



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 08:03 PM GMT

PDB ID : 1IKF
Title : A CONFORMATION OF CYCLOSPORIN A IN AQUEOUS ENVIRONMENT REVEALED BY THE X-RAY STRUCTURE OF A CYCLOSPORIN-FAB COMPLEX
Authors : Vix, O.; Altschuh, D.; Rees, B.; Thierry, J.-C.
Deposited on : 1993-12-09
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
<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

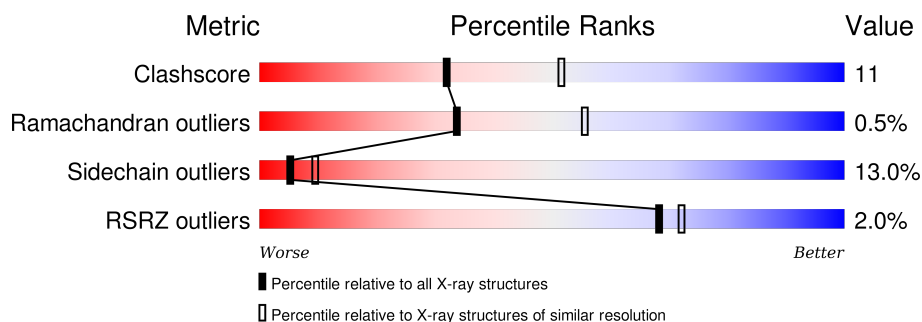
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (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 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	L	214	<div> <div></div> <div> <div></div> <div>65%</div> <div>30%</div> <div>• •</div> </div> </div>
2	H	228	<div> <div>3%</div> <div> <div></div> <div>65%</div> <div>29%</div> <div>5% •</div> </div> </div>
3	C	11	<div> <div>27%</div> <div>73%</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3651 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 R45-45-11 FAB (LIGHT CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	214	Total	C	N	O	S	0	0	0
			1652	1024	281	340	7			

- Molecule 2 is a protein called IGG1-KAPPA R45-45-11 FAB (HEAVY CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	228	Total	C	N	O	S	0	0	0
			1742	1107	288	338	9			

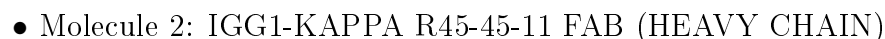
- Molecule 3 is a protein called CYCLOSPORIN A.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	11	Total	C	N	O	0	0	0
			85	62	11	12			

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	7	Total	O	0	0
			7	7		
4	H	69	Total	O	0	0
			69	69		
4	L	96	Total	O	0	0
			96	96		

- Molecule 1: IGG1-KAPPA R45-45-11 FAB (LIGHT CHAIN)



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	52.60 Å 70.20 Å 118.40 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.50 17.36 – 2.39	Depositor EDS
% Data completeness (in resolution range)	(Not available) (8.00-2.50) 78.3 (17.36-2.39)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$	-	Xtriage
Refinement program	X-PLOR	Depositor
R, R_{free}	0.164 , (Not available) 0.170 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	21.0	Xtriage
Anisotropy	0.413	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 66.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ¹	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Outliers	0 of 14219 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3651	wwPDB-VP
Average B, all atoms (Å ²)	12.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.02% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Theoretical 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

Bond lengths and bond angles in the following residue types are not validated in this section: ABA, MLE, DAL, MVA, BMT, SAR

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	0.86	0/1685	1.67	30/2286 (1.3%)
2	H	0.89	0/1788	1.76	36/2439 (1.5%)
3	C	1.05	0/10	2.04	0/11
All	All	0.88	0/3483	1.72	66/4736 (1.4%)

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
2	H	0	1

There are no bond length outliers.

All (66) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	201	ARG	NE-CZ-NH1	13.53	127.06	120.30
1	L	148	TRP	CD1-CG-CD2	10.73	114.88	106.30
1	L	32	TYR	CB-CG-CD2	-10.56	114.66	121.00
2	H	48	VAL	CG1-CB-CG2	-10.25	94.50	110.90
2	H	36	TRP	CD1-CG-CD2	8.88	113.40	106.30
2	H	47	TRP	CD1-CG-CD2	8.81	113.35	106.30
1	L	188	ARG	NE-CZ-NH1	8.69	124.65	120.30
1	L	148	TRP	CE2-CD2-CG	-8.58	100.44	107.30
1	L	108	ARG	NE-CZ-NH2	-8.31	116.15	120.30
1	L	35	TRP	CD1-CG-CD2	8.09	112.77	106.30
2	H	36	TRP	CE2-CD2-CG	-7.98	100.92	107.30
2	H	47	TRP	CE2-CD2-CG	-7.77	101.08	107.30
2	H	112	TRP	CD1-CG-CD2	7.68	112.45	106.30

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	112	TRP	CE2-CD2-CG	-7.68	101.16	107.30
2	H	167	TRP	CD1-CG-CD2	7.64	112.41	106.30
1	L	163	TRP	CE2-CD2-CG	-7.52	101.28	107.30
1	L	35	TRP	CE2-CD2-CG	-7.47	101.33	107.30
1	L	32	TYR	CB-CG-CD1	7.44	125.47	121.00
2	H	79	LEU	CA-CB-CG	7.41	132.33	115.30
2	H	112	TRP	CG-CD2-CE3	7.28	140.46	133.90
2	H	154	LEU	CA-CB-CG	7.24	131.95	115.30
1	L	163	TRP	CD1-CG-CD2	7.24	112.09	106.30
1	L	24	ARG	NE-CZ-NH2	-7.12	116.74	120.30
1	L	136	LEU	CA-CB-CG	7.08	131.59	115.30
1	L	108	ARG	NE-CZ-NH1	7.01	123.81	120.30
2	H	167	TRP	CE2-CD2-CG	-6.88	101.80	107.30
2	H	4	LEU	CA-CB-CG	6.75	130.82	115.30
2	H	190	LEU	CA-CB-CG	6.73	130.78	115.30
1	L	148	TRP	CG-CD1-NE1	-6.68	103.42	110.10
2	H	116	TRP	CD1-CG-CD2	6.52	111.52	106.30
2	H	36	TRP	CG-CD2-CE3	6.45	139.71	133.90
2	H	36	TRP	CB-CG-CD1	-6.39	118.69	127.00
1	L	31	THR	CA-CB-CG2	6.37	121.32	112.40
1	L	179	LEU	CA-CB-CG	6.36	129.92	115.30
2	H	116	TRP	CE2-CD2-CG	-6.36	102.22	107.30
1	L	163	TRP	CB-CG-CD1	-6.34	118.76	127.00
1	L	35	TRP	CG-CD2-CE3	6.26	139.54	133.90
2	H	149	VAL	CG1-CB-CG2	-6.26	100.88	110.90
2	H	48	VAL	CA-CB-CG2	6.11	120.07	110.90
1	L	18	ARG	NE-CZ-NH2	-6.11	117.25	120.30
2	H	85	ARG	NE-CZ-NH1	6.07	123.34	120.30
1	L	186	TYR	CB-CG-CD2	-6.02	117.39	121.00
1	L	33	LEU	CA-CB-CG	5.95	128.98	115.30
2	H	201	ARG	NE-CZ-NH2	-5.87	117.37	120.30
2	H	48	VAL	CA-CB-CG1	-5.81	102.19	110.90
1	L	172	THR	N-CA-CB	-5.69	99.50	110.30
1	L	163	TRP	CG-CD2-CE3	5.65	138.99	133.90
2	H	36	TRP	CG-CD1-NE1	-5.65	104.45	110.10
2	H	47	TRP	CG-CD1-NE1	-5.55	104.55	110.10
2	H	168	ASN	CB-CG-ND2	5.52	129.95	116.70
2	H	220	ASP	CB-CG-OD1	5.42	123.17	118.30
2	H	167	TRP	CG-CD1-NE1	-5.40	104.70	110.10
2	H	165	VAL	CA-C-N	5.38	129.04	117.20
1	L	213	ALA	CA-C-N	-5.36	105.41	117.20
2	H	135	TYR	CB-CG-CD1	-5.33	117.80	121.00

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	86	LEU	CA-CB-CG	5.21	127.28	115.30
1	L	188	ARG	CA-CB-CG	5.16	124.76	113.40
2	H	203	SER	N-CA-CB	-5.15	102.78	110.50
2	H	47	TRP	CG-CD2-CE3	5.15	138.53	133.90
1	L	24	ARG	NE-CZ-NH1	5.13	122.87	120.30
1	L	71	TYR	CB-CG-CD1	-5.13	117.92	121.00
1	L	35	TRP	CG-CD1-NE1	-5.08	105.02	110.10
2	H	112	TRP	CG-CD1-NE1	-5.08	105.02	110.10
1	L	82	ASP	CB-CG-OD1	5.08	122.87	118.30
2	H	111	VAL	CG1-CB-CG2	-5.01	102.88	110.90
1	L	35	TRP	CB-CG-CD1	-5.01	120.49	127.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	H	35	TYR	Sidechain

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	1652	0	1594	33	0
2	H	1742	0	1699	35	0
3	C	85	0	109	8	0
4	C	7	0	0	0	0
4	H	69	0	0	0	0
4	L	96	0	0	4	0
All	All	3651	0	3402	73	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 11.

All (73) 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:201:ARG:HG2	2:H:202:PRO:HA	1.66	0.77
1:L:106:ILE:H	1:L:166:GLN:HE22	1.36	0.73
2:H:53:ASN:HA	2:H:72:ARG:NH1	2.03	0.73
2:H:201:ARG:HH11	2:H:201:ARG:HB3	1.55	0.70
1:L:138:ASN:HD22	1:L:172:THR:HG21	1.61	0.66
1:L:34:ASN:HD22	1:L:89:GLN:HE22	1.45	0.64
1:L:31:THR:HG23	1:L:51:THR:CG2	2.31	0.61
2:H:91:THR:HG23	2:H:123:THR:HA	1.83	0.60
3:C:501:DAL:C	3:C:503:MLE:HN1	2.33	0.59
2:H:160:PRO:O	2:H:212:HIS:HE1	1.86	0.59
1:L:138:ASN:HA	1:L:172:THR:HG23	1.86	0.58
1:L:125:LEU:HD22	1:L:183:LYS:HG3	1.83	0.58
2:H:165:VAL:HG21	2:H:190:LEU:HD11	1.85	0.58
1:L:48:ILE:HG22	4:L:2031:HOH:O	2.01	0.58
2:H:24:THR:HB	2:H:27:PHE:CZ	2.39	0.58
3:C:504:MVA:HN3	3:C:505:BMT:HG2	1.87	0.57
1:L:12:SER:HA	1:L:105:GLU:O	2.06	0.56
2:H:12:VAL:HG21	2:H:86:LEU:HD23	1.89	0.55
2:H:38:ARG:HD3	2:H:48:VAL:HG11	1.90	0.54
2:H:104:THR:HB	2:H:107:GLY:O	2.07	0.53
2:H:47:TRP:HH2	2:H:59:PHE:HD1	1.56	0.53
1:L:30:SER:O	1:L:31:THR:HB	2.09	0.53
2:H:6:GLU:HG3	2:H:96:CYS:SG	2.50	0.52
2:H:60:TYR:OH	2:H:70:ILE:HG22	2.10	0.52
1:L:31:THR:HG22	1:L:50:TYR:CE1	2.46	0.50
1:L:34:ASN:HD22	1:L:89:GLN:NE2	2.10	0.50
1:L:165:ASP:HB3	4:L:2088:HOH:O	2.10	0.50
2:H:38:ARG:NH2	2:H:63:ILE:HD12	2.27	0.50
1:L:52:SER:HA	1:L:64:GLY:O	2.12	0.50
1:L:78:LEU:HD21	1:L:104:LEU:HD21	1.94	0.50
1:L:115:VAL:HA	1:L:135:PHE:O	2.12	0.49
2:H:4:LEU:HD13	2:H:24:THR:HG22	1.94	0.49
1:L:134:CYS:HB2	1:L:148:TRP:CH2	2.47	0.49
2:H:38:ARG:HG2	2:H:46:GLU:HB3	1.96	0.47
1:L:198:HIS:HD2	1:L:200:THR:OG1	1.96	0.47
1:L:31:THR:HG23	1:L:51:THR:HG22	1.96	0.47
1:L:196:ALA:HB3	1:L:205:ILE:HB	1.95	0.47
1:L:27:GLN:O	1:L:69:THR:HG22	2.16	0.46
1:L:15:LEU:HD22	1:L:15:LEU:H	1.80	0.46
3:C:501:DAL:HA	3:C:502:MLE:HN1	1.70	0.46
2:H:30:SER:O	2:H:53:ASN:HB2	2.15	0.45
2:H:47:TRP:HH2	2:H:59:PHE:CD1	2.35	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:161:ASN:ND2	1:L:177:SER:OG	2.50	0.45
3:C:502:MLE:N	3:C:503:MLE:HN1	2.32	0.45
2:H:2:VAL:HG13	2:H:27:PHE:CD1	2.51	0.45
2:H:107:GLY:HA2	3:C:510:MLE:HD22	1.99	0.45
3:C:509:VAL:HA	3:C:510:MLE:HN1	1.80	0.45
1:L:135:PHE:HE1	1:L:176:SER:HG	1.64	0.45
2:H:36:TRP:CD1	2:H:81:LEU:HG	2.53	0.44
1:L:29:ILE:HD11	1:L:71:TYR:CE2	2.53	0.44
1:L:13:ALA:HB3	1:L:78:LEU:HD11	1.98	0.44
1:L:36:TYR:HE2	1:L:89:GLN:HE21	1.64	0.44
2:H:104:THR:HG21	3:C:508:MLE:HD21	2.00	0.43
2:H:5:VAL:O	2:H:22:CYS:HA	2.17	0.43
1:L:138:ASN:HA	1:L:172:THR:CG2	2.47	0.43
4:L:2067:HOH:O	2:H:139:PRO:HA	2.19	0.43
1:L:53:ARG:HH12	2:H:104:THR:HA	1.84	0.43
2:H:109:TYR:HA	2:H:110:PRO:HD3	1.85	0.43
2:H:34:MET:HB3	2:H:79:LEU:HG	2.00	0.43
1:L:144:ILE:HG13	1:L:198:HIS:HB2	2.01	0.43
1:L:150:ILE:HD11	1:L:179:LEU:HD11	2.01	0.43
2:H:2:VAL:HG13	2:H:27:PHE:CE1	2.54	0.42
1:L:10:SER:HA	1:L:103:LYS:O	2.19	0.42
2:H:17:SER:HA	2:H:83:MET:O	2.19	0.42
2:H:132:PRO:HB3	2:H:158:TYR:HB3	2.02	0.42
2:H:144:GLN:HB3	2:H:146:ASN:OD1	2.19	0.41
2:H:144:GLN:HE21	2:H:226:ARG:NH2	2.18	0.41
2:H:177:HIS:O	2:H:192:SER:HA	2.20	0.41
1:L:31:THR:O	1:L:51:THR:HG22	2.20	0.41
2:H:97:THR:HA	2:H:115:ASP:O	2.21	0.41
1:L:155:ARG:HA	4:L:2081:HOH:O	2.19	0.41
3:C:503:MLE:HA	3:C:504:MVA:HN1	1.76	0.40
2:H:60:TYR:HB2	2:H:65:LYS:HG3	2.04	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%)	12 (6%)	1 (0%)	34	55
2	H	226/228 (99%)	217 (96%)	8 (4%)	1 (0%)	39	61
3	C	1/11 (9%)	1 (100%)	0	0	100	100
All	All	439/453 (97%)	417 (95%)	20 (5%)	2 (0%)	34	55

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	51	THR
2	H	143	ALA

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	190/190 (100%)	166 (87%)	24 (13%)	5	10
2	H	195/195 (100%)	169 (87%)	26 (13%)	5	9
3	C	1/1 (100%)	1 (100%)	0	100	100
All	All	386/386 (100%)	336 (87%)	50 (13%)	5	9

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	1	ASP
1	L	3	GLN
1	L	7	THR
1	L	11	LEU
1	L	22	SER
1	L	46	LEU
1	L	50	TYR
1	L	51	THR

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	L	55	ARG
1	L	59	PRO
1	L	67	SER
1	L	81	GLU
1	L	89	GLN
1	L	136	LEU
1	L	157	ASN
1	L	169	LYS
1	L	172	THR
1	L	175	MET
1	L	179	LEU
1	L	181	LEU
1	L	183	LYS
1	L	201	SER
1	L	202	THR
1	L	214	CYS
2	H	11	LEU
2	H	13	GLN
2	H	21	SER
2	H	30	SER
2	H	79	LEU
2	H	81	LEU
2	H	86	LEU
2	H	93	MET
2	H	99	HIS
2	H	103	ASP
2	H	129	THR
2	H	137	LEU
2	H	147	SER
2	H	151	LEU
2	H	154	LEU
2	H	166	THR
2	H	168	ASN
2	H	173	SER
2	H	196	VAL
2	H	199	SER
2	H	201	ARG
2	H	203	SER
2	H	210	VAL
2	H	223	ILE
2	H	226	ARG
2	H	228	CYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	89	GLN
1	L	138	ASN
1	L	157	ASN
1	L	161	ASN
1	L	166	GLN
1	L	198	HIS
2	H	82	GLN
2	H	212	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

9 non-standard protein/DNA/RNA residues are 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	DAL	C	501	3	3,4,5	0.83	0	0,4,6	0.00	-
3	MLE	C	502	3	7,8,9	0.79	0	4,9,11	1.53	1 (25%)
3	MLE	C	503	3	7,8,9	0.69	0	4,9,11	1.12	0
3	MVA	C	504	3	6,7,8	0.80	0	6,8,10	1.66	2 (33%)
3	BMT	C	505	3	11,12,13	1.82	3 (27%)	11,14,16	2.15	4 (36%)
3	ABA	C	506	3	4,5,6	0.57	0	3,5,7	2.52	2 (66%)
3	SAR	C	507	3	4,4,5	0.73	0	2,3,5	1.14	0
3	MLE	C	508	3	7,8,9	0.74	0	4,9,11	0.62	0
3	MLE	C	510	3	7,8,9	0.76	0	4,9,11	1.30	1 (25%)

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	DAL	C	501	3	-	0/0/2/4	0/0/0/0
3	MLE	C	502	3	-	0/4/8/10	0/0/0/0
3	MLE	C	503	3	-	0/4/8/10	0/0/0/0
3	MVA	C	504	3	-	0/5/8/10	0/0/0/0
3	BMT	C	505	3	-	1/13/16/18	0/0/0/0
3	ABA	C	506	3	-	0/2/4/6	0/0/0/0
3	SAR	C	507	3	-	0/1/2/3	0/0/0/0
3	MLE	C	508	3	-	0/4/8/10	0/0/0/0
3	MLE	C	510	3	-	0/4/8/10	0/0/0/0

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	505	BMT	CA-N	-2.28	1.43	1.47
3	C	505	BMT	CG2-CB	3.19	1.59	1.53
3	C	505	BMT	CE-CZ	3.58	1.56	1.29

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	505	BMT	O-C-CA	-3.68	114.44	125.74
3	C	506	ABA	O-C-CA	-3.63	116.04	125.49
3	C	505	BMT	CD1-CG2-CD2	-3.54	105.49	110.86
3	C	502	MLE	O-C-CA	-2.92	117.74	125.44
3	C	510	MLE	O-C-CA	-2.51	118.80	125.44
3	C	504	MVA	CG1-CB-CA	-2.24	108.37	111.68
3	C	505	BMT	OG1-CB-CA	-2.11	105.41	109.46
3	C	504	MVA	CN-N-CA	2.29	120.71	113.65
3	C	506	ABA	CG-CB-CA	2.41	119.65	113.44
3	C	505	BMT	CG2-CD2-CE	3.35	118.86	113.81

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	505	BMT	CD2-CE-CZ-CH

There are no ring outliers.

7 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	501	DAL	2	0
3	C	502	MLE	2	0
3	C	503	MLE	3	0
3	C	504	MVA	2	0
3	C	505	BMT	1	0
3	C	508	MLE	1	0
3	C	510	MLE	2	0

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 ⓘ

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	L	214/214 (100%)	-0.81	2 (0%) 85 88	2, 7, 26, 59	0
2	H	228/228 (100%)	-0.68	7 (3%) 52 57	2, 8, 32, 58	0
3	C	2/11 (18%)	-0.72	0 100 100	18, 18, 18, 30	0
All	All	444/453 (98%)	-0.74	9 (2%) 68 72	2, 8, 30, 59	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	L	214	CYS	5.9
2	H	143	ALA	5.7
2	H	145	THR	3.6
1	L	213	ALA	3.0
2	H	142	ALA	3.0
2	H	228	CYS	2.9
2	H	141	SER	2.8
2	H	144	GLN	2.6
2	H	140	GLY	2.4

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	DAL	C	501	5/6	0.91	0.16	-	25,25,28,28	0
3	MLE	C	502	9/10	0.91	0.17	-	18,19,22,23	0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	BMT	C	505	13/14	0.89	0.17	-	5,13,19,19	0
3	SAR	C	507	5/6	0.97	0.10	-	10,14,15,17	0
3	MLE	C	508	9/10	0.96	0.10	-	7,15,19,21	0
3	MLE	C	510	9/10	0.93	0.10	-	16,21,25,25	0
3	MVA	C	504	8/9	0.95	0.15	-	2,9,12,14	0
3	ABA	C	506	6/7	0.97	0.09	-	5,7,8,10	0
3	MLE	C	503	9/10	0.90	0.14	-	13,16,19,20	0

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.