



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:02 AM GMT

PDB ID : 2Z93  
Title : Crystal structure of Fab fragment of anti-ciguatoxin antibody 10C9 in complex with CTX3C-ABCD  
Authors : Ui, M.; Tanaka, Y.; Tsumoto, K.  
Deposited on : 2007-09-14  
Resolution : 2.40 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](mailto: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.

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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

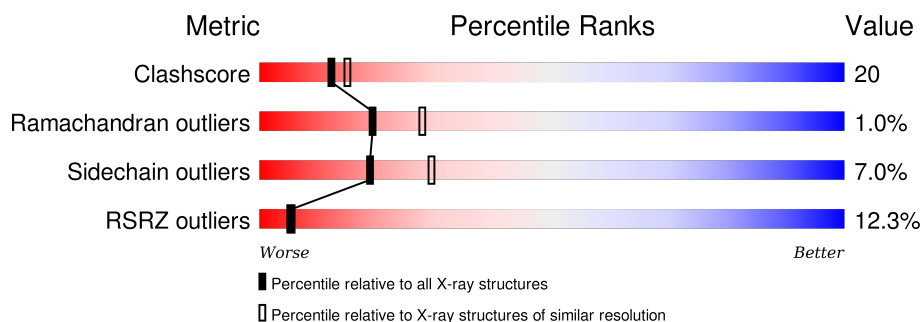
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 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	3407 (2.40-2.40)
Ramachandran outliers	100387	3351 (2.40-2.40)
Sidechain outliers	100360	3352 (2.40-2.40)
RSRZ outliers	91569	2928 (2.40-2.40)

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  $\geq 3$ , 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	A	218	
1	C	218	
2	B	213	
2	D	213	

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4307 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.

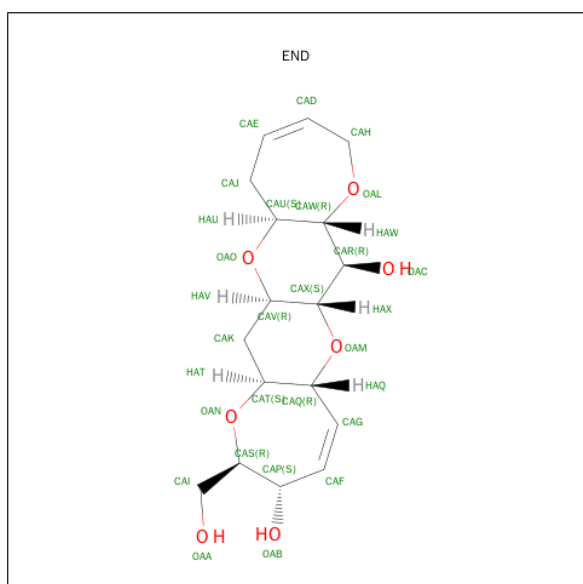
- Molecule 1 is a protein called Anti-ciguatoxin antibody 10C9 Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	81	Total	C	N	O	S	0	0	0
			594	383	93	116	2			
1	C	201	Total	C	N	O	S	0	0	0
			1565	998	254	307	6			

- Molecule 2 is a protein called Anti-ciguatoxin antibody 10C9 Fab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	80	Total	C	N	O	S	0	0	0
			623	389	106	125	3			
2	D	184	Total	C	N	O	S	0	0	0
			1380	859	226	286	9			

- Molecule 3 is 1,6:5,9:8,12:11,16-TETRAANHYDRO-2,3,4,10,13,14-HEXADEOXY-D-GLYCERO-D-ALLO-D-GULO-HEPTADEC-2,13-DIENITOL (three-letter code: END) (formula: C<sub>17</sub>H<sub>24</sub>O<sub>7</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	C	1	Total	C	O	0	0
			24	17	7		

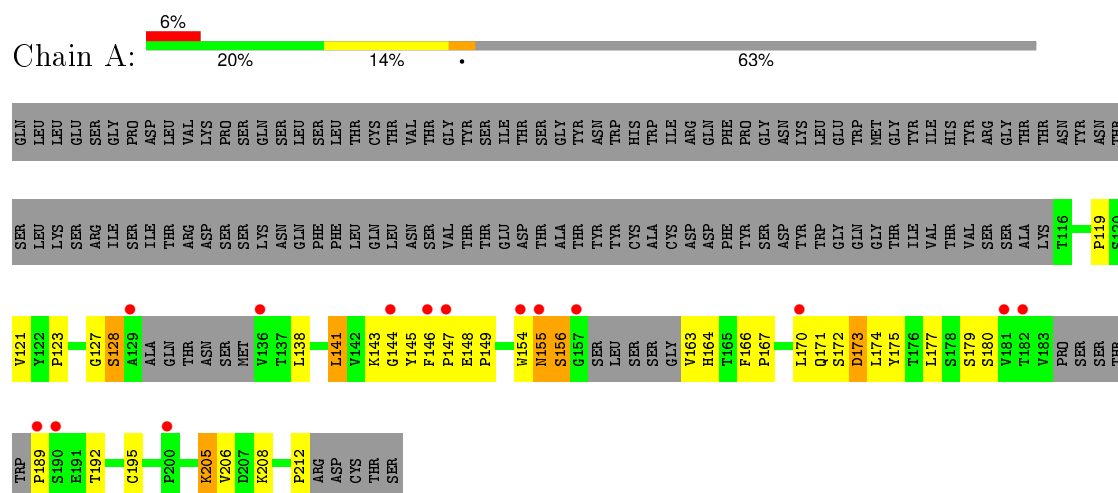
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	8	Total	O	0	0
			8	8		
4	B	18	Total	O	0	0
			18	18		
4	C	61	Total	O	0	0
			61	61		
4	D	34	Total	O	0	0
			34	34		

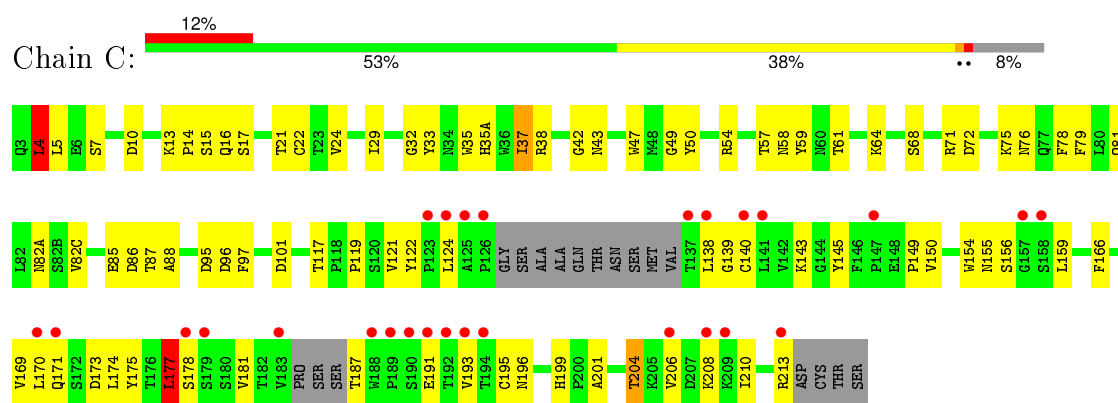
### 3 Residue-property plots

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.

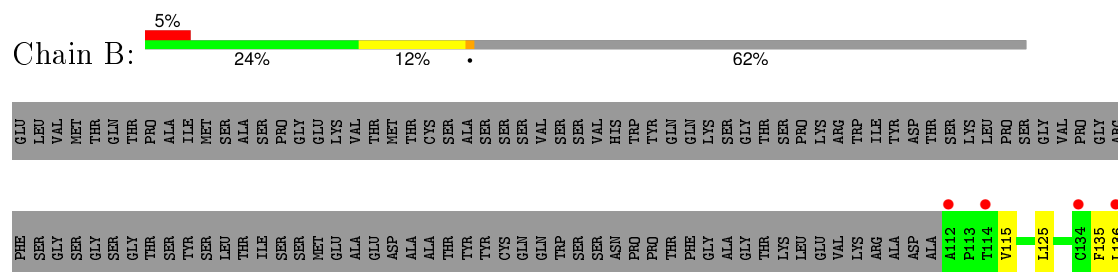
- Molecule 1: Anti-ciguatoxin antibody 10C9 Fab heavy chain

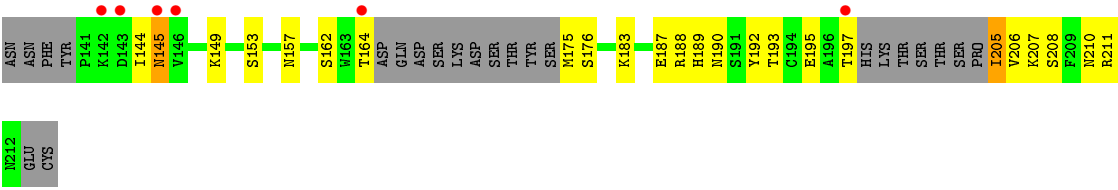


- Molecule 1: Anti-ciguatoxin antibody 10C9 Fab heavy chain

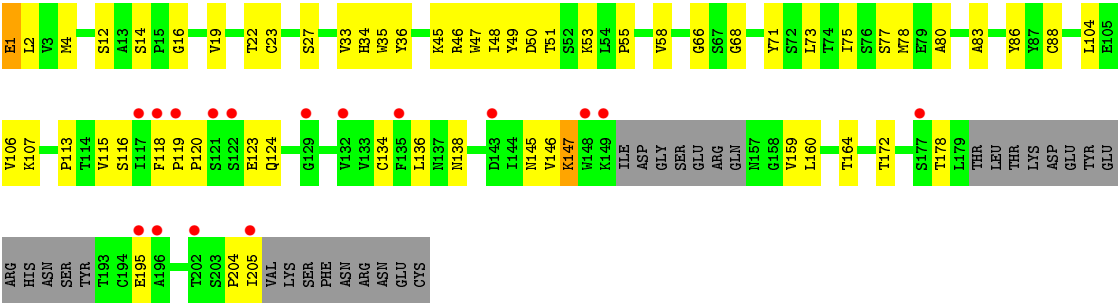


- Molecule 2: Anti-ciguatoxin antibody 10C9 Fab light chain





• Molecule 2: Anti-ciguatoxin antibody 10C9 Fab light chain



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	80.72Å 45.41Å 122.37Å 90.00° 98.90° 90.00°	Depositor
Resolution (Å)	9.99 – 2.40 42.51 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.4 (9.99-2.40) 99.5 (42.51-2.40)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.06	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.17 (at 2.39Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.297 , 0.340 (Not available) , (Not available)	Depositor DCC
$R_{free}$ test set	NotAvailable	DCC
Wilson B-factor (Å <sup>2</sup> )	34.5	Xtriage
Anisotropy	0.932	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 49.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 34703 reflections	Xtriage
$F_o, F_c$ correlation	0.85	EDS
Total number of atoms	4307	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.48% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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 [i](#)

### 5.1 Standard geometry [i](#)

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

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	0.37	0/609	0.72	0/838
1	C	0.37	0/1609	0.72	2/2204 (0.1%)
2	B	0.35	0/633	0.63	0/854
2	D	0.41	0/1413	0.67	0/1922
All	All	0.38	0/4264	0.69	2/5818 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	4	LEU	CA-CB-CG	5.52	127.99	115.30
1	C	177	LEU	CA-CB-CG	5.28	127.44	115.30

There are no chirality outliers.

There are no planarity outliers.

### 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	594	0	594	31	0
1	C	1565	0	1510	67	0
2	B	623	0	603	27	0
2	D	1380	0	1320	46	0
3	C	24	0	24	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	8	0	0	2	0
4	B	18	0	0	2	0
4	C	61	0	0	2	0
4	D	34	0	0	3	0
All	All	4307	0	4051	162	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 20.

All (162) 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:C:13:LYS:H	1:C:16:GLN:NE2	1.58	1.00
2:D:83:ALA:HB2	2:D:106:VAL:HG23	1.42	0.98
1:A:119:PRO:HB3	1:A:145:TYR:HB3	1.45	0.97
1:C:13:LYS:H	1:C:16:GLN:HE21	1.18	0.89
2:D:159:VAL:HA	2:D:178:THR:O	1.74	0.87
2:D:46:ARG:HD3	2:D:49:TYR:HB3	1.56	0.87
1:C:138:LEU:HD22	1:C:193:VAL:HG11	1.57	0.85
2:D:78:MET:HE1	2:D:104:LEU:HD21	1.59	0.84
1:C:124:LEU:HB2	1:C:139:GLY:HA3	1.58	0.83
2:D:46:ARG:HD2	4:D:216:HOH:O	1.82	0.80
1:C:187:THR:O	1:C:191:GLU:HB2	1.82	0.79
1:C:156:SER:H	1:C:196:ASN:HD21	1.31	0.76
1:C:177:LEU:HD23	1:C:178:SER:N	1.99	0.76
1:C:121:VAL:O	1:C:208:LYS:HE3	1.86	0.75
2:D:34:HIS:HD2	2:D:50:ASP:H	1.34	0.75
1:C:171:GLN:HG3	2:D:160:LEU:HD11	1.71	0.73
2:D:34:HIS:CD2	2:D:50:ASP:H	2.07	0.73
2:D:83:ALA:HB2	2:D:106:VAL:CG2	2.19	0.71
1:C:154:TRP:CZ3	1:C:195:CYS:HB3	2.26	0.70
1:C:121:VAL:HG21	1:C:206:VAL:HB	1.73	0.70
2:D:1:GLU:HG2	2:D:2:LEU:N	2.05	0.70
1:A:144:GLY:HA2	1:A:174:LEU:HB3	1.72	0.70
1:C:13:LYS:N	1:C:16:GLN:HE21	1.88	0.69
1:C:199:HIS:CE1	1:C:201:ALA:HB3	2.26	0.69
2:D:1:GLU:HA	4:D:231:HOH:O	1.93	0.69
1:C:169:VAL:HG11	2:D:160:LEU:HD22	1.75	0.68
2:D:55:PRO:HD2	2:D:58:VAL:HG21	1.76	0.68
1:C:61:THR:HA	1:C:64:LYS:HE2	1.77	0.67
1:C:119:PRO:HB3	1:C:145:TYR:HB3	1.77	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:213:ARG:HG3	1:C:213:ARG:HH11	1.61	0.65
2:D:14:SER:OG	2:D:107:LYS:HD2	1.97	0.65
1:A:141:LEU:CD1	1:A:143:LYS:HG3	2.28	0.64
2:B:189:HIS:O	2:B:211:ARG:HD3	1.96	0.64
1:A:141:LEU:HD11	1:A:143:LYS:HG3	1.80	0.63
1:A:166:PHE:HB3	2:B:162:SER:OG	1.99	0.62
1:C:124:LEU:HB2	1:C:139:GLY:CA	2.30	0.61
1:A:170:LEU:HD12	1:A:175:TYR:CZ	2.36	0.60
1:C:71:ARG:HA	1:C:78:PHE:HA	1.83	0.60
2:D:1:GLU:CD	2:D:2:LEU:HD13	2.22	0.60
2:B:195:GLU:HG3	2:B:206:VAL:HG22	1.84	0.60
2:D:73:LEU:HD23	2:D:73:LEU:C	2.22	0.59
2:B:149:LYS:HB2	2:B:193:THR:HB	1.84	0.59
1:C:29:ILE:H	1:C:76:ASN:HD21	1.50	0.59
2:D:136:LEU:N	2:D:136:LEU:HD12	2.17	0.59
2:B:205:ILE:HD13	2:B:205:ILE:N	2.18	0.58
1:C:119:PRO:HD2	1:C:204:THR:HG21	1.84	0.58
2:D:1:GLU:CG	2:D:2:LEU:N	2.67	0.58
1:A:127:GLY:O	1:A:128:SER:C	2.41	0.57
1:C:170:LEU:HB2	1:C:175:TYR:CE2	2.39	0.57
1:C:37:ILE:HG13	1:C:38:ARG:N	2.18	0.57
1:A:170:LEU:HD12	1:A:175:TYR:OH	2.06	0.56
1:C:122:TYR:CE2	2:D:124:GLN:HG3	2.41	0.56
1:C:173:ASP:O	1:C:174:LEU:HD23	2.05	0.56
1:C:170:LEU:HD12	1:C:174:LEU:O	2.06	0.55
4:C:278:HOH:O	2:D:34:HIS:HE1	1.88	0.55
1:A:166:PHE:CD1	2:B:164:THR:HG23	2.42	0.55
1:A:119:PRO:CB	1:A:145:TYR:HB3	2.26	0.54
1:C:13:LYS:N	1:C:16:GLN:NE2	2.42	0.54
2:D:116:SER:O	2:D:134:CYS:HA	2.07	0.54
1:A:170:LEU:C	1:A:170:LEU:HD23	2.29	0.54
1:A:148:GLU:OE1	1:A:149:PRO:HA	2.08	0.54
2:B:205:ILE:HA	4:B:215:HOH:O	2.08	0.53
2:B:175:MET:HG2	2:B:176:SER:N	2.24	0.53
1:C:156:SER:N	1:C:196:ASN:HD21	2.04	0.53
1:C:14:PRO:O	1:C:15:SER:HB2	2.07	0.53
1:C:139:GLY:O	1:C:154:TRP:HH2	1.91	0.53
1:C:154:TRP:CH2	1:C:195:CYS:HB3	2.43	0.53
1:A:167:PRO:HD3	2:B:164:THR:HG22	1.91	0.52
2:B:136:LEU:HD12	2:B:136:LEU:N	2.24	0.52
1:A:121:VAL:HB	1:A:206:VAL:HG11	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:140:CYS:HB2	1:C:154:TRP:CH2	2.45	0.52
1:C:138:LEU:N	1:C:138:LEU:HD12	2.25	0.51
1:C:199:HIS:HB3	1:C:204:THR:HG22	1.93	0.51
1:C:170:LEU:HD13	1:C:175:TYR:CZ	2.46	0.50
2:D:136:LEU:HD21	2:D:146:VAL:HG12	1.93	0.50
1:A:189:PRO:HG3	1:A:212:PRO:HG3	1.94	0.50
1:C:5:LEU:O	1:C:22:CYS:HA	2.12	0.50
1:C:150:VAL:HG23	1:C:199:HIS:HD2	1.77	0.49
2:D:80:ALA:HA	2:D:106:VAL:HG21	1.93	0.49
1:C:101:ASP:O	2:D:45:LYS:NZ	2.40	0.49
1:C:29:ILE:HA	1:C:35:TRP:CZ2	2.48	0.49
2:D:4:MET:HE2	2:D:88:CYS:SG	2.53	0.49
2:D:19:VAL:HG21	2:D:78:MET:CE	2.42	0.49
1:C:87:THR:O	1:C:88:ALA:HB2	2.12	0.49
2:D:78:MET:HB2	2:D:78:MET:HE3	1.58	0.49
2:D:147:LYS:HB3	2:D:195:GLU:HB3	1.95	0.49
1:C:42:GLY:O	1:C:43:ASN:HB2	2.13	0.49
1:A:180:SER:HB3	2:B:135:PHE:CE2	2.48	0.48
2:D:19:VAL:HG21	2:D:78:MET:HE3	1.93	0.48
2:B:190:ASN:O	2:B:210:ASN:HA	2.13	0.48
1:C:143:LYS:HG3	4:C:248:HOH:O	2.13	0.48
2:B:145:ASN:O	2:B:197:THR:N	2.32	0.48
1:A:163:VAL:HA	1:A:180:SER:O	2.14	0.48
1:C:138:LEU:HD23	1:C:210:ILE:HG21	1.96	0.48
1:A:147:PRO:HB2	4:A:221:HOH:O	2.14	0.47
1:C:17:SER:HA	1:C:82(A):ASN:HA	1.96	0.47
2:D:27:SER:HB2	4:D:227:HOH:O	2.15	0.47
1:C:47:TRP:CZ2	1:C:49:GLY:HA2	2.49	0.47
1:C:82(C):VAL:HA	1:C:86:ASP:OD2	2.13	0.47
2:B:190:ASN:HD21	2:B:210:ASN:HB3	1.80	0.47
2:B:193:THR:CG2	2:B:206:VAL:HG13	2.44	0.47
1:C:170:LEU:HD13	1:C:175:TYR:CE2	2.50	0.47
1:C:4:LEU:HD13	1:C:24:VAL:HG22	1.97	0.47
1:A:154:TRP:CZ3	1:A:195:CYS:HB3	2.50	0.46
2:D:75:ILE:HD12	2:D:78:MET:CE	2.46	0.46
1:C:121:VAL:HB	1:C:206:VAL:HG11	1.98	0.46
1:C:13:LYS:HB2	1:C:16:GLN:HE21	1.81	0.46
2:D:115:VAL:HG21	2:D:205:ILE:HG23	1.98	0.46
1:A:175:TYR:HE2	4:A:218:HOH:O	1.98	0.46
1:C:154:TRP:HB3	1:C:159:LEU:HD23	1.98	0.45
2:D:22:THR:HG22	2:D:23:CYS:N	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:145:ASN:H	2:B:197:THR:CG2	2.29	0.45
1:C:7:SER:OG	1:C:21:THR:HB	2.16	0.45
2:D:46:ARG:HD3	2:D:49:TYR:CB	2.40	0.45
1:C:213:ARG:CG	1:C:213:ARG:HH11	2.26	0.45
2:B:157:ASN:HA	4:B:225:HOH:O	2.17	0.45
2:D:66:GLY:HA3	2:D:71:TYR:HA	1.99	0.45
2:B:144:ILE:HD12	2:B:197:THR:HG23	1.98	0.45
2:D:35:TRP:HB2	2:D:48:ILE:HB	1.98	0.45
2:B:190:ASN:ND2	2:B:210:ASN:HB3	2.32	0.44
2:B:192:TYR:O	2:B:208:SER:HB2	2.18	0.44
1:A:172:SER:O	1:A:173:ASP:CB	2.65	0.44
2:B:125:LEU:O	2:B:183:LYS:HD2	2.18	0.44
2:B:115:VAL:HA	2:B:135:PHE:O	2.18	0.43
2:D:36:TYR:O	2:D:86:TYR:HA	2.18	0.43
1:C:35(A):HIS:ND1	1:C:50:TYR:HB3	2.32	0.43
2:B:145:ASN:H	2:B:197:THR:HG23	1.83	0.43
1:A:164:HIS:O	1:A:179:SER:HA	2.18	0.43
1:A:154:TRP:O	1:A:156:SER:N	2.51	0.43
1:C:138:LEU:HB2	1:C:181:VAL:CG1	2.48	0.43
2:D:50:ASP:HB3	2:D:53:LYS:CE	2.49	0.43
2:D:19:VAL:CG2	2:D:78:MET:HE3	2.49	0.42
2:D:4:MET:HE3	2:D:23:CYS:SG	2.58	0.42
1:C:13:LYS:CB	1:C:16:GLN:HE21	2.32	0.42
1:C:154:TRP:CE3	1:C:195:CYS:HB3	2.54	0.42
2:B:144:ILE:HA	2:B:197:THR:CG2	2.49	0.42
1:A:172:SER:O	1:A:173:ASP:HB2	2.20	0.42
2:D:113:PRO:O	2:D:205:ILE:HD13	2.19	0.42
1:A:146:PHE:CD1	1:A:147:PRO:HA	2.55	0.42
2:D:138:ASN:HA	2:D:172:THR:OG1	2.20	0.42
1:A:123:PRO:HD3	1:A:208:LYS:HD2	2.02	0.42
2:D:16:GLY:HA2	2:D:77:SER:OG	2.19	0.42
2:D:50:ASP:O	2:D:51:THR:HB	2.20	0.41
1:C:35:TRP:N	1:C:35:TRP:CD1	2.87	0.41
1:A:205:LYS:O	1:A:206:VAL:HG23	2.20	0.41
2:D:118:PHE:HA	2:D:119:PRO:HD3	1.92	0.41
1:C:117:THR:O	1:C:145:TYR:HA	2.20	0.41
1:C:29:ILE:H	1:C:76:ASN:ND2	2.16	0.41
1:C:59:TYR:HB2	1:C:64:LYS:HD3	2.01	0.41
1:A:154:TRP:CH2	1:A:195:CYS:HB3	2.56	0.41
1:C:213:ARG:CG	1:C:213:ARG:NH1	2.82	0.41
1:C:96:ASP:O	1:C:97:PHE:HB3	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:141:LEU:HD13	1:A:143:LYS:HG3	2.00	0.41
1:A:154:TRP:C	1:A:155:ASN:OD1	2.58	0.41
2:B:187:GLU:OE2	2:B:211:ARG:NH2	2.54	0.40
2:B:144:ILE:HG13	2:B:145:ASN:N	2.37	0.40
1:A:154:TRP:C	1:A:156:SER:N	2.74	0.40
1:C:32:GLY:O	1:C:33:TYR:HB2	2.21	0.40
1:C:155:ASN:OD1	1:C:193:VAL:HA	2.21	0.40
2:B:136:LEU:O	2:B:175:MET:N	2.55	0.40
1:C:166:PHE:CD1	2:D:164:THR:HG23	2.57	0.40
1:C:72:ASP:OD2	1:C:75:LYS:HE2	2.20	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	A	73/218 (34%)	64 (88%)	7 (10%)	2 (3%)	6	6
1	C	195/218 (89%)	180 (92%)	15 (8%)	0	100	100
2	B	72/213 (34%)	71 (99%)	1 (1%)	0	100	100
2	D	178/213 (84%)	170 (96%)	5 (3%)	3 (2%)	11	14
All	All	518/862 (60%)	485 (94%)	28 (5%)	5 (1%)	19	28

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	128	SER
1	A	155	ASN
2	D	204	PRO
2	D	120	PRO
2	D	68	GLY

### 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	A	72/196 (37%)	64 (89%)	8 (11%)	8	10
1	C	182/196 (93%)	168 (92%)	14 (8%)	16	24
2	B	73/188 (39%)	68 (93%)	5 (7%)	20	31
2	D	160/188 (85%)	153 (96%)	7 (4%)	35	53
All	All	487/768 (63%)	453 (93%)	34 (7%)	19	29

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	138	LEU
1	A	141	LEU
1	A	156	SER
1	A	171	GLN
1	A	173	ASP
1	A	177	LEU
1	A	192	THR
1	A	205	LYS
2	B	145	ASN
2	B	153	SER
2	B	188	ARG
2	B	205	ILE
2	B	207	LYS
1	C	4	LEU
1	C	10	ASP
1	C	37	ILE
1	C	54	ARG
1	C	57	THR
1	C	58	ASN
1	C	68	SER
1	C	79	PHE
1	C	81	GLN
1	C	85	GLU
1	C	95	ASP

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Mol	Chain	Res	Type
1	C	149	PRO
1	C	177	LEU
1	C	204	THR
2	D	1	GLU
2	D	12	SER
2	D	33	VAL
2	D	47	TRP
2	D	123	GLU
2	D	145	ASN
2	D	147	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	171	GLN
1	A	196	ASN
2	B	145	ASN
2	B	156	GLN
2	B	212	ASN
1	C	16	GLN
1	C	76	ASN
1	C	77	GLN
1	C	81	GLN
1	C	196	ASN
2	D	34	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

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	END	C	1	-	24,27,27	1.92	4 (16%)	24,39,39	2.51	9 (37%)

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	END	C	1	-	-	0/2/55/55	0/3/4/4

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	1	END	CAQ-CAG	-6.65	1.40	1.49
3	C	1	END	CAH-CAD	-3.58	1.40	1.48
3	C	1	END	CAJ-CAE	-3.48	1.40	1.49
3	C	1	END	CAG-CAF	3.85	1.38	1.32

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1	END	CAS-OAN-CAT	-4.74	107.29	114.96
3	C	1	END	CAJ-CAU-CAW	-2.45	110.68	114.14
3	C	1	END	OAN-CAT-CAK	2.06	110.66	107.42
3	C	1	END	CAW-CAR-CAX	2.15	113.39	109.06
3	C	1	END	CAH-CAD-CAE	2.30	131.24	127.24
3	C	1	END	OAL-CAW-CAU	2.37	115.38	110.70
3	C	1	END	OAB-CAP-CAS	2.40	111.81	108.48
3	C	1	END	OAM-CAQ-CAT	5.61	113.82	109.00
3	C	1	END	CAH-OAL-CAW	7.66	122.06	114.94



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 ⓘ

### 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	81/218 (37%)	1.22	14 (17%) 2 2	42, 54, 75, 83	0
1	C	201/218 (92%)	0.73	27 (13%) 4 4	18, 37, 69, 77	0
2	B	80/213 (37%)	0.81	10 (12%) 5 5	28, 43, 59, 63	0
2	D	184/213 (86%)	0.61	16 (8%) 13 12	13, 29, 71, 76	0
All	All	546/862 (63%)	0.77	67 (12%) 5 5	13, 41, 70, 83	0

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	140	CYS	6.0
1	C	189	PRO	4.9
1	C	193	VAL	4.8
2	D	205	ILE	4.5
1	A	190	SER	4.3
1	C	125	ALA	4.2
1	C	137	THR	4.2
2	B	164	THR	3.9
2	B	145	ASN	3.9
1	A	181	VAL	3.9
1	C	183	VAL	3.9
1	C	126	PRO	3.6
1	A	154	TRP	3.6
1	C	191	GLU	3.6
2	B	197	THR	3.6
2	D	121	SER	3.5
1	C	171	GLN	3.5
2	D	195	GLU	3.5
2	B	136	LEU	3.4
2	D	122	SER	3.4
2	D	132	VAL	3.3

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Mol	Chain	Res	Type	RSRZ
1	A	170	LEU	3.3
1	C	190	SER	3.3
2	D	196	ALA	3.2
1	C	192	THR	3.2
1	A	136	VAL	3.2
1	C	194	THR	3.1
1	A	182	THR	3.1
1	A	155	ASN	3.0
2	D	117	ILE	2.9
2	B	143	ASP	2.9
1	C	124	LEU	2.9
1	A	189	PRO	2.9
1	A	129	ALA	2.8
1	C	141	LEU	2.8
2	D	119	PRO	2.8
2	D	202	THR	2.7
1	C	123	PRO	2.7
2	B	114	THR	2.7
1	C	188	TRP	2.7
1	A	144	GLY	2.6
1	A	200	PRO	2.6
1	C	157	GLY	2.6
2	B	112	ALA	2.6
1	A	146	PHE	2.5
1	A	147	PRO	2.5
2	D	177	SER	2.5
1	C	179	SER	2.5
2	D	143	ASP	2.5
1	C	138	LEU	2.5
1	C	213	ARG	2.5
1	C	178	SER	2.4
1	C	170	LEU	2.4
2	D	129	GLY	2.4
2	B	142	LYS	2.3
1	C	158	SER	2.3
2	D	135	PHE	2.3
2	B	146	VAL	2.3
2	D	148	TRP	2.2
1	C	147	PRO	2.2
1	C	208	LYS	2.2
2	D	149	LYS	2.2
1	A	157	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
2	D	118	PHE	2.2
1	C	209	LYS	2.2
1	C	206	VAL	2.1
2	B	134	CYS	2.1

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

There are no non-standard protein/DNA/RNA residues in this entry.

## 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

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, 95<sup>th</sup> 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( $\text{\AA}^2$ )	Q<0.9
3	END	C	1	24/24	0.94	0.17	-0.32	16,19,24,26	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.