1	A	136	LEU
1	A	138	ARG
1	A	152	GLN
1	A	174	THR
1	A	177	VAL
1	A	188	GLN
1	A	191	THR
1	A	193	ASP
1	A	244	ILE
1	A	247	LEU
1	A	260	LYS
1	A	280	ILE

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Mol	Chain	Res	Type
1	A	284	ILE
1	A	331	THR
1	A	334	LEU
1	A	368	GLN
1	B	25	VAL
1	B	72	LEU
1	B	80	ILE
1	B	84	ASP
1	B	104	LEU
1	B	108	THR
1	B	136	LEU
1	B	138	ARG
1	B	174	THR
1	B	178	ASP
1	B	185	SER
1	B	188	GLN
1	B	191	THR
1	B	193	ASP
1	B	244	ILE
1	B	247	LEU
1	B	260	LYS
1	B	275	ASN
1	B	280	ILE
1	B	284	ILE
1	B	295	SER
1	B	331	THR
1	B	334	LEU
1	B	354	SER
1	B	368	GLN
1	C	2	ASN
1	C	25	VAL
1	C	28	GLU
1	C	72	LEU
1	C	80	ILE
1	C	104	LEU
1	C	108	THR
1	C	122	THR
1	C	136	LEU
1	C	152	GLN
1	C	154	VAL
1	C	158	ASP
1	C	174	THR

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Mol	Chain	Res	Type
1	C	188	GLN
1	C	191	THR
1	C	192	HIS
1	C	193	ASP
1	C	230	ILE
1	C	242	GLU
1	C	244	ILE
1	C	247	LEU
1	C	260	LYS
1	C	270	GLU
1	C	280	ILE
1	C	284	ILE
1	C	331	THR
1	C	334	LEU
1	C	337	ARG
1	C	354	SER
1	C	368	GLN
1	D	25	VAL
1	D	28	GLU
1	D	54	LEU
1	D	62	ARG
1	D	72	LEU
1	D	80	ILE
1	D	108	THR
1	D	125	THR
1	D	136	LEU
1	D	154	VAL
1	D	155	SER
1	D	174	THR
1	D	177	VAL
1	D	185	SER
1	D	188	GLN
1	D	191	THR
1	D	230	ILE
1	D	242	GLU
1	D	244	ILE
1	D	247	LEU
1	D	255	ASN
1	D	270	GLU
1	D	284	ILE
1	D	295	SER
1	D	316	ASP

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Mol	Chain	Res	Type
1	D	324	ASN
1	D	331	THR
1	D	334	LEU
1	D	342	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	346	GLN
1	C	101	ASN

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](#)

6 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	D	401	1	15,15,16	6.52	10 (66%)	20,22,23	1.44	3 (15%)
2	PLP	B	401	1	15,15,16	6.37	9 (60%)	20,22,23	1.59	3 (15%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	PLP	C	401	1	15,15,16	6.50	8 (53%)	20,22,23	1.30	3 (15%)
2	PLP	A	401	1	15,15,16	6.41	8 (53%)	20,22,23	1.50	3 (15%)
2	PLP	D	402	1	15,15,16	6.76	10 (66%)	20,22,23	1.67	5 (25%)
2	PLP	B	402	1	15,15,16	6.58	8 (53%)	20,22,23	2.12	6 (30%)

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	D	401	1	-	0/6/6/8	0/1/1/1
2	PLP	B	401	1	-	2/6/6/8	0/1/1/1
2	PLP	C	401	1	-	1/6/6/8	0/1/1/1
2	PLP	A	401	1	-	2/6/6/8	0/1/1/1
2	PLP	D	402	1	-	0/6/6/8	0/1/1/1
2	PLP	B	402	1	-	0/6/6/8	0/1/1/1

All (53) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	402	PLP	C3-C2	15.03	1.55	1.40
2	D	402	PLP	C3-C2	14.79	1.55	1.40
2	D	401	PLP	C3-C2	14.45	1.55	1.40
2	B	401	PLP	C3-C2	14.44	1.55	1.40
2	A	401	PLP	C3-C2	14.44	1.55	1.40
2	C	401	PLP	C3-C2	14.37	1.55	1.40
2	D	402	PLP	C5-C4	12.18	1.54	1.40
2	C	401	PLP	C5-C4	11.26	1.53	1.40
2	D	401	PLP	C5-C4	11.16	1.52	1.40
2	B	402	PLP	C5-C4	11.05	1.52	1.40
2	A	401	PLP	C5-C4	11.00	1.52	1.40
2	B	401	PLP	C5-C4	10.73	1.52	1.40
2	D	402	PLP	C2-N1	9.69	1.52	1.33
2	D	401	PLP	C2-N1	9.60	1.51	1.33
2	B	402	PLP	C2-N1	9.57	1.51	1.33
2	A	401	PLP	C2-N1	9.54	1.51	1.33
2	C	401	PLP	C2-N1	9.40	1.51	1.33
2	B	401	PLP	C2-N1	9.37	1.51	1.33
2	D	402	PLP	C6-N1	8.76	1.52	1.34
2	D	401	PLP	C6-N1	8.66	1.52	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	PLP	C6-N1	8.63	1.52	1.34
2	A	401	PLP	C6-N1	8.55	1.52	1.34
2	C	401	PLP	C6-N1	8.49	1.52	1.34
2	B	402	PLP	C6-N1	8.40	1.52	1.34
2	D	402	PLP	C6-C5	8.24	1.54	1.37
2	B	402	PLP	C6-C5	8.09	1.54	1.37
2	A	401	PLP	C6-C5	7.91	1.54	1.37
2	D	401	PLP	C6-C5	7.83	1.54	1.37
2	C	401	PLP	C6-C5	7.79	1.54	1.37
2	B	401	PLP	C6-C5	7.77	1.53	1.37
2	B	401	PLP	C3-C4	5.98	1.52	1.40
2	C	401	PLP	C3-C4	5.95	1.52	1.40
2	D	402	PLP	C3-C4	5.80	1.52	1.40
2	A	401	PLP	C3-C4	5.73	1.52	1.40
2	D	401	PLP	C3-C4	5.73	1.52	1.40
2	B	402	PLP	C3-C4	5.66	1.51	1.40
2	D	402	PLP	P-O1P	4.70	1.65	1.50
2	C	401	PLP	P-O1P	4.63	1.65	1.50
2	D	401	PLP	P-O1P	4.57	1.65	1.50
2	B	402	PLP	P-O1P	4.38	1.64	1.50
2	B	402	PLP	P-O3P	2.89	1.66	1.54
2	D	401	PLP	P-O3P	2.82	1.65	1.54
2	C	401	PLP	P-O3P	2.79	1.65	1.54
2	B	401	PLP	P-O3P	2.76	1.65	1.54
2	A	401	PLP	P-O2P	2.74	1.65	1.54
2	A	401	PLP	P-O3P	2.73	1.65	1.54
2	D	402	PLP	P-O3P	2.70	1.65	1.54
2	B	401	PLP	P-O2P	2.61	1.64	1.54
2	B	401	PLP	C2A-C2	2.14	1.54	1.50
2	D	401	PLP	C2A-C2	2.11	1.53	1.50
2	D	402	PLP	C2A-C2	2.06	1.53	1.50
2	D	402	PLP	P-O2P	-2.03	1.47	1.54
2	D	401	PLP	P-O2P	-2.00	1.47	1.54

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	402	PLP	C6-C5-C4	5.51	122.50	118.16
2	A	401	PLP	O4P-C5A-C5	-4.07	101.59	109.35
2	B	401	PLP	O4P-C5A-C5	-4.03	101.67	109.35
2	D	402	PLP	C4A-C4-C5	4.01	125.06	120.94
2	B	402	PLP	O4P-C5A-C5	-3.79	102.12	109.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	PLP	C6-C5-C4	3.66	121.04	118.16
2	D	401	PLP	O4P-C5A-C5	-3.58	102.53	109.35
2	A	401	PLP	C6-C5-C4	3.45	120.88	118.16
2	B	402	PLP	C4A-C4-C5	3.37	124.41	120.94
2	B	401	PLP	C5-C6-N1	-3.08	118.69	123.82
2	C	401	PLP	O4P-C5A-C5	-3.06	103.52	109.35
2	B	402	PLP	C5A-C5-C6	-3.00	114.43	119.37
2	D	401	PLP	C6-C5-C4	2.94	120.47	118.16
2	D	401	PLP	C5-C6-N1	-2.88	119.03	123.82
2	C	401	PLP	C6-C5-C4	2.82	120.37	118.16
2	A	401	PLP	C5-C6-N1	-2.74	119.25	123.82
2	C	401	PLP	C5-C6-N1	-2.71	119.31	123.82
2	D	402	PLP	O4P-C5A-C5	-2.61	104.37	109.35
2	D	402	PLP	C6-C5-C4	2.61	120.21	118.16
2	D	402	PLP	C5A-C5-C6	-2.58	115.12	119.37
2	B	402	PLP	O2P-P-O4P	2.57	113.58	106.73
2	D	402	PLP	C4A-C4-C3	-2.52	116.23	120.50
2	B	402	PLP	C5-C6-N1	-2.16	120.22	123.82

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	PLP	C4-C5-C5A-O4P
2	B	401	PLP	C6-C5-C5A-O4P
2	C	401	PLP	C6-C5-C5A-O4P
2	A	401	PLP	C6-C5-C5A-O4P
2	B	401	PLP	C4-C5-C5A-O4P

There are no ring outliers.

3 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	401	PLP	1	0
2	D	402	PLP	3	0
2	B	402	PLP	9	1

5.7 Other polymers ⓘ

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	372/381 (97%)	-0.08	2 (0%) 91 92	7, 12, 21, 38	0
1	B	372/381 (97%)	-0.08	4 (1%) 80 84	7, 12, 20, 36	0
1	C	372/381 (97%)	-0.08	3 (0%) 86 88	7, 12, 21, 38	0
1	D	372/381 (97%)	-0.18	0 100 100	6, 11, 19, 31	0
All	All	1488/1524 (97%)	-0.11	9 (0%) 89 91	6, 12, 21, 38	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	373	SER	3.5
1	B	31	HIS	3.0
1	C	31	HIS	2.9
1	A	193	ASP	2.6
1	B	177	VAL	2.5
1	B	340	ASP	2.3
1	C	373	SER	2.2
1	A	31	HIS	2.1
1	C	192	HIS	2.1

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	PLP	B	402	15/16	0.78	0.27	20,25,33,33	0
2	PLP	D	402	15/16	0.91	0.14	13,17,24,27	0
2	PLP	D	401	15/16	0.95	0.14	7,10,12,15	0
2	PLP	A	401	15/16	0.96	0.13	7,9,12,17	0
2	PLP	B	401	15/16	0.96	0.15	7,9,13,14	0
2	PLP	C	401	15/16	0.96	0.13	7,10,12,13	0

6.5 Other polymers [i](#)

There are no such residues in this entry.