



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:28 PM GMT

PDB ID : 1TTP
Title : TRYPTOPHAN SYNTHASE (E.C.4.2.1.20) IN THE PRESENCE OF CESIUM, ROOM TEMPERATURE
Authors : Rhee, S.; Parris, K.; Ahmed, S.; Miles, E.W.; Davies, D.R.
Deposited on : 1995-10-11
Resolution : 2.30 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

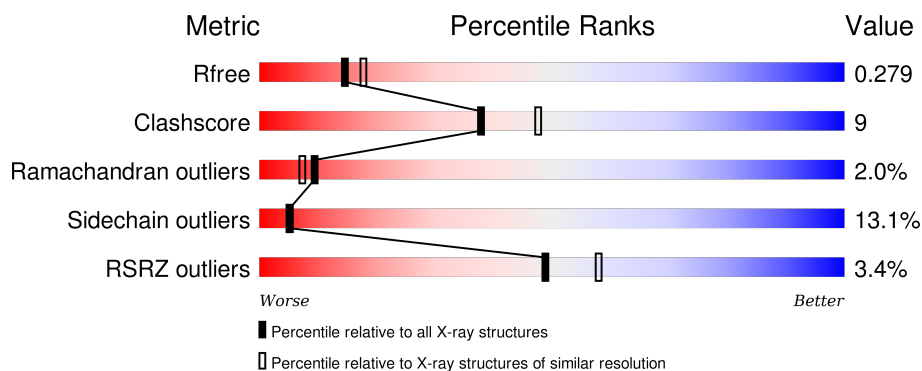
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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}	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	268	
2	B	397	

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 4918 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 TRYPTOPHAN SYNTHASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	256	Total	C	N	O	S	0	0	0
			1937	1236	333	360	8			

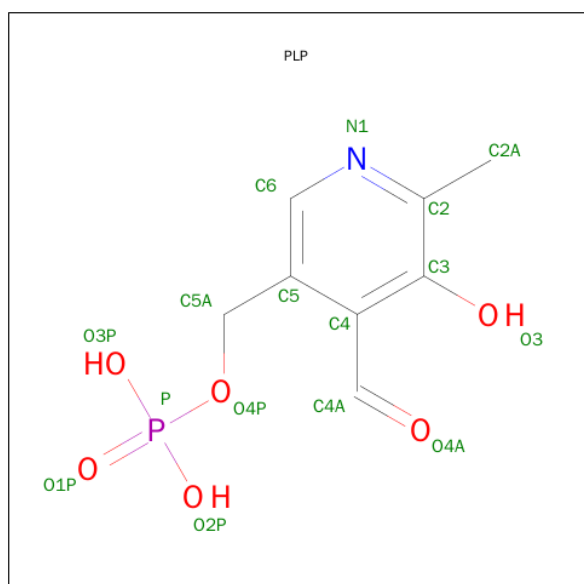
- Molecule 2 is a protein called TRYPTOPHAN SYNTHASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	389	Total	C	N	O	S	0	0	0
			2950	1855	518	558	19			

- Molecule 3 is CESIUM ION (three-letter code: CS) (formula: Cs).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	2	Total	Cs	0	0
			2	2		

- Molecule 4 is PYRIDOXAL-5'-PHOSPHATE (three-letter code: PLP) (formula: C₈H₁₀NO₆P).



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

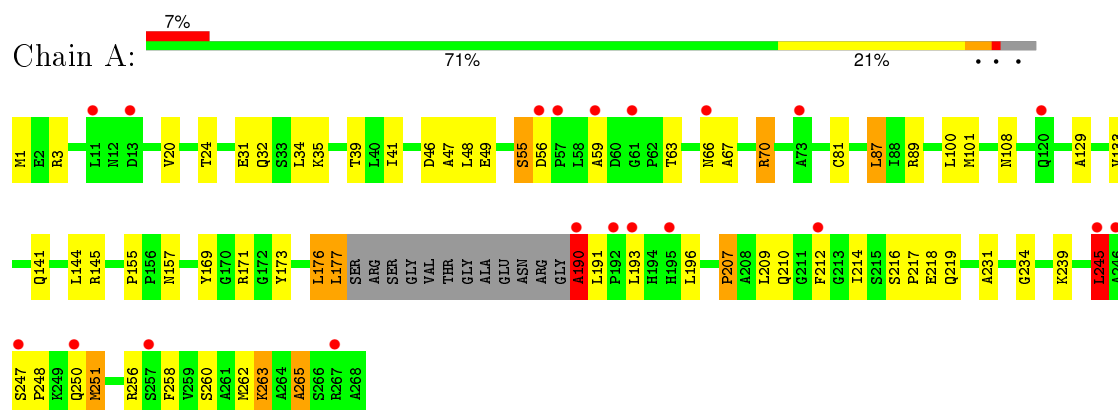
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	6	Total	O	0	0
			6	6		
5	B	8	Total	O	0	0
			8	8		

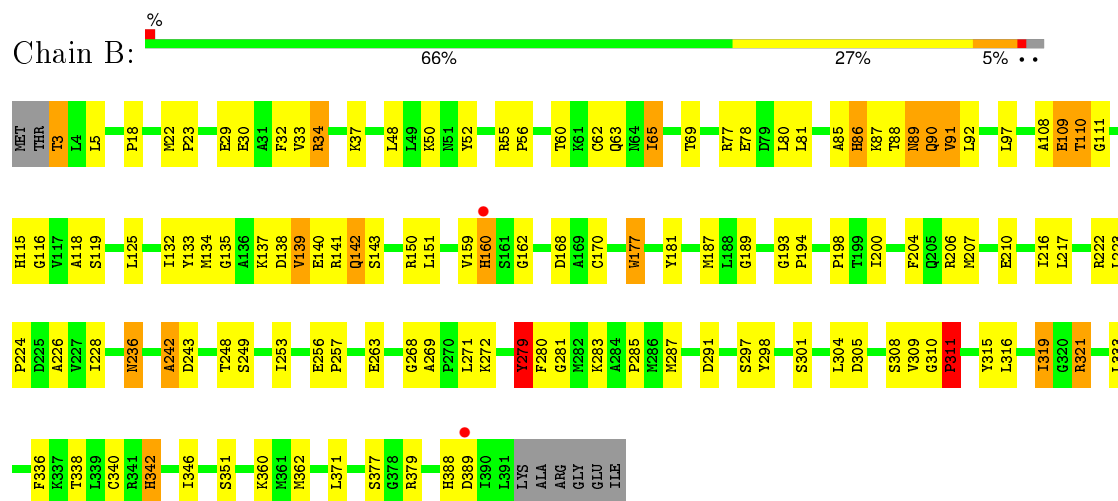
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: TRYPTOPHAN SYNTHASE



• Molecule 2: TRYPTOPHAN SYNTHASE



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	185.30 Å 61.30 Å 68.10 Å 90.00° 94.80° 90.00°	Depositor
Resolution (Å)	8.00 – 2.30 8.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	67.2 (8.00-2.30) 67.4 (8.00-2.30)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.39 (at 2.30 Å)	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.217 , 0.302 0.199 , 0.279	Depositor DCC
R_{free} test set	2230 reflections (9.95%)	DCC
Wilson B-factor (Å ²)	24.5	Xtriage
Anisotropy	0.929	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 82.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Outliers	0 of 22409 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4918	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.39% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: CS, 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	2.15	9/1976 (0.5%)	1.71	27/2684 (1.0%)
2	B	0.94	1/3008 (0.0%)	1.49	26/4064 (0.6%)
All	All	1.54	10/4984 (0.2%)	1.58	53/6748 (0.8%)

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	3

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	190	ALA	N-CA	60.83	2.68	1.46
1	A	190	ALA	C-O	42.89	2.04	1.23
1	A	177	LEU	N-CA	28.91	2.04	1.46
1	A	177	LEU	CA-CB	25.72	2.12	1.53
1	A	177	LEU	CG-CD2	21.31	2.30	1.51
1	A	190	ALA	CA-CB	14.57	1.83	1.52
1	A	190	ALA	C-N	12.59	1.63	1.34
1	A	177	LEU	CB-CG	-9.66	1.24	1.52
1	A	177	LEU	C-O	-6.83	1.10	1.23
2	B	87	LYS	CB-CG	6.66	1.70	1.52

All (53) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	177	LEU	CB-CG-CD2	22.18	148.72	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	176	LEU	O-C-N	19.39	153.73	122.70
1	A	177	LEU	CA-CB-CG	-18.91	71.81	115.30
1	A	176	LEU	C-N-CA	-18.56	75.31	121.70
1	A	177	LEU	CD1-CG-CD2	-16.71	60.37	110.50
1	A	176	LEU	CA-C-N	-16.55	80.79	117.20
1	A	190	ALA	O-C-N	-11.18	104.82	122.70
1	A	190	ALA	N-CA-C	10.85	140.29	111.00
1	A	190	ALA	CA-C-N	10.81	140.99	117.20
1	A	190	ALA	CA-C-O	-10.72	97.59	120.10
1	A	177	LEU	CA-C-O	-10.45	98.16	120.10
2	B	279	TYR	CB-CG-CD2	-10.16	114.91	121.00
1	A	190	ALA	N-CA-CB	-10.07	96.00	110.10
1	A	169	TYR	CB-CG-CD2	-8.61	115.83	121.00
2	B	379	ARG	NE-CZ-NH2	-8.50	116.05	120.30
2	B	222	ARG	NE-CZ-NH2	-8.46	116.07	120.30
2	B	279	TYR	CB-CG-CD1	8.23	125.94	121.00
1	A	177	LEU	CB-CG-CD1	-8.16	97.12	111.00
2	B	222	ARG	NE-CZ-NH1	8.03	124.32	120.30
2	B	34	ARG	NE-CZ-NH1	8.01	124.30	120.30
1	A	177	LEU	CB-CA-C	-7.95	95.09	110.20
2	B	379	ARG	NE-CZ-NH1	7.77	124.19	120.30
1	A	3	ARG	NE-CZ-NH1	7.72	124.16	120.30
1	A	190	ALA	C-N-CA	7.68	140.91	121.70
2	B	91	VAL	CG1-CB-CG2	-6.92	99.83	110.90
1	A	171	ARG	NE-CZ-NH2	-6.91	116.85	120.30
1	A	171	ARG	NE-CZ-NH1	6.66	123.63	120.30
1	A	145	ARG	NE-CZ-NH1	6.55	123.58	120.30
2	B	177	TRP	CE2-CD2-CG	-6.51	102.09	107.30
2	B	315	TYR	CB-CG-CD2	-6.49	117.11	121.00
2	B	177	TRP	CD1-CG-CD2	6.41	111.43	106.30
1	A	245	LEU	CA-CB-CG	6.37	129.95	115.30
2	B	3	THR	CA-CB-CG2	6.37	121.31	112.40
1	A	190	ALA	CB-CA-C	-6.23	100.75	110.10
2	B	80	LEU	CA-CB-CG	6.21	129.57	115.30
2	B	150	ARG	NE-CZ-NH1	6.05	123.32	120.30
2	B	88	THR	CA-C-N	5.85	130.08	117.20
2	B	287	MET	CA-CB-CG	-5.80	103.44	113.30
2	B	150	ARG	CA-CB-CG	5.79	126.14	113.40
2	B	55	ARG	NE-CZ-NH2	-5.71	117.44	120.30
2	B	55	ARG	NE-CZ-NH1	5.66	123.13	120.30
1	A	87	LEU	CB-CG-CD2	-5.41	101.80	111.00
2	B	389	ASP	CA-CB-CG	5.40	125.27	113.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	251	MET	CA-CB-CG	5.26	122.25	113.30
2	B	56	PRO	CA-C-N	5.25	128.76	117.20
2	B	177	TRP	CG-CD2-CE3	5.18	138.56	133.90
2	B	206	ARG	NE-CZ-NH1	5.14	122.87	120.30
1	A	133	VAL	CA-CB-CG2	-5.12	103.22	110.90
1	A	81	CYS	CA-CB-SG	-5.12	104.78	114.00
1	A	70	ARG	NE-CZ-NH2	-5.12	117.74	120.30
2	B	3	THR	CA-CB-OG1	-5.11	98.26	109.00
2	B	298	TYR	CB-CG-CD2	-5.08	117.95	121.00
2	B	177	TRP	CB-CG-CD1	-5.08	120.40	127.00

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	1	MET	Mainchain
1	A	190	ALA	Mainchain
1	A	247	SER	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	1937	0	1949	36	0
2	B	2950	0	2924	59	0
3	B	2	0	0	0	1
4	B	15	0	7	3	0
5	A	6	0	0	0	0
5	B	8	0	0	1	0
All	All	4918	0	4880	90	1

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 9.

All (90) 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:A:177:LEU:HG	1:A:177:LEU:CA	1.32	1.54
1:A:190:ALA:CA	1:A:190:ALA:CB	1.83	1.52
1:A:177:LEU:HD11	1:A:177:LEU:CD2	1.47	1.41
1:A:177:LEU:CD2	1:A:177:LEU:CD1	2.06	1.31
1:A:177:LEU:CA	1:A:177:LEU:CG	2.10	1.28
1:A:177:LEU:CA	1:A:177:LEU:CB	2.12	1.27
1:A:176:LEU:C	1:A:177:LEU:CA	2.09	1.20
1:A:177:LEU:CA	1:A:177:LEU:N	2.04	1.19
1:A:177:LEU:CG	1:A:177:LEU:CD2	2.30	1.07
1:A:177:LEU:HG	1:A:177:LEU:HA	1.39	1.01
1:A:177:LEU:HD11	1:A:177:LEU:HD22	1.02	0.98
1:A:190:ALA:O	1:A:190:ALA:C	2.04	0.95
1:A:176:LEU:C	1:A:177:LEU:HA	1.85	0.94
1:A:176:LEU:O	1:A:177:LEU:HA	1.68	0.92
1:A:177:LEU:HD22	1:A:177:LEU:CD1	1.90	0.87
2:B:89:ASN:HD22	2:B:89:ASN:H	1.25	0.85
2:B:216:ILE:HG21	2:B:224:PRO:HD3	1.59	0.85
2:B:194:PRO:HG3	2:B:281:GLY:HA3	1.61	0.83
2:B:89:ASN:H	2:B:89:ASN:ND2	1.92	0.68
1:A:190:ALA:C	1:A:190:ALA:CB	2.61	0.67
1:A:214:ILE:HG12	1:A:231:ALA:HB1	1.77	0.67
2:B:256:GLU:HG2	2:B:269:ALA:HA	1.78	0.64
2:B:236:ASN:HD22	4:B:402:PLP:H6	1.66	0.61
2:B:90:GLN:HA	2:B:204:PHE:HB3	1.82	0.60
1:A:177:LEU:C	1:A:177:LEU:CB	2.70	0.60
2:B:108:ALA:HB3	2:B:132:ILE:HD13	1.83	0.60
1:A:39:THR:HG23	1:A:256:ARG:NH1	2.17	0.59
1:A:108:ASN:HD21	2:B:291:ASP:HA	1.67	0.59
2:B:338:THR:O	2:B:342:HIS:HB2	2.03	0.57
1:A:190:ALA:CA	1:A:190:ALA:N	2.68	0.57
2:B:62:CYS:SG	2:B:65:ILE:HD11	2.44	0.57
2:B:111:GLY:O	2:B:138:ASP:HB3	2.06	0.55
2:B:60:THR:OG1	2:B:77:ARG:HD2	2.06	0.55
2:B:271:LEU:HD12	2:B:309:VAL:HG11	1.88	0.55
2:B:48:LEU:HB3	2:B:92:LEU:HD13	1.88	0.55
2:B:340:CYS:SG	2:B:346:ILE:HG12	2.47	0.54
2:B:137:LYS:O	2:B:141:ARG:HG2	2.07	0.54
2:B:116:GLY:HA2	2:B:132:ILE:HD12	1.90	0.53
2:B:3:THR:HG22	2:B:5:LEU:O	2.08	0.53
2:B:89:ASN:HD22	2:B:89:ASN:N	2.02	0.53
1:A:67:ALA:HA	1:A:70:ARG:HD3	1.90	0.53
2:B:110:THR:HG23	2:B:115:HIS:HB3	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:39:THR:HG23	1:A:256:ARG:HH11	1.74	0.52
2:B:52:TYR:CE1	2:B:125:LEU:HD13	2.45	0.51
2:B:226:ALA:HB1	2:B:253:ILE:CD1	2.41	0.51
2:B:236:ASN:HB3	4:B:402:PLP:O3P	2.10	0.51
2:B:91:VAL:HG11	2:B:118:ALA:O	2.10	0.51
1:A:59:ALA:HB2	2:B:18:PRO:HG3	1.93	0.50
2:B:301:SER:HB2	2:B:304:LEU:HB2	1.92	0.50
2:B:242:ALA:HB3	5:B:412:HOH:O	2.12	0.50
2:B:134:MET:SD	2:B:139:VAL:HG23	2.51	0.50
2:B:142:GLN:HA	2:B:142:GLN:HE21	1.78	0.49
1:A:56:ASP:HB2	2:B:279:TYR:OH	2.12	0.49
1:A:157:ASN:ND2	2:B:23:PRO:HG3	2.29	0.48
2:B:30:GLU:O	2:B:34:ARG:HG3	2.13	0.48
2:B:177:TRP:O	2:B:181:TYR:HB3	2.13	0.48
2:B:63:GLN:HA	2:B:63:GLN:HE21	1.79	0.47
2:B:193:GLY:HA2	2:B:280:PHE:O	2.14	0.47
2:B:109:GLU:HA	2:B:133:TYR:O	2.14	0.47
2:B:377:SER:C	4:B:402:PLP:H2A2	2.35	0.46
1:A:248:PRO:O	1:A:251:MET:HB3	2.15	0.46
2:B:91:VAL:HG13	2:B:187:MET:SD	2.56	0.46
2:B:125:LEU:HG	2:B:125:LEU:O	2.16	0.45
1:A:258:PHE:O	1:A:262:MET:HG2	2.17	0.45
2:B:115:HIS:CE1	2:B:189:GLY:HA2	2.52	0.45
2:B:69:THR:HG21	2:B:362:MET:HB2	1.98	0.45
2:B:29:GLU:O	2:B:33:VAL:HG23	2.18	0.44
2:B:135:GLY:HA2	2:B:159:VAL:HB	1.99	0.44
1:A:173:TYR:HB2	1:A:209:LEU:HD13	1.98	0.44
2:B:32:PHE:CD1	2:B:200:ILE:HG12	2.53	0.44
1:A:196:LEU:HD23	1:A:196:LEU:HA	1.85	0.44
2:B:321:ARG:O	2:B:321:ARG:HD3	2.18	0.43
2:B:162:GLY:HA3	2:B:168:ASP:OD1	2.18	0.43
2:B:78:GLU:HB2	2:B:377:SER:HA	1.99	0.43
2:B:316:LEU:HA	2:B:319:ILE:CD1	2.49	0.43
1:A:41:ILE:HD11	1:A:48:LEU:HD11	2.01	0.42
2:B:81:LEU:HD22	2:B:86:HIS:HA	2.02	0.41
2:B:109:GLU:HG2	2:B:170:CYS:HA	2.03	0.41
2:B:257:PRO:HG2	2:B:268:GLY:O	2.20	0.41
2:B:272:LYS:HA	2:B:272:LYS:HD3	1.89	0.41
2:B:333:LEU:O	2:B:336:PHE:HB3	2.20	0.41
1:A:217:PRO:HB3	1:A:265:ALA:HB2	2.03	0.41
1:A:46:ASP:OD2	1:A:263:LYS:HE3	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:89:ARG:HH11	1:A:89:ARG:HD3	1.75	0.41
2:B:216:ILE:HD12	2:B:216:ILE:HA	1.73	0.40
2:B:90:GLN:HE21	2:B:90:GLN:HB2	1.62	0.40
2:B:310:GLY:HA2	2:B:311:PRO:HD2	1.78	0.40
1:A:108:ASN:ND2	2:B:291:ASP:HA	2.36	0.40
1:A:20:VAL:HG22	1:A:47:ALA:HB3	2.04	0.40
2:B:22:MET:N	2:B:23:PRO:HD2	2.37	0.40

All (1) 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:401:CS:CS	3:B:401:CS:CS[2_655]	1.88	0.32

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	252/268 (94%)	233 (92%)	12 (5%)	7 (3%)	6	4
2	B	387/397 (98%)	349 (90%)	32 (8%)	6 (2%)	12	11
All	All	639/665 (96%)	582 (91%)	44 (7%)	13 (2%)	9	7

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	55	SER
2	B	86	HIS
1	A	129	ALA
1	A	245	LEU
2	B	160	HIS
1	A	191	LEU

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Mol	Chain	Res	Type
1	A	207	PRO
2	B	85	ALA
2	B	242	ALA
2	B	311	PRO
1	A	265	ALA
1	A	234	GLY
2	B	297	SER

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	200/208 (96%)	173 (86%)	27 (14%)	5	4
2	B	305/311 (98%)	266 (87%)	39 (13%)	5	5
All	All	505/519 (97%)	439 (87%)	66 (13%)	5	5

All (66) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	THR
1	A	31	GLU
1	A	32	GLN
1	A	34	LEU
1	A	35	LYS
1	A	49	GLU
1	A	55	SER
1	A	63	THR
1	A	66	ASN
1	A	87	LEU
1	A	100	LEU
1	A	101	MET
1	A	141	GLN
1	A	144	LEU
1	A	155	PRO
1	A	193	LEU

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Mol	Chain	Res	Type
1	A	207	PRO
1	A	210	GLN
1	A	212	PHE
1	A	216	SER
1	A	218	GLU
1	A	219	GLN
1	A	239	LYS
1	A	245	LEU
1	A	250	GLN
1	A	260	SER
1	A	263	LYS
2	B	37	LYS
2	B	50	LYS
2	B	65	ILE
2	B	89	ASN
2	B	90	GLN
2	B	97	LEU
2	B	109	GLU
2	B	110	THR
2	B	119	SER
2	B	139	VAL
2	B	140	GLU
2	B	142	GLN
2	B	143	SER
2	B	151	LEU
2	B	160	HIS
2	B	198	PRO
2	B	207	MET
2	B	210	GLU
2	B	217	LEU
2	B	223	LEU
2	B	228	ILE
2	B	236	ASN
2	B	243	ASP
2	B	248	THR
2	B	249	SER
2	B	263	GLU
2	B	279	TYR
2	B	283	LYS
2	B	285	PRO
2	B	305	ASP
2	B	308	SER

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Mol	Chain	Res	Type
2	B	311	PRO
2	B	319	ILE
2	B	321	ARG
2	B	342	HIS
2	B	351	SER
2	B	360	LYS
2	B	371	LEU
2	B	388	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	10	GLN
2	B	26	ASN
2	B	42	GLN
2	B	63	GLN
2	B	89	ASN
2	B	90	GLN
2	B	94	GLN
2	B	142	GLN
2	B	195	HIS

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](#)

Of 3 ligands modelled in this entry, 2 are monoatomic - leaving 1 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	PLP	B	402	2	15,15,16	4.93	10 (66%)	21,22,23	2.45	7 (33%)

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
4	PLP	B	402	2	-	0/6/6/8	0/1/1/1

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	402	PLP	P-O3P	-5.64	1.34	1.54
4	B	402	PLP	O4P-C5A	-2.72	1.33	1.44
4	B	402	PLP	P-O4P	-2.59	1.51	1.60
4	B	402	PLP	C2-N1	2.72	1.39	1.34
4	B	402	PLP	C6-N1	3.71	1.42	1.34
4	B	402	PLP	C6-C5	5.07	1.48	1.37
4	B	402	PLP	C3-C4	6.93	1.57	1.40
4	B	402	PLP	C4A-C4	7.27	1.66	1.51
4	B	402	PLP	C5-C4	7.97	1.49	1.40
4	B	402	PLP	C3-C2	10.11	1.47	1.40

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	402	PLP	C3-C2-N1	-3.10	116.33	120.61
4	B	402	PLP	C6-C5-C4	-2.72	115.84	118.15
4	B	402	PLP	C5A-C5-C4	3.18	125.86	121.65
4	B	402	PLP	C6-N1-C2	3.72	126.86	119.28
4	B	402	PLP	O4P-C5A-C5	4.70	116.76	108.99
4	B	402	PLP	C2A-C2-C3	4.75	126.76	121.04
4	B	402	PLP	O3P-P-O4P	5.10	121.24	106.56

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	402	PLP	3	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	256/268 (95%)	0.19	20 (7%) 16 22	9, 38, 79, 86	4 (1%)
2	B	389/397 (97%)	-0.44	2 (0%) 91 94	3, 20, 56, 75	3 (0%)
All	All	645/665 (96%)	-0.19	22 (3%) 49 58	3, 26, 68, 86	7 (1%)

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	56	ASP	4.2
1	A	250	GLN	4.0
1	A	257	SER	4.0
2	B	389	ASP	3.9
1	A	190	ALA	3.5
1	A	245	LEU	3.3
1	A	192	PRO	3.2
1	A	195	HIS	2.9
1	A	59	ALA	2.9
1	A	193	LEU	2.9
1	A	66	ASN	2.9
1	A	73	ALA	2.8
1	A	212	PHE	2.6
1	A	61	GLY	2.6
2	B	160	HIS	2.5
1	A	247	SER	2.5
1	A	13	ASP	2.4
1	A	246	ALA	2.4
1	A	11	LEU	2.2
1	A	57	PRO	2.1
1	A	267	ARG	2.1
1	A	120	GLN	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	PLP	B	402	15/16	0.97	0.11	0.06	17,27,34,36	0
3	CS	B	400	1/1	0.98	0.03	-2.92	60,60,60,60	0
3	CS	B	401	1/1	0.88	0.18	-	116,116,116,116	0

6.5 Other polymers [i](#)

There are no such residues in this entry.