



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 06:57 PM GMT

PDB ID : 1DBB
Title : THREE-DIMENSIONAL STRUCTURE OF AN ANTI-STEROID FAB' AND
PROGESTERONE-FAB' COMPLEX
Authors : Arevalo, J.H.; Wilson, I.A.
Deposited on : 1992-11-11
Resolution : 2.70 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

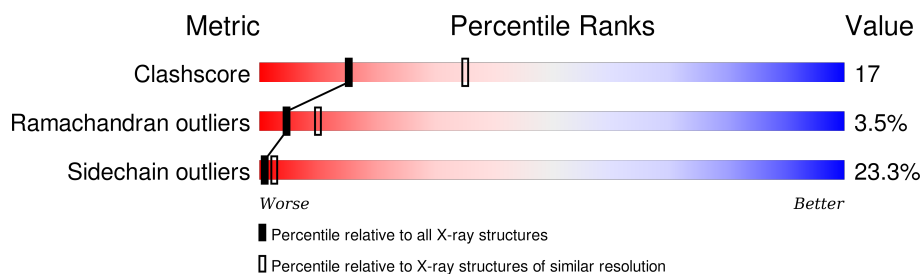
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)

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	216	
2	H	219	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3377 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 DB3 FAB (LIGHT CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	216	Total	C	N	O	S	0	0	0
			1679	1051	286	335	7			

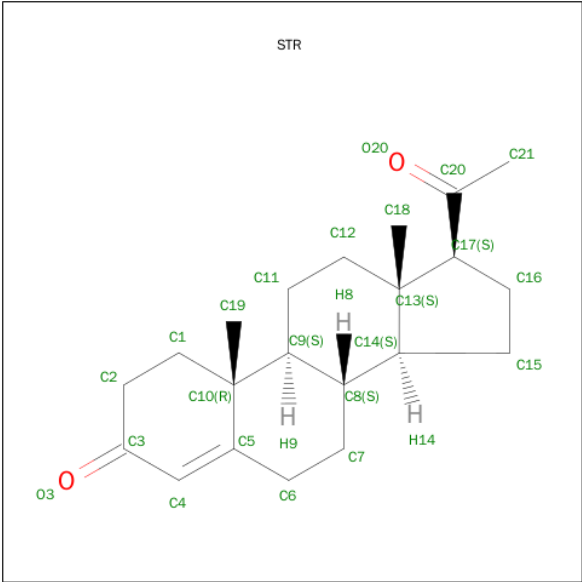
There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
L	2	VAL	ILE	CONFLICT	GB 1589925
L	7	ILE	SER	CONFLICT	GB 1589925
L	14	ASN	SER	CONFLICT	GB 1589925
L	27B	LEU	VAL	CONFLICT	GB 1589925
L	27C	ILE	VAL	CONFLICT	GB 1589925
L	34	HIS	GLU	CONFLICT	GB 1589925
L	36	TYR	PHE	CONFLICT	GB 1589925
L	48	MET	ILE	CONFLICT	GB 1589925
L	56	TYR	SER	CONFLICT	GB 1589925
L	85	ILE	VAL	CONFLICT	GB 1589925
L	87	PHE	TYR	CONFLICT	GB 1589925
L	89	SER	PHE	CONFLICT	GB 1589925
L	91	SER	ALA	CONFLICT	GB 1589925
L	96	PRO	TRP	CONFLICT	GB 1589925

- Molecule 2 is a protein called IGG1-KAPPA DB3 FAB (HEAVY CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	219	Total	C	N	O	S	0	0	0
			1675	1071	270	328	6			

- Molecule 3 is PROGESTERONE (three-letter code: STR) (formula: C₂₁H₃₀O₂).



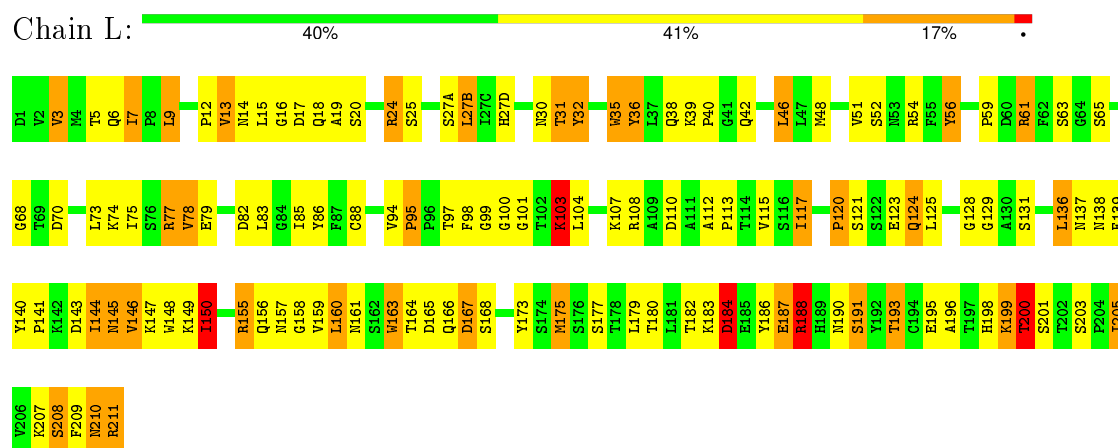
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	H	1	Total	C	O	0	0
			23	21	2		

3 Residue-property plots

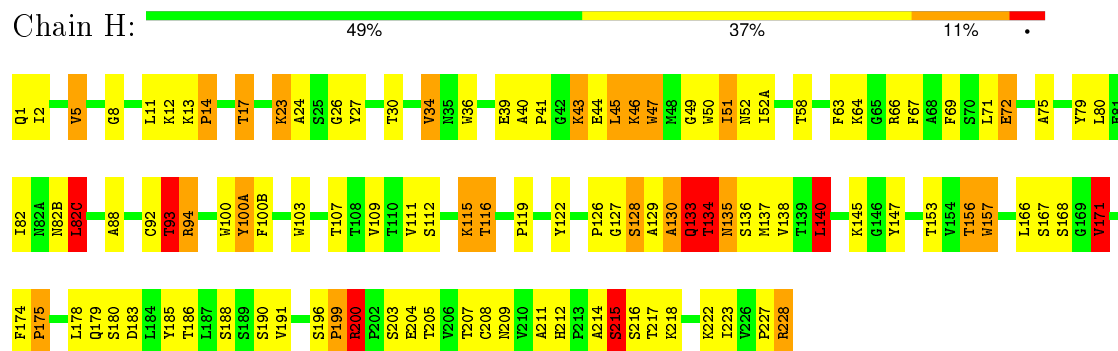
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: IGG1-KAPPA DB3 FAB (LIGHT CHAIN)



• Molecule 2: IGG1-KAPPA DB3 FAB (HEAVY CHAIN)



4 Data and refinement statistics

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

Property	Value	Source
Space group	P 64 2 2	Depositor
Cell constants a, b, c, α , β , γ	134.76Å 134.76Å 124.21Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	(Not available) – 2.70	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-2.70)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.210 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3377	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: STR

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	0/1719	1.89	40/2331 (1.7%)
2	H	1.03	0/1723	1.90	52/2358 (2.2%)
All	All	1.01	0/3442	1.90	92/4689 (2.0%)

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	L	0	1
2	H	0	3
All	All	0	4

There are no bond length outliers.

All (92) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	200	ARG	NE-CZ-NH1	12.56	126.58	120.30
2	H	100(A)	TYR	CB-CG-CD2	-12.19	113.69	121.00
1	L	155	ARG	NE-CZ-NH2	-11.87	114.37	120.30
1	L	211	ARG	NE-CZ-NH1	11.73	126.16	120.30
1	L	155	ARG	NE-CZ-NH1	10.60	125.60	120.30
2	H	66	ARG	NE-CZ-NH1	10.18	125.39	120.30
1	L	35	TRP	CD1-CG-CD2	8.90	113.42	106.30
2	H	103	TRP	CD1-CG-CD2	8.46	113.07	106.30
1	L	163	TRP	CD1-CG-CD2	8.33	112.96	106.30
2	H	47	TRP	CD1-CG-CD2	8.30	112.94	106.30
1	L	173	TYR	CB-CG-CD2	-8.02	116.19	121.00
2	H	47	TRP	CE2-CD2-CG	-7.87	101.01	107.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	171	VAL	CG1-CB-CG2	-7.79	98.43	110.90
2	H	103	TRP	CE2-CD2-CG	-7.47	101.33	107.30
2	H	100(A)	TYR	CB-CG-CD1	7.44	125.46	121.00
1	L	148	TRP	CD1-CG-CD2	7.39	112.21	106.30
2	H	140	LEU	CA-CB-CG	7.28	132.04	115.30
2	H	199	PRO	CA-C-N	7.27	133.20	117.20
1	L	163	TRP	CE2-CD2-CG	-7.25	101.50	107.30
1	L	211	ARG	NE-CZ-NH2	-7.11	116.75	120.30
1	L	188	ARG	NE-CZ-NH1	7.10	123.85	120.30
1	L	32	TYR	CB-CG-CD1	-7.10	116.74	121.00
2	H	228	ARG	NE-CZ-NH2	-7.07	116.77	120.30
2	H	103	TRP	CG-CD2-CE3	7.06	140.25	133.90
2	H	82(C)	LEU	CA-CB-CG	7.05	131.51	115.30
1	L	35	TRP	CE2-CD2-CG	-7.04	101.67	107.30
2	H	227	PRO	O-C-N	6.97	133.85	122.70
1	L	148	TRP	CE2-CD2-CG	-6.96	101.73	107.30
2	H	45	LEU	N-CA-C	-6.80	92.64	111.00
2	H	93	THR	CA-CB-CG2	-6.58	103.19	112.40
1	L	46	LEU	CA-CB-CG	6.55	130.36	115.30
1	L	36	TYR	CB-CG-CD2	-6.51	117.09	121.00
2	H	157	TRP	CE2-CD2-CE3	6.43	126.42	118.70
1	L	156	GLN	O-C-N	-6.41	112.45	122.70
2	H	157	TRP	CE2-CD2-CG	-6.40	102.18	107.30
1	L	184	ASP	CA-CB-CG	-6.36	99.41	113.40
2	H	128	SER	CA-C-N	-6.25	103.45	117.20
1	L	163	TRP	CG-CD1-NE1	-6.25	103.85	110.10
2	H	47	TRP	CG-CD2-CE3	6.19	139.47	133.90
2	H	94	ARG	NE-CZ-NH2	6.15	123.38	120.30
2	H	228	ARG	N-CA-C	-6.15	94.40	111.00
1	L	94	VAL	CG1-CB-CG2	-6.15	101.07	110.90
1	L	103	LYS	CA-CB-CG	6.13	126.89	113.40
1	L	167	ASP	CB-CG-OD1	6.13	123.81	118.30
2	H	36	TRP	CE2-CD2-CG	-6.10	102.42	107.30
1	L	144	ILE	CB-CA-C	-6.10	99.40	111.60
2	H	47	TRP	CG-CD1-NE1	-6.10	104.00	110.10
2	H	36	TRP	CD1-CG-CD2	6.05	111.14	106.30
2	H	11	LEU	CA-CB-CG	6.05	129.21	115.30
2	H	50	TRP	CE2-CD2-CG	-6.04	102.47	107.30
1	L	35	TRP	CG-CD1-NE1	-6.03	104.07	110.10
2	H	227	PRO	CA-C-N	-5.98	104.05	117.20
1	L	54	ARG	NE-CZ-NH2	-5.93	117.33	120.30
2	H	109	VAL	CA-CB-CG2	-5.85	102.12	110.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	186	TYR	CB-CG-CD1	-5.80	117.52	121.00
2	H	157	TRP	CD1-CG-CD2	5.79	110.93	106.30
1	L	95	PRO	N-CD-CG	-5.68	94.68	103.20
2	H	47	TRP	CB-CG-CD1	-5.65	119.66	127.00
1	L	3	VAL	N-CA-CB	-5.63	99.11	111.50
2	H	50	TRP	NE1-CE2-CZ2	-5.62	124.22	130.40
1	L	188	ARG	NE-CZ-NH2	-5.58	117.51	120.30
1	L	13	VAL	CG1-CB-CG2	-5.58	101.97	110.90
2	H	199	PRO	CA-N-CD	-5.56	103.72	111.50
2	H	100	TRP	CE2-CD2-CG	-5.55	102.86	107.30
2	H	103	TRP	CG-CD1-NE1	-5.47	104.63	110.10
2	H	216	SER	N-CA-CB	-5.47	102.29	110.50
2	H	175	PRO	N-CD-CG	-5.43	95.06	103.20
1	L	48	MET	CA-CB-CG	5.42	122.52	113.30
1	L	110	ASP	CB-CG-OD1	5.42	123.18	118.30
2	H	228	ARG	NE-CZ-NH1	5.41	123.01	120.30
1	L	3	VAL	CA-CB-CG2	-5.40	102.80	110.90
2	H	100	TRP	CD1-CG-CD2	5.36	110.59	106.30
1	L	56	TYR	CB-CG-CD2	-5.35	117.79	121.00
2	H	175	PRO	O-C-N	5.33	131.22	122.70
2	H	185	TYR	CB-CG-CD2	-5.28	117.83	121.00
2	H	128	SER	CA-C-O	5.27	131.17	120.10
2	H	153	THR	N-CA-CB	-5.27	100.29	110.30
2	H	208	CYS	CA-CB-SG	5.25	123.44	114.00
2	H	175	PRO	CA-C-N	-5.21	105.73	117.20
1	L	120	PRO	CA-C-N	-5.18	105.80	117.20
1	L	156	GLN	CA-C-N	5.18	128.60	117.20
1	L	144	ILE	N-CA-CB	5.18	122.70	110.80
2	H	92	CYS	CA-CB-SG	5.15	123.27	114.00
2	H	66	ARG	NE-CZ-NH2	-5.14	117.73	120.30
1	L	205	ILE	CA-C-N	-5.12	105.93	117.20
1	L	9	LEU	CA-CB-CG	5.10	127.03	115.30
1	L	24	ARG	NE-CZ-NH2	-5.07	117.76	120.30
2	H	100(A)	TYR	CA-CB-CG	5.07	123.03	113.40
2	H	103	TRP	CB-CG-CD1	-5.06	120.42	127.00
1	L	136	LEU	N-CA-C	-5.01	97.47	111.00
2	H	228	ARG	CA-CB-CG	5.01	124.42	113.40
2	H	58	THR	CA-CB-CG2	5.00	119.41	112.40

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	H	130	ALA	Peptide
2	H	134	THR	Peptide
2	H	200	ARG	Peptide
1	L	150	ILE	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	1679	0	1625	68	0
2	H	1675	0	1633	53	0
3	H	23	0	29	0	0
All	All	3377	0	3287	116	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 17.

All (116) 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:L:150:ILE:HD12	1:L:155:ARG:HG3	1.30	1.06
2:H:93:THR:HG21	2:H:100(B):PHE:HB3	1.53	0.88
1:L:149:LYS:HB2	1:L:193:THR:HG23	1.61	0.80
2:H:34:VAL:HG13	2:H:51:ILE:HG23	1.66	0.76
1:L:193:THR:HB	1:L:208:SER:OG	1.87	0.74
1:L:196:ALA:HB3	1:L:205:ILE:HG23	1.70	0.74
1:L:12:PRO:HB2	1:L:107:LYS:HB2	1.72	0.71
1:L:13:VAL:HG11	1:L:78:VAL:HG11	1.75	0.69
2:H:130:ALA:O	2:H:134:THR:N	2.26	0.68
1:L:199:LYS:HD3	1:L:199:LYS:H	1.58	0.68
2:H:156:THR:HG23	2:H:209:ASN:HB2	1.77	0.67
1:L:27(B):LEU:O	1:L:31:THR:HA	1.97	0.64
2:H:40:ALA:HB3	2:H:43:LYS:HB2	1.80	0.64
1:L:144:ILE:HG22	1:L:198:HIS:CD2	2.34	0.63
1:L:107:LYS:HA	1:L:140:TYR:OH	2.02	0.60
1:L:123:GLU:HG3	2:H:122:TYR:CD1	2.36	0.60
1:L:128:GLY:HA2	1:L:183:LYS:HE2	1.84	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:136:SER:O	2:H:137:MET:SD	2.60	0.59
1:L:144:ILE:HD11	1:L:175:MET:SD	2.42	0.59
1:L:161:ASN:OD1	1:L:177:SER:HB3	2.03	0.59
1:L:86:TYR:O	1:L:101:GLY:HA2	2.03	0.58
1:L:85:ILE:HD12	1:L:103:LYS:HE3	1.85	0.58
1:L:83:LEU:HD23	1:L:104:LEU:O	2.02	0.58
2:H:47:TRP:HE1	2:H:100(B):PHE:HZ	1.51	0.58
2:H:196:SER:O	2:H:203:SER:HB3	2.04	0.58
1:L:164:THR:HG23	2:H:174:PHE:HD1	1.69	0.57
1:L:155:ARG:NH2	1:L:157:ASN:HB3	2.19	0.57
1:L:144:ILE:HG22	1:L:198:HIS:HB2	1.87	0.57
1:L:63:SER:O	1:L:73:LEU:HD12	2.04	0.57
1:L:129:GLY:H	1:L:183:LYS:H	1.53	0.56
2:H:119:PRO:HB3	2:H:147:TYR:HB3	1.87	0.56
1:L:187:GLU:HA	1:L:211:ARG:NE	2.20	0.56
1:L:187:GLU:HA	1:L:211:ARG:CZ	2.35	0.56
2:H:49:GLY:HA3	2:H:69:PHE:HE2	1.71	0.56
1:L:95:PRO:O	1:L:97:THR:HG23	2.06	0.56
2:H:179:GLN:HE22	2:H:186:THR:HG21	1.71	0.56
2:H:5:VAL:HG13	2:H:23:LYS:HG2	1.87	0.56
1:L:124:GLN:HG2	2:H:122:TYR:CD2	2.42	0.55
2:H:171:VAL:HA	2:H:190:SER:O	2.07	0.55
1:L:35:TRP:CZ3	1:L:88:CYS:HB3	2.41	0.55
1:L:143:ASP:HB2	1:L:199:LYS:NZ	2.23	0.54
2:H:93:THR:CG2	2:H:100(B):PHE:HB3	2.33	0.54
1:L:184:ASP:O	1:L:188:ARG:HG3	2.08	0.54
1:L:112:ALA:HB2	1:L:200:THR:HB	1.90	0.53
2:H:212:HIS:ND1	2:H:215:SER:HB3	2.24	0.53
1:L:157:ASN:OD1	1:L:158:GLY:N	2.43	0.52
2:H:8:GLY:O	2:H:107:THR:HG23	2.10	0.52
2:H:52(A):ILE:HD12	2:H:71:LEU:HD11	1.91	0.52
1:L:199:LYS:HD3	1:L:199:LYS:N	2.24	0.51
2:H:69:PHE:HA	2:H:79:TYR:O	2.11	0.51
1:L:145:ASN:N	1:L:145:ASN:HD22	2.07	0.51
2:H:12:LYS:O	2:H:111:VAL:HA	2.10	0.51
1:L:164:THR:HG23	2:H:174:PHE:CD1	2.45	0.51
2:H:166:LEU:HD23	2:H:191:VAL:HG21	1.93	0.51
2:H:2:ILE:HD12	2:H:26:GLY:O	2.11	0.50
1:L:143:ASP:HB2	1:L:199:LYS:HZ1	1.76	0.50
2:H:24:ALA:HB1	2:H:27:TYR:CE2	2.46	0.50
2:H:207:THR:HG23	2:H:222:LYS:HD2	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:49:GLY:HA3	2:H:69:PHE:CE2	2.46	0.49
2:H:34:VAL:HG12	2:H:52(A):ILE:HD11	1.95	0.49
1:L:27(A):SER:HA	1:L:68:GLY:O	2.13	0.49
1:L:38:GLN:NE2	1:L:42:GLN:O	2.46	0.49
1:L:6:GLN:O	1:L:7:ILE:HD12	2.13	0.49
2:H:39:GLU:O	2:H:88:ALA:HB1	2.13	0.49
1:L:158:GLY:O	1:L:160:LEU:HD23	2.13	0.48
1:L:14:ASN:HB2	1:L:17:ASP:CG	2.33	0.48
2:H:17:THR:HA	2:H:82:ILE:O	2.14	0.48
1:L:190:ASN:HB3	1:L:210:ASN:OD1	2.13	0.48
2:H:211:ALA:HB2	2:H:218:LYS:HD3	1.95	0.48
1:L:83:LEU:H	1:L:83:LEU:HD12	1.78	0.48
1:L:36:TYR:CE1	1:L:46:LEU:HD23	2.48	0.48
1:L:121:SER:O	1:L:125:LEU:HD23	2.14	0.47
2:H:82(C):LEU:HD22	2:H:82(C):LEU:H	1.79	0.47
1:L:98:PHE:HB2	2:H:45:LEU:O	2.15	0.47
2:H:180:SER:O	2:H:183:ASP:HB2	2.14	0.47
1:L:27(D):HIS:N	1:L:30:ASN:O	2.48	0.46
1:L:85:ILE:HG23	1:L:103:LYS:HE3	1.97	0.46
1:L:150:ILE:HD12	1:L:155:ARG:CG	2.23	0.46
2:H:2:ILE:CG2	2:H:27:TYR:HB3	2.45	0.46
2:H:72:GLU:O	2:H:75:ALA:HB3	2.16	0.46
2:H:115:LYS:HD3	2:H:116:THR:H	1.81	0.46
1:L:184:ASP:HA	1:L:187:GLU:H	1.81	0.45
1:L:79:GLU:HB2	1:L:82:ASP:OD1	2.16	0.45
2:H:133:GLN:HE21	2:H:133:GLN:HB2	1.59	0.45
2:H:41:PRO:O	2:H:43:LYS:HD2	2.16	0.44
2:H:130:ALA:O	2:H:134:THR:HG23	2.18	0.44
1:L:39:LYS:HB3	1:L:40:PRO:HD2	1.99	0.43
2:H:82(B):ASN:HD22	2:H:82(B):ASN:N	2.16	0.43
1:L:85:ILE:CD1	1:L:103:LYS:HG3	2.48	0.43
2:H:5:VAL:HG13	2:H:23:LYS:CG	2.47	0.43
1:L:14:ASN:O	1:L:17:ASP:HB2	2.18	0.43
2:H:179:GLN:NE2	2:H:186:THR:HG21	2.33	0.43
2:H:46:LYS:HD3	2:H:47:TRP:N	2.34	0.43
1:L:6:GLN:HG3	1:L:99:GLY:O	2.19	0.43
1:L:61:ARG:O	1:L:75:ILE:HA	2.19	0.43
2:H:14:PRO:HA	2:H:82(C):LEU:O	2.19	0.42
2:H:127:GLY:O	2:H:129:ALA:N	2.52	0.42
1:L:144:ILE:CG2	1:L:198:HIS:CD2	3.02	0.42
2:H:167:SER:O	2:H:171:VAL:HG12	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:16:GLY:O	1:L:77:ARG:NH1	2.53	0.42
2:H:140:LEU:HG	2:H:223:ILE:HG21	2.02	0.42
1:L:199:LYS:O	1:L:201:SER:N	2.53	0.41
2:H:133:GLN:O	2:H:135:ASN:N	2.53	0.41
1:L:120:PRO:HB3	1:L:131:SER:O	2.20	0.41
1:L:144:ILE:HG21	1:L:144:ILE:HD13	1.84	0.41
1:L:117:ILE:HD12	1:L:209:PHE:HD1	1.84	0.41
1:L:149:LYS:N	1:L:193:THR:O	2.54	0.41
1:L:75:ILE:HD12	1:L:75:ILE:N	2.35	0.41
1:L:146:VAL:HA	1:L:195:GLU:O	2.20	0.41
1:L:18:GLN:HG2	1:L:19:ALA:N	2.35	0.41
1:L:113:PRO:HB3	1:L:139:PHE:CD2	2.55	0.41
2:H:63:PHE:O	2:H:67:PHE:HB2	2.21	0.41
1:L:32:TYR:CD1	1:L:32:TYR:N	2.89	0.41
2:H:13:LYS:HD3	2:H:112:SER:O	2.20	0.41
2:H:136:SER:C	2:H:137:MET:SD	3.00	0.40
1:L:61:ARG:HD2	1:L:77:ARG:O	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	214/216 (99%)	170 (79%)	36 (17%)	8 (4%)	4	9
2	H	217/219 (99%)	184 (85%)	26 (12%)	7 (3%)	5	12
All	All	431/435 (99%)	354 (82%)	62 (14%)	15 (4%)	4	10

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	159	VAL

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Mol	Chain	Res	Type
1	L	191	SER
1	L	200	THR
2	H	128	SER
2	H	133	GLN
1	L	100	GLY
1	L	187	GLU
2	H	175	PRO
1	L	167	ASP
2	H	214	ALA
2	H	215	SER
2	H	135	ASN
1	L	138	ASN
2	H	14	PRO
1	L	51	VAL

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	194/194 (100%)	144 (74%)	50 (26%)	0	2
2	H	188/188 (100%)	149 (79%)	39 (21%)	1	4
All	All	382/382 (100%)	293 (77%)	89 (23%)	1	2

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

Mol	Chain	Res	Type
1	L	3	VAL
1	L	5	THR
1	L	7	ILE
1	L	9	LEU
1	L	15	LEU
1	L	20	SER
1	L	24	ARG
1	L	25	SER
1	L	27(B)	LEU

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Mol	Chain	Res	Type
1	L	31	THR
1	L	52	SER
1	L	56	TYR
1	L	59	PRO
1	L	61	ARG
1	L	65	SER
1	L	70	ASP
1	L	74	LYS
1	L	77	ARG
1	L	78	VAL
1	L	103	LYS
1	L	108	ARG
1	L	115	VAL
1	L	117	ILE
1	L	124	GLN
1	L	136	LEU
1	L	137	ASN
1	L	141	PRO
1	L	145	ASN
1	L	146	VAL
1	L	147	LYS
1	L	150	ILE
1	L	160	LEU
1	L	163	TRP
1	L	165	ASP
1	L	166	GLN
1	L	168	SER
1	L	175	MET
1	L	179	LEU
1	L	180	THR
1	L	182	THR
1	L	184	ASP
1	L	188	ARG
1	L	191	SER
1	L	193	THR
1	L	199	LYS
1	L	200	THR
1	L	203	SER
1	L	207	LYS
1	L	208	SER
1	L	210	ASN
2	H	1	GLN

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Mol	Chain	Res	Type
2	H	5	VAL
2	H	17	THR
2	H	23	LYS
2	H	30	THR
2	H	34	VAL
2	H	43	LYS
2	H	44	GLU
2	H	46	LYS
2	H	51	ILE
2	H	52	ASN
2	H	64	LYS
2	H	72	GLU
2	H	80	LEU
2	H	82(C)	LEU
2	H	93	THR
2	H	94	ARG
2	H	100(A)	TYR
2	H	115	LYS
2	H	116	THR
2	H	126	PRO
2	H	133	GLN
2	H	134	THR
2	H	138	VAL
2	H	140	LEU
2	H	145	LYS
2	H	156	THR
2	H	157	TRP
2	H	168	SER
2	H	171	VAL
2	H	178	LEU
2	H	188	SER
2	H	199	PRO
2	H	200	ARG
2	H	204	GLU
2	H	205	THR
2	H	215	SER
2	H	217	THR
2	H	228	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	14	ASN
1	L	34	HIS
1	L	38	GLN
1	L	145	ASN
1	L	166	GLN
2	H	35	ASN
2	H	82(A)	ASN
2	H	82(B)	ASN
2	H	133	GLN

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](#)

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	$\# Z > 2$
3	STR	H	229	-	26,26,26	2.07	11 (42%)	42,42,42	2.43	16 (38%)

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
3	STR	H	229	-	-	0/4/62/62	0/4/4/4

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	H	229	STR	C4-C3	-4.29	1.36	1.45
3	H	229	STR	C7-C6	-2.84	1.46	1.52
3	H	229	STR	C11-C9	-2.69	1.49	1.53
3	H	229	STR	C12-C11	-2.65	1.47	1.53
3	H	229	STR	C13-C14	-2.51	1.49	1.55
3	H	229	STR	C13-C17	-2.40	1.51	1.56
3	H	229	STR	C7-C8	-2.37	1.49	1.53
3	H	229	STR	C10-C9	-2.35	1.51	1.56
3	H	229	STR	C17-C20	-2.05	1.48	1.51
3	H	229	STR	C12-C13	2.56	1.59	1.54
3	H	229	STR	C4-C5	5.34	1.41	1.34

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	H	229	STR	C6-C5-C10	-3.49	112.74	116.71
3	H	229	STR	C1-C2-C3	-3.04	105.47	111.64
3	H	229	STR	C13-C14-C8	-2.78	110.02	114.37
3	H	229	STR	C10-C9-C8	-2.64	108.78	112.67
3	H	229	STR	C15-C14-C13	-2.57	100.48	103.82
3	H	229	STR	C5-C4-C3	-2.55	120.23	123.75
3	H	229	STR	C12-C11-C9	-2.37	109.10	113.10
3	H	229	STR	C7-C8-C14	-2.22	108.34	112.02
3	H	229	STR	C18-C13-C14	-2.16	107.52	111.75
3	H	229	STR	C11-C9-C10	-2.07	110.36	113.11
3	H	229	STR	C21-C20-C17	2.13	120.51	117.53
3	H	229	STR	C11-C12-C13	2.51	117.32	112.84
3	H	229	STR	C19-C10-C9	2.75	115.17	111.67
3	H	229	STR	C13-C17-C20	4.99	121.42	115.13
3	H	229	STR	C17-C13-C14	7.12	107.15	99.74
3	H	229	STR	C15-C14-C8	7.66	131.13	119.03

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

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

6.4 Ligands [i](#)

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

6.5 Other polymers [i](#)

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