



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 06:20 PM GMT

PDB ID : 1ABJ  
Title : STRUCTURE OF THE HIRULOG 3-THROMBIN COMPLEX AND NATURE OF THE S' SUBSITES OF SUBSTRATES AND INHIBITORS  
Authors : Qiu, X.; Tulinsky, A.  
Deposited on : 1992-08-24  
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) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

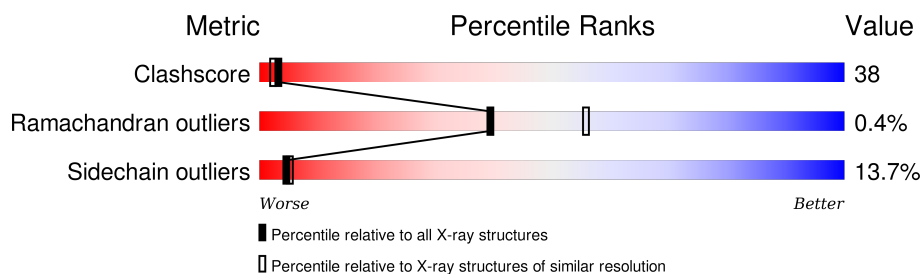
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 Å.

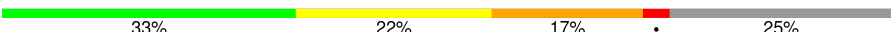

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



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)

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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	L	36	
2	H	259	

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 criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	OG6	H	1	X	-	-	-

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 2531 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.

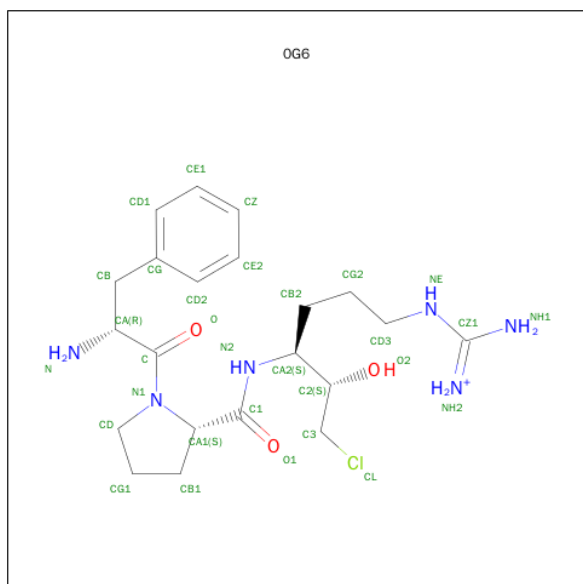
- Molecule 1 is a protein called ALPHA-THROMBIN (SMALL SUBUNIT).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	27	Total	C	N	O	S	0	0	0
			222	140	36	45	1			

- Molecule 2 is a protein called ALPHA-THROMBIN (LARGE SUBUNIT).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	258	Total	C	N	O	S	0	0	0
			2083	1329	369	371	14			

- Molecule 3 is D-PHENYLALANYL-N-[(2S,3S)-6-{[AMINO(IMINIO)METHYL]AMINO}-1-CHLORO-2-HYDROXYHEXAN-3-YL]-L-PROLINAMIDE (three-letter code: 0G6) (formula: C<sub>21</sub>H<sub>34</sub>ClN<sub>6</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	H	1	Total	C	N	O	0	0
			30	21	6	3		

- Molecule 4 is water.


Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	L	14	Total 14	O 14	0	0
4	H	182	Total 182	O 182	0	0

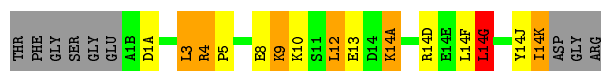
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

Note EDS was not executed.

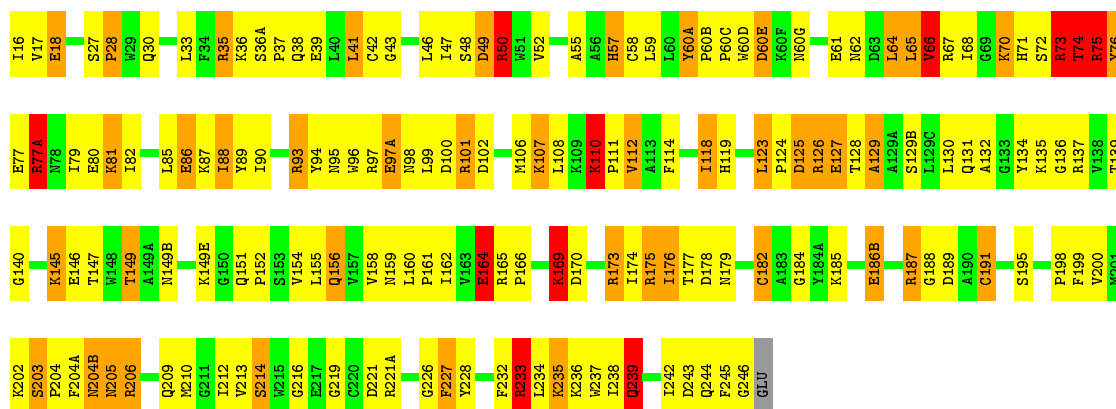
#### • Molecule 1: ALPHA-THROMBIN (SMALL SUBUNIT)

Chain L: 



#### • Molecule 2: ALPHA-THROMBIN (LARGE SUBUNIT)

Chain H: 



## 4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	88.26Å 68.02Å 61.22Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	7.00 – 2.40	Depositor
% Data completeness (in resolution range)	(Not available) (7.00-2.40)	Depositor
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	PROLSQ	Depositor
R, $R_{free}$	0.144 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2531	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	28.0	wwPDB-VP

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 0G6

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	0.97	0/224	2.12	7/298 (2.3%)
2	H	1.03	1/2138 (0.0%)	2.14	77/2891 (2.7%)
All	All	1.03	1/2362 (0.0%)	2.14	84/3189 (2.6%)

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	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	214	SER	CA-CB	6.75	1.63	1.52

All (84) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	137	ARG	NE-CZ-NH1	17.37	128.99	120.30
2	H	221(A)	ARG	NE-CZ-NH2	-17.18	111.71	120.30
2	H	187	ARG	NE-CZ-NH1	15.71	128.15	120.30
2	H	187	ARG	NE-CZ-NH2	-13.77	113.42	120.30
2	H	35	ARG	NE-CZ-NH1	13.15	126.88	120.30
2	H	73	ARG	NE-CZ-NH2	-12.87	113.87	120.30
2	H	125	ASP	CB-CG-OD1	12.36	129.43	118.30
1	L	4	ARG	NE-CZ-NH1	11.82	126.21	120.30
2	H	233	ARG	NE-CZ-NH1	11.24	125.92	120.30
2	H	101	ARG	NE-CZ-NH2	-10.59	115.01	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	50	ARG	NE-CZ-NH1	10.38	125.49	120.30
2	H	35	ARG	NE-CZ-NH2	-10.18	115.21	120.30
2	H	235	LYS	CA-CB-CG	10.15	135.74	113.40
2	H	125	ASP	CB-CG-OD2	-9.66	109.61	118.30
2	H	243	ASP	CB-CG-OD2	9.46	126.82	118.30
2	H	67	ARG	NE-CZ-NH2	9.26	124.93	120.30
2	H	233	ARG	CD-NE-CZ	8.99	136.18	123.60
2	H	66	VAL	CB-CA-C	8.13	126.86	111.40
2	H	77(A)	ARG	NE-CZ-NH2	-8.04	116.28	120.30
2	H	149	THR	N-CA-CB	7.93	125.37	110.30
2	H	101	ARG	CD-NE-CZ	-7.55	113.03	123.60
1	L	1(A)	ASP	CB-CG-OD1	-7.44	111.60	118.30
2	H	175	ARG	NE-CZ-NH1	-7.27	116.66	120.30
2	H	57	HIS	N-CA-CB	7.27	123.69	110.60
2	H	73	ARG	NE-CZ-NH1	7.10	123.85	120.30
2	H	110	LYS	N-CA-CB	7.05	123.28	110.60
2	H	127	GLU	OE1-CD-OE2	7.05	131.76	123.30
2	H	137	ARG	NE-CZ-NH2	-6.81	116.89	120.30
2	H	27	SER	CB-CA-C	6.76	122.94	110.10
2	H	164	GLU	OE1-CD-OE2	-6.67	115.30	123.30
2	H	221(A)	ARG	NH1-CZ-NH2	6.66	126.72	119.40
2	H	49	ASP	CB-CG-OD1	-6.66	112.31	118.30
1	L	3	LEU	CA-CB-CG	6.50	130.25	115.30
1	L	14(D)	ARG	NE-CZ-NH2	-6.42	117.09	120.30
2	H	60(E)	ASP	CB-CG-OD1	-6.42	112.53	118.30
2	H	60(A)	TYR	CB-CG-CD1	-6.37	117.18	121.00
2	H	227	PHE	C-N-CA	6.19	137.18	121.70
1	L	14(G)	LEU	CB-CA-C	6.10	121.79	110.20
2	H	186(B)	GLU	OE1-CD-OE2	-6.03	116.07	123.30
2	H	137	ARG	NH1-CZ-NH2	-5.98	112.83	119.40
2	H	170	ASP	CB-CG-OD1	5.91	123.62	118.30
2	H	175	ARG	CD-NE-CZ	-5.91	115.33	123.60
2	H	191	CYS	CA-CB-SG	5.91	124.63	114.00
2	H	242	ILE	CA-CB-CG2	5.84	122.58	110.90
2	H	118	ILE	O-C-N	5.81	132.00	122.70
2	H	38	GLN	OE1-CD-NE2	5.79	135.22	121.90
2	H	97(A)	GLU	OE1-CD-OE2	5.79	130.25	123.30
2	H	60(A)	TYR	CB-CG-CD2	5.73	124.44	121.00
2	H	129	ALA	C-N-CA	5.71	135.96	121.70
2	H	43	GLY	N-CA-C	-5.70	98.84	113.10
2	H	206	ARG	CD-NE-CZ	5.65	131.51	123.60
2	H	73	ARG	CA-C-O	5.63	131.91	120.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	234	LEU	CB-CA-C	5.62	120.88	110.20
2	H	137	ARG	CD-NE-CZ	5.58	131.41	123.60
2	H	76	TYR	CB-CG-CD1	-5.54	117.68	121.00
2	H	41	LEU	CA-CB-CG	5.50	127.95	115.30
2	H	35	ARG	CB-CA-C	-5.47	99.46	110.40
2	H	74	THR	N-CA-CB	-5.44	99.96	110.30
2	H	246	GLY	CA-C-O	-5.42	110.84	120.60
2	H	86	GLU	CG-CD-OE2	-5.42	107.47	118.30
2	H	147	THR	N-CA-CB	-5.41	100.03	110.30
2	H	209	GLN	N-CA-CB	5.41	120.33	110.60
2	H	239	GLN	OE1-CD-NE2	5.38	134.26	121.90
2	H	149(E)	LYS	CA-CB-CG	5.37	125.21	113.40
2	H	123	LEU	CB-CA-C	5.36	120.38	110.20
2	H	137	ARG	O-C-N	5.35	131.26	122.70
2	H	158	VAL	CB-CA-C	5.33	121.54	111.40
2	H	75	ARG	NE-CZ-NH1	5.32	122.96	120.30
2	H	203	SER	N-CA-CB	5.30	118.46	110.50
2	H	169	LYS	O-C-N	5.30	131.18	122.70
2	H	97(A)	GLU	CG-CD-OE2	-5.28	107.73	118.30
2	H	62	ASN	CA-CB-CG	-5.28	101.78	113.40
2	H	86	GLU	CG-CD-OE1	5.28	128.86	118.30
1	L	3	LEU	N-CA-CB	5.23	120.86	110.40
2	H	64	LEU	CA-C-N	-5.21	105.75	117.20
2	H	93	ARG	NE-CZ-NH2	-5.20	117.70	120.30
1	L	14(D)	ARG	NE-CZ-NH1	5.20	122.90	120.30
2	H	156	GLN	CA-CB-CG	5.10	124.62	113.40
2	H	18	GLU	O-C-N	5.07	131.81	123.20
2	H	97	ARG	NE-CZ-NH1	5.06	122.83	120.30
2	H	173	ARG	NE-CZ-NH2	5.04	122.82	120.30
2	H	88	ILE	O-C-N	5.02	130.73	122.70
2	H	233	ARG	NH1-CZ-NH2	-5.01	113.88	119.40
2	H	67	ARG	CB-CA-C	5.00	120.41	110.40

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	H	126	ARG	Sidechain
2	H	73	ARG	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	222	0	225	19	0
2	H	2083	0	2058	160	0
3	H	30	0	30	5	0
4	H	182	0	0	12	0
4	L	14	0	0	2	0
All	All	2531	0	2313	173	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 38.

All (173) 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:5:PRO:HA	1:L:9:LYS:HB2	1.19	1.18
2:H:195:SER:HG	3:H:1:0G6:C2	1.62	1.07
2:H:50:ARG:NH1	2:H:108:LEU:O	1.92	1.01
2:H:195:SER:OG	3:H:1:0G6:C3	2.09	1.01
2:H:126:ARG:NH1	2:H:126:ARG:HB3	1.84	0.93
2:H:60(D):TRP:O	2:H:60(E):ASP:HB2	1.70	0.92
2:H:73:ARG:HD3	2:H:152:PRO:O	1.69	0.91
2:H:50:ARG:NH1	2:H:107:LYS:HE2	1.87	0.90
1:L:5:PRO:HA	1:L:9:LYS:CB	2.01	0.90
2:H:49:ASP:HB3	2:H:114:PHE:CZ	2.08	0.89
2:H:74:THR:HG23	2:H:75:ARG:HD3	1.56	0.88
2:H:205:ASN:O	2:H:205:ASN:ND2	2.07	0.88
2:H:60(G):ASN:ND2	4:H:528:HOH:O	2.05	0.86
2:H:205:ASN:HD22	2:H:205:ASN:C	1.72	0.86
1:L:10:LYS:HB3	1:L:12:LEU:HD22	1.57	0.86
1:L:5:PRO:CA	1:L:9:LYS:HB2	2.06	0.86
2:H:57:HIS:CD2	3:H:1:0G6:C3	2.58	0.85
2:H:93:ARG:O	2:H:101:ARG:HD2	1.78	0.83
1:L:10:LYS:CB	1:L:12:LEU:HD22	2.10	0.82
2:H:136:GLY:HA3	2:H:199:PHE:CZ	2.14	0.81
2:H:126:ARG:CZ	2:H:126:ARG:HB3	2.09	0.81
1:L:8:GLU:OE2	2:H:202:LYS:HE3	1.81	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:75:ARG:HG2	2:H:75:ARG:NH1	1.98	0.78
2:H:75:ARG:HG2	2:H:75:ARG:HH11	1.47	0.77
1:L:14(A):LYS:NZ	1:L:14(A):LYS:HB2	2.00	0.77
2:H:125:ASP:OD1	2:H:128:THR:HB	1.84	0.77
2:H:205:ASN:ND2	2:H:205:ASN:C	2.38	0.76
2:H:164:GLU:OE1	2:H:164:GLU:N	2.13	0.76
2:H:169:LYS:HD3	4:H:559:HOH:O	1.87	0.75
2:H:74:THR:CG2	2:H:75:ARG:HD3	2.17	0.75
2:H:76:TYR:CE2	2:H:77(A):ARG:HA	2.23	0.73
2:H:87:LYS:HG2	2:H:89:TYR:CZ	2.24	0.73
2:H:239:GLN:NE2	4:H:471:HOH:O	2.22	0.73
2:H:49:ASP:HB3	2:H:114:PHE:CE1	2.24	0.72
2:H:74:THR:HG23	2:H:75:ARG:CD	2.20	0.71
2:H:60(A):TYR:CZ	2:H:60(C):PRO:HG2	2.27	0.70
2:H:236:LYS:HB2	4:H:436:HOH:O	1.91	0.69
2:H:126:ARG:CB	2:H:126:ARG:NH1	2.55	0.69
1:L:3:LEU:HD12	1:L:9:LYS:NZ	2.09	0.67
2:H:71:HIS:CE1	2:H:154:VAL:HG13	2.30	0.67
2:H:195:SER:CB	3:H:1:0G6:C3	2.73	0.66
2:H:187:ARG:NH2	2:H:221:ASP:O	2.27	0.66
2:H:165:ARG:HB2	2:H:166:PRO:HD3	1.78	0.66
2:H:126:ARG:HG2	2:H:232:PHE:HZ	1.60	0.66
2:H:80:GLU:O	2:H:81:LYS:HD2	1.98	0.64
1:L:14(A):LYS:HZ3	1:L:14(A):LYS:HB2	1.59	0.64
1:L:14(F):LEU:HD11	2:H:159:ASN:HB3	1.80	0.64
2:H:188:GLY:O	2:H:189:ASP:HB2	1.98	0.63
2:H:70:LYS:HE3	2:H:72:SER:O	1.98	0.62
2:H:50:ARG:O	2:H:107:LYS:HA	1.99	0.62
2:H:75:ARG:CG	2:H:75:ARG:HH11	2.12	0.62
2:H:136:GLY:HA3	2:H:199:PHE:CE1	2.33	0.62
1:L:3:LEU:HD12	1:L:9:LYS:HZ1	1.63	0.62
2:H:94:TYR:HA	2:H:101:ARG:HB2	1.82	0.61
2:H:125:ASP:OD2	2:H:127:GLU:HB3	2.01	0.61
2:H:203:SER:HB3	2:H:204(B):ASN:ND2	2.15	0.61
2:H:164:GLU:C	2:H:166:PRO:HD2	2.22	0.60
2:H:30:GLN:O	4:H:419:HOH:O	2.15	0.59
2:H:90:ILE:HD12	4:H:525:HOH:O	2.01	0.59
2:H:49:ASP:N	2:H:49:ASP:OD1	2.33	0.58
2:H:216:GLY:O	3:H:1:0G6:N	2.37	0.58
2:H:127:GLU:HA	2:H:127:GLU:OE1	2.03	0.58
2:H:135:LYS:HA	2:H:161:PRO:HA	1.85	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:176:ILE:N	2:H:176:ILE:HD13	2.18	0.57
2:H:213:VAL:HG13	4:H:409:HOH:O	2.04	0.57
1:L:13:GLU:OE1	4:L:442:HOH:O	2.17	0.57
1:L:8:GLU:OE2	2:H:202:LYS:CE	2.50	0.57
2:H:47:ILE:O	2:H:48:SER:HB3	2.04	0.57
2:H:55:ALA:O	2:H:58:CYS:HB2	2.03	0.57
1:L:14(J):TYR:O	1:L:14(K):ILE:C	2.44	0.56
2:H:85:LEU:CD2	2:H:106:MET:HB3	2.36	0.56
1:L:9:LYS:HB3	4:L:484:HOH:O	2.05	0.56
2:H:165:ARG:N	2:H:166:PRO:CD	2.68	0.56
2:H:17:VAL:O	2:H:18:GLU:HB2	2.06	0.55
2:H:49:ASP:HB2	2:H:112:VAL:O	2.07	0.55
2:H:60(B):PRO:N	2:H:60(C):PRO:HD2	2.22	0.55
2:H:39:GLU:HG3	2:H:41:LEU:CD1	2.37	0.55
2:H:176:ILE:HG22	2:H:177:THR:N	2.22	0.54
2:H:130:LEU:HG	2:H:210:MET:HE2	1.89	0.54
2:H:33:LEU:HD23	2:H:66:VAL:HG23	1.88	0.54
2:H:59:LEU:HD21	2:H:106:MET:HE1	1.90	0.54
2:H:160:LEU:HD22	2:H:184:GLY:HA2	1.90	0.54
2:H:77(A):ARG:HG2	2:H:77(A):ARG:HH11	1.73	0.54
2:H:198:PRO:HB2	2:H:200:VAL:HG13	1.90	0.54
2:H:146:GLU:HG3	4:H:521:HOH:O	2.08	0.53
2:H:185:LYS:H	2:H:186(B):GLU:HG3	1.74	0.53
2:H:164:GLU:CD	2:H:164:GLU:H	2.07	0.53
2:H:140:GLY:O	2:H:155:LEU:HD12	2.09	0.53
2:H:60(A):TYR:CE1	2:H:60(C):PRO:HG2	2.45	0.52
2:H:100:ASP:O	2:H:101:ARG:HB2	2.09	0.52
2:H:85:LEU:HD22	2:H:106:MET:HB3	1.91	0.52
2:H:49:ASP:O	2:H:111:PRO:HA	2.09	0.51
2:H:17:VAL:O	2:H:188:GLY:HA2	2.10	0.51
2:H:16:ILE:N	4:H:403:HOH:O	2.44	0.51
2:H:60(B):PRO:N	2:H:60(C):PRO:CD	2.74	0.50
2:H:87:LYS:HG3	2:H:88:ILE:N	2.27	0.49
2:H:60(B):PRO:HG2	2:H:96:TRP:CE2	2.48	0.49
2:H:139:THR:HA	2:H:156:GLN:O	2.13	0.49
2:H:71:HIS:CE1	2:H:154:VAL:CG1	2.96	0.49
2:H:165:ARG:NH2	2:H:177:THR:O	2.37	0.49
2:H:178:ASP:O	2:H:233:ARG:HD2	2.13	0.49
2:H:244:GLN:HB3	2:H:245:PHE:CD1	2.48	0.48
2:H:97(A):GLU:HG2	2:H:98:ASN:N	2.28	0.48
2:H:124:PRO:HB3	2:H:128:THR:HG22	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:36(A):SER:HA	2:H:37:PRO:C	2.32	0.48
2:H:179:ASN:OD1	2:H:233:ARG:HD3	2.13	0.48
2:H:182:CYS:HA	2:H:226:GLY:O	2.13	0.48
1:L:14(G):LEU:HA	1:L:14(G):LEU:HD12	1.55	0.48
2:H:75:ARG:HD3	2:H:75:ARG:N	2.29	0.48
2:H:131:GLN:HB2	2:H:134:TYR:CD2	2.49	0.48
2:H:127:GLU:OE1	2:H:127:GLU:CA	2.60	0.48
2:H:28:PRO:HB2	2:H:119:HIS:HB3	1.96	0.47
2:H:126:ARG:HG2	2:H:232:PHE:CZ	2.46	0.47
2:H:145:LYS:CE	2:H:149(B):ASN:HB3	2.44	0.47
2:H:154:VAL:HG13	2:H:155:LEU:N	2.30	0.47
2:H:50:ARG:HD3	2:H:111:PRO:N	2.30	0.46
1:L:14(F):LEU:CD1	2:H:159:ASN:HB3	2.42	0.46
2:H:50:ARG:HH12	2:H:107:LYS:HE2	1.73	0.46
2:H:165:ARG:N	2:H:166:PRO:HD2	2.30	0.46
2:H:204:PRO:HG2	2:H:204(A):PHE:CZ	2.50	0.46
2:H:126:ARG:CB	2:H:126:ARG:HH11	2.29	0.46
2:H:129:ALA:HA	2:H:210:MET:HE1	1.96	0.46
2:H:50:ARG:NH1	2:H:107:LYS:CE	2.71	0.46
2:H:125:ASP:OD1	2:H:128:THR:CB	2.58	0.46
2:H:46:LEU:HD11	2:H:48:SER:O	2.16	0.46
1:L:4:ARG:HG2	2:H:28:PRO:CG	2.46	0.46
2:H:42:CYS:HB3	2:H:195:SER:O	2.15	0.45
2:H:57:HIS:ND1	2:H:102:ASP:OD2	2.49	0.45
2:H:70:LYS:HB3	2:H:70:LYS:HE3	1.56	0.45
2:H:33:LEU:HD23	2:H:33:LEU:HA	1.60	0.45
2:H:66:VAL:O	2:H:82:ILE:HA	2.16	0.45
2:H:95:ASN:HB3	2:H:100:ASP:HB3	1.99	0.45
1:L:14(J):TYR:HE2	2:H:202:LYS:O	1.99	0.45
2:H:80:GLU:C	2:H:81:LYS:HD2	2.36	0.45
2:H:236:LYS:HD2	2:H:236:LYS:N	2.32	0.45
2:H:237:TRP:O	2:H:238:ILE:C	2.51	0.45
2:H:97(A):GLU:CD	2:H:175:ARG:HH21	2.21	0.45
2:H:210:MET:HB3	2:H:210:MET:HE2	1.76	0.44
2:H:77(A):ARG:HG2	2:H:77(A):ARG:NH1	2.32	0.44
2:H:100:ASP:O	2:H:101:ARG:CB	2.66	0.44
2:H:204(B):ASN:ND2	2:H:206:ARG:HB2	2.33	0.44
2:H:132:ALA:HA	2:H:162:ILE:HG22	1.99	0.44
2:H:107:LYS:HE2	2:H:107:LYS:HB2	1.67	0.43
2:H:114:PHE:CD1	2:H:114:PHE:N	2.82	0.43
2:H:35:ARG:NH2	4:H:524:HOH:O	2.34	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:152:PRO:HG3	2:H:156:GLN:HG2	1.99	0.43
2:H:101:ARG:HD2	2:H:101:ARG:HH11	1.56	0.43
2:H:87:LYS:HG3	2:H:88:ILE:H	1.83	0.43
2:H:235:LYS:HA	2:H:238:ILE:HB	2.00	0.43
2:H:174:ILE:O	2:H:176:ILE:HD13	2.19	0.43
2:H:57:HIS:CE1	2:H:214:SER:O	2.72	0.43
2:H:145:LYS:HE2	2:H:149(B):ASN:HB3	2.01	0.43
2:H:219:GLY:HA2	4:H:443:HOH:O	2.19	0.43
2:H:98:ASN:N	2:H:98:ASN:OD1	2.49	0.43
2:H:160:LEU:HA	2:H:161:PRO:HD3	1.97	0.42
2:H:61:GLU:HG2	2:H:86:GLU:O	2.18	0.42
2:H:214:SER:HB3	2:H:227:PHE:O	2.18	0.42
2:H:146:GLU:O	2:H:149(B):ASN:ND2	2.51	0.42
2:H:68:ILE:CG2	2:H:118:ILE:HG23	2.49	0.42
2:H:136:GLY:CA	2:H:199:PHE:CZ	2.97	0.42
2:H:154:VAL:CG1	2:H:155:LEU:N	2.82	0.42
2:H:129(B):SER:O	2:H:131:GLN:NE2	2.53	0.42
2:H:123:LEU:HA	2:H:123:LEU:HD23	1.83	0.42
2:H:114:PHE:HA	2:H:118:ILE:O	2.19	0.42
2:H:52:VAL:HB	2:H:106:MET:HB2	2.03	0.41
2:H:131:GLN:O	2:H:134:TYR:HB2	2.20	0.41
2:H:169:LYS:HB3	4:H:559:HOH:O	2.21	0.41
2:H:212:ILE:O	2:H:228:TYR:HA	2.21	0.41
2:H:176:ILE:CG2	2:H:177:THR:N	2.84	0.40
2:H:60(B):PRO:HG2	2:H:96:TRP:CD2	2.56	0.40
2:H:65:LEU:HA	2:H:65:LEU:HD12	1.69	0.40
2:H:110:LYS:CB	2:H:111:PRO:HD2	2.52	0.40
2:H:70:LYS:NZ	2:H:77:GLU:OE1	2.36	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	25/36 (69%)	21 (84%)	4 (16%)	0	100	100
2	H	256/259 (99%)	237 (93%)	18 (7%)	1 (0%)	39	56
All	All	281/295 (95%)	258 (92%)	22 (8%)	1 (0%)	39	56

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	H	77(A)	ARG

### 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	25/31 (81%)	20 (80%)	5 (20%)	1	1
2	H	224/225 (100%)	195 (87%)	29 (13%)	5	6
All	All	249/256 (97%)	215 (86%)	34 (14%)	4	5

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

Mol	Chain	Res	Type
1	L	9	LYS
1	L	12	LEU
1	L	14(A)	LYS
1	L	14(G)	LEU
1	L	14(K)	ILE
2	H	28	PRO
2	H	36	LYS
2	H	50	ARG
2	H	64	LEU
2	H	65	LEU
2	H	66	VAL
2	H	70	LYS
2	H	73	ARG
2	H	74	THR
2	H	75	ARG

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Mol	Chain	Res	Type
2	H	79	ILE
2	H	81	LYS
2	H	99	LEU
2	H	107	LYS
2	H	110	LYS
2	H	112	VAL
2	H	145	LYS
2	H	149	THR
2	H	151	GLN
2	H	164	GLU
2	H	169	LYS
2	H	173	ARG
2	H	176	ILE
2	H	182	CYS
2	H	191	CYS
2	H	204(B)	ASN
2	H	205	ASN
2	H	233	ARG
2	H	239	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
2	H	71	HIS
2	H	78	ASN
2	H	131	GLN
2	H	204(B)	ASN
2	H	205	ASN
2	H	239	GLN

### 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 [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
3	0G6	H	1	2	27,31,32	2.43	2 (7%)	35,41,42	1.57	7 (20%)

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	0G6	H	1	2	1/1/8/10	0/29/41/43	0/2/2/2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	H	1	0G6	C3-C2	-9.73	1.27	1.51
3	H	1	0G6	O2-C2	-6.68	1.28	1.43

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	H	1	0G6	CE1-CD1-CG	-2.48	116.70	120.65
3	H	1	0G6	CZ-CE2-CD2	-2.20	116.96	120.19
3	H	1	0G6	CB1-CA1-N1	2.02	105.88	102.99
3	H	1	0G6	CE2-CD2-CG	2.20	124.15	120.65
3	H	1	0G6	O2-C2-C3	2.95	118.14	109.61
3	H	1	0G6	CB-CA-C	3.89	117.52	108.80
3	H	1	0G6	O2-C2-CA2	4.45	118.60	108.58

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	H	1	0G6	C2

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	H	1	0G6	5	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 ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

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

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

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

### 6.3 Carbohydrates ⓘ

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

### 6.4 Ligands ⓘ

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

### 6.5 Other polymers ⓘ

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