



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 08:01 PM GMT

PDB ID : 1IC5
Title : CRYSTAL STRUCTURE OF HYHEL-10 FV MUTANT(HD99A)-HEN
LYSOZYME COMPLEX
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Deposited on : 2001-03-30
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) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

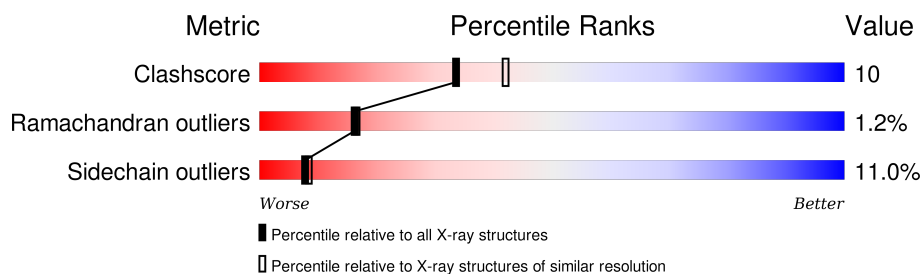
1 Overall quality at a glance

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))
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	L	107	
2	H	114	
3	Y	129	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 2834 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 LYSOZYME BINDING IG KAPPA CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	107	Total	C	N	O	S	0	0	0
			818	511	138	166	3			

- Molecule 2 is a protein called IGG1 FAB CHAIN H.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	114	Total	C	N	O	S	0	0	0
			899	566	144	186	3			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
H	99	ALA	ASP	ENGINEERED	UNP P01823

- Molecule 3 is a protein called LYSOZYME C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	Y	129	Total	C	N	O	S	0	0	0
			1001	613	193	185	10			

- Molecule 4 is water.

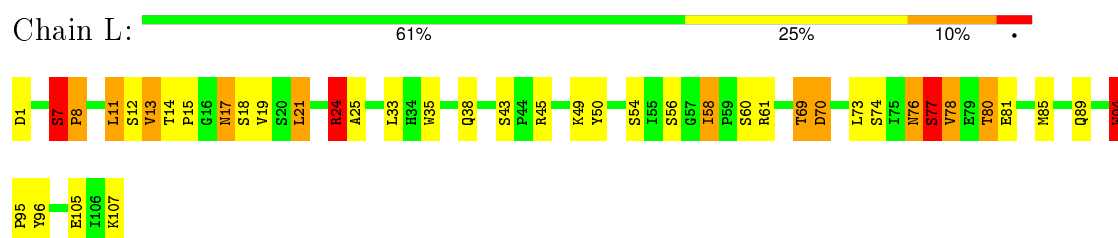
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	H	37	Total	O	0	0
			37	37		
4	L	35	Total	O	0	0
			35	35		
4	Y	44	Total	O	0	0
			44	44		

3 Residue-property plots

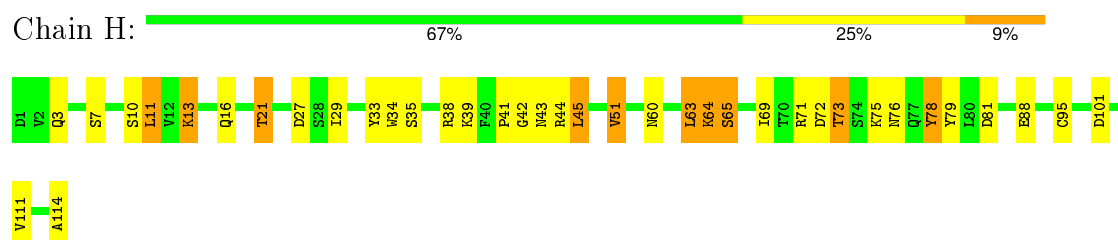
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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 ($RSRZ > 2$). 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.

Note EDS was not executed.

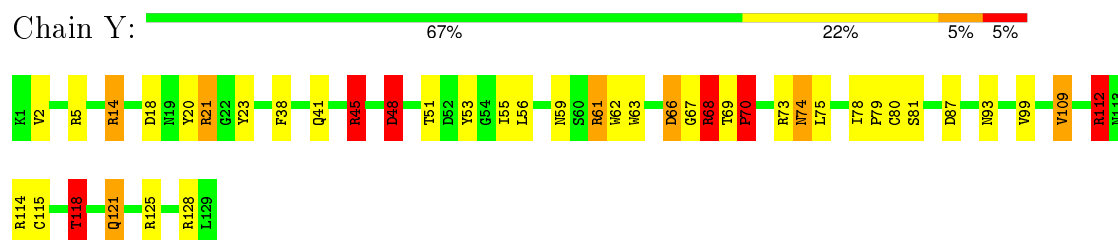
• Molecule 1: LYSOZYME BINDING IG KAPPA CHAIN



• Molecule 2: IGG1 FAB CHAIN H



• Molecule 3: LYSOZYME C



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	57.00 Å 57.00 Å 238.09 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.30	Depositor
% Data completeness (in resolution range)	(Not available) (8.00-2.30)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	REFMAC	Depositor
R, R_{free}	0.165 , 0.252	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2834	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

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	L	0.99	1/837 (0.1%)	2.11	33/1133 (2.9%)
2	H	0.99	0/922	2.02	28/1259 (2.2%)
3	Y	0.96	0/1021	1.99	35/1379 (2.5%)
All	All	0.98	1/2780 (0.0%)	2.04	96/3771 (2.5%)

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	L	0	5

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	L	18	SER	CB-OG	5.33	1.49	1.42

All (96) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	71	ARG	CD-NE-CZ	16.48	146.68	123.60
1	L	1	ASP	CB-CG-OD2	-13.64	106.03	118.30
2	H	81	ASP	CB-CG-OD1	12.60	129.64	118.30
1	L	49	LYS	CA-CB-CG	11.44	138.57	113.40
2	H	38	ARG	NE-CZ-NH2	-11.14	114.73	120.30
1	L	61	ARG	NE-CZ-NH2	10.30	125.45	120.30
3	Y	128	ARG	NE-CZ-NH1	-10.24	115.18	120.30
3	Y	21	ARG	NE-CZ-NH2	-10.22	115.19	120.30
2	H	71	ARG	NE-CZ-NH1	-10.04	115.28	120.30
2	H	41	PRO	CA-C-N	9.83	135.86	116.20
3	Y	66	ASP	CB-CG-OD2	9.81	127.13	118.30
1	L	96	TYR	CB-CG-CD1	9.33	126.59	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	Y	23	TYR	CB-CG-CD1	-9.07	115.56	121.00
3	Y	125	ARG	NE-CZ-NH1	9.01	124.80	120.30
1	L	45	ARG	NE-CZ-NH2	8.69	124.64	120.30
1	L	69	THR	N-CA-CB	-8.38	94.38	110.30
2	H	33	TYR	CB-CG-CD1	8.30	125.98	121.00
1	L	24	ARG	NE-CZ-NH2	8.28	124.44	120.30
2	H	44	ARG	CD-NE-CZ	8.20	135.08	123.60
3	Y	68	ARG	NE-CZ-NH1	-7.96	116.32	120.30
3	Y	87	ASP	CB-CG-OD1	7.91	125.41	118.30
1	L	7	SER	CA-C-O	-7.83	103.66	120.10
3	Y	114	ARG	NE-CZ-NH2	-7.82	116.39	120.30
1	L	50	TYR	CB-CG-CD1	-7.54	116.47	121.00
2	H	7	SER	N-CA-CB	7.51	121.76	110.50
1	L	50	TYR	CB-CG-CD2	7.42	125.45	121.00
2	H	51	VAL	CB-CA-C	-7.35	97.44	111.40
3	Y	112	ARG	CD-NE-CZ	7.24	133.73	123.60
1	L	94	TRP	CA-C-O	-7.08	105.24	120.10
1	L	45	ARG	CD-NE-CZ	7.05	133.47	123.60
2	H	71	ARG	NE-CZ-NH2	6.95	123.78	120.30
1	L	8	PRO	CA-N-CD	-6.94	101.78	111.50
3	Y	121	GLN	CB-CG-CD	6.86	129.44	111.60
3	Y	114	ARG	NE-CZ-NH1	-6.75	116.92	120.30
3	Y	114	ARG	NH1-CZ-NH2	6.62	126.68	119.40
3	Y	18	ASP	CB-CG-OD2	-6.57	112.39	118.30
1	L	1	ASP	OD1-CG-OD2	6.55	135.75	123.30
2	H	33	TYR	CB-CG-CD2	-6.46	117.12	121.00
3	Y	128	ARG	NE-CZ-NH2	6.46	123.53	120.30
1	L	54	SER	N-CA-CB	-6.35	100.97	110.50
3	Y	23	TYR	CB-CG-CD2	6.34	124.80	121.00
1	L	56	SER	CB-CA-C	-6.31	98.11	110.10
3	Y	21	ARG	NE-CZ-NH1	6.25	123.42	120.30
1	L	70	ASP	CB-CG-OD2	-6.18	112.73	118.30
2	H	111	VAL	O-C-N	6.18	132.59	122.70
3	Y	2	VAL	N-CA-CB	6.17	125.08	111.50
2	H	101	ASP	CB-CG-OD1	6.14	123.83	118.30
3	Y	93	ASN	CB-CG-OD1	6.13	133.86	121.60
3	Y	41	GLN	CG-CD-OE1	-5.94	109.72	121.60
1	L	96	TYR	CB-CG-CD2	-5.93	117.44	121.00
1	L	81	GLU	CA-CB-CG	5.89	126.36	113.40
2	H	81	ASP	OD1-CG-OD2	-5.87	112.14	123.30
2	H	51	VAL	CG1-CB-CG2	5.86	120.28	110.90
2	H	45	LEU	CA-CB-CG	5.83	128.71	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	70	ASP	CB-CG-OD1	-5.77	113.10	118.30
3	Y	66	ASP	OD1-CG-OD2	-5.77	112.34	123.30
1	L	69	THR	OG1-CB-CG2	5.77	123.26	110.00
1	L	60	SER	CB-CA-C	-5.76	99.15	110.10
3	Y	45	ARG	CB-CG-CD	5.76	126.58	111.60
2	H	73	THR	N-CA-CB	-5.67	99.53	110.30
2	H	78	TYR	CB-CG-CD1	5.66	124.40	121.00
1	L	54	SER	CA-CB-OG	-5.65	95.94	111.20
1	L	43	SER	N-CA-CB	-5.62	102.07	110.50
1	L	70	ASP	OD1-CG-OD2	5.60	133.95	123.30
1	L	81	GLU	OE1-CD-OE2	-5.60	116.58	123.30
1	L	80	THR	N-CA-CB	5.59	120.92	110.30
3	Y	118	THR	N-CA-CB	-5.56	99.74	110.30
3	Y	125	ARG	NH1-CZ-NH2	-5.55	113.30	119.40
2	H	35	SER	N-CA-CB	-5.52	102.21	110.50
3	Y	14	ARG	NE-CZ-NH2	-5.52	117.54	120.30
1	L	8	PRO	N-CA-CB	5.51	109.91	103.30
2	H	11	LEU	CA-CB-CG	5.44	127.82	115.30
2	H	95	CYS	N-CA-CB	5.43	120.37	110.60
1	L	21	LEU	CB-CG-CD2	5.41	120.20	111.00
3	Y	45	ARG	CB-CA-C	5.41	121.23	110.40
3	Y	109	VAL	CA-CB-CG1	5.38	118.97	110.90
2	H	76	ASN	O-C-N	-5.38	114.10	122.70
3	Y	109	VAL	CA-CB-CG2	-5.38	102.84	110.90
2	H	41	PRO	O-C-N	-5.34	114.11	123.20
2	H	45	LEU	CB-CG-CD1	5.34	120.07	111.00
1	L	54	SER	O-C-N	-5.31	114.20	122.70
3	Y	38	PHE	CB-CG-CD1	5.29	124.50	120.80
3	Y	48	ASP	CA-C-N	5.26	126.71	116.20
2	H	27	ASP	CB-CG-OD1	5.25	123.02	118.30
2	H	64	LYS	CA-CB-CG	-5.24	101.86	113.40
3	Y	5	ARG	NE-CZ-NH1	5.20	122.90	120.30
1	L	58	ILE	CG1-CB-CG2	-5.20	99.95	111.40
3	Y	5	ARG	CD-NE-CZ	5.19	130.87	123.60
2	H	10	SER	N-CA-CB	-5.18	102.73	110.50
2	H	3	GLN	O-C-N	-5.18	114.42	122.70
1	L	58	ILE	CA-CB-CG1	-5.17	101.19	111.00
3	Y	114	ARG	CB-CG-CD	-5.13	98.25	111.60
1	L	70	ASP	CA-CB-CG	-5.11	102.17	113.40
3	Y	61	ARG	O-C-N	-5.10	114.54	122.70
3	Y	87	ASP	CB-CG-OD2	-5.08	113.73	118.30
3	Y	20	TYR	CB-CG-CD1	-5.07	117.96	121.00

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	L	33	LEU	Mainchain
1	L	7	SER	Mainchain,Peptide
1	L	94	TRP	Mainchain,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	L	818	0	784	16	0
2	H	899	0	846	13	0
3	Y	1001	0	959	26	0
4	H	37	0	0	1	0
4	L	35	0	0	0	0
4	Y	44	0	0	2	0
All	All	2834	0	2589	54	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 10.

All (54) 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)
3:Y:61:ARG:HH12	3:Y:70:PRO:HB2	1.25	1.02
3:Y:74:ASN:HD21	3:Y:78:ILE:H	1.14	0.94
1:L:12:SER:HB2	1:L:107:LYS:HB2	1.61	0.81
3:Y:59:ASN:ND2	3:Y:61:ARG:H	1.85	0.75
3:Y:78:ILE:HG13	3:Y:79:PRO:HD2	1.73	0.71
2:H:42:GLY:O	2:H:43:ASN:HB2	1.91	0.69
3:Y:109:VAL:HG22	3:Y:112:ARG:NH2	2.08	0.69
3:Y:45:ARG:HD3	3:Y:68:ARG:HH22	1.62	0.65
1:L:38:GLN:OE1	2:H:39:LYS:HE3	1.98	0.63
3:Y:61:ARG:NH1	3:Y:70:PRO:HB2	2.08	0.60
3:Y:45:ARG:HG3	3:Y:51:THR:OG1	2.02	0.59
3:Y:74:ASN:HD21	3:Y:78:ILE:N	1.93	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:Y:118:THR:CG2	4:Y:134:HOH:O	2.51	0.58
2:H:34:TRP:HB3	2:H:78:TYR:CZ	2.38	0.57
3:Y:66:ASP:OD1	3:Y:69:THR:OG1	2.21	0.57
3:Y:118:THR:HG21	4:Y:134:HOH:O	2.03	0.57
2:H:13:LYS:HE3	2:H:114:ALA:H	1.69	0.57
1:L:12:SER:HA	1:L:105:GLU:O	2.09	0.53
3:Y:74:ASN:ND2	3:Y:78:ILE:H	1.96	0.53
3:Y:45:ARG:HD3	3:Y:68:ARG:NH2	2.23	0.53
1:L:17:ASN:O	1:L:78:VAL:HG13	2.10	0.51
3:Y:115:CYS:O	3:Y:118:THR:HB	2.10	0.51
2:H:64:LYS:O	2:H:65:SER:HB2	2.11	0.51
2:H:51:VAL:HG22	2:H:69:ILE:HG21	1.92	0.51
1:L:94:TRP:CD2	1:L:95:PRO:HA	2.47	0.50
2:H:51:VAL:HG22	2:H:69:ILE:CG2	2.43	0.48
1:L:7:SER:OG	1:L:8:PRO:N	2.45	0.48
2:H:60:ASN:HB3	2:H:63:LEU:HD22	1.95	0.48
1:L:13:VAL:HG21	1:L:19:VAL:CG1	2.44	0.48
2:H:72:ASP:OD2	2:H:75:LYS:HE3	2.13	0.48
3:Y:66:ASP:O	3:Y:68:ARG:N	2.46	0.48
1:L:24:ARG:HG2	1:L:25:ALA:N	2.28	0.47
1:L:76:ASN:HD22	1:L:77:SER:N	2.13	0.47
3:Y:68:ARG:HB2	3:Y:68:ARG:HH11	1.80	0.46
1:L:85:MET:HG3	4:H:137:HOH:O	2.15	0.46
3:Y:14:ARG:HD2	3:Y:14:ARG:HH11	1.58	0.46
3:Y:62:TRP:HZ3	3:Y:73:ARG:NH1	2.13	0.46
1:L:14:THR:HG22	1:L:15:PRO:O	2.17	0.44
3:Y:53:TYR:CD2	3:Y:80:CYS:HB3	2.53	0.43
1:L:13:VAL:HG11	1:L:78:VAL:HG21	1.99	0.43
3:Y:48:ASP:O	3:Y:61:ARG:NH2	2.51	0.43
3:Y:68:ARG:CB	3:Y:68:ARG:HH11	2.32	0.43
1:L:35:TRP:CE2	1:L:73:LEU:HB2	2.55	0.42
3:Y:21:ARG:HA	3:Y:21:ARG:HD2	1.84	0.42
3:Y:59:ASN:ND2	3:Y:61:ARG:N	2.62	0.41
3:Y:63:TRP:CE3	3:Y:75:LEU:HB2	2.55	0.41
1:L:11:LEU:HD23	1:L:11:LEU:HA	1.91	0.41
3:Y:56:LEU:HD23	3:Y:56:LEU:HA	1.97	0.41
2:H:64:LYS:O	2:H:65:SER:CB	2.69	0.41
2:H:13:LYS:O	2:H:16:GLN:HG2	2.19	0.41
1:L:76:ASN:HD22	1:L:76:ASN:C	2.25	0.40
2:H:29:ILE:HD13	2:H:29:ILE:HG21	1.87	0.40
1:L:14:THR:O	1:L:15:PRO:C	2.58	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:21:THR:HB	2:H:79:TYR:CD1	2.57	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	L	105/107 (98%)	100 (95%)	4 (4%)	1 (1%)	19	21
2	H	112/114 (98%)	110 (98%)	2 (2%)	0	100	100
3	Y	127/129 (98%)	120 (94%)	4 (3%)	3 (2%)	7	5
All	All	344/350 (98%)	330 (96%)	10 (3%)	4 (1%)	16	16

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	Y	70	PRO
3	Y	48	ASP
3	Y	67	GLY
1	L	77	SER

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	L	94/94 (100%)	79 (84%)	15 (16%)	3	3
2	H	101/101 (100%)	93 (92%)	8 (8%)	15	19
3	Y	105/105 (100%)	95 (90%)	10 (10%)	11	12
All	All	300/300 (100%)	267 (89%)	33 (11%)	8	8

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	7	SER
1	L	11	LEU
1	L	13	VAL
1	L	17	ASN
1	L	21	LEU
1	L	24	ARG
1	L	58	ILE
1	L	69	THR
1	L	70	ASP
1	L	74	SER
1	L	76	ASN
1	L	77	SER
1	L	78	VAL
1	L	80	THR
1	L	89	GLN
2	H	11	LEU
2	H	13	LYS
2	H	21	THR
2	H	45	LEU
2	H	63	LEU
2	H	65	SER
2	H	73	THR
2	H	88	GLU
3	Y	45	ARG
3	Y	55	ILE
3	Y	68	ARG
3	Y	70	PRO
3	Y	74	ASN
3	Y	81	SER
3	Y	99	VAL
3	Y	112	ARG
3	Y	118	THR
3	Y	121	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	17	ASN
1	L	76	ASN
1	L	89	GLN
2	H	5	GLN
2	H	77	GLN
3	Y	59	ASN
3	Y	74	ASN

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](#)

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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.