Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:20 AM GMT

PDB ID	:	3AFF
Title	:	Crystal structure of the HsaA monooxygenase from M. tuberculosis
Authors	:	D'Angelo, I.; Lin, L.Y.; Dresen, C.; Tocheva, E.I.; Strynadka, N.; Eltis, L.D.
Deposited on	:	2010-02-28
Resolution	:	2.00 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:

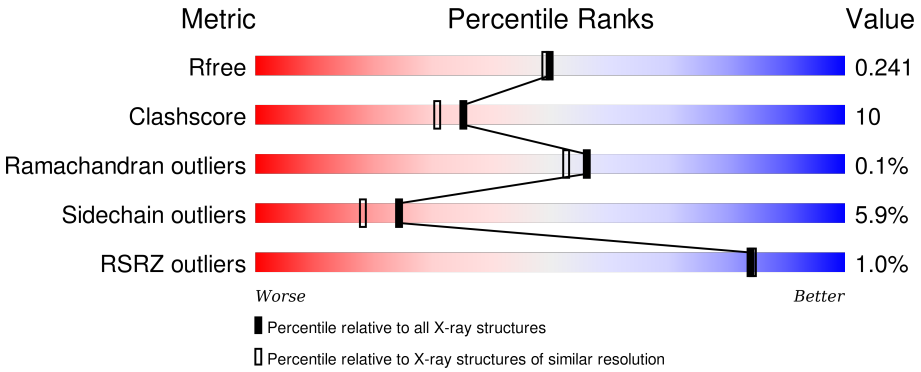
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)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

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

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	394	% 77% 12% • 7%
1	B	394	% 76% 13% • 8%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 6321 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 Hydroxylase, putative.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	366	Total	C	N	O	S	0	5	0
			2860	1806	512	533	9			
1	B	364	Total	C	N	O	S	0	4	0
			2839	1792	510	528	9			

- Molecule 2 is water.

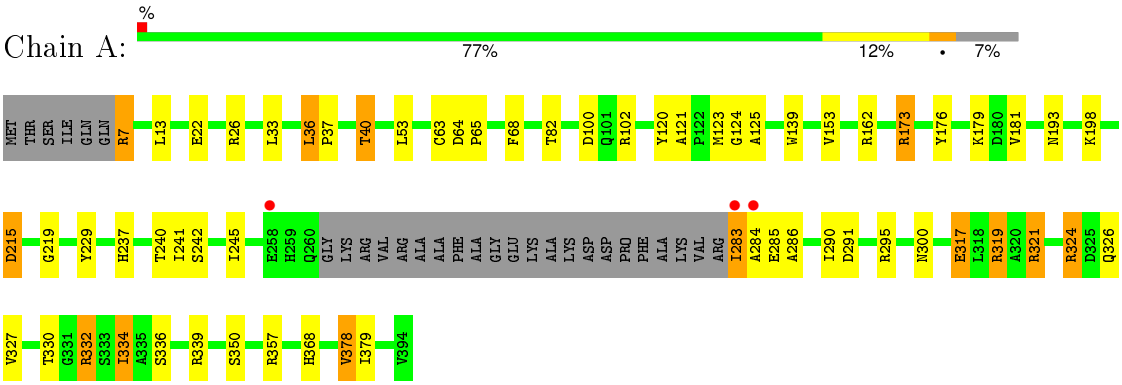
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	318	Total	O	0	0
			318	318		
2	B	304	Total	O	0	0
			304	304		

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.

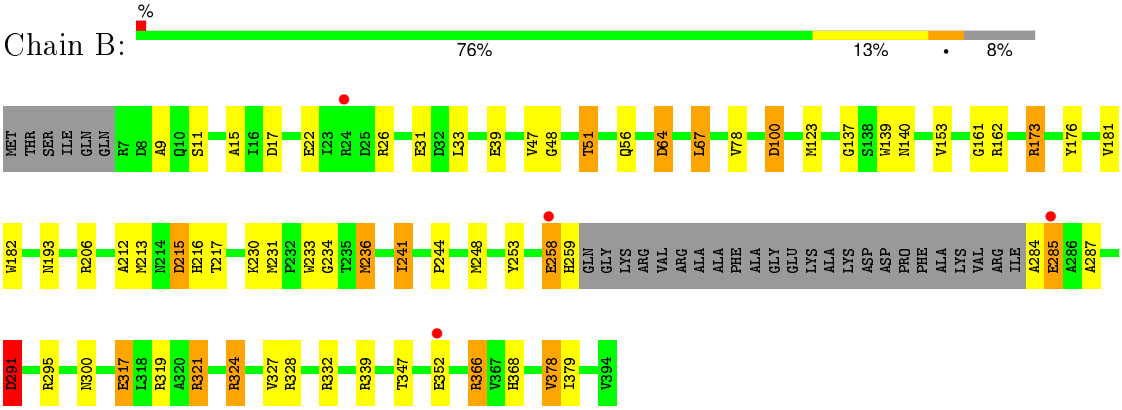
- Molecule 1: Hydroxylase, putative

Chain A:



- Molecule 1: Hydroxylase, putative

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α, β, γ	69.14Å 175.77Å 179.49Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.76 – 2.00 19.76 – 2.00	Depositor EDS
% Data completeness (in resolution range)	100.0 (19.76-2.00) 98.0 (19.76-2.00)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$< I/\sigma(I) >$ [1]	3.78 (at 2.01Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.198 , 0.239 0.200 , 0.241	Depositor DCC
R_{free} test set	3631 reflections (5.26%)	DCC
Wilson B-factor (Å^2)	26.0	Xtriage
Anisotropy	0.009	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.38 , 49.1	EDS
Estimated twinning fraction	0.019 for -h,-l,-k	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Outliers	0 of 72618 reflections	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	6321	wwPDB-VP
Average B, all atoms (Å^2)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.32% 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

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	1.05	3/2951 (0.1%)	0.98	12/4020 (0.3%)
1	B	1.03	4/2926 (0.1%)	0.98	16/3985 (0.4%)
All	All	1.04	7/5877 (0.1%)	0.98	28/8005 (0.3%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	317	GLU	CG-CD	6.87	1.62	1.51
1	B	206	ARG	CZ-NH2	6.56	1.41	1.33
1	B	317	GLU	CB-CG	5.79	1.63	1.52
1	A	120	TYR	CD1-CE1	5.43	1.47	1.39
1	B	215	ASP	CB-CG	-5.34	1.40	1.51
1	A	215	ASP	CB-CG	-5.13	1.41	1.51
1	A	22	GLU	CG-CD	5.01	1.59	1.51

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	319	ARG	NE-CZ-NH2	-12.02	114.29	120.30
1	A	215	ASP	CB-CG-OD1	-11.39	108.05	118.30
1	B	236	MET	CG-SD-CE	-10.25	83.80	100.20
1	B	319	ARG	NE-CZ-NH1	9.77	125.18	120.30
1	B	206	ARG	NE-CZ-NH1	-9.54	115.53	120.30
1	A	319	ARG	NE-CZ-NH2	-8.76	115.92	120.30
1	B	215	ASP	CB-CG-OD1	-8.52	110.63	118.30
1	B	321	ARG	NE-CZ-NH2	-8.19	116.20	120.30
1	A	319	ARG	NE-CZ-NH1	8.11	124.36	120.30
1	A	321	ARG	NE-CZ-NH1	-6.42	117.09	120.30
1	A	173	ARG	NE-CZ-NH2	-6.14	117.23	120.30
1	B	366	ARG	NE-CZ-NH2	5.82	123.21	120.30
1	B	241	ILE	CG1-CB-CG2	-5.67	98.92	111.40
1	B	378	VAL	CG1-CB-CG2	5.56	119.79	110.90
1	A	378	VAL	CG1-CB-CG2	5.54	119.76	110.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	64	ASP	CB-CG-OD2	-5.52	113.33	118.30
1	A	334	ILE	CG1-CB-CG2	-5.46	99.38	111.40
1	B	100[A]	ASP	CB-CG-OD1	-5.41	113.43	118.30
1	B	100[B]	ASP	CB-CG-OD1	-5.41	113.43	118.30
1	B	321	ARG	NE-CZ-NH1	5.36	122.98	120.30
1	A	215	ASP	CB-CG-OD2	5.35	123.11	118.30
1	A	7	ARG	NE-CZ-NH2	-5.30	117.65	120.30
1	A	64	ASP	CB-CG-OD2	-5.29	113.54	118.30
1	A	321	ARG	NE-CZ-NH2	5.29	122.94	120.30
1	B	319	ARG	CD-NE-CZ	5.24	130.94	123.60
1	B	291	ASP	CB-CG-OD2	-5.10	113.71	118.30
1	A	215	ASP	CB-CA-C	-5.10	100.21	110.40
1	B	215	ASP	CB-CA-C	-5.05	100.30	110.40

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2860	0	2749	62	0
1	B	2839	0	2732	53	1
2	A	318	0	0	28	1
2	B	304	0	0	20	1
All	All	6321	0	5481	110	2

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 10.

All (110) 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:237:HIS:O	1:A:240:THR:HG22	1.24	1.25
1:A:40:THR:HG21	2:A:570:HOH:O	1.55	1.05
1:B:140:ASN:HB3	2:B:444:HOH:O	1.57	1.04

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:231:MET:HB3	1:B:236:MET:HE1	1.43	0.95
1:A:173:ARG:NE	2:A:493:HOH:O	1.91	0.95
1:A:237:HIS:O	1:A:240:THR:CG2	2.16	0.94
1:A:339:ARG:HG2	2:A:705:HOH:O	1.71	0.90
1:B:231:MET:CB	1:B:236:MET:HE1	2.01	0.89
1:B:284:ALA:HB1	2:B:634:HOH:O	1.73	0.88
1:A:173:ARG:NH2	2:A:493:HOH:O	2.06	0.88
1:A:283:ILE:HG22	2:A:564:HOH:O	1.78	0.82
1:A:245:ILE:HG21	1:A:330:THR:HG23	1.67	0.77
1:A:63:CYS:SG	2:A:629:HOH:O	2.44	0.76
1:A:283:ILE:CG2	2:A:564:HOH:O	2.31	0.76
1:B:368:HIS:HE1	2:B:457:HOH:O	1.69	0.76
1:B:291:ASP:OD2	2:B:481:HOH:O	2.04	0.76
1:A:368:HIS:HD2	2:A:455:HOH:O	1.69	0.75
1:A:283:ILE:CG2	1:A:284:ALA:H	1.99	0.74
1:A:300:ASN:HD21	1:A:321:ARG:HH12	1.36	0.71
1:B:368:HIS:HD2	2:B:451:HOH:O	1.76	0.69
1:B:285:GLU:OE1	2:B:443:HOH:O	2.11	0.68
1:B:324:ARG:HD2	1:B:328:ARG:HH21	1.59	0.68
1:A:283:ILE:CG2	1:A:284:ALA:N	2.57	0.68
1:A:100[A]:ASP:HB2	2:A:580:HOH:O	1.94	0.67
1:A:121:ALA:HB3	1:A:123:MET:HE2	1.76	0.67
1:A:332:ARG:HG2	1:A:332:ARG:HH11	1.60	0.67
1:A:324:ARG:HD3	2:A:583:HOH:O	1.93	0.67
1:A:102:ARG:HD3	2:A:461:HOH:O	1.94	0.67
1:A:121:ALA:HB3	1:A:123:MET:CE	2.25	0.66
1:A:173:ARG:CZ	2:A:493:HOH:O	2.27	0.65
1:B:259:HIS:HD2	2:B:657:HOH:O	1.80	0.65
1:B:317:GLU:OE2	2:B:527:HOH:O	2.15	0.64
1:A:286:ALA:HB1	2:A:495:HOH:O	1.96	0.64
1:B:366:ARG:HD3	2:B:570:HOH:O	1.98	0.62
1:A:283:ILE:HG23	1:A:284:ALA:H	1.64	0.62
1:A:300:ASN:HD21	1:A:321:ARG:NH1	1.97	0.62
1:B:15:ALA:HB1	1:B:47:VAL:HG22	1.82	0.61
1:B:216:HIS:HD2	2:B:689:HOH:O	1.84	0.60
1:B:300:ASN:ND2	1:B:321:ARG:HH11	2.01	0.59
1:A:181:VAL:H	1:A:193:ASN:ND2	2.01	0.58
1:B:231:MET:HB2	1:B:236:MET:HE1	1.84	0.58
1:B:213:MET:HE2	1:B:233:TRP:HB2	1.85	0.58
1:B:100[A]:ASP:HB2	2:B:446:HOH:O	2.04	0.58
1:B:123:MET:HE1	1:B:140:ASN:H	1.69	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:48:GLY:O	1:B:51:THR:HB	2.03	0.57
1:A:283:ILE:HG22	1:A:284:ALA:H	1.68	0.57
1:B:285:GLU:CD	1:B:285:GLU:H	2.09	0.55
1:A:336:SER:HB3	2:A:495:HOH:O	2.06	0.55
1:B:26:ARG:NH2	1:B:39:GLU:OE1	2.38	0.54
1:B:162[B]:ARG:NE	2:B:600:HOH:O	2.40	0.54
1:A:339:ARG:NH2	2:A:674:HOH:O	2.39	0.54
1:A:65:PRO:HA	2:A:629:HOH:O	2.09	0.53
1:B:181:VAL:H	1:B:193:ASN:ND2	2.08	0.52
1:A:7:ARG:N	2:A:573:HOH:O	2.42	0.52
1:B:9:ALA:HB1	1:B:67:LEU:HD23	1.91	0.52
1:B:216:HIS:CD2	2:B:689:HOH:O	2.62	0.51
1:A:290:ILE:HD11	2:A:495:HOH:O	2.10	0.51
1:A:162[B]:ARG:CD	1:B:137:GLY:HA2	2.41	0.51
1:B:332:ARG:NE	2:B:694:HOH:O	2.44	0.50
1:A:283:ILE:HG22	1:A:284:ALA:N	2.24	0.50
1:B:259:HIS:CD2	2:B:657:HOH:O	2.60	0.50
1:B:182:TRP:HB2	1:B:193:ASN:ND2	2.26	0.50
1:A:317:GLU:HG2	2:A:539:HOH:O	2.11	0.49
1:A:357:ARG:HD2	2:A:589:HOH:O	2.12	0.49
1:A:321:ARG:NE	2:A:648:HOH:O	2.31	0.49
1:A:181:VAL:H	1:A:193:ASN:HD22	1.60	0.49
1:B:22:GLU:O	1:B:26:ARG:HG2	2.14	0.48
1:A:339:ARG:NH2	2:A:575:HOH:O	2.30	0.48
1:A:179:LYS:HG2	1:A:181:VAL:HG13	1.94	0.48
1:B:216:HIS:CD2	2:B:573:HOH:O	2.66	0.48
1:B:328:ARG:HD3	2:B:550:HOH:O	2.13	0.47
1:B:258:GLU:O	1:B:258:GLU:HG3	2.14	0.47
1:A:124:GLY:HA3	1:A:139:TRP:CH2	2.49	0.47
1:B:230:LYS:NZ	2:B:417:HOH:O	2.36	0.47
1:B:253:TYR:CZ	1:B:287:ALA:HB2	2.50	0.47
1:A:125:ALA:HB3	1:B:161:GLY:O	2.16	0.46
1:B:258:GLU:O	1:B:259:HIS:HB2	2.14	0.46
1:A:68:PHE:HB2	2:A:629:HOH:O	2.16	0.46
1:B:291:ASP:OD1	1:B:295:ARG:NH1	2.49	0.46
1:A:139:TRP:CE2	1:A:153:VAL:HB	2.51	0.46
1:A:295:ARG:NE	2:A:436:HOH:O	2.24	0.46
1:A:229:TYR:HD1	2:A:683:HOH:O	1.97	0.46
1:A:26:ARG:HH11	1:A:40:THR:HG22	1.81	0.45
1:A:162[B]:ARG:HD2	1:B:137:GLY:HA2	1.99	0.45
1:B:284:ALA:HB3	2:B:443:HOH:O	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:36:LEU:HD11	1:A:82:THR:HA	2.00	0.44
1:B:215:ASP:HB3	1:B:217:THR:HG23	2.00	0.44
1:B:173[B]:ARG:HA	1:B:176:TYR:CE2	2.54	0.43
1:A:379:ILE:CD1	1:B:379:ILE:CD1	2.97	0.43
1:A:121:ALA:HB3	1:A:123:MET:HE3	1.98	0.43
1:A:327:VAL:HG13	2:A:610:HOH:O	2.18	0.43
1:B:139:TRP:CE2	1:B:153:VAL:HB	2.54	0.43
1:B:300:ASN:HD21	1:B:321:ARG:HH11	1.65	0.43
1:B:173[A]:ARG:HA	1:B:176:TYR:CE2	2.54	0.43
1:B:212:ALA:HA	1:B:215:ASP:HB2	2.01	0.43
1:B:234:GLY:HA2	2:B:587:HOH:O	2.19	0.43
1:A:379:ILE:HD12	1:B:379:ILE:HD12	2.02	0.42
1:A:283:ILE:HG23	1:A:284:ALA:N	2.29	0.42
1:A:37:PRO:HD2	1:A:40:THR:HG23	2.02	0.42
1:B:64:ASP:HB3	1:B:67:LEU:HD22	2.01	0.42
1:B:244:PRO:O	1:B:248:MET:HG3	2.20	0.42
1:A:53:LEU:HD21	2:A:629:HOH:O	2.18	0.42
1:A:100[B]:ASP:OD1	1:A:219:GLY:HA3	2.19	0.42
1:A:332:ARG:HG2	1:A:332:ARG:NH1	2.32	0.41
1:A:283:ILE:HD12	1:A:283:ILE:HA	1.84	0.41
1:A:237:HIS:HB3	2:A:695:HOH:O	2.20	0.41
1:B:9:ALA:CB	1:B:67:LEU:HD23	2.51	0.41
1:A:242:SER:OG	1:A:326:GLN:NE2	2.54	0.40
1:A:173:ARG:HA	1:A:176:TYR:CE2	2.57	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:317:GLU:CG	2:A:650:HOH:O[3_655]	2.18	0.02
2:B:553:HOH:O	2:B:553:HOH:O[2_755]	2.19	0.01

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	367/394 (93%)	364 (99%)	3 (1%)	0	100 100
1	B	364/394 (92%)	359 (99%)	4 (1%)	1 (0%)	46 41
All	All	731/788 (93%)	723 (99%)	7 (1%)	1 (0%)	56 53

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	258	GLU

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	294/310 (95%)	277 (94%)	17 (6%)	25 19
1	B	291/310 (94%)	272 (94%)	19 (6%)	21 15
All	All	585/620 (94%)	549 (94%)	36 (6%)	24 16

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	LEU
1	A	33	LEU
1	A	36	LEU
1	A	40	THR
1	A	198	LYS
1	A	215	ASP
1	A	241	ILE
1	A	283	ILE
1	A	285	GLU
1	A	291[A]	ASP
1	A	291[B]	ASP
1	A	317	GLU

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Mol	Chain	Res	Type
1	A	324	ARG
1	A	332	ARG
1	A	334	ILE
1	A	350	SER
1	A	378	VAL
1	B	11	SER
1	B	17	ASP
1	B	31	GLU
1	B	33	LEU
1	B	51	THR
1	B	56	GLN
1	B	67	LEU
1	B	78	VAL
1	B	173[A]	ARG
1	B	173[B]	ARG
1	B	241	ILE
1	B	285	GLU
1	B	291	ASP
1	B	324	ARG
1	B	327	VAL
1	B	339	ARG
1	B	347	THR
1	B	352	GLU
1	B	378	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	193	ASN
1	A	237	HIS
1	A	300	ASN
1	A	326	GLN
1	A	368	HIS
1	A	371	ASN
1	A	382	ASN
1	B	193	ASN
1	B	300	ASN
1	B	326	GLN
1	B	368	HIS
1	B	371	ASN
1	B	382	ASN

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

There are no ligands in this entry.

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

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	366/394 (92%)	-0.39	3 (0%)	87	88	15, 24, 41, 57	0
1	B	364/394 (92%)	-0.32	4 (1%)	82	83	15, 26, 44, 72	0
All	All	730/788 (92%)	-0.35	7 (0%)	84	84	15, 25, 43, 72	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	258	GLU	3.7
1	B	352	GLU	3.7
1	A	258	GLU	3.1
1	A	283	ILE	2.7
1	A	284	ALA	2.5
1	B	285	GLU	2.4
1	B	24	ARG	2.1

6.2 Non-standard residues in protein, DNA, RNA chains

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

6.3 Carbohydrates

There are no carbohydrates in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.