



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:23 PM UTC

PDB ID : 1FDL / pdb_00001fdl
Title : CRYSTALLOGRAPHIC REFINEMENT OF THE THREE-DIMENSIONAL
STRUCTURE OF THE FAB D1.3-LYSOZYME COMPLEX AT 2.5-
ANGSTROMS RESOLUTION
Authors : Fischmann, T.O.; Poljak, R.J.
Deposited on : 1990-08-27
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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

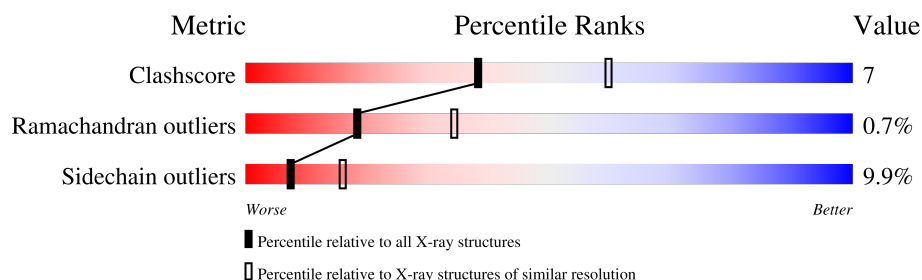
1 Overall quality at a glance

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	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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 $\leq 5\%$

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	L	214	
2	H	218	
3	Y	129	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4309 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 IGG1-KAPPA D1.3 FAB (LIGHT CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	214	Total	C	N	O	S	0	0	0
			1665	1037	282	339	7			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
L	50	TYR	ASN	conflict	GB 2072141
L	51	THR	ALA	conflict	GB 2072141
L	52	THR	LYS	conflict	GB 2072141
L	85	SER	THR	conflict	GB 2072141
L	89	GLN	HIS	conflict	GB 2072141
L	96	ARG	TRP	conflict	GB 2072141
L	106	ILE	VAL	conflict	GB 2072141
L	118	PHE	LEU	conflict	GB 2072141

- Molecule 2 is a protein called IGG1-KAPPA D1.3 FAB (HEAVY CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	218	Total	C	N	O	S	0	0	0
			1643	1030	279	326	8			

- Molecule 3 is a protein called HEN EGG WHITE LYSOZYME.

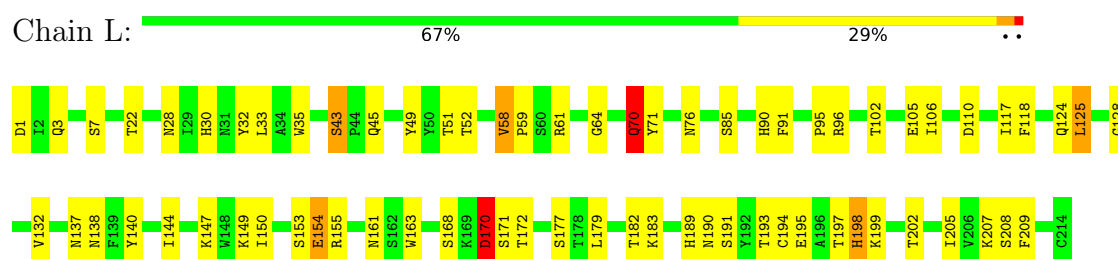
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	Y	129	Total	C	N	O	S	0	0	0
			1001	613	193	185	10			

3 Residue-property plots

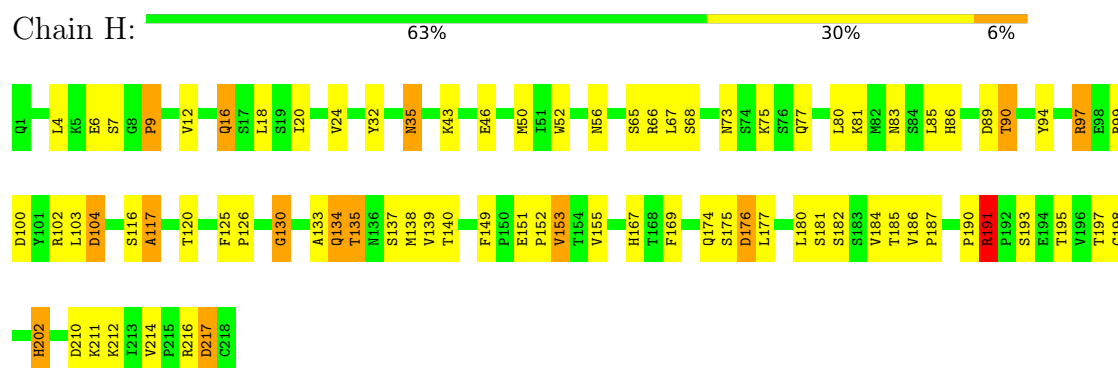
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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.

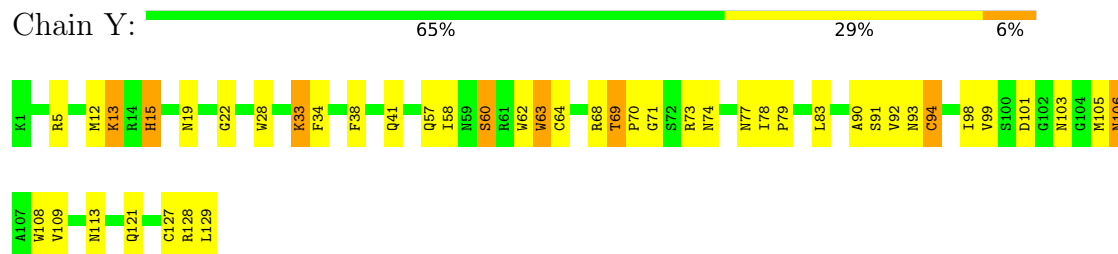
• Molecule 1: IGG1-KAPPA D1.3 FAB (LIGHT CHAIN)



• Molecule 2: IGG1-KAPPA D1.3 FAB (HEAVY CHAIN)



• Molecule 3: HEN EGG WHITE LYSOZYME



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	56.00Å 143.50Å 49.30Å 90.00° 120.40° 90.00°	Depositor
Resolution (Å)	(Not available) – 2.50	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-2.50)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.184 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4309	wwPDB-VP
Average B, all atoms (Å ²)	44.0	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	0.98	5/1705 (0.3%)	1.79	31/2315 (1.3%)
2	H	1.11	4/1684 (0.2%)	2.01	47/2300 (2.0%)
3	Y	1.11	1/1021 (0.1%)	1.96	28/1379 (2.0%)
All	All	1.06	10/4410 (0.2%)	1.91	106/5994 (1.8%)

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	202	HIS	CD2-NE2	-7.04	1.30	1.37
1	L	198	HIS	CD2-NE2	-7.03	1.30	1.37
1	L	189	HIS	CD2-NE2	-6.98	1.30	1.37
2	H	9	PRO	N-CA	-6.96	1.42	1.47
1	L	90	HIS	CD2-NE2	-6.82	1.30	1.37
1	L	30	HIS	CD2-NE2	-6.81	1.30	1.37
2	H	167	HIS	CD2-NE2	-6.69	1.30	1.37
3	Y	15	HIS	CD2-NE2	-6.51	1.30	1.37
2	H	86	HIS	CD2-NE2	-5.84	1.31	1.37
1	L	198	HIS	CG-ND1	-5.54	1.32	1.38

All (106) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	170	ASP	CA-CB-CG	10.12	122.72	112.60
1	L	209	PHE	CA-CB-CG	9.81	123.61	113.80
2	H	66	ARG	NE-CZ-NH2	-9.66	110.50	119.20
2	H	176	ASP	CA-CB-CG	9.12	121.72	112.60
1	L	138	ASN	N-CA-C	8.41	122.74	111.30
1	L	137	ASN	OD1-CG-ND2	-8.26	114.34	122.60
1	L	140	TYR	O-C-N	-8.14	115.49	121.84
2	H	104	ASP	CA-CB-CG	8.08	120.68	112.60
2	H	130	GLY	N-CA-C	-7.90	102.85	111.93
2	H	217	ASP	N-CA-C	7.89	122.30	109.59
2	H	90	THR	N-CA-C	-7.88	95.88	108.73

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	Y	108	TRP	CG-CD2-CE3	7.74	141.64	133.90
1	L	171	SER	N-CA-C	7.73	121.81	111.30
3	Y	101	ASP	CA-CB-CG	7.56	120.16	112.60
2	H	134	GLN	OE1-CD-NE2	-7.55	115.05	122.60
2	H	56	ASN	OD1-CG-ND2	-7.55	115.05	122.60
2	H	9	PRO	CB-CA-C	7.53	118.21	109.92
2	H	167	HIS	CB-CG-CD2	-7.48	121.47	131.20
3	Y	69	THR	N-CA-C	7.46	121.62	109.15
3	Y	108	TRP	CB-CG-CD1	-7.43	115.76	126.90
3	Y	34	PHE	CA-CB-CG	-7.39	106.41	113.80
2	H	217	ASP	CA-CB-CG	7.34	119.94	112.60
3	Y	93	ASN	OD1-CG-ND2	-7.21	115.39	122.60
1	L	30	HIS	CB-CG-CD2	-7.17	121.88	131.20
2	H	77	GLN	N-CA-C	6.98	119.79	108.41
3	Y	121	GLN	OE1-CD-NE2	-6.92	115.68	122.60
2	H	83	ASN	OD1-CG-ND2	-6.76	115.84	122.60
3	Y	109	VAL	N-CA-C	6.74	117.50	110.62
3	Y	103	ASN	N-CA-C	6.63	121.04	113.15
1	L	194	CYS	N-CA-C	-6.62	97.48	108.34
1	L	189	HIS	CB-CG-CD2	-6.57	122.66	131.20
2	H	135	THR	N-CA-C	-6.46	100.69	109.54
1	L	35	TRP	CG-CD2-CE3	6.45	140.35	133.90
1	L	70	GLN	OE1-CD-NE2	-6.45	116.15	122.60
3	Y	12	MET	N-CA-C	-6.34	104.36	111.28
2	H	117	ALA	N-CA-C	6.27	118.88	110.35
1	L	7	SER	O-C-N	-6.10	117.58	121.85
3	Y	33	LYS	CA-C-O	-6.08	114.44	120.82
2	H	83	ASN	CA-CB-CG	6.07	118.67	112.60
2	H	202	HIS	CB-CG-CD2	-6.03	123.36	131.20
2	H	56	ASN	CA-CB-CG	6.03	118.63	112.60
2	H	202	HIS	CA-C-N	6.00	125.49	119.24
2	H	202	HIS	C-N-CA	6.00	125.49	119.24
1	L	171	SER	N-CA-CB	-6.00	103.04	111.62
3	Y	113	ASN	CB-CG-ND2	5.99	125.38	116.40
3	Y	113	ASN	OD1-CG-ND2	-5.98	116.62	122.60
3	Y	74	ASN	OD1-CG-ND2	-5.93	116.67	122.60
1	L	58	VAL	CA-C-N	5.88	125.78	119.85
1	L	58	VAL	C-N-CA	5.88	125.78	119.85
1	L	28	ASN	OD1-CG-ND2	-5.81	116.79	122.60
3	Y	74	ASN	N-CA-C	5.74	119.10	111.30
2	H	56	ASN	CB-CG-ND2	5.73	125.00	116.40
2	H	16	GLN	OE1-CD-NE2	-5.73	116.87	122.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	90	HIS	CA-CB-CG	5.72	119.52	113.80
2	H	86	HIS	CB-CG-CD2	-5.69	123.80	131.20
1	L	35	TRP	CB-CG-CD1	-5.69	118.37	126.90
1	L	161	ASN	OD1-CG-ND2	-5.68	116.92	122.60
2	H	191	ARG	NE-CZ-NH1	5.66	127.16	121.50
2	H	174	GLN	O-C-N	-5.64	116.65	120.83
2	H	191	ARG	CA-CB-CG	5.59	125.27	114.10
1	L	28	ASN	N-CA-C	-5.58	101.59	109.96
2	H	73	ASN	N-CA-C	5.55	117.01	111.07
3	Y	15	HIS	CB-CG-CD2	-5.55	123.98	131.20
2	H	197	THR	CA-C-O	-5.53	114.30	120.66
3	Y	60	SER	N-CA-C	5.52	119.11	112.38
1	L	3	GLN	CB-CG-CD	-5.52	103.22	112.60
3	Y	77	ASN	CA-CB-CG	5.50	118.10	112.60
2	H	90	THR	O-C-N	5.47	129.46	123.22
2	H	169	PHE	CA-CB-CG	-5.47	108.33	113.80
2	H	20	ILE	O-C-N	-5.46	117.45	123.18
1	L	43	SER	CA-C-O	-5.45	114.50	120.17
3	Y	128	ARG	N-CA-C	-5.45	100.30	108.96
3	Y	57	GLN	N-CA-C	5.43	118.89	111.39
1	L	153	SER	CA-CB-OG	5.43	121.96	111.10
3	Y	68	ARG	CA-CB-CG	5.43	124.96	114.10
3	Y	108	TRP	CE2-CD2-CG	-5.41	100.70	107.20
3	Y	90	ALA	N-CA-C	5.39	119.12	112.54
1	L	132	VAL	O-C-N	-5.38	116.81	122.95
1	L	118	PHE	CA-C-N	5.36	125.90	120.38
1	L	118	PHE	C-N-CA	5.36	125.90	120.38
1	L	163	TRP	CE2-CD2-CG	-5.36	100.77	107.20
1	L	30	HIS	CB-CG-ND1	5.32	130.69	122.70
2	H	6	GLU	CA-CB-CG	5.25	124.59	114.10
1	L	205	ILE	N-CA-C	-5.24	101.03	108.58
2	H	176	ASP	N-CA-CB	5.24	119.35	110.49
2	H	94	TYR	CB-CA-C	-5.24	99.78	109.37
2	H	16	GLN	CB-CG-CD	5.24	121.50	112.60
3	Y	28	TRP	CE2-CD2-CG	-5.21	100.95	107.20
2	H	176	ASP	CB-CA-C	-5.20	100.07	110.42
2	H	100	ASP	N-CA-C	5.19	117.51	111.02
2	H	52	TRP	CG-CD2-CE3	5.18	139.08	133.90
2	H	97	ARG	NE-CZ-NH2	-5.18	114.53	119.20
2	H	52	TRP	CE2-CD2-CG	-5.18	100.99	107.20
2	H	167	HIS	CB-CG-ND1	5.18	130.47	122.70
2	H	212	LYS	O-C-N	-5.18	116.98	123.04

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	Y	13	LYS	CA-C-O	-5.17	115.39	120.82
1	L	106	ILE	CB-CA-C	-5.16	105.66	111.23
2	H	77	GLN	CA-CB-CG	-5.16	103.79	114.10
2	H	52	TRP	CB-CG-CD1	-5.14	119.19	126.90
3	Y	63	TRP	N-CA-C	5.14	119.68	113.41
2	H	185	THR	CA-CB-OG1	-5.12	101.91	109.60
3	Y	105	MET	CA-CB-CG	-5.11	103.88	114.10
1	L	163	TRP	CD1-CG-CD2	5.10	114.47	106.30
2	H	120	THR	CA-C-N	5.08	125.62	120.38
2	H	120	THR	C-N-CA	5.08	125.62	120.38
3	Y	94	CYS	N-CA-C	-5.00	105.93	112.23

There are no chirality outliers.

There are no planarity outliers.

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	1665	0	1583	20	0
2	H	1643	0	1611	26	0
3	Y	1001	0	959	15	0
All	All	4309	0	4153	58	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 7.

All (58) 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)
2:H:191:ARG:HH11	2:H:191:ARG:HG2	1.31	0.94
2:H:191:ARG:HG2	2:H:191:ARG:NH1	1.97	0.75
1:L:61:ARG:HB2	1:L:76:ASN:O	1.95	0.67
1:L:144:ILE:HG13	1:L:198:HIS:HB2	1.78	0.66
1:L:155:ARG:HH11	1:L:179:LEU:HD11	1.62	0.63
2:H:130:GLY:HA2	2:H:216:ARG:HD2	1.82	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:117:ALA:HB2	2:H:176:ASP:HB3	1.86	0.56
3:Y:15:HIS:HB3	3:Y:92:VAL:HG11	1.87	0.56
1:L:32:TYR:HB3	1:L:91:PHE:HB2	1.87	0.56
2:H:139:VAL:HG23	2:H:186:VAL:HG23	1.86	0.55
1:L:96:ARG:NH1	2:H:50:MET:SD	2.80	0.54
2:H:35:ASN:HD22	2:H:35:ASN:N	2.06	0.54
1:L:147:LYS:HG2	1:L:195:GLU:HB3	1.88	0.54
1:L:124:GLN:HG3	2:H:125:PHE:CE1	2.47	0.50
1:L:150:ILE:HD11	1:L:179:LEU:HD21	1.93	0.50
3:Y:19:ASN:OD1	3:Y:22:GLY:HA2	2.13	0.49
1:L:150:ILE:HD12	1:L:155:ARG:HD2	1.95	0.48
3:Y:33:LYS:HB2	3:Y:38:PHE:CE2	2.49	0.48
1:L:85:SER:HA	1:L:102:THR:O	2.13	0.48
1:L:149:LYS:HG2	1:L:154:GLU:HA	1.96	0.48
1:L:51:THR:HG23	1:L:71:TYR:HD2	1.79	0.47
2:H:155:VAL:HG11	2:H:182:SER:OG	2.14	0.47
2:H:18:LEU:O	2:H:81:LYS:HA	2.14	0.47
2:H:198:CYS:O	2:H:210:ASP:HA	2.14	0.47
2:H:138:MET:HA	2:H:187:PRO:HA	1.96	0.47
2:H:180:LEU:HD23	2:H:181:SER:N	2.30	0.47
3:Y:13:LYS:HD3	3:Y:129:LEU:HD22	1.98	0.46
3:Y:62:TRP:HB2	3:Y:63:TRP:CD1	2.49	0.46
3:Y:94:CYS:O	3:Y:98:ILE:HG13	2.16	0.46
3:Y:71:GLY:O	3:Y:73:ARG:HD2	2.16	0.46
3:Y:33:LYS:HB2	3:Y:38:PHE:CZ	2.51	0.46
2:H:97:ARG:O	2:H:103:LEU:HA	2.16	0.45
1:L:125:LEU:O	1:L:183:LYS:HD3	2.15	0.45
3:Y:69:THR:HA	3:Y:70:PRO:HD3	1.76	0.45
2:H:67:LEU:CD2	2:H:80:LEU:HD11	2.47	0.45
2:H:35:ASN:CG	2:H:103:LEU:HD21	2.43	0.44
2:H:140:THR:HA	2:H:184:VAL:O	2.18	0.44
2:H:126:PRO:HD3	2:H:211:LYS:HD2	1.99	0.44
3:Y:60:SER:HA	3:Y:64:CYS:SG	2.57	0.44
1:L:22:THR:HG23	1:L:70:GLN:NE2	2.32	0.44
2:H:32:TYR:HE2	2:H:99:ARG:HG3	1.83	0.44
2:H:149:PHE:HB2	2:H:177:LEU:HD23	2.00	0.44
2:H:67:LEU:HD22	2:H:80:LEU:HD11	2.00	0.44
2:H:85:LEU:HD23	2:H:89:ASP:OD1	2.18	0.43
3:Y:106:ASN:HD22	3:Y:106:ASN:H	1.64	0.43
3:Y:78:ILE:HD12	3:Y:78:ILE:HA	1.87	0.43
3:Y:78:ILE:HA	3:Y:79:PRO:HD2	1.79	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:58:VAL:HA	1:L:59:PRO:HD3	1.79	0.42
3:Y:5:ARG:NH1	3:Y:127:CYS:SG	2.92	0.42
1:L:33:LEU:HD22	1:L:71:TYR:CB	2.51	0.41
1:L:52:THR:HG22	1:L:64:GLY:O	2.20	0.41
2:H:4:LEU:HG	2:H:24:VAL:HG12	2.01	0.41
3:Y:58:ILE:HB	3:Y:83:LEU:HD13	2.02	0.41
1:L:49:TYR:CG	2:H:102:ARG:HD2	2.56	0.41
1:L:117:ILE:HG22	1:L:207:LYS:HG3	2.03	0.40
2:H:153:VAL:HG23	2:H:202:HIS:HD2	1.86	0.40
2:H:97:ARG:HH21	2:H:104:ASP:CG	2.28	0.40
1:L:128:GLY:HA2	1:L:183:LYS:HB2	2.02	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	212/214 (99%)	199 (94%)	11 (5%)	2 (1%)	14	27
2	H	216/218 (99%)	201 (93%)	13 (6%)	2 (1%)	14	27
3	Y	127/129 (98%)	115 (91%)	12 (9%)	0	100	100
All	All	555/561 (99%)	515 (93%)	36 (6%)	4 (1%)	18	34

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	170	ASP
1	L	199	LYS
2	H	133	ALA
2	H	137	SER

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	L	189/189 (100%)	169 (89%)	20 (11%)	6	14
2	H	190/190 (100%)	166 (87%)	24 (13%)	4	9
3	Y	105/105 (100%)	101 (96%)	4 (4%)	29	56
All	All	484/484 (100%)	436 (90%)	48 (10%)	7	16

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	1	ASP
1	L	43	SER
1	L	45	GLN
1	L	70	GLN
1	L	95	PRO
1	L	105	GLU
1	L	110	ASP
1	L	125	LEU
1	L	154	GLU
1	L	168	SER
1	L	170	ASP
1	L	172	THR
1	L	177	SER
1	L	182	THR
1	L	190	ASN
1	L	191	SER
1	L	193	THR
1	L	197	THR
1	L	202	THR
1	L	208	SER
2	H	7	SER
2	H	9	PRO
2	H	12	VAL
2	H	16	GLN
2	H	35	ASN
2	H	43	LYS

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Mol	Chain	Res	Type
2	H	46	GLU
2	H	65	SER
2	H	68	SER
2	H	75	LYS
2	H	90	THR
2	H	116	SER
2	H	134	GLN
2	H	135	THR
2	H	151	GLU
2	H	152	PRO
2	H	153	VAL
2	H	175	SER
2	H	190	PRO
2	H	191	ARG
2	H	193	SER
2	H	195	THR
2	H	214	VAL
2	H	217	ASP
3	Y	41	GLN
3	Y	91	SER
3	Y	99	VAL
3	Y	106	ASN

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

Mol	Chain	Res	Type
1	L	30	HIS
1	L	76	ASN
1	L	189	HIS
3	Y	44	ASN
3	Y	57	GLN
3	Y	65	ASN
3	Y	106	ASN

5.3.3 RNA ⓘ

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 oligosaccharides 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 is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.