Full wwPDB X-ray Structure Validation Report ⓘ

May 26, 2020 – 01:07 pm BST

PDB ID	:	1CFQ
Title	:	ANTI-P24 (HIV-1) FAB FRAGMENT CB41
Authors	:	Keitel, T.; Kramer, A.; Wessner, H.; Scholz, C.; Schneider-Mergener, J.; Hoehne, W.
Deposited on	:	1999-03-19
Resolution	:	2.80 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 following versions of software and data (see references ⓘ) were used in the production of this report:

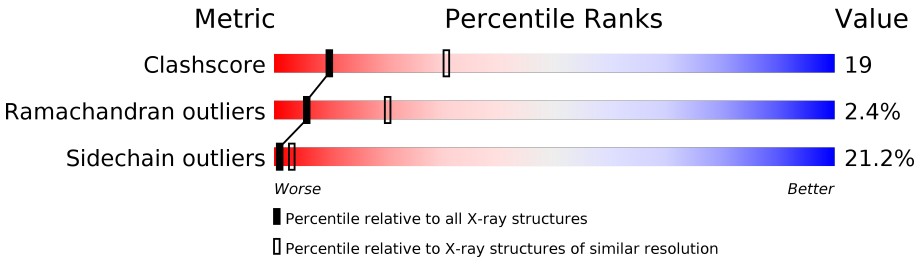
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

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 2.80 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	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)

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 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 <=5%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	214	54% 35% 10% .
2	B	213	54% 35% 9% .

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3303 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 PROTEIN (IGG2A KAPPA ANTIBODY CB41 (LIGHT CHAIN)).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	214	1678	1052	276	340	10	0	0	0

- Molecule 2 is a protein called PROTEIN (IGG2A KAPPA ANTIBODY CB41 (HEAVY CHAIN)).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	213	1595	1011	263	315	6	0	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	12	12	12	0	0
3	B	18	Total	O	0	0
			18	18		

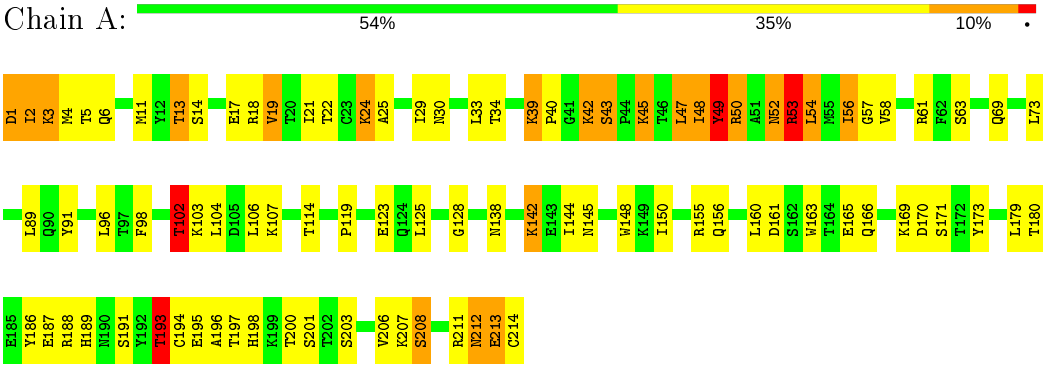
3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

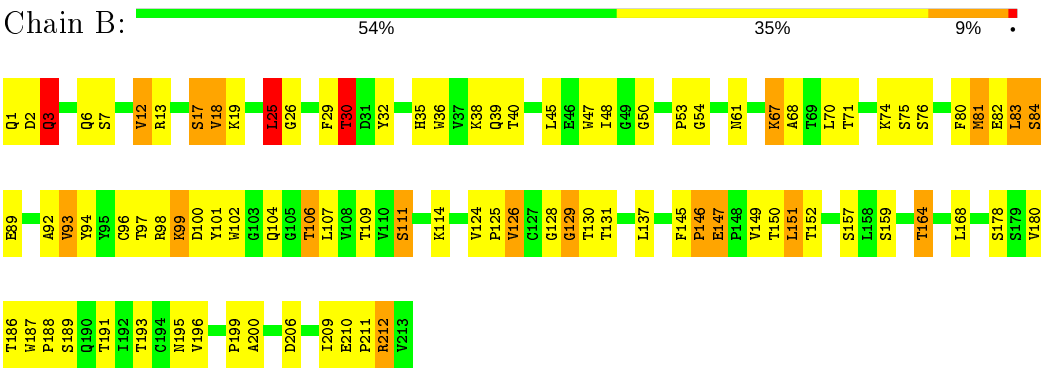
- Molecule 1: PROTEIN (IGG2A KAPPA ANTIBODY CB41 (LIGHT CHAIN))

Chain A:



- Molecule 2: PROTEIN (IGG2A KAPPA ANTIBODY CB41 (HEAVY CHAIN))

Chain B:



4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	147.34Å 55.60Å 75.67Å 90.00° 118.91° 90.00°	Depositor
Resolution (Å)	65.00 – 2.80	Depositor
% Data completeness (in resolution range)	84.0 (65.00-2.80)	Depositor
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CCP4	Depositor
R, R_{free}	0.221 , 0.312	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3303	wwPDB-VP
Average B, all atoms (Å^2)	35.0	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	A	0.43	0/1715	1.25	10/2321 (0.4%)
2	B	0.46	0/1635	1.30	11/2233 (0.5%)
All	All	0.45	0/3350	1.27	21/4554 (0.5%)

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	A	0	2
2	B	0	2
All	All	0	4

There are no bond length outliers.

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	147	GLU	CA-C-O	-11.07	96.85	120.10
1	A	1	ASP	CA-C-O	-8.20	102.89	120.10
2	B	212	ARG	NE-CZ-NH1	-7.64	116.48	120.30
1	A	1	ASP	CA-C-N	7.22	133.07	117.20
1	A	188	ARG	NE-CZ-NH2	-7.05	116.78	120.30
2	B	129	GLY	N-CA-C	6.84	130.19	113.10
1	A	102	THR	N-CA-CB	6.83	123.28	110.30
2	B	71	THR	N-CA-CB	6.30	122.28	110.30
2	B	30	THR	N-CA-CB	6.11	121.90	110.30
1	A	53	ARG	NE-CZ-NH2	-5.98	117.31	120.30
1	A	49	TYR	CA-CB-CG	5.77	124.36	113.40
2	B	25	LEU	CA-CB-CG	5.61	128.20	115.30
1	A	188	ARG	NE-CZ-NH1	5.60	123.10	120.30
2	B	100	ASP	CB-CG-OD1	5.55	123.30	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	2	ILE	CB-CA-C	-5.36	100.88	111.60
2	B	40	THR	N-CA-CB	5.30	120.37	110.30
2	B	212	ARG	NH1-CZ-NH2	5.29	125.22	119.40
1	A	193	THR	CA-CB-CG2	5.25	119.75	112.40
2	B	67	LYS	CA-CB-CG	5.21	124.86	113.40
1	A	114	THR	N-CA-CB	5.18	120.14	110.30
2	B	83	LEU	CA-CB-CG	5.06	126.95	115.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	138	ASN	Mainchain
1	A	2	ILE	Mainchain
2	B	147	GLU	Mainchain
2	B	164	THR	Mainchain

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	1678	0	1625	68	3
2	B	1595	0	1576	62	0
3	A	12	0	0	2	0
3	B	18	0	0	1	0
All	All	3303	0	3201	124	3

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 19.

All (124) 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:B:124:VAL:HG22	2:B:209:ILE:HG23	1.39	1.03
1:A:106:LEU:H	1:A:166:GLN:HE22	1.12	0.92

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1:ASP:N	1:A:3:LYS:NZ	2.26	0.83
1:A:119:PRO:HD2	2:B:212:ARG:HH12	1.45	0.81
1:A:1:ASP:H3	1:A:3:LYS:NZ	1.79	0.81
1:A:39:LYS:HB2	1:A:42:LYS:HE3	1.63	0.80
1:A:1:ASP:N	1:A:3:LYS:HZ1	1.80	0.80
1:A:150:ILE:HD11	1:A:179:LEU:HD21	1.68	0.76
1:A:6:GLN:HE21	1:A:102:THR:HG23	1.52	0.75
1:A:56:ILE:HG13	1:A:57:GLY:H	1.53	0.74
1:A:150:ILE:HG21	1:A:155:ARG:NH2	2.04	0.73
2:B:151:LEU:HD22	2:B:196:VAL:HG22	1.69	0.73
2:B:193:THR:HG21	2:B:206:ASP:HB3	1.70	0.73
2:B:183:THR:HG23	2:B:185:ASN:H	1.55	0.71
1:A:193:THR:CG2	1:A:208:SER:HB3	2.20	0.71
1:A:106:LEU:N	1:A:166:GLN:HE22	1.86	0.69
1:A:144:ILE:HG22	1:A:163:TRP:CH2	2.29	0.67
2:B:98:ARG:O	2:B:99:LYS:HB2	1.93	0.67
1:A:193:THR:HG23	1:A:208:SER:HB3	1.75	0.66
2:B:17:SER:OG	2:B:84:SER:HA	1.98	0.63
2:B:124:VAL:HG21	2:B:210:GLU:O	1.99	0.63
1:A:119:PRO:HD2	2:B:212:ARG:NH1	2.14	0.62
2:B:18:VAL:HG13	2:B:86:LEU:HD11	1.82	0.62
1:A:89:LEU:HD13	1:A:98:PHE:CE2	2.34	0.61
2:B:39:GLN:HB3	2:B:93:VAL:HG23	1.80	0.61
1:A:145:ASN:O	1:A:196:ALA:HA	2.01	0.61
2:B:39:GLN:O	2:B:92:ALA:HB1	2.01	0.61
1:A:119:PRO:CD	2:B:212:ARG:HH12	2.13	0.60
2:B:137:LEU:HB2	2:B:180:VAL:HG12	1.81	0.60
2:B:38:LYS:HD2	2:B:94:TYR:CE2	2.37	0.60
2:B:1:GLN:HB3	2:B:26:GLY:HA3	1.83	0.59
2:B:3:GLN:HB3	2:B:25:LEU:HG	1.85	0.59
1:A:6:GLN:NE2	1:A:102:THR:HG23	2.16	0.59
1:A:144:ILE:HG22	1:A:163:TRP:HH2	1.67	0.59
1:A:1:ASP:N	1:A:3:LYS:HZ3	1.99	0.59
2:B:30:THR:HA	2:B:53:PRO:HB2	1.83	0.59
2:B:183:THR:HG22	2:B:186:THR:OG1	2.03	0.59
1:A:56:ILE:HG13	1:A:57:GLY:N	2.18	0.58
1:A:1:ASP:H2	1:A:3:LYS:NZ	2.00	0.57
2:B:124:VAL:HG12	2:B:125:PRO:HD2	1.86	0.57
2:B:2:ASP:HB3	2:B:101:TYR:CD1	2.40	0.57
1:A:21:ILE:HG23	1:A:102:THR:HG21	1.86	0.56
1:A:4:MET:SD	1:A:25:ALA:HB2	2.44	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:198:HIS:HD2	1:A:200:THR:H	1.51	0.56
2:B:19:LYS:HG3	2:B:82:GLU:HB2	1.87	0.56
1:A:13:THR:HG21	1:A:19:VAL:HG12	1.88	0.56
1:A:13:THR:HG22	1:A:17:GLU:OE2	2.06	0.55
1:A:161:ASP:O	2:B:168:LEU:HD21	2.06	0.54
1:A:165:GLU:HB3	3:A:219:HOH:O	2.07	0.54
1:A:150:ILE:HD13	1:A:155:ARG:CD	2.38	0.54
2:B:164:THR:HG23	2:B:178:SER:HB2	1.89	0.54
1:A:33:LEU:HA	1:A:89:LEU:O	2.08	0.54
1:A:182:THR:HG22	1:A:184:ASP:H	1.73	0.53
2:B:187:TRP:CG	2:B:188:PRO:HA	2.44	0.53
2:B:12:VAL:HG11	2:B:18:VAL:CG1	2.38	0.53
2:B:152:THR:HG23	2:B:195:ASN:HB2	1.91	0.53
1:A:1:ASP:H3	1:A:3:LYS:HZ1	1.44	0.53
2:B:39:GLN:HG3	2:B:45:LEU:HD13	1.90	0.52
2:B:38:LYS:HB2	2:B:48:ILE:HD11	1.92	0.52
2:B:36:TRP:CD1	2:B:70:LEU:HD22	2.45	0.52
2:B:6:GLN:NE2	2:B:106:THR:HG22	2.25	0.51
1:A:1:ASP:O	1:A:3:LYS:NZ	2.35	0.51
1:A:198:HIS:CD2	1:A:200:THR:H	2.27	0.51
1:A:49:TYR:CD1	1:A:53:ARG:HB2	2.46	0.51
2:B:124:VAL:HG11	2:B:211:PRO:HA	1.94	0.50
2:B:128:GLY:O	2:B:130:THR:N	2.40	0.50
1:A:212:ASN:O	1:A:213:GLU:HB2	2.12	0.50
1:A:42:LYS:HD2	1:A:43:SER:O	2.13	0.49
1:A:160:LEU:HB3	2:B:168:LEU:CD1	2.43	0.48
2:B:2:ASP:HB3	2:B:101:TYR:CE1	2.48	0.48
2:B:97:THR:HB	2:B:102:TRP:CE3	2.49	0.48
2:B:1:GLN:NE2	2:B:25:LEU:HD12	2.28	0.48
2:B:29:PHE:CE1	2:B:53:PRO:HB3	2.49	0.47
2:B:68:ALA:HB1	2:B:81:MET:CE	2.45	0.47
1:A:24:LYS:HA	1:A:69:GLN:O	2.14	0.47
2:B:212:ARG:NE	3:B:217:HOH:O	2.47	0.47
1:A:19:VAL:HG21	1:A:104:LEU:HD21	1.96	0.47
1:A:183:LYS:O	1:A:187:GLU:HG3	2.15	0.47
2:B:124:VAL:O	2:B:212:ARG:NH2	2.48	0.46
1:A:150:ILE:HD11	1:A:179:LEU:CD2	2.40	0.46
2:B:126:VAL:HA	2:B:212:ARG:HH11	1.80	0.46
2:B:199:PRO:O	2:B:200:ALA:C	2.54	0.46
1:A:54:LEU:HD21	1:A:58:VAL:O	2.16	0.46
1:A:6:GLN:HA	1:A:22:THR:O	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:152:THR:CG2	2:B:195:ASN:HB2	2.46	0.46
1:A:142:LYS:HG3	1:A:173:TYR:CE1	2.51	0.45
1:A:6:GLN:HE21	1:A:102:THR:CG2	2.25	0.45
1:A:150:ILE:HD13	1:A:155:ARG:HD2	1.98	0.45
2:B:137:LEU:HD22	2:B:209:ILE:HG21	1.98	0.45
1:A:142:LYS:HE3	1:A:173:TYR:CE1	2.52	0.45
2:B:13:ARG:HD3	2:B:111:SER:O	2.16	0.45
2:B:47:TRP:CE3	2:B:61:ASN:HB2	2.51	0.44
1:A:163:TRP:HZ2	3:A:221:HOH:O	1.99	0.44
1:A:91:TYR:HA	1:A:96:LEU:HD22	1.99	0.44
1:A:125:LEU:O	1:A:183:LYS:HD2	2.17	0.44
2:B:68:ALA:HB1	2:B:81:MET:HE2	2.00	0.43
1:A:142:LYS:HE3	1:A:173:TYR:CZ	2.54	0.43
2:B:145:PHE:HA	2:B:146:PRO:HA	1.72	0.43
2:B:19:LYS:HE3	2:B:80:PHE:CD2	2.54	0.43
1:A:45:LYS:HB2	1:A:45:LYS:HE2	1.84	0.43
1:A:18:ARG:HA	1:A:75:ILE:O	2.19	0.43
2:B:124:VAL:HG13	2:B:187:TRP:CZ2	2.54	0.43
2:B:124:VAL:HG13	2:B:187:TRP:CH2	2.54	0.43
1:A:144:ILE:HG13	1:A:198:HIS:HB2	2.00	0.43
1:A:50:ARG:O	1:A:50:ARG:HG3	2.19	0.43
1:A:52:ASN:HD22	1:A:53:ARG:HG2	1.83	0.43
1:A:128:GLY:O	1:A:183:LYS:N	2.50	0.42
1:A:106:LEU:H	1:A:166:GLN:NE2	1.95	0.42
2:B:12:VAL:HG21	2:B:86:LEU:CD1	2.49	0.42
1:A:155:ARG:HH12	1:A:181:LEU:CD1	2.33	0.42
2:B:12:VAL:HG11	2:B:18:VAL:HG11	2.00	0.42
2:B:1:GLN:CB	2:B:26:GLY:HA3	2.48	0.42
2:B:29:PHE:O	2:B:30:THR:HG23	2.19	0.42
1:A:47:LEU:O	1:A:48:ILE:HD12	2.20	0.41
1:A:193:THR:HG22	1:A:208:SER:HB3	1.99	0.41
2:B:93:VAL:HG13	2:B:107:LEU:HD12	2.03	0.41
2:B:1:GLN:HE21	2:B:3:GLN:NE2	2.19	0.41
1:A:119:PRO:HD2	2:B:212:ARG:HH22	1.84	0.41
2:B:193:THR:CG2	2:B:206:ASP:HB3	2.45	0.41
1:A:195:GLU:HG3	1:A:206:VAL:HG22	2.02	0.41
1:A:170:ASP:O	1:A:171:SER:HB2	2.19	0.41
1:A:186:TYR:OH	1:A:211:ARG:HG3	2.20	0.41
2:B:1:GLN:HE22	2:B:25:LEU:HD12	1.85	0.40
2:B:35:HIS:ND1	2:B:50:GLY:HA3	2.36	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1:ASP:N	1:A:189:HIS:NE2[2_555]	1.72	0.48
1:A:1:ASP:O	1:A:189:HIS:CE1[2_555]	1.90	0.30
1:A:1:ASP:O	1:A:189:HIS:NE2[2_555]	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	212/214 (99%)	191 (90%)	18 (8%)	3 (1%)	11 34
2	B	211/213 (99%)	186 (88%)	18 (8%)	7 (3%)	4 13
All	All	423/427 (99%)	377 (89%)	36 (8%)	10 (2%)	6 20

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	56	ILE
2	B	3	GLN
2	B	30	THR
2	B	129	GLY
1	A	54	LEU
2	B	99	LYS
2	B	89	GLU
2	B	189	SER
1	A	40	PRO
2	B	54	GLY

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	193/193 (100%)	148 (77%)	45 (23%)	1 2
2	B	180/180 (100%)	146 (81%)	34 (19%)	1 5
All	All	373/373 (100%)	294 (79%)	79 (21%)	1 3

All (79) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LYS
1	A	5	THR
1	A	11	MET
1	A	13	THR
1	A	14	SER
1	A	19	VAL
1	A	24	LYS
1	A	29	ILE
1	A	30	ASN
1	A	34	THR
1	A	39	LYS
1	A	42	LYS
1	A	43	SER
1	A	45	LYS
1	A	47	LEU
1	A	48	ILE
1	A	49	TYR
1	A	50	ARG
1	A	52	ASN
1	A	53	ARG
1	A	61	ARG
1	A	63	SER
1	A	73	LEU
1	A	79	GLU
1	A	102	THR
1	A	103	LYS
1	A	107	LYS
1	A	123	GLU
1	A	142	LYS
1	A	148	TRP
1	A	156	GLN
1	A	169	LYS
1	A	180	THR

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Mol	Chain	Res	Type
1	A	183	LYS
1	A	191	SER
1	A	193	THR
1	A	194	CYS
1	A	197	THR
1	A	201	SER
1	A	203	SER
1	A	207	LYS
1	A	208	SER
1	A	212	ASN
1	A	213	GLU
1	A	214	CYS
2	B	3	GLN
2	B	7	SER
2	B	12	VAL
2	B	17	SER
2	B	18	VAL
2	B	25	LEU
2	B	30	THR
2	B	32	TYR
2	B	67	LYS
2	B	74	LYS
2	B	75	SER
2	B	76	SER
2	B	81	MET
2	B	83	LEU
2	B	84	SER
2	B	88	SER
2	B	93	VAL
2	B	96	CYS
2	B	104	GLN
2	B	106	THR
2	B	109	THR
2	B	111	SER
2	B	114	LYS
2	B	126	VAL
2	B	131	THR
2	B	146	PRO
2	B	149	VAL
2	B	150	THR
2	B	151	LEU
2	B	157	SER

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Mol	Chain	Res	Type
2	B	159	SER
2	B	183	THR
2	B	184	SER
2	B	191	THR

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

Mol	Chain	Res	Type
1	A	6	GLN
1	A	52	ASN
1	A	69	GLN
1	A	124	GLN
1	A	166	GLN
1	A	198	HIS
2	B	1	GLN
2	B	163	HIS
2	B	198	HIS

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

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.