



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 06:58 am GMT

PDB ID : 1JVQ  
Title : Crystal structure at 2.6Å of the ternary complex between antithrombin, a P14-P8 reactive loop peptide, and an exogenous tetrapeptide  
Authors : Zhou, A.; Huntington, J.A.; Lomas, D.A.; Carrell, R.W.; Stein, P.E.  
Deposited on : 2001-08-31  
Resolution : 2.60 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 : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : trunk28620  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
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) : recalc28949

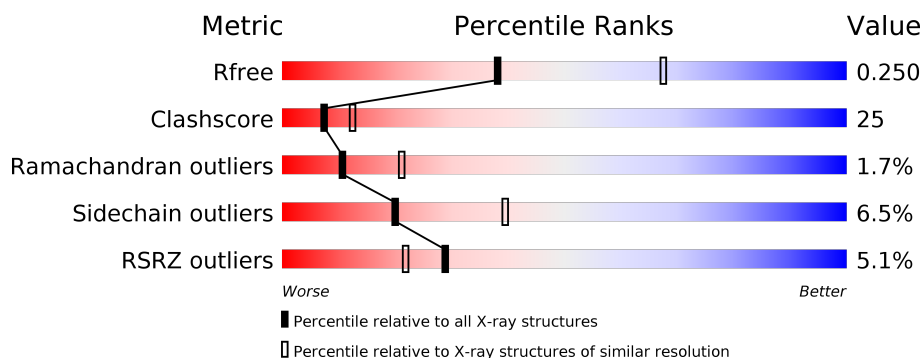
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.60 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))
$R_{free}$	100719	2542 (2.60-2.60)
Clashscore	112137	2895 (2.60-2.60)
Ramachandran outliers	110173	2848 (2.60-2.60)
Sidechain outliers	110143	2848 (2.60-2.60)
RSRZ outliers	101464	2550 (2.60-2.60)

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	I	432	<div> <div>2%</div> <div> <div></div> <div>59%</div> <div>32%</div> <div>• 5%</div> </div> </div>
1	L	432	<div> <div>8%</div> <div> <div></div> <div>52%</div> <div>36%</div> <div>5% • 6%</div> </div> </div>
2	C	8	<div> <div></div> <div> <div></div> <div>75%</div> <div>25%</div> </div> </div>
3	D	5	<div> <div></div> <div> <div></div> <div>40%</div> <div>60%</div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	NDG	L	841	-	-	X	-
5	NAG	I	841	-	-	X	-
5	NAG	I	861	-	-	X	-

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6801 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 ANTITHROMBIN-III.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	406	Total	C	N	O	S	0	0	0
			3249	2073	547	611	18			
1	I	410	Total	C	N	O	S	0	0	0
			3284	2093	554	619	18			

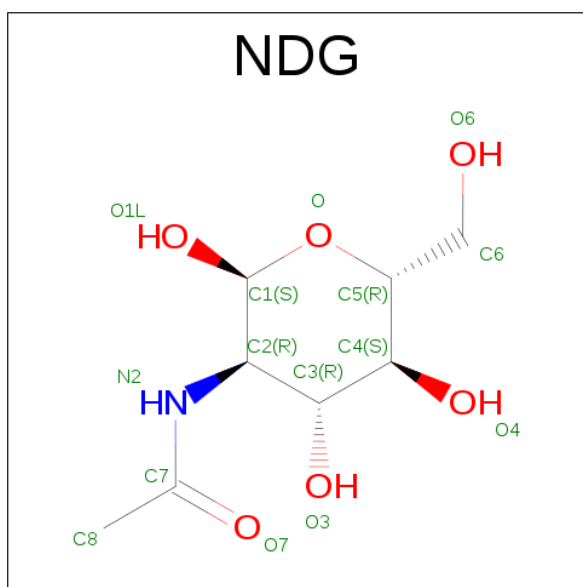
- Molecule 2 is a protein called P14-P8 reactive loop peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	8	Total	C	N	O	0	0	0
			47	26	7	14			

- Molecule 3 is a protein called exogenous Cholecystokinin tetrapeptide.

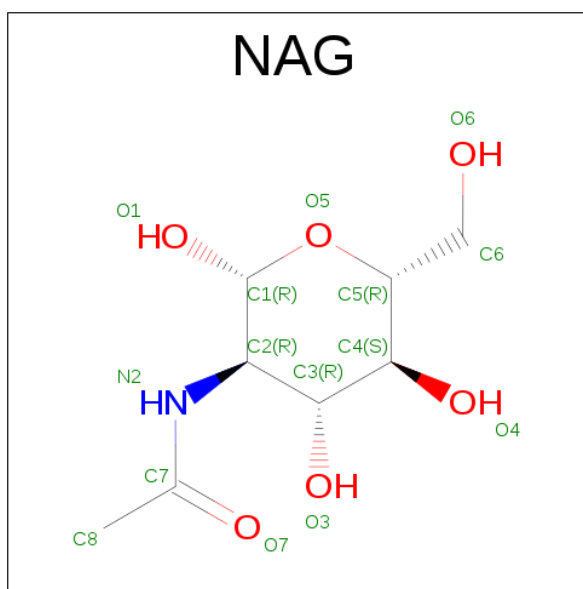
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	5	Total	C	N	O	S	0	0	1
			42	29	6	6	1			

- Molecule 4 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NDG) (formula:  $C_8H_{15}NO_6$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	L	1	Total	C	N	O	0	0
			15	8	1	6		
4	L	1	Total	C	N	O	0	0
			15	8	1	6		
4	L	1	Total	C	N	O	0	0
			15	8	1	6		
4	L	1	Total	C	N	O	0	0
			15	8	1	6		
4	I	1	Total	C	N	O	0	0
			15	8	1	6		
4	I	1	Total	C	N	O	0	0
			15	8	1	6		
4	I	1	Total	C	N	O	0	0
			15	8	1	6		

- Molecule 5 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula:  $C_8H_{15}NO_6$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	I	1	Total	C	N	O	0	0
			15	8	1	6		
5	I	1	Total	C	N	O	0	0
			15	8	1	6		

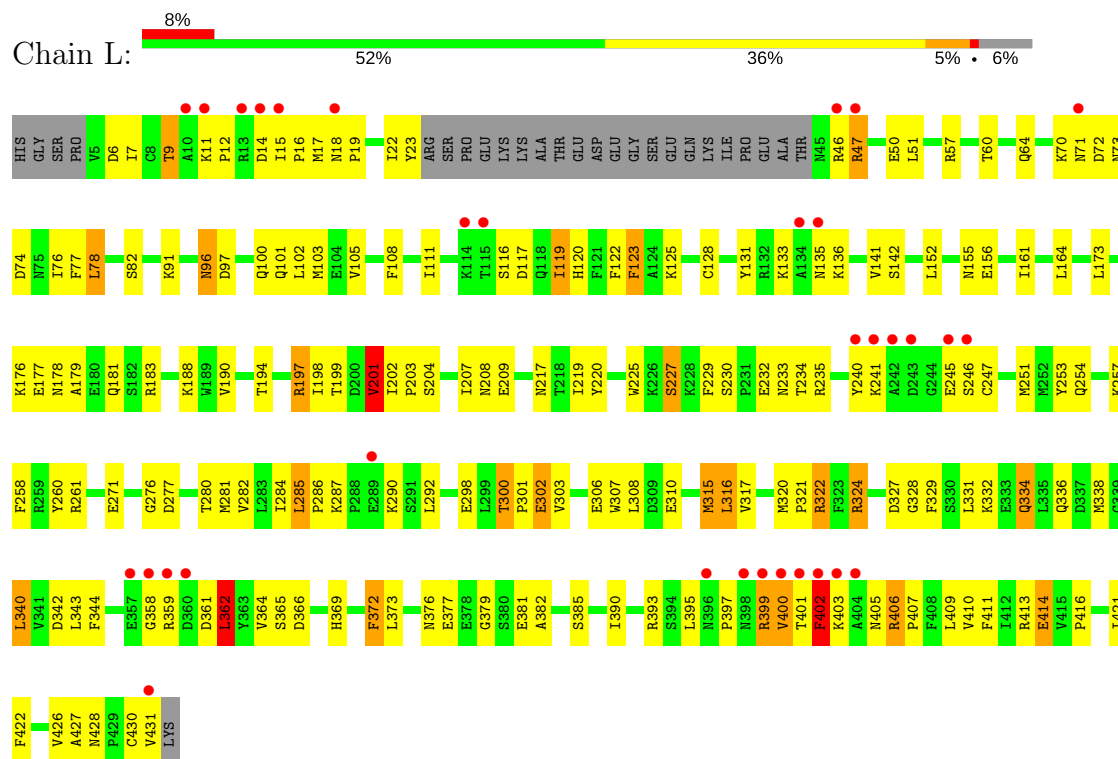
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	I	30	Total	O	0	0
			30	30		
6	L	14	Total	O	0	0
			14	14		

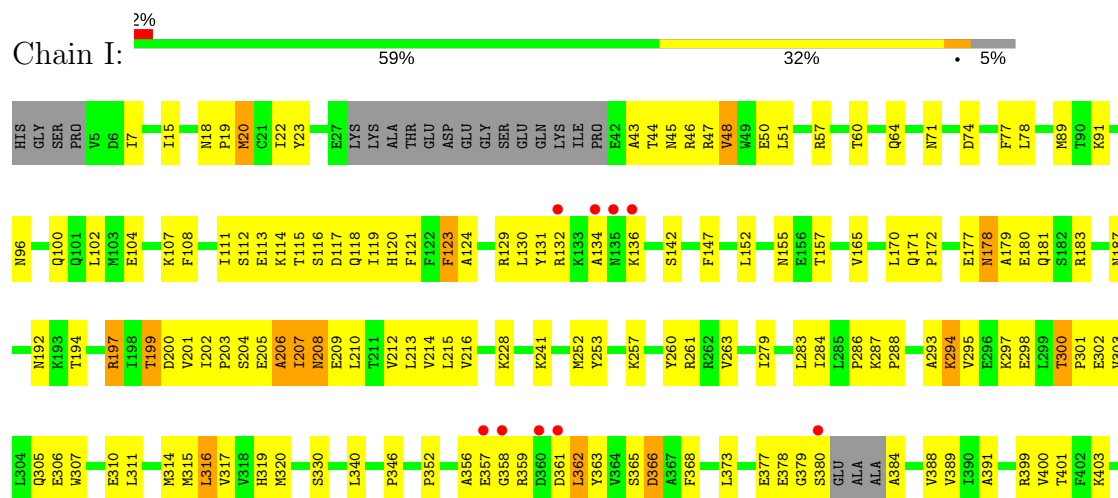
### 3 Residue-property plots

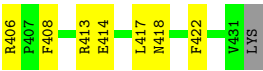
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are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



#### • Molecule 1: ANTITHROMBIN-III





- Molecule 2: P14-P8 reactive loop peptide



- Molecule 3: exogenous Cholecystokinin tetrapeptide





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	69.14Å 100.62Å 87.16Å 90.00° 104.42° 90.00°	Depositor
Resolution (Å)	38.63 – 2.60 40.55 – 2.60	Depositor EDS
% Data completeness (in resolution range)	96.4 (38.63-2.60) 96.4 (40.55-2.60)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	0.09	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.94 (at 2.61Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.204 , 0.258 0.205 , 0.250	Depositor DCC
$R_{free}$ test set	1029 reflections (3.09%)	DCC
Wilson B-factor (Å <sup>2</sup> )	56.9	Xtriage
Anisotropy	0.020	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 54.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	6801	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.31% 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, 0.333 respectively for untwinned datasets, and 0.375, 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: NH2, NDG, ACE, NAG

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	I	0.54	0/3348	0.74	0/4518
1	L	0.52	0/3313	0.72	2/4472 (0.0%)
2	C	0.84	0/44	0.76	0/58
3	D	0.61	0/43	0.67	0/57
All	All	0.53	0/6748	0.73	2/9105 (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	L	362	LEU	CA-CB-CG	5.36	127.63	115.30
1	L	414	GLU	N-CA-C	-5.16	97.08	111.00

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	I	3284	0	3284	152	0
1	L	3249	0	3252	168	0
2	C	47	0	41	3	0
3	D	42	0	31	8	0
4	I	45	0	45	11	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	L	60	0	60	13	0
5	I	30	0	30	15	0
6	I	30	0	0	1	0
6	L	14	0	0	0	0
All	All	6801	0	6743	333	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 25.

All (333) 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:155:ASN:HD21	4:L:841:NDG:C1	1.63	1.12
1:I:316:LEU:HB3	1:I:400:VAL:HG13	1.44	0.98
1:L:7:ILE:HD11	1:L:164:LEU:HG	1.47	0.96
1:I:208:ASN:HD22	1:I:209:GLU:H	0.95	0.94
1:L:91:LYS:NZ	1:L:120:HIS:NE2	2.15	0.94
1:L:71:ASN:HD21	1:L:403:LYS:HE2	1.35	0.92
1:I:208:ASN:HD22	1:I:210:LEU:H	1.18	0.91
4:L:841:NDG:H4	4:L:842:NDG:C1	2.01	0.91
1:L:230:SER:HB3	1:L:232:GLU:OE2	1.72	0.88
1:I:187:ASN:HD21	1:I:202:ILE:H	1.20	0.88
1:I:208:ASN:ND2	1:I:209:GLU:H	1.73	0.87
1:L:155:ASN:ND2	4:L:841:NDG:C1	2.37	0.86
1:L:300:THR:HG22	1:L:303:VAL:H	1.38	0.86
1:I:300:THR:CG2	1:I:302:GLU:HG2	2.06	0.86
4:L:841:NDG:H4	4:L:842:NDG:H1	1.58	0.84
1:I:152:LEU:HD11	1:I:212:VAL:HB	1.57	0.84
1:I:20:MET:HE3	1:I:352:PRO:HB2	1.59	0.82
1:I:283:LEU:HD11	1:I:320:MET:CE	2.10	0.82
1:I:178:ASN:N	1:I:178:ASN:HD22	1.75	0.81
1:I:208:ASN:ND2	1:I:210:LEU:H	1.76	0.81
4:L:841:NDG:H4	4:L:842:NDG:O	1.80	0.81
1:L:271:GLU:OE2	1:L:280:THR:HG21	1.81	0.80
1:I:22:ILE:HG23	1:I:115:THR:HG23	1.62	0.80
1:L:332:LYS:O	1:L:336:GLN:HG3	1.82	0.80
1:I:208:ASN:HD22	1:I:209:GLU:N	1.79	0.80
1:I:183:ARG:NE	1:I:204:SER:HA	1.97	0.79
1:L:208:ASN:HD22	1:L:393:ARG:NH1	1.81	0.78
4:L:841:NDG:C4	4:L:842:NDG:H1	2.13	0.78
1:L:155:ASN:ND2	4:L:841:NDG:O1L	2.14	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:155:ASN:ND2	4:L:841:NDG:H1	1.97	0.78
1:I:199:THR:HG22	1:I:200:ASP:N	1.99	0.78
3:D:9:TRP:CZ3	3:D:11:ASP:HB2	2.19	0.77
1:I:60:THR:O	1:I:64:GLN:HG3	1.84	0.77
1:I:152:LEU:CD1	1:I:212:VAL:HB	2.14	0.76
1:I:192:ASN:ND2	5:I:861:NAG:C5	2.49	0.76
1:I:294:LYS:O	1:I:298:GLU:HG3	1.86	0.75
1:L:57:ARG:HG2	1:L:301:PRO:HG2	1.68	0.75
1:L:17:MET:C	1:L:19:PRO:HD3	2.07	0.75
1:I:192:ASN:HD21	5:I:861:NAG:C5	1.99	0.74
1:I:22:ILE:HG23	1:I:115:THR:CG2	2.18	0.73
1:I:177:GLU:C	1:I:178:ASN:HD22	1.91	0.73
1:I:300:THR:HG21	1:I:302:GLU:HG2	1.69	0.73
1:I:300:THR:HG22	1:I:303:VAL:H	1.54	0.72
1:L:257:LYS:HG3	1:L:315:MET:HE2	1.71	0.72
1:L:96:ASN:ND2	4:L:801:NDG:O1L	2.22	0.72
1:L:71:ASN:HB3	1:L:74:ASP:OD2	1.89	0.72
1:I:192:ASN:HD21	5:I:861:NAG:H5	1.54	0.71
1:L:284:ILE:HD13	1:L:307:TRP:CZ3	2.25	0.71
5:I:841:NAG:O4	4:I:842:NDG:H1	1.90	0.71
1:L:183:ARG:CZ	1:L:204:SER:HA	2.20	0.71
1:L:133:LYS:O	1:L:136:LYS:HG3	1.91	0.71
1:L:188:LYS:HE2	4:L:861:NDG:H6C1	1.71	0.70
1:I:283:LEU:HD11	1:I:320:MET:HE3	1.73	0.70
1:L:241:LYS:HE3	1:L:247:CYS:SG	2.33	0.69
1:L:203:PRO:HB2	1:L:395:LEU:CD1	2.22	0.69
1:I:192:ASN:ND2	5:I:861:NAG:O5	2.26	0.69
1:L:209:GLU:CD	1:L:209:GLU:H	1.94	0.68
1:I:96:ASN:HD21	4:I:801:NDG:C5	2.06	0.68
1:I:44:THR:HG22	1:I:45:ASN:H	1.59	0.68
1:I:129:ARG:HB3	1:I:417:LEU:HD11	1.76	0.67
1:L:285:LEU:HD23	1:L:285:LEU:N	2.09	0.67
1:L:47:ARG:HG2	1:L:122:PHE:CZ	2.30	0.67
1:I:208:ASN:HD22	1:I:210:LEU:N	1.90	0.67
1:I:155:ASN:ND2	5:I:841:NAG:O1	2.28	0.67
1:I:319:HIS:HB2	1:I:403:LYS:HA	1.75	0.67
1:L:15:ILE:HG23	1:L:164:LEU:HD21	1.76	0.67
1:L:71:ASN:ND2	1:L:403:LYS:HE2	2.08	0.67
1:L:300:THR:CG2	1:L:302:GLU:HG2	2.25	0.67
1:L:7:ILE:CD1	1:L:164:LEU:HG	2.23	0.66
1:I:91:LYS:NZ	1:I:120:HIS:NE2	2.32	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:190:VAL:HG11	1:L:201:VAL:CG2	2.26	0.66
1:L:96:ASN:N	1:L:96:ASN:HD22	1.94	0.65
1:I:187:ASN:ND2	1:I:202:ILE:H	1.94	0.65
1:I:294:LYS:HA	1:I:294:LYS:HE3	1.78	0.65
1:I:96:ASN:HD21	4:I:801:NDG:H5	1.60	0.65
1:I:130:LEU:HD23	1:I:414:GLU:HG3	1.78	0.64
1:L:116:SER:O	1:L:119:ILE:HG22	1.97	0.64
1:I:356:ALA:HB1	5:I:841:NAG:H81	1.79	0.64
1:L:190:VAL:HG11	1:L:201:VAL:HG21	1.79	0.64
1:I:178:ASN:N	1:I:178:ASN:ND2	2.46	0.63
5:I:861:NAG:C4	4:I:862:NDG:H1L	2.10	0.63
1:I:215:LEU:HD23	1:I:216:VAL:N	2.13	0.63
1:I:115:THR:H	1:I:118:GLN:NE2	1.97	0.63
1:L:102:LEU:HD23	1:L:340:LEU:HD21	1.81	0.62
1:I:89:MET:HB3	1:I:215:LEU:HD11	1.81	0.61
1:L:401:THR:HG22	1:L:402:PHE:N	2.14	0.61
1:I:194:THR:O	1:I:197:ARG:HG2	2.01	0.61
1:L:362:LEU:HB3	1:L:390:ILE:CG2	2.31	0.61
1:L:70:LYS:HD2	1:L:76:ILE:HG12	1.83	0.60
1:I:378:GLU:OE1	1:I:384:ALA:HB3	2.01	0.60
4:L:841:NDG:C3	4:L:842:NDG:H1	2.32	0.60
1:L:284:ILE:HB	1:L:409:LEU:HB2	1.82	0.60
1:L:287:LYS:HG2	1:L:290:LYS:HB2	1.84	0.60
1:I:208:ASN:HD21	1:I:210:LEU:HB2	1.66	0.59
1:I:215:LEU:HD23	1:I:215:LEU:C	2.23	0.59
1:I:306:GLU:O	1:I:310:GLU:HG3	2.03	0.59
1:L:428:ASN:O	1:L:431:VAL:HG13	2.02	0.59
1:L:334:GLN:CD	1:L:334:GLN:H	2.06	0.59
1:I:208:ASN:ND2	1:I:210:LEU:N	2.47	0.59
1:L:108:PHE:HB3	1:L:119:ILE:HG13	1.84	0.58
1:L:257:LYS:HG3	1:L:315:MET:CE	2.33	0.58
1:I:205:GLU:OE2	1:I:206:ALA:HB2	2.04	0.58
1:I:45:ASN:HB3	1:I:48:VAL:HG13	1.85	0.58
1:I:155:ASN:CG	5:I:841:NAG:O1	2.42	0.58
1:L:91:LYS:HE2	1:L:103:MET:HE2	1.84	0.58
1:I:346:PRO:HG3	1:I:363:TYR:CE1	2.38	0.58
1:I:100:GLN:O	1:I:104:GLU:HG3	2.04	0.58
4:I:801:NDG:O1L	4:I:801:NDG:C7	2.51	0.58
1:L:232:GLU:H	1:L:232:GLU:CD	2.05	0.58
1:L:230:SER:HB2	1:L:233:ASN:ND2	2.19	0.58
1:L:284:ILE:C	1:L:285:LEU:HD23	2.24	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:19:PRO:HG3	1:I:22:ILE:HD11	1.86	0.58
1:I:115:THR:HG22	1:I:117:ASP:H	1.69	0.57
1:L:18:ASN:N	1:L:19:PRO:HD3	2.19	0.57
1:L:230:SER:HB2	1:L:233:ASN:HD22	1.69	0.57
1:L:19:PRO:HG2	1:L:117:ASP:HB2	1.86	0.57
1:L:6:ASP:HB3	1:L:9:THR:OG1	2.05	0.57
1:L:372:PHE:O	1:L:382:ALA:HA	2.05	0.57
1:I:180:GLU:OE2	1:I:183:ARG:NH1	2.38	0.57
1:I:44:THR:HG22	1:I:45:ASN:N	2.20	0.57
1:I:47:ARG:NH1	1:I:50:GLU:OE1	2.38	0.57
1:I:7:ILE:HG13	1:I:15:ILE:CD1	2.34	0.57
1:L:281:MET:HA	1:L:411:PHE:O	2.05	0.57
1:I:210:LEU:N	1:I:210:LEU:HD12	2.20	0.56
1:L:14:ASP:O	1:L:16:PRO:HD3	2.05	0.56
1:L:421:ILE:HG22	1:L:422:PHE:CD1	2.40	0.56
1:L:71:ASN:O	1:L:73:ASN:N	2.38	0.56
1:L:96:ASN:N	1:L:96:ASN:ND2	2.53	0.56
1:I:111:ILE:HG22	1:I:114:LYS:HG3	1.87	0.55
1:I:213:LEU:HB3	3:D:12:PHE:HB2	1.89	0.55
1:L:260:TYR:CG	1:L:261:ARG:N	2.74	0.55
1:I:253:TYR:HD1	1:I:319:HIS:CE1	2.24	0.55
1:L:23:TYR:CZ	1:L:100:GLN:HG3	2.42	0.55
1:I:183:ARG:HE	1:I:204:SER:HA	1.71	0.55
1:I:18:ASN:OD1	4:I:842:NDG:H8C2	2.06	0.55
1:L:258:PHE:CD1	1:L:316:LEU:HD21	2.42	0.55
1:L:203:PRO:HB2	1:L:395:LEU:HD11	1.89	0.55
1:I:208:ASN:ND2	1:I:210:LEU:HD13	2.22	0.54
1:I:283:LEU:HD11	1:I:320:MET:HE1	1.87	0.54
1:L:18:ASN:H	1:L:161:ILE:HD11	1.72	0.54
1:L:372:PHE:CD1	1:L:372:PHE:C	2.80	0.54
1:I:201:VAL:HG23	1:I:202:ILE:HG13	1.90	0.54
1:L:324:ARG:O	1:L:431:VAL:HG11	2.08	0.54
1:L:321:PRO:HD3	1:I:391:ALA:O	2.08	0.53
1:I:108:PHE:HB3	1:I:119:ILE:HG12	1.89	0.53
1:I:155:ASN:OD1	1:I:157:THR:N	2.37	0.53
1:I:179:ALA:HB1	1:I:207:ILE:HG22	1.89	0.53
1:L:401:THR:CG2	1:L:402:PHE:N	2.70	0.53
3:D:9:TRP:HZ3	3:D:11:ASP:HB2	1.66	0.53
1:L:229:PHE:O	1:L:377:GLU:HB3	2.08	0.53
1:I:22:ILE:HD12	1:I:115:THR:CG2	2.39	0.52
1:I:132:ARG:HG3	1:I:134:ALA:H	1.74	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:403:LYS:HB3	1:L:403:LYS:NZ	2.24	0.52
1:L:406:ARG:H	1:L:406:ARG:HE	1.58	0.52
1:I:300:THR:HG23	1:I:302:GLU:OE2	2.09	0.52
1:L:233:ASN:HB3	1:L:253:TYR:HB3	1.92	0.52
1:L:7:ILE:HD11	1:L:164:LEU:O	2.09	0.52
5:I:841:NAG:H62	4:I:842:NDG:H8C3	1.92	0.52
1:I:365:SER:HB3	3:D:11:ASP:OD2	2.10	0.52
1:L:300:THR:HG22	1:L:303:VAL:HG23	1.92	0.51
1:L:225:TRP:CD1	1:L:379:GLY:HA2	2.45	0.51
1:L:328:GLY:O	1:L:329:PHE:HB3	2.10	0.51
1:L:176:LYS:HE3	1:L:359:ARG:CZ	2.39	0.51
1:L:414:GLU:CD	1:L:416:PRO:HD2	2.31	0.51
1:L:286:PRO:HD3	1:L:292:LEU:HD13	1.93	0.51
1:L:399:ARG:HG2	1:L:400:VAL:N	2.22	0.51
1:L:241:LYS:HE2	1:L:402:PHE:HZ	1.75	0.51
1:L:97:ASP:HB2	1:L:342:ASP:OD2	2.11	0.51
1:L:300:THR:HG21	1:L:302:GLU:HG2	1.93	0.51
4:I:801:NDG:O1L	4:I:801:NDG:C8	2.58	0.51
1:L:407:PRO:HB3	1:L:427:ALA:HB2	1.92	0.51
1:I:102:LEU:HD23	1:I:340:LEU:HD11	1.93	0.51
1:L:91:LYS:HD2	1:L:103:MET:HE3	1.93	0.51
1:L:15:ILE:CG2	1:L:164:LEU:HD21	2.39	0.50
4:L:801:NDG:O7	4:L:801:NDG:H3	2.10	0.50
1:I:20:MET:CE	1:I:352:PRO:HB2	2.36	0.50
1:L:198:ILE:HG22	1:L:201:VAL:HG22	1.93	0.50
1:I:132:ARG:HD2	1:I:136:LYS:HG3	1.94	0.50
1:L:234:THR:HA	1:L:251:MET:O	2.12	0.50
1:I:46:ARG:O	1:I:50:GLU:HG3	2.12	0.50
1:L:276:GLY:O	1:L:277:ASP:HB2	2.12	0.50
1:I:293:ALA:O	1:I:297:LYS:HG3	2.12	0.50
1:I:206:ALA:HB1	3:D:9:TRP:HH2	1.77	0.50
1:I:124:ALA:HB2	1:I:165:VAL:HG13	1.94	0.49
1:I:210:LEU:H	1:I:210:LEU:HD12	1.75	0.49
1:I:346:PRO:HG3	1:I:363:TYR:CZ	2.48	0.49
1:L:324:ARG:HA	1:L:373:LEU:O	2.12	0.49
1:I:199:THR:HG22	1:I:200:ASP:H	1.76	0.49
1:I:178:ASN:HB3	1:I:181:GLN:HB3	1.94	0.49
1:L:421:ILE:HG22	1:L:422:PHE:CE1	2.48	0.48
1:I:15:ILE:HD12	1:I:121:PHE:HE1	1.78	0.48
1:L:125:LYS:O	1:L:128:CYS:HB2	2.13	0.48
1:L:302:GLU:CD	1:L:302:GLU:H	2.15	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:131:TYR:CE2	1:I:142:SER:HB2	2.48	0.48
1:I:45:ASN:CG	1:I:48:VAL:HG13	2.33	0.48
1:L:82:SER:HA	1:L:217:ASN:ND2	2.29	0.48
1:L:306:GLU:O	1:L:310:GLU:HG3	2.13	0.48
1:L:285:LEU:HD11	1:I:389:VAL:HG13	1.95	0.48
1:L:51:LEU:CD2	1:L:123:PHE:HA	2.44	0.48
1:I:47:ARG:NH1	1:I:114:LYS:HE3	2.28	0.48
1:L:23:TYR:CE2	1:L:100:GLN:HG3	2.48	0.48
1:I:286:PRO:HB3	1:I:295:VAL:HG21	1.96	0.48
1:L:317:VAL:HG21	1:I:228:LYS:HG3	1.96	0.48
1:I:260:TYR:CG	1:I:261:ARG:N	2.82	0.48
1:L:77:PHE:CE2	1:L:373:LEU:HB2	2.48	0.48
1:I:257:LYS:HG3	1:I:315:MET:CE	2.43	0.47
1:L:410:VAL:HG12	1:L:411:PHE:N	2.30	0.47
1:I:183:ARG:HE	1:I:203:PRO:C	2.17	0.47
1:I:252:MET:SD	1:I:377:GLU:HG3	2.54	0.47
1:L:300:THR:CG2	1:L:303:VAL:HG23	2.44	0.47
1:I:301:PRO:HD2	1:I:302:GLU:OE2	2.14	0.47
1:I:317:VAL:HB	1:I:401:THR:HG22	1.95	0.47
1:L:234:THR:O	1:L:235:ARG:HD2	2.14	0.47
1:L:334:GLN:O	1:L:338:MET:HG3	2.15	0.47
1:I:261:ARG:HB3	1:I:311:LEU:HD23	1.96	0.47
1:L:11:LYS:HB3	1:L:12:PRO:HD2	1.96	0.47
1:L:208:ASN:HD22	1:L:393:ARG:HH12	1.61	0.47
1:L:334:GLN:N	1:L:334:GLN:CD	2.68	0.47
1:L:343:LEU:HG	1:L:364:VAL:CG2	2.45	0.47
1:L:298:GLU:O	1:L:303:VAL:HG21	2.14	0.47
1:L:178:ASN:HB3	1:L:181:GLN:HB2	1.96	0.47
1:L:292:LEU:CD2	1:L:407:PRO:HG2	2.44	0.47
1:L:101:GLN:O	1:L:105:VAL:HG23	2.15	0.47
1:L:183:ARG:NE	1:L:204:SER:HA	2.29	0.47
1:I:147:PHE:HB2	1:I:214:VAL:HB	1.97	0.47
1:L:141:VAL:CG2	1:L:220:TYR:HB3	2.45	0.47
1:L:22:ILE:HG22	1:L:23:TYR:N	2.30	0.47
1:L:227:SER:HB2	1:L:254:GLN:HE22	1.79	0.47
1:I:202:ILE:HD13	1:I:207:ILE:HD11	1.96	0.46
1:I:60:THR:OG1	1:I:301:PRO:HG3	2.16	0.46
1:L:152:LEU:HD22	1:L:358:GLY:HA3	1.97	0.46
1:I:183:ARG:HE	1:I:204:SER:N	2.14	0.46
1:L:77:PHE:CZ	1:L:373:LEU:HB2	2.50	0.46
4:L:841:NDG:H6C2	4:L:842:NDG:H6C1	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:129:ARG:CB	1:I:417:LEU:HD11	2.45	0.46
1:L:60:THR:O	1:L:64:GLN:HG3	2.16	0.46
1:I:45:ASN:CB	1:I:48:VAL:HG13	2.45	0.46
1:L:336:GLN:HA	1:L:340:LEU:O	2.16	0.46
1:I:22:ILE:HD11	1:I:117:ASP:HB3	1.96	0.46
1:L:332:LYS:HD3	1:L:344:PHE:CD2	2.51	0.46
1:I:22:ILE:CD1	1:I:117:ASP:HB3	2.46	0.46
1:I:210:LEU:H	1:I:210:LEU:CD1	2.29	0.46
1:L:91:LYS:HE2	1:L:103:MET:CE	2.46	0.46
1:L:179:ALA:HB1	1:L:207:ILE:HG22	1.98	0.46
1:L:7:ILE:CD1	1:L:164:LEU:O	2.64	0.46
1:I:132:ARG:HD2	1:I:136:LYS:HE3	1.98	0.45
1:I:192:ASN:ND2	5:I:861:NAG:H61	2.30	0.45
1:L:51:LEU:HA	1:L:111:ILE:HD11	1.98	0.45
1:L:7:ILE:HD11	1:L:164:LEU:C	2.37	0.45
1:I:22:ILE:HD12	1:I:115:THR:HG21	1.96	0.45
1:L:91:LYS:HE3	1:L:120:HIS:CE1	2.51	0.45
1:I:192:ASN:ND2	5:I:861:NAG:C6	2.79	0.45
1:L:284:ILE:HD12	1:L:411:PHE:HE1	1.82	0.45
1:L:414:GLU:OE1	1:L:416:PRO:HD2	2.16	0.45
1:I:23:TYR:CE1	1:I:100:GLN:HG3	2.51	0.45
1:I:358:GLY:O	1:I:359:ARG:HB3	2.16	0.45
1:L:300:THR:HG23	1:L:302:GLU:HG2	1.98	0.45
1:I:379:GLY:O	1:I:380:SER:C	2.55	0.45
1:L:285:LEU:HD11	1:I:389:VAL:CG1	2.47	0.45
1:L:194:THR:O	1:L:197:ARG:HB2	2.16	0.45
1:I:51:LEU:HA	1:I:111:ILE:HD11	1.98	0.44
5:I:861:NAG:C4	4:I:862:NDG:O1L	2.64	0.44
1:L:258:PHE:CD1	1:L:316:LEU:CD2	3.00	0.44
1:I:155:ASN:ND2	5:I:841:NAG:C1	2.80	0.44
1:L:197:ARG:HD3	1:L:381:GLU:OE2	2.18	0.44
1:L:430:CYS:O	1:L:431:VAL:C	2.55	0.44
1:I:284:ILE:O	1:I:408:PHE:HB2	2.18	0.44
1:L:131:TYR:CE2	1:L:142:SER:HB2	2.53	0.44
1:I:206:ALA:O	1:I:207:ILE:HB	2.18	0.44
1:I:57:ARG:HG2	1:I:107:LYS:HG3	2.00	0.44
1:L:217:ASN:HD21	1:L:219:ILE:HG13	1.82	0.44
1:I:77:PHE:CZ	1:I:422:PHE:HB3	2.52	0.44
1:L:197:ARG:NH1	1:L:381:GLU:OE1	2.51	0.44
1:I:155:ASN:OD1	1:I:155:ASN:C	2.56	0.44
1:L:176:LYS:HE3	1:L:359:ARG:NH1	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:365:SER:O	1:L:366:ASP:HB2	2.17	0.43
1:L:229:PHE:HB2	1:L:377:GLU:HA	2.00	0.43
1:L:22:ILE:HD13	1:L:117:ASP:OD2	2.17	0.43
1:L:329:PHE:HE1	1:L:331:LEU:HD23	1.83	0.43
1:I:15:ILE:HD12	1:I:121:PHE:CE1	2.53	0.43
1:I:187:ASN:HD21	1:I:202:ILE:N	2.01	0.43
1:L:241:LYS:HE2	1:L:402:PHE:CZ	2.53	0.43
1:I:287:LYS:HB2	1:I:288:PRO:HD2	2.01	0.43
1:L:198:ILE:CG2	1:L:201:VAL:HG22	2.48	0.43
5:I:841:NAG:O4	4:I:842:NDG:O7	2.35	0.43
1:I:366:ASP:HB3	1:I:368:PHE:CE2	2.54	0.43
2:C:8:THR:OXT	3:D:9:TRP:CA	2.67	0.43
1:L:343:LEU:HG	1:L:364:VAL:HG23	2.01	0.43
1:I:77:PHE:CE2	1:I:373:LEU:HB2	2.54	0.42
1:I:257:LYS:HA	1:I:314:MET:O	2.19	0.42
1:L:258:PHE:HD1	1:L:316:LEU:CD2	2.33	0.42
1:I:96:ASN:ND2	4:I:801:NDG:C5	2.78	0.42
1:L:46:ARG:O	1:L:50:GLU:HG3	2.20	0.42
1:L:71:ASN:C	1:L:73:ASN:H	2.23	0.42
1:I:253:TYR:CD2	1:I:253:TYR:C	2.93	0.42
1:I:359:ARG:O	1:I:359:ARG:CG	2.68	0.42
1:L:202:ILE:HA	1:L:203:PRO:HD3	1.89	0.42
1:L:91:LYS:CE	1:L:103:MET:CE	2.97	0.42
1:I:23:TYR:HB3	1:I:116:SER:HB3	2.01	0.41
1:I:399:ARG:HH11	1:I:399:ARG:HG2	1.85	0.41
1:I:406:ARG:HD2	6:I:873:HOH:O	2.19	0.41
1:L:131:TYR:CZ	1:L:142:SER:HB2	2.55	0.41
1:I:314:MET:SD	1:I:400:VAL:HG11	2.60	0.41
1:I:51:LEU:O	1:I:51:LEU:HG	2.20	0.41
1:I:356:ALA:O	1:I:357:GLU:CB	2.69	0.41
1:I:183:ARG:HE	1:I:204:SER:CA	2.33	0.41
1:L:240:TYR:HA	1:L:246:SER:OG	2.20	0.41
1:I:51:LEU:CD2	1:I:123:PHE:HA	2.51	0.41
1:L:287:LYS:HG2	1:L:290:LYS:HG3	2.03	0.41
1:L:406:ARG:HE	1:L:406:ARG:N	2.18	0.41
1:L:76:ILE:HA	1:L:327:ASP:OD2	2.21	0.41
1:I:362:LEU:HD21	3:D:12:PHE:HB3	2.02	0.41
1:L:317:VAL:HB	1:I:388:VAL:HG22	2.02	0.41
1:I:279:ILE:HA	1:I:413:ARG:O	2.20	0.41
1:I:71:ASN:HB3	1:I:74:ASP:OD2	2.21	0.41
1:L:119:ILE:HD13	1:L:119:ILE:C	2.41	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:171:GLN:HA	1:I:172:PRO:HD3	1.90	0.41
1:L:119:ILE:HG23	1:L:120:HIS:N	2.36	0.41
1:L:308:LEU:HD13	1:L:413:ARG:NH2	2.35	0.41
1:I:205:GLU:CD	1:I:206:ALA:N	2.74	0.41
1:I:330:SER:HB2	1:I:368:PHE:CE1	2.56	0.41
1:L:369:HIS:HA	1:L:385:SER:O	2.21	0.41
1:I:197:ARG:HD3	2:C:3:GLU:OE2	2.20	0.40
1:L:271:GLU:HG3	1:L:282:VAL:HG22	2.03	0.40
1:L:287:LYS:HG2	1:L:290:LYS:CB	2.48	0.40
1:L:410:VAL:HG21	1:L:426:VAL:HG23	2.03	0.40
1:L:406:ARG:HE	1:L:406:ARG:CA	2.34	0.40
1:L:78:LEU:HD22	1:L:78:LEU:C	2.41	0.40
2:C:8:THR:OXT	3:D:9:TRP:N	2.54	0.40
1:I:201:VAL:HG23	1:I:202:ILE:N	2.36	0.40
1:L:322:ARG:NH1	1:L:376:ASN:HB2	2.36	0.40
1:L:197:ARG:HG3	1:L:372:PHE:CZ	2.56	0.40
1:L:91:LYS:HE3	1:L:91:LYS:HB3	1.81	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	I	404/432 (94%)	378 (94%)	18 (4%)	8 (2%)	9	17
1	L	402/432 (93%)	363 (90%)	33 (8%)	6 (2%)	12	24
2	C	6/8 (75%)	6 (100%)	0	0	100	100
3	D	3/5 (60%)	3 (100%)	0	0	100	100
All	All	815/877 (93%)	750 (92%)	51 (6%)	14 (2%)	11	21

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	I	43	ALA
1	I	361	ASP
1	L	72	ASP
1	L	402	PHE
1	I	112	SER
1	L	397	PRO
1	I	206	ALA
1	L	135	ASN
1	L	400	VAL
1	I	113	GLU
1	I	199	THR
1	I	207	ILE
1	I	263	VAL
1	L	201	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	I	366/383 (96%)	349 (95%)	17 (5%)	31	58
1	L	361/383 (94%)	330 (91%)	31 (9%)	12	23
2	C	4/4 (100%)	4 (100%)	0	100	100
3	D	4/4 (100%)	4 (100%)	0	100	100
All	All	735/774 (95%)	687 (94%)	48 (6%)	20	39

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

Mol	Chain	Res	Type
1	L	9	THR
1	L	47	ARG
1	L	78	LEU
1	L	96	ASN
1	L	119	ILE
1	L	123	PHE
1	L	156	GLU
1	L	173	LEU

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Mol	Chain	Res	Type
1	L	177	GLU
1	L	197	ARG
1	L	199	THR
1	L	201	VAL
1	L	227	SER
1	L	245	GLU
1	L	285	LEU
1	L	300	THR
1	L	302	GLU
1	L	315	MET
1	L	316	LEU
1	L	320	MET
1	L	322	ARG
1	L	324	ARG
1	L	334	GLN
1	L	340	LEU
1	L	361	ASP
1	L	362	LEU
1	L	372	PHE
1	L	399	ARG
1	L	402	PHE
1	L	405	ASN
1	L	406	ARG
1	I	20	MET
1	I	48	VAL
1	I	78	LEU
1	I	123	PHE
1	I	170	LEU
1	I	178	ASN
1	I	197	ARG
1	I	208	ASN
1	I	241	LYS
1	I	294	LYS
1	I	300	THR
1	I	305	GLN
1	I	307	TRP
1	I	316	LEU
1	I	362	LEU
1	I	366	ASP
1	I	418	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	45	ASN
1	L	65	HIS
1	L	96	ASN
1	L	135	ASN
1	L	155	ASN
1	L	208	ASN
1	L	217	ASN
1	L	233	ASN
1	L	254	GLN
1	L	336	GLN
1	L	405	ASN
1	L	428	ASN
1	I	96	ASN
1	I	118	GLN
1	I	127	ASN
1	I	178	ASN
1	I	187	ASN
1	I	192	ASN
1	I	208	ASN
1	I	217	ASN
1	I	254	GLN
1	I	305	GLN
1	I	319	HIS
1	I	336	GLN
1	I	418	ASN

### 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](#)

9 ligands 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	$\# Z  > 2$
4	NDG	I	801	-	15,15,15	0.86	1 (6%)	21,21,21	1.42	2 (9%)
5	NAG	I	841	-	15,15,15	0.31	0	21,21,21	0.70	0
4	NDG	I	842	-	15,15,15	0.47	0	21,21,21	0.98	2 (9%)
5	NAG	I	861	-	15,15,15	0.36	0	21,21,21	0.89	0
4	NDG	I	862	-	15,15,15	0.35	0	21,21,21	0.66	0
4	NDG	L	801	-	15,15,15	0.60	0	21,21,21	0.63	0
4	NDG	L	841	-	15,15,15	0.67	0	21,21,21	1.11	2 (9%)
4	NDG	L	842	-	15,15,15	0.45	0	21,21,21	0.70	0
4	NDG	L	861	-	15,15,15	0.46	0	21,21,21	0.58	0

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
4	NDG	I	801	-	-	0/6/26/26	0/1/1/1
5	NAG	I	841	-	-	0/6/26/26	0/1/1/1
4	NDG	I	842	-	-	0/6/26/26	0/1/1/1
5	NAG	I	861	-	-	0/6/26/26	0/1/1/1
4	NDG	I	862	-	-	0/6/26/26	0/1/1/1
4	NDG	L	801	-	-	0/6/26/26	0/1/1/1
4	NDG	L	841	-	-	0/6/26/26	0/1/1/1
4	NDG	L	842	-	-	0/6/26/26	0/1/1/1
4	NDG	L	861	-	-	0/6/26/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	I	801	NDG	C1-C2	2.61	1.56	1.52

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	I	842	NDG	C1-C2-N2	-2.06	108.34	110.73
4	I	842	NDG	O-C1-C2	2.32	111.85	109.52
4	L	841	NDG	C3-C4-C5	2.71	115.00	110.22
4	L	841	NDG	C4-C3-C2	3.09	114.91	110.33
4	I	801	NDG	C1-C2-C3	3.47	115.28	110.54
4	I	801	NDG	O-C1-C2	4.24	113.77	109.52

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

9 monomers are involved in 34 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	I	801	NDG	5	0
5	I	841	NAG	7	0
4	I	842	NDG	4	0
5	I	861	NAG	8	0
4	I	862	NDG	2	0
4	L	801	NDG	2	0
4	L	841	NDG	10	0
4	L	842	NDG	6	0
4	L	861	NDG	1	0

## 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 [i](#)

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

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	I	410/432 (94%)	-0.09	9 (2%) 62 56	28, 48, 95, 199	0
1	L	406/432 (93%)	0.24	33 (8%) 13 8	30, 55, 120, 168	0
2	C	7/8 (87%)	-0.21	0 100 100	27, 37, 50, 51	0
3	D	4/5 (80%)	-0.15	0 100 100	49, 54, 56, 60	0
All	All	827/877 (94%)	0.07	42 (5%) 29 22	27, 52, 107, 199	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	I	358	GLY	8.6
1	L	402	PHE	5.9
1	L	242	ALA	5.4
1	I	360	ASP	5.3
1	L	404	ALA	4.6
1	I	380	SER	4.3
1	L	358	GLY	4.2
1	L	240	TYR	4.1
1	L	14	ASP	4.0
1	L	431	VAL	4.0
1	L	114	LYS	3.9
1	L	359	ARG	3.8
1	L	357	GLU	3.7
1	I	132	ARG	3.7
1	L	243	ASP	3.5
1	L	401	THR	3.4
1	L	13	ARG	3.3
1	I	357	GLU	3.2
1	L	18	ASN	3.0
1	L	400	VAL	3.0
1	L	15	ILE	3.0

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Mol	Chain	Res	Type	RSRZ
1	I	136	LYS	2.9
1	L	360	ASP	2.8
1	L	10	ALA	2.8
1	L	398	ASN	2.7
1	L	135	ASN	2.6
1	L	71	ASN	2.6
1	I	135	ASN	2.5
1	L	403	LYS	2.5
1	L	46	ARG	2.4
1	L	134	ALA	2.3
1	L	115	THR	2.3
1	I	134	ALA	2.3
1	L	11	LYS	2.3
1	I	361	ASP	2.3
1	L	47	ARG	2.2
1	L	396	ASN	2.2
1	L	399	ARG	2.2
1	L	246	SER	2.1
1	L	289	GLU	2.1
1	L	245	GLU	2.1
1	L	241	LYS	2.0

## 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(Å <sup>2</sup> )	Q<0.9
4	NDG	L	841	15/15	0.86	0.35	1.66	62,90,100,109	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
5	NAG	I	841	15/15	0.94	0.15	-0.49	44,60,74,82	0
4	NDG	L	861	15/15	0.58	0.42	-	117,151,163,173	0
4	NDG	L	842	15/15	0.82	0.41	-	77,129,132,135	0
4	NDG	L	801	15/15	0.73	0.32	-	64,114,124,133	0
5	NAG	I	861	15/15	0.75	0.28	-	40,112,119,127	0
4	NDG	I	862	15/15	0.87	0.41	-	61,151,155,156	0
4	NDG	I	842	15/15	0.59	0.34	-	88,131,135,136	0
4	NDG	I	801	15/15	0.72	0.38	-	61,115,124,131	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.