Full wwPDB X-ray Structure Validation Report (i)

Jan 31, 2016 – 10:57 PM GMT

PDB ID	:	1VZC
Title	:	L. CASEI THYMIDYLATE SYNTHASE MUTANT E60Q BINARY COMPLEX WITH FDUMP
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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

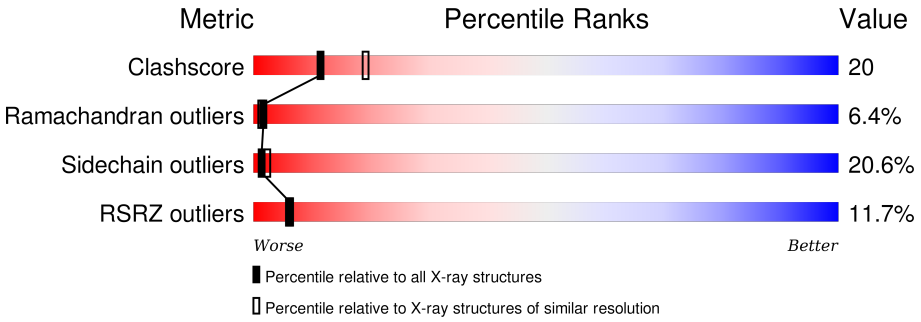
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)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

1 Overall quality at a glance (i)

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 <=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	12% 39% 41% 16% •

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 criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	UFP	A	317	X	-	-	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3267 atoms, of which 615 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.

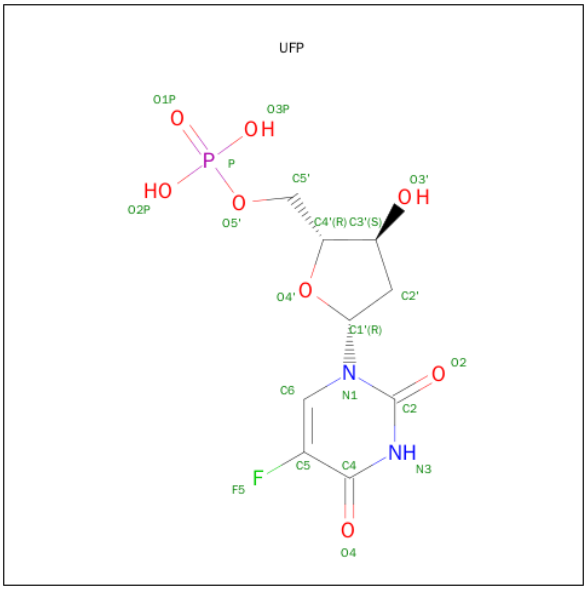
- Molecule 1 is a protein called THYMIDYLATE SYNTHASE.

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 5-FLUORO-2'-DEOXYURIDINE-5'-MONOPHOSPHATE (three-letter code: UFP) (formula: $C_9H_{12}FN_2O_8P$).



Mol	Chain	Residues	Atoms							ZeroOcc	AltConf
			Total	C	F	H	N	O	P		
2	A	1	22	9	1	1	2	8	1	0	0

- Molecule 3 is water.

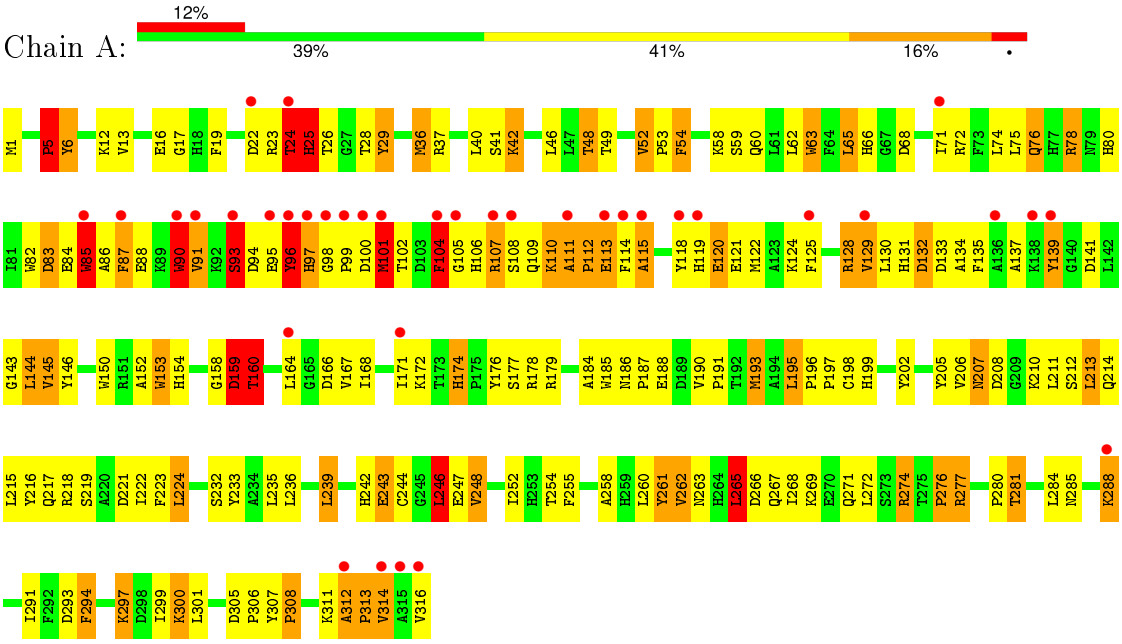
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	41	Total	H	O	0	0
			123	82	41		

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 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: THYMIDYLATE SYNTHASE

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α, β, γ	79.50Å 79.50Å 229.10Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	15.00 – 2.50 59.01 – 2.05	Depositor EDS
% Data completeness (in resolution range)	(Not available) (15.00-2.50) 41.8 (59.01-2.05)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.51 (at 2.05Å)	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.201 , (Not available) 0.307 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å^2)	15.0	Xtriage
Anisotropy	0.337	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 347.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.46$, $< L^2 > = 0.29$	Xtriage
Outliers	0 of 11633 reflections	Xtriage
F_o,F_c correlation	0.76	EDS
Total number of atoms	3267	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 4.62% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: UFP

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	1.47	11/2674 (0.4%)	2.20	110/3634 (3.0%)

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	8

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	150	TRP	CG-CD2	-6.01	1.33	1.43
1	A	23	ARG	NE-CZ	5.91	1.40	1.33
1	A	63	TRP	CA-CB	-5.46	1.42	1.53
1	A	177	SER	CA-CB	-5.42	1.44	1.52
1	A	212	SER	CA-CB	-5.42	1.44	1.52
1	A	63	TRP	CD1-NE1	-5.42	1.28	1.38
1	A	150	TRP	NE1-CE2	-5.39	1.30	1.37
1	A	120	GLU	CG-CD	5.32	1.59	1.51
1	A	146	TYR	CA-CB	-5.28	1.42	1.53
1	A	185	TRP	CG-CD2	-5.21	1.34	1.43
1	A	54	PHE	CA-CB	-5.19	1.42	1.53

All (110) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	179	ARG	NE-CZ-NH1	15.68	128.14	120.30
1	A	274	ARG	NE-CZ-NH1	14.08	127.34	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	179	ARG	NE-CZ-NH2	-11.32	114.64	120.30
1	A	78	ARG	NE-CZ-NH2	-11.23	114.68	120.30
1	A	176	TYR	CB-CG-CD2	-10.87	114.48	121.00
1	A	63	TRP	CD1-CG-CD2	10.01	114.31	106.30
1	A	25	HIS	N-CA-C	9.56	136.81	111.00
1	A	85	TRP	CB-CG-CD1	-9.50	114.66	127.00
1	A	82	TRP	CD1-CG-CD2	9.23	113.69	106.30
1	A	202	TYR	CB-CG-CD1	-9.07	115.56	121.00
1	A	23	ARG	NE-CZ-NH1	9.05	124.83	120.30
1	A	90	TRP	CE2-CD2-CG	-8.62	100.41	107.30
1	A	85	TRP	CE2-CD2-CG	-8.45	100.54	107.30
1	A	96	TYR	CA-C-N	-8.44	98.64	117.20
1	A	63	TRP	CE2-CD2-CG	-8.27	100.69	107.30
1	A	66	HIS	CB-CA-C	-8.23	93.95	110.40
1	A	82	TRP	CE2-CD2-CG	-8.21	100.73	107.30
1	A	85	TRP	CG-CD2-CE3	8.21	141.29	133.90
1	A	101	MET	CA-CB-CG	-8.01	99.69	113.30
1	A	246	LEU	CA-CB-CG	8.00	133.70	115.30
1	A	248	VAL	CG1-CB-CG2	-7.92	98.22	110.90
1	A	176	TYR	CB-CG-CD1	7.74	125.64	121.00
1	A	174	HIS	CA-CB-CG	7.61	126.54	113.60
1	A	277	ARG	NE-CZ-NH2	-7.50	116.55	120.30
1	A	85	TRP	CD1-CG-CD2	7.50	112.30	106.30
1	A	146	TYR	CB-CG-CD1	-7.49	116.50	121.00
1	A	185	TRP	CD1-CG-CD2	7.48	112.28	106.30
1	A	93	SER	CA-C-N	-7.44	100.83	117.20
1	A	153	TRP	CD1-CG-CD2	7.37	112.20	106.30
1	A	90	TRP	CD1-CG-CD2	7.21	112.07	106.30
1	A	233	TYR	CB-CG-CD2	-7.12	116.72	121.00
1	A	110	LYS	N-CA-C	6.96	129.80	111.00
1	A	111	ALA	N-CA-C	6.95	129.76	111.00
1	A	141	ASP	CB-CG-OD2	-6.95	112.04	118.30
1	A	221	ASP	CB-CG-OD2	-6.94	112.05	118.30
1	A	144	LEU	CA-CB-CG	6.93	131.24	115.30
1	A	5	PRO	CA-N-CD	-6.74	102.06	111.50
1	A	221	ASP	CB-CG-OD1	6.71	124.33	118.30
1	A	202	TYR	CB-CG-CD2	6.65	124.99	121.00
1	A	281	THR	CA-CB-CG2	6.65	121.70	112.40
1	A	185	TRP	CE2-CD2-CG	-6.64	101.99	107.30
1	A	300	LYS	CB-CG-CD	6.61	128.78	111.60
1	A	62	LEU	CA-CB-CG	6.56	130.40	115.30
1	A	150	TRP	CD1-CG-CD2	6.56	111.55	106.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	109	GLN	CB-CG-CD	6.54	128.59	111.60
1	A	153	TRP	CE2-CD2-CG	-6.53	102.08	107.30
1	A	66	HIS	CA-CB-CG	6.46	124.58	113.60
1	A	29	TYR	CB-CG-CD1	-6.44	117.13	121.00
1	A	107	ARG	NE-CZ-NH2	-6.39	117.10	120.30
1	A	242	HIS	CA-C-N	6.24	130.92	117.20
1	A	150	TRP	CE2-CD2-CG	-6.19	102.35	107.30
1	A	104	PHE	CA-CB-CG	6.18	128.74	113.90
1	A	255	PHE	CB-CG-CD2	-6.10	116.53	120.80
1	A	281	THR	CA-CB-OG1	-6.10	96.19	109.00
1	A	78	ARG	CA-CB-CG	6.09	126.81	113.40
1	A	90	TRP	CG-CD2-CE3	6.09	139.38	133.90
1	A	37	ARG	NE-CZ-NH2	-6.08	117.26	120.30
1	A	308	PRO	N-CD-CG	-5.97	94.24	103.20
1	A	193	MET	CG-SD-CE	5.90	109.64	100.20
1	A	96	TYR	CB-CG-CD2	-5.89	117.47	121.00
1	A	96	TYR	N-CA-CB	5.89	121.20	110.60
1	A	274	ARG	NH1-CZ-NH2	-5.89	112.92	119.40
1	A	19	PHE	CB-CG-CD1	5.88	124.92	120.80
1	A	87	PHE	CB-CG-CD2	-5.86	116.70	120.80
1	A	294	PHE	CB-CG-CD2	-5.84	116.71	120.80
1	A	97	HIS	N-CA-C	-5.84	95.23	111.00
1	A	172	LYS	CB-CG-CD	5.82	126.73	111.60
1	A	129	VAL	CB-CA-C	-5.79	100.39	111.40
1	A	66	HIS	N-CA-CB	5.73	120.92	110.60
1	A	88	GLU	CA-CB-CG	5.69	125.92	113.40
1	A	97	HIS	CA-CB-CG	5.67	123.25	113.60
1	A	84	GLU	CA-CB-CG	5.67	125.86	113.40
1	A	96	TYR	O-C-N	5.66	131.76	122.70
1	A	104	PHE	N-CA-C	5.66	126.29	111.00
1	A	24	THR	N-CA-CB	-5.63	99.61	110.30
1	A	72	ARG	NE-CZ-NH2	-5.62	117.49	120.30
1	A	166	ASP	CB-CG-OD2	5.61	123.35	118.30
1	A	297	LYS	CA-CB-CG	-5.60	101.09	113.40
1	A	146	TYR	CD1-CG-CD2	5.55	124.00	117.90
1	A	48	THR	O-C-N	-5.51	113.89	122.70
1	A	6	TYR	CA-C-N	5.49	129.29	117.20
1	A	36	MET	CB-CG-SD	-5.45	96.04	112.40
1	A	146	TYR	CA-CB-CG	-5.45	103.05	113.40
1	A	22	ASP	CB-CG-OD2	-5.43	113.41	118.30
1	A	294	PHE	CA-C-N	-5.43	105.25	117.20
1	A	274	ARG	CG-CD-NE	5.42	123.19	111.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	265	LEU	CA-CB-CG	5.38	127.67	115.30
1	A	78	ARG	NE-CZ-NH1	5.35	122.98	120.30
1	A	145	VAL	N-CA-C	-5.32	96.63	111.00
1	A	145	VAL	CA-C-N	-5.30	105.53	117.20
1	A	199	HIS	CA-C-N	-5.30	105.55	117.20
1	A	199	HIS	ND1-CG-CD2	5.28	116.19	108.80
1	A	12	LYS	CA-C-N	5.26	128.77	117.20
1	A	160	THR	CA-CB-CG2	5.25	119.76	112.40
1	A	160	THR	CA-CB-OG1	-5.25	97.97	109.00
1	A	153	TRP	CG-CD1-NE1	-5.24	104.86	110.10
1	A	166	ASP	CB-CG-OD1	-5.23	113.60	118.30
1	A	254	THR	CA-C-N	-5.23	105.70	117.20
1	A	110	LYS	C-N-CA	5.20	134.69	121.70
1	A	36	MET	CG-SD-CE	5.17	108.48	100.20
1	A	128	ARG	NE-CZ-NH1	5.16	122.88	120.30
1	A	159	ASP	CB-CG-OD1	-5.15	113.67	118.30
1	A	300	LYS	CG-CD-CE	5.13	127.28	111.90
1	A	199	HIS	O-C-N	5.11	130.88	122.70
1	A	90	TRP	CB-CG-CD1	-5.11	120.36	127.00
1	A	210	LYS	N-CA-C	5.08	124.72	111.00
1	A	23	ARG	CA-CB-CG	5.05	124.51	113.40
1	A	262	VAL	CA-CB-CG2	-5.05	103.33	110.90
1	A	65	LEU	N-CA-CB	-5.03	100.34	110.40
1	A	139	TYR	CB-CG-CD2	-5.01	118.00	121.00

There are no chirality outliers.

All (8) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	139	TYR	Sidechain
1	A	216	TYR	Sidechain
1	A	261	TYR	Sidechain
1	A	29	TYR	Sidechain
1	A	312	ALA	Peptide
1	A	6	TYR	Sidechain
1	A	96	TYR	Sidechain
1	A	98	GLY	Peptide

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	2590	532	2501	105	3
2	A	21	1	10	0	0
3	A	41	82	0	2	0
All	All	2652	615	2511	105	3

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 20.

All (105) 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:243:GLU:HG3	1:A:291:ILE:HG23	1.53	0.89
1:A:76:GLN:HB3	1:A:130:LEU:HD11	1.64	0.79
1:A:195:LEU:CD1	3:A:359:HOH:O	2.30	0.79
1:A:104:PHE:HB2	1:A:118:TYR:CD2	2.22	0.74
1:A:195:LEU:HD11	3:A:359:HOH:O	1.87	0.73
1:A:101:MET:HE1	1:A:118:TYR:HB2	1.73	0.71
1:A:244:CYS:HB2	1:A:246:LEU:HD22	1.72	0.70
1:A:108:SER:HB2	1:A:119:HIS:HE1	1.57	0.70
1:A:152:ALA:HA	1:A:160:THR:HG23	1.73	0.70
1:A:223:PHE:HD1	1:A:268:ILE:HD13	1.57	0.69
1:A:71:ILE:HG22	1:A:129:VAL:HG11	1.75	0.69
1:A:288:LYS:HZ1	1:A:293:ASP:HB3	1.59	0.67
1:A:265:LEU:O	1:A:269:LYS:HG3	1.94	0.67
1:A:63:TRP:CD1	1:A:68:ASP:HB3	2.30	0.67
1:A:87:PHE:HE1	1:A:91:VAL:HG23	1.60	0.66
1:A:87:PHE:CE1	1:A:91:VAL:HG23	2.31	0.66
1:A:101:MET:CE	1:A:101:MET:HA	2.24	0.66
1:A:205:TYR:CE2	1:A:207:ASN:HB3	2.32	0.65
1:A:90:TRP:O	1:A:93:SER:HB2	1.97	0.63
1:A:65:LEU:HD21	1:A:291:ILE:HD12	1.80	0.63
1:A:71:ILE:HG13	1:A:86:ALA:HB2	1.80	0.63
1:A:105:GLY:HA2	1:A:108:SER:OG	1.99	0.62
1:A:125:PHE:O	1:A:129:VAL:HG23	1.98	0.62
1:A:1:MET:HG3	1:A:276:PRO:HB2	1.83	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:288:LYS:NZ	1:A:293:ASP:HB3	2.16	0.60
1:A:90:TRP:HA	1:A:90:TRP:CE3	2.35	0.60
1:A:101:MET:SD	1:A:118:TYR:HA	2.42	0.60
1:A:288:LYS:HZ2	1:A:288:LYS:HB3	1.68	0.58
1:A:297:LYS:NZ	1:A:300:LYS:HB2	2.18	0.58
1:A:85:TRP:HZ2	1:A:195:LEU:HB2	1.68	0.57
1:A:87:PHE:CD2	1:A:122:MET:SD	2.98	0.57
1:A:274:ARG:HG2	1:A:307:TYR:CD2	2.40	0.56
1:A:153:TRP:O	1:A:160:THR:HA	2.04	0.56
1:A:48:THR:HB	1:A:277:ARG:HB2	1.86	0.56
1:A:13:VAL:O	1:A:17:GLY:HA3	2.06	0.55
1:A:265:LEU:O	1:A:268:ILE:HG22	2.06	0.54
1:A:196:PRO:HG2	1:A:218:ARG:HH22	1.72	0.54
1:A:104:PHE:HD2	1:A:118:TYR:HE2	1.56	0.53
1:A:184:ALA:O	1:A:197:PRO:HG2	2.08	0.53
1:A:128:ARG:HB3	1:A:135:PHE:CG	2.43	0.53
1:A:284:LEU:HD22	1:A:299:ILE:HG13	1.89	0.53
1:A:86:ALA:HB3	1:A:125:PHE:HE2	1.73	0.52
1:A:224:LEU:HD11	1:A:312:ALA:HB2	1.92	0.52
1:A:223:PHE:HD2	1:A:224:LEU:HD13	1.75	0.51
1:A:277:ARG:HD2	1:A:305:ASP:O	2.10	0.51
1:A:213:LEU:O	1:A:252:ILE:HG12	2.10	0.51
1:A:108:SER:HB2	1:A:119:HIS:CE1	2.42	0.51
1:A:104:PHE:HB2	1:A:118:TYR:CE2	2.45	0.51
1:A:24:THR:HG22	1:A:316:VAL:HA	1.92	0.51
1:A:46:LEU:HG	1:A:52:VAL:HG13	1.92	0.51
1:A:281:THR:O	1:A:301:LEU:HD22	2.11	0.51
1:A:243:GLU:HG3	1:A:291:ILE:CG2	2.34	0.51
1:A:85:TRP:CZ2	1:A:195:LEU:HB2	2.46	0.50
1:A:104:PHE:O	1:A:108:SER:HB3	2.11	0.50
1:A:110:LYS:O	1:A:112:PRO:HD2	2.12	0.50
1:A:223:PHE:CD1	1:A:268:ILE:HD13	2.43	0.50
1:A:90:TRP:HA	1:A:90:TRP:HE3	1.77	0.50
1:A:101:MET:CE	1:A:118:TYR:HB2	2.39	0.50
1:A:5:PRO:HB3	1:A:36:MET:HG3	1.94	0.49
1:A:40:LEU:HB2	1:A:248:VAL:CG1	2.43	0.49
1:A:297:LYS:HZ3	1:A:300:LYS:HB2	1.77	0.49
1:A:214:GLN:HA	1:A:252:ILE:HB	1.94	0.49
1:A:167:VAL:O	1:A:171:ILE:HG13	2.13	0.48
1:A:134:ALA:O	1:A:137:ALA:HB3	2.13	0.48
1:A:285:ASN:ND2	1:A:288:LYS:HG3	2.28	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:26:THR:HG22	1:A:262:VAL:HB	1.96	0.48
1:A:107:ARG:HB2	1:A:114:PHE:CD2	2.49	0.47
1:A:187:PRO:O	1:A:191:PRO:HD2	2.14	0.47
1:A:153:TRP:CZ2	1:A:186:ASN:HB2	2.50	0.46
1:A:186:ASN:O	1:A:190:VAL:HG23	2.14	0.46
1:A:198:CYS:O	1:A:217:GLN:HG3	2.15	0.46
1:A:54:PHE:CE2	1:A:301:LEU:HB2	2.50	0.46
1:A:26:THR:HG21	1:A:263:ASN:H	1.80	0.46
1:A:314:VAL:HB	1:A:316:VAL:HG22	1.96	0.46
1:A:291:ILE:HA	1:A:294:PHE:CD1	2.51	0.45
1:A:108:SER:HA	1:A:115:ALA:HA	1.98	0.45
1:A:260:LEU:HD21	1:A:268:ILE:HG21	1.98	0.45
1:A:40:LEU:HB2	1:A:248:VAL:HG12	1.98	0.45
1:A:63:TRP:CZ2	1:A:74:LEU:HD11	2.52	0.45
1:A:118:TYR:HD1	1:A:119:HIS:ND1	2.15	0.44
1:A:131:HIS:O	1:A:132:ASP:HB2	2.18	0.44
1:A:260:LEU:CD2	1:A:265:LEU:HD12	2.48	0.44
1:A:284:LEU:HD22	1:A:299:ILE:CG1	2.47	0.44
1:A:80:HIS:O	1:A:83:ASP:HB2	2.18	0.44
1:A:235:LEU:HG	1:A:239:LEU:CD2	2.48	0.43
1:A:101:MET:SD	1:A:118:TYR:HB2	2.59	0.42
1:A:168:ILE:HA	1:A:171:ILE:HD12	2.00	0.42
1:A:87:PHE:HZ	1:A:101:MET:HG2	1.84	0.42
1:A:267:GLN:N	1:A:267:GLN:OE1	2.50	0.42
1:A:26:THR:CG2	1:A:262:VAL:HB	2.50	0.42
1:A:41:SER:OG	1:A:42:LYS:NZ	2.52	0.42
1:A:239:LEU:O	1:A:291:ILE:HG21	2.20	0.42
1:A:222:ILE:HD11	1:A:258:ALA:C	2.39	0.42
1:A:288:LYS:NZ	1:A:288:LYS:HB3	2.35	0.41
1:A:75:LEU:HA	1:A:75:LEU:HD23	1.89	0.41
1:A:104:PHE:HD2	1:A:118:TYR:CE2	2.37	0.41
1:A:71:ILE:HA	1:A:71:ILE:HD13	1.69	0.41
1:A:154:HIS:HE1	1:A:188:GLU:OE2	2.04	0.41
1:A:153:TRP:CH2	1:A:186:ASN:HB2	2.56	0.41
1:A:268:ILE:O	1:A:272:LEU:HD12	2.21	0.41
1:A:90:TRP:HH2	1:A:128:ARG:NH2	2.19	0.41
1:A:213:LEU:HA	1:A:213:LEU:HD23	1.91	0.40
1:A:28:THR:HG22	1:A:261:TYR:HA	2.03	0.40
1:A:205:TYR:HE2	1:A:207:ASN:HB3	1.79	0.40
1:A:269:LYS:O	1:A:272:LEU:HB2	2.22	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:107:ARG:HH21	1:A:107:ARG:HH21[11_654]	0.98	0.62
1:A:178:ARG:HH11	1:A:219:SER:HG[11_554]	1.35	0.25
1:A:107:ARG:NH2	1:A:107:ARG:NH2[11_654]	2.16	0.04

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	314/316 (99%)	248 (79%)	46 (15%)	20 (6%)	2 1

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	24	THR
1	A	25	HIS
1	A	102	THR
1	A	104	PHE
1	A	111	ALA
1	A	112	PRO
1	A	113	GLU
1	A	115	ALA
1	A	132	ASP
1	A	208	ASP
1	A	313	PRO
1	A	99	PRO
1	A	143	GLY
1	A	308	PRO
1	A	5	PRO
1	A	53	PRO
1	A	106	HIS
1	A	159	ASP
1	A	49	THR

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Mol	Chain	Res	Type
1	A	158	GLY

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%)	220 (79%)	57 (21%)	1 2

All (57) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	GLU
1	A	24	THR
1	A	25	HIS
1	A	42	LYS
1	A	52	VAL
1	A	58	LYS
1	A	59	SER
1	A	60	GLN
1	A	76	GLN
1	A	78	ARG
1	A	83	ASP
1	A	85	TRP
1	A	90	TRP
1	A	91	VAL
1	A	93	SER
1	A	94	ASP
1	A	95	GLU
1	A	96	TYR
1	A	97	HIS
1	A	100	ASP
1	A	101	MET
1	A	104	PHE
1	A	113	GLU
1	A	120	GLU
1	A	121	GLU

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Mol	Chain	Res	Type
1	A	124	LYS
1	A	133	ASP
1	A	144	LEU
1	A	145	VAL
1	A	159	ASP
1	A	160	THR
1	A	164	LEU
1	A	174	HIS
1	A	193	MET
1	A	195	LEU
1	A	206	VAL
1	A	207	ASN
1	A	211	LEU
1	A	213	LEU
1	A	215	LEU
1	A	224	LEU
1	A	232	SER
1	A	236	LEU
1	A	239	LEU
1	A	243	GLU
1	A	246	LEU
1	A	247	GLU
1	A	265	LEU
1	A	266	ASP
1	A	271	GLN
1	A	276	PRO
1	A	280	PRO
1	A	288	LYS
1	A	306	PRO
1	A	311	LYS
1	A	313	PRO
1	A	314	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	76	GLN
1	A	77	HIS
1	A	154	HIS
1	A	186	ASN
1	A	207	ASN
1	A	289	HIS

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	UFP	A	317	-	18,22,22	2.38	4 (22%)	19,33,33	2.10	4 (21%)

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	UFP	A	317	-	1/1/4/4	0/6/22/22	0/2/2/2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	317	UFP	F5-C5	-2.08	1.32	1.35
2	A	317	UFP	O4'-C1'	3.46	1.50	1.42
2	A	317	UFP	C4-N3	3.56	1.41	1.36
2	A	317	UFP	C4-C5	7.34	1.47	1.38

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	317	UFP	C2'-C3'-C4'	-3.35	95.83	102.77
2	A	317	UFP	O5'-P-O1P	-2.87	99.84	107.14
2	A	317	UFP	O4'-C4'-C3'	-2.62	99.08	105.67
2	A	317	UFP	O4'-C1'-N1	6.30	118.62	107.72

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	317	UFP	C1'

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

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, 95^{th} 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%)	0.62	37 (11%) 6 6	3, 11, 25, 31	0

All (37) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	115	ALA	11.4
1	A	24	THR	9.4
1	A	114	PHE	9.3
1	A	315	ALA	8.5
1	A	91	VAL	8.4
1	A	101	MET	7.3
1	A	136	ALA	6.3
1	A	104	PHE	4.6
1	A	316	VAL	4.5
1	A	95	GLU	4.4
1	A	90	TRP	4.3
1	A	93	SER	4.0
1	A	118	TYR	3.9
1	A	98	GLY	3.7
1	A	108	SER	3.4
1	A	87	PHE	3.3
1	A	139	TYR	3.2
1	A	125	PHE	3.2
1	A	113	GLU	3.2
1	A	107	ARG	3.0
1	A	100	ASP	2.9
1	A	105	GLY	2.8
1	A	22	ASP	2.8
1	A	138	LYS	2.8
1	A	96	TYR	2.7
1	A	119	HIS	2.7
1	A	129	VAL	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	71	ILE	2.6
1	A	97	HIS	2.6
1	A	85	TRP	2.5
1	A	312	ALA	2.3
1	A	99	PRO	2.2
1	A	314	VAL	2.2
1	A	164	LEU	2.1
1	A	111	ALA	2.1
1	A	171	ILE	2.1
1	A	288	LYS	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, 95^{th} 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	UFP	A	317	21/21	0.83	0.24	0.16	13,24,27,28	0

6.5 Other polymers (i)

There are no such residues in this entry.