



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:57 PM GMT

PDB ID : 1VZB
Title : L. CASEI THYMIDYLATE SYNTHASE MUTANT E60Q BINARY COMPLEX WITH DUMP
Authors : Birdsall, D.L.; Finer-Moore, J.; Stroud, R.M.
Deposited on : 1996-09-18
Resolution : 2.50 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

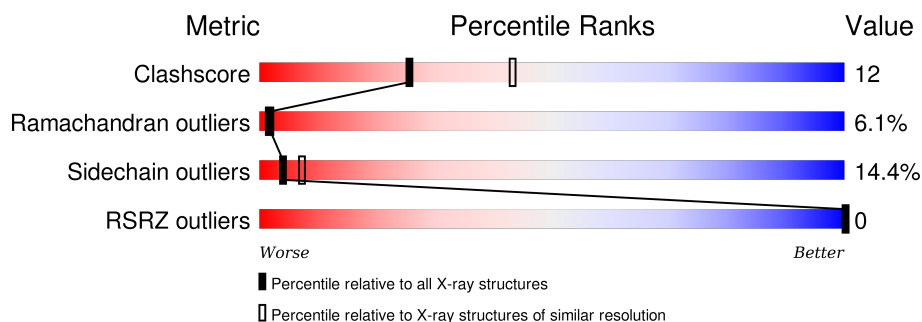
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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))
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

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	316	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 3275 atoms, of which 621 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 THYMIDYLATE SYNTHASE.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	316	3122	1677	532	439	466	8	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	60	GLN	GLU	ENGINEERED	UNP P00469
A	111	ALA	ASP	CONFLICT	UNP P00469

- Molecule 2 is 2'-DEOXYURIDINE 5'-MONOPHOSPHATE (three-letter code: UMP) (formula: $C_9H_{13}N_2O_8P$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	C	H	N	O	P		
2	A	1	21	9	1	2	8	1	0	0

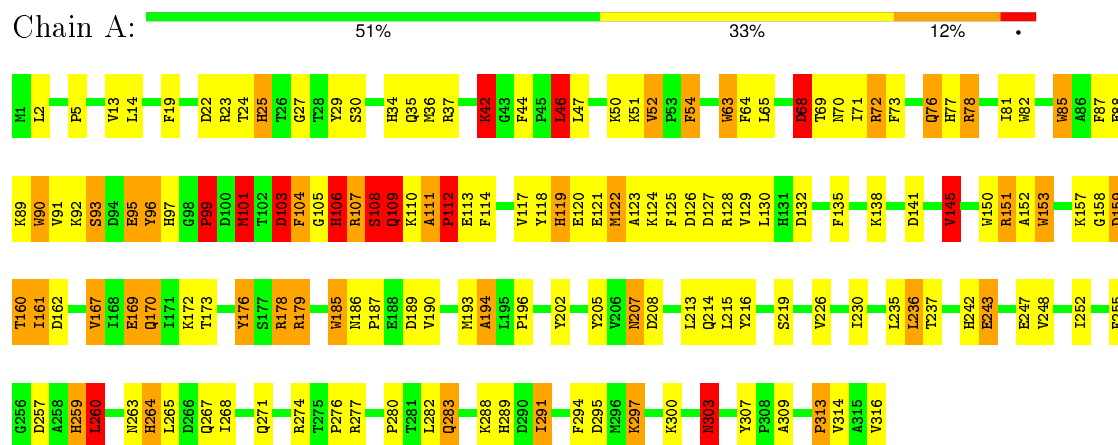
- Molecule 3 is water.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	44	Total	H	O	0	0
			132	88	44		

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 ($\text{RSRZ} > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: THYMIDYLATE SYNTHASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	78.30 Å 78.30 Å 242.30 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	15.00 – 2.50 39.15 – 2.20	Depositor EDS
% Data completeness (in resolution range)	(Not available) (15.00-2.50) 50.6 (39.15-2.20)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.74 (at 2.20 Å)	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.215 , (Not available) 0.218 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	40.3	Xtriage
Anisotropy	0.069	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 74.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 11787 reflections (0.008%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3275	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.86% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: UMP

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	1.24	3/2674 (0.1%)	2.10	97/3634 (2.7%)

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	4

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	219	SER	CA-CB	-8.29	1.40	1.52
1	A	95	GLU	CB-CG	7.31	1.66	1.52
1	A	82	TRP	NE1-CE2	-5.21	1.30	1.37

All (97) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	72	ARG	NE-CZ-NH2	-14.13	113.24	120.30
1	A	90	TRP	CD1-CG-CD2	10.29	114.53	106.30
1	A	63	TRP	CD1-CG-CD2	9.78	114.12	106.30
1	A	150	TRP	CD1-CG-CD2	9.65	114.02	106.30
1	A	277	ARG	NE-CZ-NH2	-9.54	115.53	120.30
1	A	13	VAL	CA-CB-CG2	-9.09	97.27	110.90
1	A	179	ARG	NE-CZ-NH1	8.98	124.79	120.30
1	A	189	ASP	CB-CG-OD1	8.79	126.21	118.30
1	A	153	TRP	CD1-CG-CD2	8.62	113.19	106.30
1	A	36	MET	CA-CB-CG	8.41	127.60	113.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	90	TRP	CE2-CD2-CG	-8.29	100.67	107.30
1	A	167	VAL	CG1-CB-CG2	-8.13	97.89	110.90
1	A	185	TRP	CD1-CG-CD2	7.86	112.59	106.30
1	A	107	ARG	NE-CZ-NH2	-7.70	116.45	120.30
1	A	63	TRP	CE2-CD2-CG	-7.58	101.24	107.30
1	A	150	TRP	CE2-CD2-CG	-7.48	101.31	107.30
1	A	193	MET	CA-CB-CG	7.45	125.97	113.30
1	A	205	TYR	CB-CG-CD2	-7.39	116.57	121.00
1	A	173	THR	O-C-N	-7.24	111.12	122.70
1	A	257	ASP	CA-C-N	7.19	133.02	117.20
1	A	153	TRP	CE2-CD2-CG	-7.13	101.59	107.30
1	A	189	ASP	CA-CB-CG	7.09	129.01	113.40
1	A	260	LEU	CA-CB-CG	6.96	131.30	115.30
1	A	150	TRP	CG-CD1-NE1	-6.93	103.17	110.10
1	A	216	TYR	CB-CG-CD1	-6.89	116.87	121.00
1	A	185	TRP	CE2-CD2-CG	-6.88	101.80	107.30
1	A	112	PRO	CA-N-CD	-6.87	101.88	111.50
1	A	37	ARG	NE-CZ-NH1	6.86	123.73	120.30
1	A	108	SER	N-CA-C	-6.86	92.49	111.00
1	A	128	ARG	NE-CZ-NH2	-6.85	116.88	120.30
1	A	82	TRP	CE2-CD2-CG	-6.84	101.83	107.30
1	A	162	ASP	CB-CG-OD1	6.81	124.43	118.30
1	A	112	PRO	N-CA-C	6.77	129.69	112.10
1	A	173	THR	CA-C-N	6.72	131.98	117.20
1	A	141	ASP	CA-C-N	-6.67	102.52	117.20
1	A	90	TRP	CG-CD1-NE1	-6.53	103.57	110.10
1	A	158	GLY	CA-C-N	-6.50	102.91	117.20
1	A	248	VAL	CG1-CB-CG2	6.49	121.29	110.90
1	A	179	ARG	NE-CZ-NH2	-6.48	117.06	120.30
1	A	52	VAL	CG1-CB-CG2	-6.46	100.57	110.90
1	A	291	ILE	CA-CB-CG2	-6.45	98.00	110.90
1	A	90	TRP	CB-CG-CD1	-6.44	118.62	127.00
1	A	37	ARG	N-CA-CB	-6.42	99.05	110.60
1	A	13	VAL	CA-CB-CG1	6.38	120.47	110.90
1	A	248	VAL	N-CA-CB	-6.35	97.52	111.50
1	A	242	HIS	CA-CB-CG	6.23	124.19	113.60
1	A	257	ASP	O-C-N	-6.13	112.90	122.70
1	A	190	VAL	CA-CB-CG2	-6.01	101.88	110.90
1	A	42	LYS	CA-CB-CG	6.00	126.61	113.40
1	A	68	ASP	CA-C-N	-6.00	104.01	117.20
1	A	288	LYS	CA-CB-CG	-5.91	100.41	113.40
1	A	314	VAL	N-CA-C	5.89	126.90	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	178	ARG	CA-CB-CG	5.84	126.25	113.40
1	A	90	TRP	CG-CD2-CE3	5.81	139.13	133.90
1	A	50	LYS	CA-CB-CG	5.79	126.15	113.40
1	A	85	TRP	CE2-CD2-CG	-5.79	102.67	107.30
1	A	255	PHE	CB-CG-CD2	-5.79	116.75	120.80
1	A	114	PHE	O-C-N	-5.73	113.53	122.70
1	A	29	TYR	CB-CG-CD1	5.68	124.41	121.00
1	A	47	LEU	CB-CG-CD1	-5.68	101.35	111.00
1	A	158	GLY	C-N-CA	5.61	135.72	121.70
1	A	101	MET	N-CA-C	5.57	126.04	111.00
1	A	283	GLN	CA-CB-CG	5.57	125.65	113.40
1	A	150	TRP	CG-CD2-CE3	5.55	138.89	133.90
1	A	151	ARG	CA-CB-CG	-5.51	101.27	113.40
1	A	128	ARG	NE-CZ-NH1	5.50	123.05	120.30
1	A	189	ASP	CB-CG-OD2	-5.48	113.37	118.30
1	A	90	TRP	N-CA-C	-5.47	96.23	111.00
1	A	236	LEU	CA-CB-CG	5.44	127.81	115.30
1	A	103	ASP	CA-C-N	-5.43	105.25	117.20
1	A	176	TYR	N-CA-C	-5.43	96.34	111.00
1	A	243	GLU	CA-CB-CG	-5.42	101.49	113.40
1	A	82	TRP	CD1-CG-CD2	5.41	110.63	106.30
1	A	109	GLN	N-CA-C	5.38	125.53	111.00
1	A	216	TYR	CB-CG-CD2	5.38	124.22	121.00
1	A	29	TYR	N-CA-C	-5.37	96.49	111.00
1	A	159	ASP	N-CA-CB	-5.36	100.95	110.60
1	A	138	LYS	CB-CG-CD	5.34	125.47	111.60
1	A	153	TRP	CG-CD1-NE1	-5.32	104.78	110.10
1	A	106	HIS	CA-C-N	5.29	128.84	117.20
1	A	113	GLU	N-CA-C	-5.28	96.75	111.00
1	A	145	VAL	CG1-CB-CG2	-5.23	102.53	110.90
1	A	161	ILE	N-CA-CB	-5.23	98.78	110.80
1	A	123	ALA	N-CA-CB	5.22	117.42	110.10
1	A	64	PHE	CB-CG-CD2	-5.21	117.15	120.80
1	A	127	ASP	CB-CG-OD1	-5.20	113.62	118.30
1	A	158	GLY	O-C-N	5.20	131.02	122.70
1	A	35	GLN	N-CA-CB	5.18	119.92	110.60
1	A	207	ASN	CA-CB-CG	5.16	124.75	113.40
1	A	78	ARG	CB-CG-CD	5.15	125.00	111.60
1	A	95	GLU	CB-CG-CD	5.13	128.06	114.20
1	A	46	LEU	CB-CG-CD1	-5.12	102.29	111.00
1	A	202	TYR	CB-CG-CD1	-5.11	117.94	121.00
1	A	119	HIS	CB-CA-C	-5.06	100.27	110.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	264	HIS	CA-CB-CG	5.05	122.19	113.60
1	A	303	ASN	CB-CG-ND2	5.03	128.77	116.70
1	A	123	ALA	CB-CA-C	-5.01	102.58	110.10

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	176	TYR	Sidechain
1	A	25	HIS	Sidechain
1	A	259	HIS	Sidechain
1	A	34	HIS	Sidechain

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	2590	532	2501	64	1
2	A	20	1	11	0	0
3	A	44	88	0	0	0
All	All	2654	621	2512	64	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 12.

All (64) 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:153:TRP:HB3	1:A:161:ILE:HG22	1.70	0.73
1:A:44:PHE:HB3	1:A:280:PRO:HG2	1.70	0.73
1:A:194:ALA:O	1:A:196:PRO:HD3	1.88	0.72
1:A:153:TRP:HB3	1:A:161:ILE:CG2	2.23	0.68
1:A:104:PHE:HD1	1:A:105:GLY:H	1.46	0.64
1:A:260:LEU:HD12	1:A:265:LEU:HD23	1.80	0.64
1:A:106:HIS:HA	1:A:109:GLN:HB2	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:108:SER:O	1:A:110:LYS:N	2.33	0.62
1:A:274:ARG:NH1	1:A:307:TYR:HB3	2.16	0.61
1:A:226:VAL:O	1:A:230:ILE:HG13	2.02	0.59
1:A:71:ILE:HG22	1:A:129:VAL:HG11	1.86	0.57
1:A:152:ALA:HB1	1:A:160:THR:HG23	1.87	0.55
1:A:111:ALA:N	1:A:112:PRO:HD2	2.22	0.55
1:A:121:GLU:O	1:A:124:LYS:HB3	2.08	0.54
1:A:213:LEU:HD13	1:A:237:THR:OG1	2.07	0.54
1:A:19:PHE:CZ	1:A:27:GLY:HA3	2.43	0.53
1:A:151:ARG:HG2	1:A:151:ARG:HH11	1.73	0.53
1:A:63:TRP:CD1	1:A:68:ASP:HB3	2.44	0.52
1:A:132:ASP:HB3	1:A:135:PHE:HB3	1.92	0.52
1:A:169:GLU:HG2	1:A:172:LYS:HE3	1.91	0.52
1:A:167:VAL:HA	1:A:170:GLN:HE21	1.76	0.51
1:A:2:LEU:O	1:A:5:PRO:HD2	2.10	0.51
1:A:106:HIS:HD1	1:A:316:VAL:C	2.13	0.50
1:A:71:ILE:CG2	1:A:129:VAL:HG11	2.41	0.50
1:A:303:ASN:OD1	1:A:303:ASN:N	2.43	0.49
1:A:72:ARG:O	1:A:76:GLN:HG3	2.13	0.49
1:A:126:ASP:O	1:A:130:LEU:HD22	2.13	0.49
1:A:69:THR:HG22	1:A:145:VAL:HG13	1.94	0.49
1:A:85:TRP:O	1:A:88:GLU:HB3	2.13	0.49
1:A:263:ASN:O	1:A:264:HIS:HD2	1.95	0.48
1:A:291:ILE:HA	1:A:294:PHE:CD1	2.49	0.48
1:A:89:LYS:O	1:A:93:SER:HB2	2.14	0.47
1:A:51:LYS:H	1:A:309:ALA:HA	1.80	0.47
1:A:265:LEU:O	1:A:268:ILE:HG22	2.14	0.47
1:A:120:GLU:O	1:A:124:LYS:N	2.45	0.47
1:A:14:LEU:HD23	1:A:14:LEU:HA	1.64	0.47
1:A:104:PHE:O	1:A:106:HIS:N	2.48	0.46
1:A:87:PHE:CD2	1:A:122:MET:SD	3.09	0.45
1:A:93:SER:O	1:A:96:TYR:HB3	2.16	0.45
1:A:297:LYS:H	1:A:297:LYS:HD3	1.82	0.45
1:A:151:ARG:HG2	1:A:151:ARG:NH1	2.32	0.44
1:A:104:PHE:O	1:A:107:ARG:N	2.49	0.44
1:A:243:GLU:HG3	1:A:289:HIS:O	2.18	0.44
1:A:30:SER:HB3	1:A:259:HIS:HB3	1.99	0.44
1:A:267:GLN:OE1	1:A:267:GLN:N	2.51	0.44
1:A:95:GLU:O	1:A:97:HIS:N	2.49	0.44
1:A:90:TRP:O	1:A:93:SER:N	2.51	0.43
1:A:70:ASN:ND2	1:A:72:ARG:H	2.17	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:124:LYS:HB2	1:A:124:LYS:HE3	1.83	0.43
1:A:99:PRO:HB2	1:A:101:MET:HE1	2.00	0.43
1:A:87:PHE:HZ	1:A:121:GLU:HB3	1.83	0.42
1:A:73:PHE:CE1	1:A:77:HIS:HE1	2.38	0.42
1:A:186:ASN:HA	1:A:187:PRO:HD3	1.79	0.42
1:A:271:GLN:HA	1:A:274:ARG:HD2	2.01	0.42
1:A:72:ARG:HD3	1:A:76:GLN:HE21	1.85	0.42
1:A:214:GLN:HG3	1:A:252:ILE:HB	2.00	0.42
1:A:46:LEU:HD22	1:A:52:VAL:HB	2.02	0.41
1:A:72:ARG:HD3	1:A:76:GLN:NE2	2.35	0.41
1:A:107:ARG:O	1:A:111:ALA:HB3	2.19	0.41
1:A:87:PHE:HB2	1:A:122:MET:SD	2.61	0.41
1:A:122:MET:O	1:A:125:PHE:HB3	2.21	0.41
1:A:268:ILE:HA	1:A:268:ILE:HD12	1.78	0.40
1:A:185:TRP:O	1:A:187:PRO:HD3	2.22	0.40
1:A:260:LEU:HD12	1:A:265:LEU:CD2	2.49	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)
1:A:23:ARG:HH11	1:A:179:ARG:HH21[11_554]	1.12	0.48

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	314/316 (99%)	246 (78%)	49 (16%)	19 (6%)	2 1

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	25	HIS
1	A	101	MET
1	A	104	PHE
1	A	109	GLN
1	A	112	PRO
1	A	159	ASP
1	A	194	ALA
1	A	96	TYR
1	A	103	ASP
1	A	111	ALA
1	A	81	ILE
1	A	42	LYS
1	A	76	GLN
1	A	118	TYR
1	A	313	PRO
1	A	54	PHE
1	A	99	PRO
1	A	276	PRO
1	A	117	VAL

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	277/277 (100%)	237 (86%)	40 (14%)	4 7

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	ASP
1	A	24	THR
1	A	42	LYS
1	A	46	LEU
1	A	54	PHE
1	A	65	LEU
1	A	68	ASP
1	A	78	ARG

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Mol	Chain	Res	Type
1	A	91	VAL
1	A	92	LYS
1	A	93	SER
1	A	99	PRO
1	A	101	MET
1	A	103	ASP
1	A	106	HIS
1	A	108	SER
1	A	109	GLN
1	A	112	PRO
1	A	119	HIS
1	A	122	MET
1	A	145	VAL
1	A	157	LYS
1	A	160	THR
1	A	169	GLU
1	A	170	GLN
1	A	178	ARG
1	A	207	ASN
1	A	208	ASP
1	A	215	LEU
1	A	235	LEU
1	A	236	LEU
1	A	247	GLU
1	A	260	LEU
1	A	282	LEU
1	A	283	GLN
1	A	295	ASP
1	A	297	LYS
1	A	300	LYS
1	A	303	ASN
1	A	313	PRO

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	76	GLN
1	A	170	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 ⓘ

1 ligand is 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	UMP	A	317	-	16,21,21	1.76	3 (18%)	20,31,31	2.04	3 (15%)

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	UMP	A	317	-	-	0/6/22/22	0/2/2/2

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	317	UMP	C6-N1	-4.39	1.29	1.35
2	A	317	UMP	C5'-C4'	2.32	1.59	1.51
2	A	317	UMP	C4-N3	3.17	1.41	1.36

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	A	317	UMP	O4'-C1'-C2'	-3.73	98.84	106.27
2	A	317	UMP	C2'-C3'-C4'	-3.28	95.98	102.77
2	A	317	UMP	O4'-C1'-N1	6.49	118.95	107.72

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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 [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	316/316 (100%)	-0.92	0 100 100	10, 23, 37, 43	0

There are no RSRZ outliers to report.

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(Å ²)	Q<0.9
2	UMP	A	317	20/20	0.97	0.09	-0.71	15,26,28,28	0

6.5 Other polymers [i](#)

There are no such residues in this entry.