



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 03:38 PM GMT

PDB ID : 4CY6
Title : apo structure of 2-hydroxybiphenyl 3-monooxygenase HbpA
Authors : Jensen, C.N.; Farrugia, J.E.; Frank, A.; Man, H.; Hart, S.; Turkenburg, J.P.; Grogan, G.
Deposited on : 2014-04-10
Resolution : 2.76 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

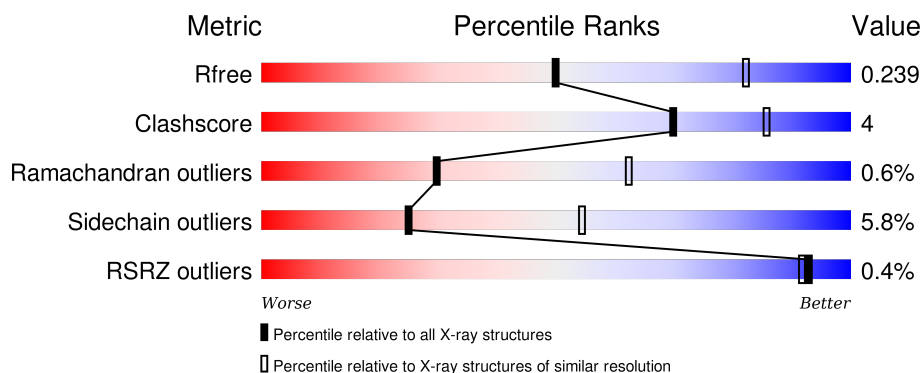
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION





The reported resolution of this entry is 2.76 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	3340 (2.80-2.72)
Clashscore	102246	3829 (2.80-2.72)
Ramachandran outliers	100387	3767 (2.80-2.72)
Sidechain outliers	100360	3770 (2.80-2.72)
RSRZ outliers	91569	3352 (2.80-2.72)

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	586	 81% 12% • •
1	B	586	 83% 10% • 5%
1	C	586	 81% 12% • • 5%
1	D	586	 82% 11% • • 6%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 16265 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 2-HYDROXYBIPHENYL-3-MONOOXYGENASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	561	Total	C	N	O	S	0	0	0
			4095	2603	723	749	20			
1	B	555	Total	C	N	O	S	0	0	0
			3982	2532	707	724	19			
1	C	554	Total	C	N	O	S	0	0	0
			4017	2545	717	736	19			
1	D	553	Total	C	N	O	S	0	0	0
			4032	2556	716	740	20			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	347	GLN	THR	ENGINEERED MUTATION	UNP O06647
B	347	GLN	THR	ENGINEERED MUTATION	UNP O06647
C	347	GLN	THR	ENGINEERED MUTATION	UNP O06647
D	347	GLN	THR	ENGINEERED MUTATION	UNP O06647

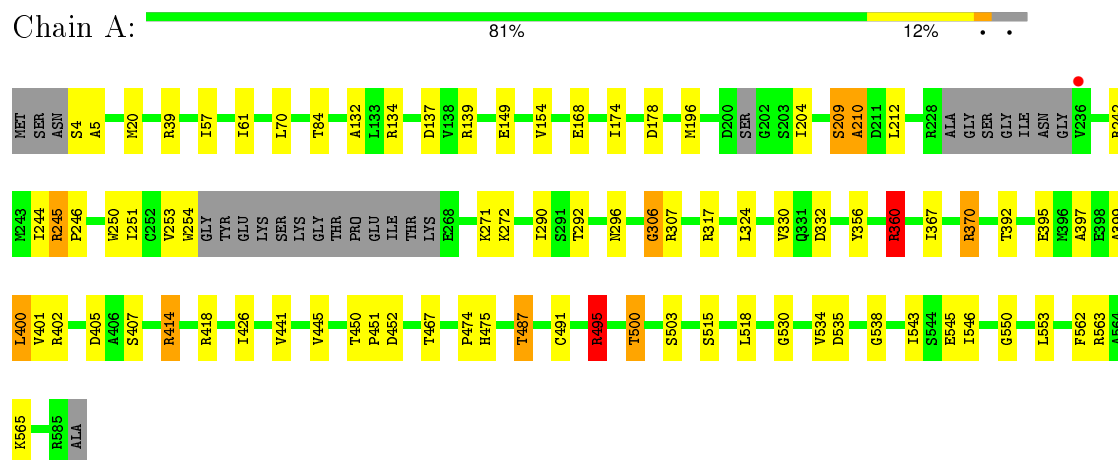
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	40	Total	O	0	0
			40	40		
2	B	35	Total	O	0	0
			35	35		
2	C	33	Total	O	0	0
			33	33		
2	D	31	Total	O	0	0
			31	31		

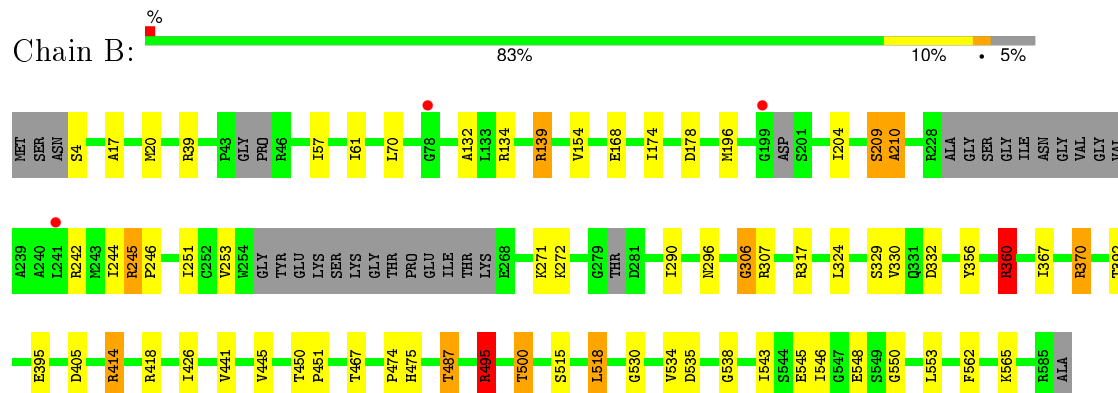
3 Residue-property plots

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.

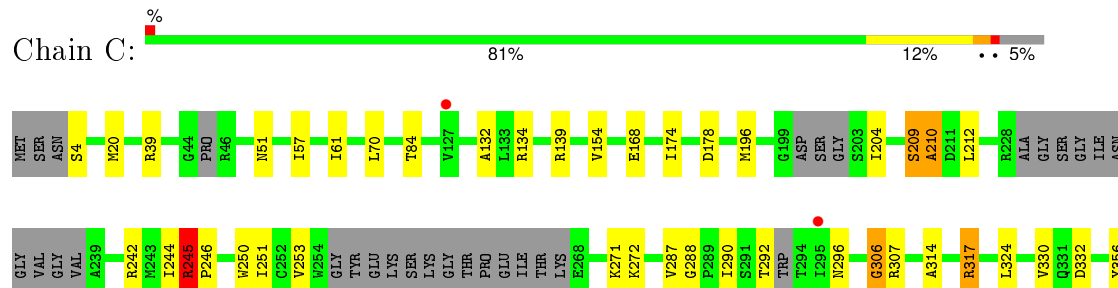
• Molecule 1: 2-HYDROXYBIPHENYL-3-MONOOXYGENASE



• Molecule 1: 2-HYDROXYBIPHENYL-3-MONOOXYGENASE



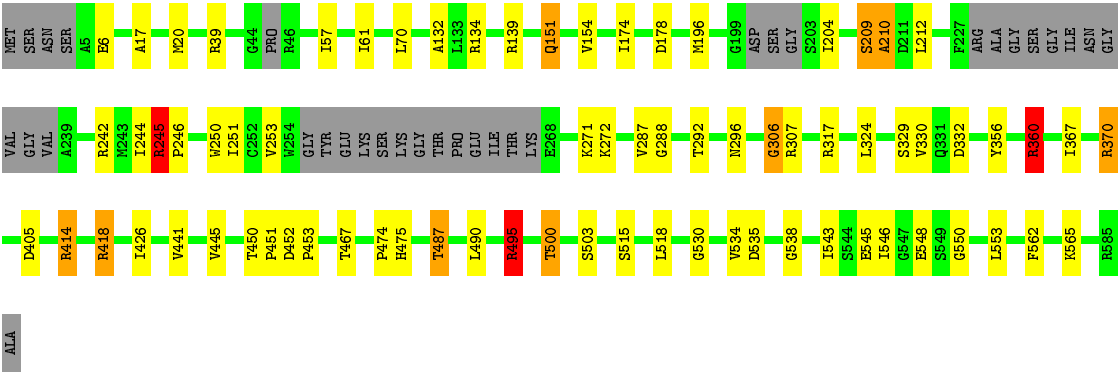
• Molecule 1: 2-HYDROXYBIPHENYL-3-MONOOXYGENASE





• Molecule 1: 2-HYDROXYBIPHENYL-3-MONOOXYGENASE

Chain D: 82% 11% 6%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	79.35Å 94.67Å 99.16Å 115.03° 96.09° 109.59°	Depositor
Resolution (Å)	85.97 – 2.76 85.97 – 2.76	Depositor EDS
% Data completeness (in resolution range)	96.8 (85.97-2.76) 79.6 (85.97-2.76)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.19 (at 2.77Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.190 , 0.241 0.195 , 0.239	Depositor DCC
R_{free} test set	2997 reflections (5.38%)	DCC
Wilson B-factor (Å ²)	52.8	Xtriage
Anisotropy	0.419	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 56.5	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	0 of 58683 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	16265	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.95% 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

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	0.67	0/4184	0.83	5/5695 (0.1%)
1	B	0.68	0/4068	0.84	4/5547 (0.1%)
1	C	0.66	0/4102	0.84	6/5587 (0.1%)
1	D	0.67	1/4119 (0.0%)	0.84	7/5610 (0.1%)
All	All	0.67	1/16473 (0.0%)	0.84	22/22439 (0.1%)

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
1	B	0	4
1	C	0	4
1	D	0	5
All	All	0	17

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	453	PRO	N-CD	5.13	1.55	1.47

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	495	ARG	NE-CZ-NH1	8.92	124.76	120.30
1	C	495	ARG	NE-CZ-NH1	8.47	124.53	120.30
1	B	495	ARG	NE-CZ-NH1	8.17	124.39	120.30
1	A	495	ARG	NE-CZ-NH1	8.03	124.31	120.30
1	A	535	ASP	CB-CG-OD1	-7.66	111.41	118.30
1	B	535	ASP	CB-CG-OD1	-7.57	111.49	118.30
1	A	360	ARG	NE-CZ-NH1	7.50	124.05	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	535	ASP	CB-CG-OD1	-7.43	111.61	118.30
1	D	360	ARG	NE-CZ-NH1	7.37	123.98	120.30
1	B	360	ARG	NE-CZ-NH1	6.70	123.65	120.30
1	D	535	ASP	CB-CG-OD1	-6.58	112.38	118.30
1	C	360	ARG	NE-CZ-NH1	6.33	123.46	120.30
1	D	414	ARG	NE-CZ-NH1	6.22	123.41	120.30
1	D	418	ARG	NE-CZ-NH1	6.04	123.32	120.30
1	D	452	ASP	C-N-CD	5.72	140.42	128.40
1	C	414	ARG	NE-CZ-NH1	5.55	123.08	120.30
1	B	414	ARG	NE-CZ-NH1	5.40	123.00	120.30
1	C	245	ARG	NE-CZ-NH1	5.35	122.98	120.30
1	A	414	ARG	NE-CZ-NH1	5.28	122.94	120.30
1	A	491	CYS	N-CA-C	5.17	124.96	111.00
1	C	491	CYS	N-CA-C	5.12	124.83	111.00
1	D	245	ARG	NE-CZ-NH1	5.02	122.81	120.30

There are no chirality outliers.

All (17) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	271	LYS	Peptide
1	A	272	LYS	Peptide
1	A	530	GLY	Peptide
1	A	545	GLU	Peptide
1	B	271	LYS	Peptide
1	B	272	LYS	Peptide
1	B	530	GLY	Peptide
1	B	545	GLU	Peptide
1	C	271	LYS	Peptide
1	C	272	LYS	Peptide
1	C	531	GLN	Peptide
1	C	545	GLU	Peptide
1	D	271	LYS	Peptide
1	D	272	LYS	Peptide
1	D	490	LEU	Peptide
1	D	530	GLY	Peptide
1	D	545	GLU	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	4095	0	3918	41	0
1	B	3982	0	3750	37	0
1	C	4017	0	3798	39	0
1	D	4032	0	3830	35	0
2	A	40	0	0	2	0
2	B	35	0	0	0	0
2	C	33	0	0	2	0
2	D	31	0	0	0	0
All	All	16265	0	15296	140	0

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 4.

All (140) 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:132:ALA:O	1:B:39:ARG:NH1	1.97	0.96
1:A:196:MET:O	1:A:296:ASN:ND2	2.07	0.86
1:D:196:MET:O	1:D:296:ASN:ND2	2.08	0.86
1:B:196:MET:O	1:B:296:ASN:ND2	2.08	0.85
1:C:196:MET:O	1:C:296:ASN:ND2	2.11	0.83
1:B:405:ASP:O	1:B:414:ARG:NH2	2.13	0.82
1:C:39:ARG:NH1	1:D:132:ALA:O	2.14	0.81
1:A:405:ASP:O	1:A:414:ARG:NH2	2.13	0.80
1:A:39:ARG:NH1	1:B:132:ALA:O	2.14	0.80
1:D:405:ASP:O	1:D:414:ARG:NH2	2.14	0.80
1:D:70:LEU:O	1:D:245:ARG:NH2	2.16	0.79
1:C:405:ASP:O	1:C:414:ARG:NH2	2.16	0.79
1:C:70:LEU:O	1:C:245:ARG:NH2	2.18	0.77
1:B:515:SER:O	1:B:518:LEU:O	2.03	0.76
1:B:70:LEU:O	1:B:245:ARG:NH2	2.17	0.76
1:C:132:ALA:O	1:D:39:ARG:NH1	2.20	0.75
1:A:70:LEU:O	1:A:245:ARG:NH2	2.19	0.75
1:C:515:SER:O	1:C:518:LEU:O	2.06	0.74
1:A:515:SER:O	1:A:518:LEU:O	2.06	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:154:VAL:HG21	1:B:174:ILE:HG13	1.71	0.73
1:D:515:SER:O	1:D:518:LEU:O	2.06	0.72
1:D:154:VAL:HG21	1:D:174:ILE:HG13	1.74	0.69
1:D:244:ILE:O	1:D:245:ARG:HD2	1.93	0.69
1:C:154:VAL:HG21	1:C:174:ILE:HG13	1.74	0.69
1:A:154:VAL:HG21	1:A:174:ILE:HG13	1.75	0.68
1:C:244:ILE:O	1:C:245:ARG:HD2	1.93	0.68
1:A:244:ILE:O	1:A:245:ARG:HD2	1.93	0.68
1:B:244:ILE:O	1:B:245:ARG:HD2	1.93	0.68
1:A:500:THR:HG23	1:A:550:GLY:O	1.96	0.65
1:B:500:THR:HG23	1:B:550:GLY:O	1.96	0.65
1:A:149:GLU:OE1	1:A:306:GLY:O	2.15	0.65
1:C:500:THR:HG23	1:C:550:GLY:O	1.98	0.64
1:D:306:GLY:O	1:D:307:ARG:HB2	1.98	0.64
1:A:306:GLY:O	1:A:307:ARG:HB2	1.97	0.63
1:B:306:GLY:O	1:B:307:ARG:HB2	1.99	0.63
1:D:500:THR:HG23	1:D:550:GLY:O	1.98	0.63
2:A:2016:HOH:O	1:B:139:ARG:HD3	1.98	0.62
1:B:414:ARG:NH1	1:D:548:GLU:OE2	2.33	0.62
1:C:306:GLY:O	1:C:307:ARG:HB2	1.98	0.61
1:A:84:THR:HG21	1:A:400:LEU:HD11	1.83	0.60
1:B:418:ARG:HD3	1:D:538:GLY:HA3	1.83	0.60
1:D:474:PRO:O	1:D:487:THR:HG21	2.02	0.59
1:B:474:PRO:O	1:B:487:THR:HG21	2.01	0.58
1:C:84:THR:HG21	1:C:400:LEU:HD11	1.85	0.58
1:A:474:PRO:O	1:A:487:THR:HG21	2.04	0.57
1:B:495:ARG:HH11	1:B:495:ARG:CG	2.18	0.57
1:A:495:ARG:CG	1:A:495:ARG:HH11	2.18	0.57
1:D:332:ASP:OD1	1:D:360:ARG:NH1	2.38	0.57
1:B:546:ILE:HD12	1:B:565:LYS:HA	1.87	0.56
1:C:474:PRO:O	1:C:487:THR:HG21	2.05	0.56
1:C:495:ARG:HH11	1:C:495:ARG:CG	2.19	0.55
1:A:495:ARG:HG3	1:A:495:ARG:HH11	1.72	0.54
1:B:332:ASP:OD1	1:B:360:ARG:NH1	2.41	0.54
1:C:332:ASP:OD1	1:C:360:ARG:NH1	2.41	0.54
1:C:546:ILE:HD12	1:C:565:LYS:HA	1.90	0.53
1:A:332:ASP:OD1	1:A:360:ARG:NH1	2.39	0.53
1:A:414:ARG:NH1	1:C:548:GLU:OE2	2.42	0.53
1:D:495:ARG:HH11	1:D:495:ARG:HG3	1.74	0.53
1:C:495:ARG:HH11	1:C:495:ARG:HG3	1.73	0.53
1:B:495:ARG:HH11	1:B:495:ARG:HG3	1.73	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:495:ARG:HH11	1:D:495:ARG:CG	2.22	0.52
1:D:450:THR:HG22	1:D:451:PRO:O	2.11	0.51
1:B:548:GLU:OE2	1:D:414:ARG:NH1	2.45	0.50
1:D:546:ILE:HD12	1:D:565:LYS:HA	1.94	0.50
1:A:553:LEU:HB3	1:A:562:PHE:HB3	1.93	0.49
1:A:5:ALA:N	2:A:2001:HOH:O	2.46	0.49
1:A:204:ILE:O	1:A:253:VAL:O	2.31	0.48
1:A:546:ILE:HD12	1:A:565:LYS:HA	1.95	0.48
1:C:553:LEU:HB3	1:C:562:PHE:HB3	1.96	0.48
1:D:204:ILE:O	1:D:253:VAL:O	2.32	0.47
1:B:553:LEU:HB3	1:B:562:PHE:HB3	1.96	0.47
1:A:137:ASP:OD1	1:B:39:ARG:HD3	2.15	0.47
1:A:209:SER:O	1:A:210:ALA:HB2	2.15	0.47
1:D:360:ARG:HG3	1:D:360:ARG:HH11	1.80	0.47
1:D:553:LEU:HB3	1:D:562:PHE:HB3	1.96	0.47
1:D:151:GLN:HA	1:D:151:GLN:HE21	1.80	0.47
1:C:209:SER:O	1:C:210:ALA:HB2	2.15	0.47
1:B:209:SER:O	1:B:210:ALA:HB2	2.15	0.46
1:B:204:ILE:O	1:B:253:VAL:O	2.33	0.46
1:B:475:HIS:HA	1:B:487:THR:HG22	1.98	0.46
1:C:399:ALA:HA	1:C:402:ARG:HD3	1.96	0.46
1:A:399:ALA:HA	1:A:402:ARG:HD3	1.98	0.45
1:B:20:MET:HG2	1:B:330:VAL:HG13	1.97	0.45
1:D:209:SER:O	1:D:210:ALA:HB2	2.16	0.45
1:D:20:MET:HG2	1:D:330:VAL:HG13	1.97	0.45
1:B:154:VAL:HG21	1:B:174:ILE:CG1	2.45	0.45
1:C:204:ILE:O	1:C:253:VAL:O	2.34	0.45
1:D:370:ARG:NH2	1:D:426:ILE:HG12	2.32	0.45
1:A:20:MET:HG2	1:A:330:VAL:HG13	1.98	0.44
1:B:370:ARG:NH2	1:B:426:ILE:HG12	2.32	0.44
1:B:242:ARG:HD3	1:B:251:ILE:HD11	2.00	0.44
1:A:405:ASP:OD1	1:A:407:SER:OG	2.32	0.44
1:C:20:MET:HG2	1:C:330:VAL:HG13	2.00	0.44
1:A:450:THR:HG22	1:A:451:PRO:O	2.17	0.43
1:C:398:GLU:O	1:C:402:ARG:HD3	2.19	0.43
1:B:450:THR:HG22	1:B:451:PRO:O	2.19	0.43
1:D:212:LEU:HD12	1:D:250:TRP:CE2	2.53	0.43
1:B:392:THR:HG22	1:B:395:GLU:OE1	2.19	0.43
1:C:450:THR:HG22	1:C:451:PRO:O	2.18	0.43
1:A:57:ILE:O	1:A:61:ILE:HG12	2.19	0.43
1:D:242:ARG:HD3	1:D:251:ILE:HD11	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:418:ARG:HD3	1:C:538:GLY:HA3	1.99	0.42
1:C:287:VAL:HG12	1:C:288:GLY:O	2.20	0.42
1:C:360:ARG:NH2	2:C:2021:HOH:O	2.41	0.42
1:A:538:GLY:HA3	1:C:418:ARG:HD3	2.00	0.42
1:A:356:TYR:O	1:A:360:ARG:HB2	2.20	0.42
1:C:370:ARG:NH2	1:C:426:ILE:HG12	2.35	0.42
1:D:61:ILE:O	1:D:61:ILE:HG22	2.20	0.42
1:A:475:HIS:HA	1:A:487:THR:HG22	2.01	0.42
1:C:4:SER:HA	1:C:168:GLU:O	2.20	0.42
1:B:495:ARG:CG	1:B:495:ARG:NH1	2.83	0.42
1:D:57:ILE:O	1:D:61:ILE:HG12	2.19	0.42
1:A:212:LEU:HD12	1:A:250:TRP:CE2	2.55	0.42
1:C:242:ARG:HD3	1:C:251:ILE:HD11	2.00	0.42
1:C:464:GLN:HG2	2:C:2031:HOH:O	2.19	0.42
1:B:4:SER:HA	1:B:168:GLU:O	2.20	0.42
1:A:360:ARG:HH11	1:A:360:ARG:HG3	1.85	0.42
1:B:61:ILE:O	1:B:61:ILE:HG22	2.20	0.42
1:A:242:ARG:HD3	1:A:251:ILE:HD11	2.02	0.42
1:A:392:THR:HG22	1:A:395:GLU:OE1	2.20	0.41
1:D:475:HIS:HA	1:D:487:THR:HG22	2.02	0.41
1:A:546:ILE:CG1	1:A:563:ARG:HD2	2.50	0.41
1:C:314:ALA:O	1:C:317:ARG:NH1	2.54	0.41
1:A:370:ARG:NH2	1:A:426:ILE:HG12	2.36	0.41
1:A:154:VAL:HG21	1:A:174:ILE:CG1	2.49	0.41
1:B:57:ILE:O	1:B:61:ILE:HG12	2.21	0.41
1:A:4:SER:HA	1:A:168:GLU:O	2.21	0.41
1:B:17:ALA:HB2	1:B:329:SER:HB3	2.03	0.41
1:C:356:TYR:O	1:C:360:ARG:HB2	2.21	0.41
1:C:360:ARG:HG3	1:C:360:ARG:HH11	1.85	0.41
1:D:287:VAL:HG12	1:D:288:GLY:O	2.21	0.41
1:B:538:GLY:HA3	1:D:418:ARG:HD3	2.02	0.41
1:D:17:ALA:HB2	1:D:329:SER:HB3	2.03	0.41
1:D:356:TYR:O	1:D:360:ARG:HB2	2.21	0.40
1:C:57:ILE:O	1:C:61:ILE:HG12	2.21	0.40
1:C:212:LEU:HD12	1:C:250:TRP:CE2	2.55	0.40
1:C:51:ASN:OD1	1:C:51:ASN:C	2.60	0.40
1:A:397:ALA:O	1:A:401:VAL:HG13	2.21	0.40
1:C:154:VAL:HG21	1:C:174:ILE:CG1	2.47	0.40
1:B:356:TYR:O	1:B:360:ARG:HB2	2.21	0.40

There are no symmetry-related clashes.

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	553/586 (94%)	521 (94%)	29 (5%)	3 (0%)	34	67
1	B	543/586 (93%)	514 (95%)	26 (5%)	3 (1%)	30	62
1	C	542/586 (92%)	510 (94%)	28 (5%)	4 (1%)	26	59
1	D	543/586 (93%)	513 (94%)	27 (5%)	3 (1%)	30	62
All	All	2181/2344 (93%)	2058 (94%)	110 (5%)	13 (1%)	30	62

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	306	GLY
1	B	306	GLY
1	C	306	GLY
1	D	306	GLY
1	A	210	ALA
1	B	210	ALA
1	C	210	ALA
1	C	533	PHE
1	D	210	ALA
1	A	246	PRO
1	B	246	PRO
1	C	246	PRO
1	D	246	PRO

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	392/467 (84%)	368 (94%)	24 (6%)	23	52
1	B	371/467 (79%)	351 (95%)	20 (5%)	27	58
1	C	379/467 (81%)	356 (94%)	23 (6%)	23	52
1	D	385/467 (82%)	363 (94%)	22 (6%)	25	55
All	All	1527/1868 (82%)	1438 (94%)	89 (6%)	25	54

All (89) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	134	ARG
1	A	139	ARG
1	A	178	ASP
1	A	209	SER
1	A	245	ARG
1	A	254	TRP
1	A	290	ILE
1	A	292	THR
1	A	317	ARG
1	A	324	LEU
1	A	360	ARG
1	A	367	ILE
1	A	370	ARG
1	A	400	LEU
1	A	441	VAL
1	A	445	VAL
1	A	452	ASP
1	A	467	THR
1	A	487	THR
1	A	495	ARG
1	A	500	THR
1	A	503	SER
1	A	534	VAL
1	A	543	ILE
1	B	134	ARG
1	B	139	ARG
1	B	178	ASP
1	B	209	SER
1	B	245	ARG
1	B	290	ILE
1	B	317	ARG
1	B	324	LEU
1	B	360	ARG

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Mol	Chain	Res	Type
1	B	367	ILE
1	B	370	ARG
1	B	441	VAL
1	B	445	VAL
1	B	467	THR
1	B	487	THR
1	B	495	ARG
1	B	500	THR
1	B	518	LEU
1	B	534	VAL
1	B	543	ILE
1	C	134	ARG
1	C	139	ARG
1	C	178	ASP
1	C	209	SER
1	C	245	ARG
1	C	290	ILE
1	C	292	THR
1	C	317	ARG
1	C	324	LEU
1	C	360	ARG
1	C	367	ILE
1	C	370	ARG
1	C	400	LEU
1	C	441	VAL
1	C	445	VAL
1	C	467	THR
1	C	487	THR
1	C	495	ARG
1	C	500	THR
1	C	503	SER
1	C	534	VAL
1	C	543	ILE
1	C	569	ARG
1	D	6	GLU
1	D	134	ARG
1	D	139	ARG
1	D	151	GLN
1	D	178	ASP
1	D	209	SER
1	D	245	ARG
1	D	292	THR

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Mol	Chain	Res	Type
1	D	317	ARG
1	D	324	LEU
1	D	360	ARG
1	D	367	ILE
1	D	370	ARG
1	D	441	VAL
1	D	445	VAL
1	D	467	THR
1	D	487	THR
1	D	495	ARG
1	D	500	THR
1	D	503	SER
1	D	534	VAL
1	D	543	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	80	HIS
1	A	107	HIS
1	A	275	HIS
1	C	80	HIS
1	C	107	HIS
1	C	275	HIS
1	C	516	GLN
1	D	107	HIS
1	D	151	GLN
1	D	275	HIS
1	D	481	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 [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	561/586 (95%)	-0.26	1 (0%) 95 95	39, 59, 81, 118	0
1	B	555/586 (94%)	-0.22	3 (0%) 91 90	39, 61, 86, 119	0
1	C	554/586 (94%)	-0.20	4 (0%) 89 86	40, 60, 87, 113	0
1	D	553/586 (94%)	-0.23	0 100 100	37, 60, 86, 117	0
All	All	2223/2344 (94%)	-0.23	8 (0%) 93 92	37, 60, 86, 119	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	241	LEU	3.1
1	B	78	GLY	2.7
1	C	127	VAL	2.3
1	C	543	ILE	2.3
1	B	199	GLY	2.3
1	A	236	VAL	2.3
1	C	295	ILE	2.2
1	C	389	PRO	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.