



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 06:32 PM GMT

PDB ID : 1BA8  
Title : THROMBIN INHIBITOR WITH A RIGID TRIPEPTIDYL ALDEHYDES  
Authors : Krishnan, R.; Zhang, E.; Hakansson, K.; Arni, R.K.; Tulinsky, A.; Lim-Wilby, M.S.L.; Levy, O.E.; Semple, J.E.; Brunck, T.K.  
Deposited on : 1998-04-23  
Resolution : 1.80 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](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

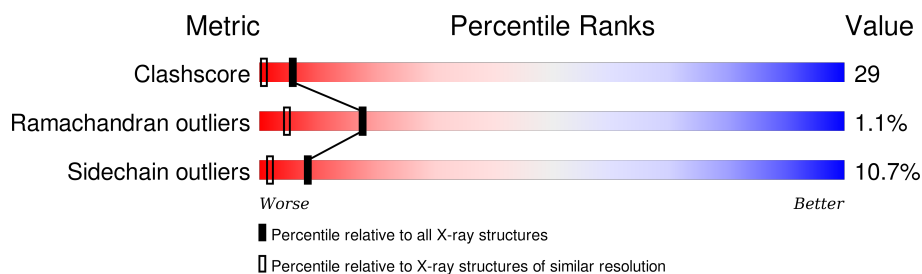
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.80 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	102246	5383 (1.80-1.80)
Ramachandran outliers	100387	5320 (1.80-1.80)
Sidechain outliers	100360	5319 (1.80-1.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	A	36	
2	B	259	
3	C	13	

## 2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 2551 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 THROMBIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	30	Total	C	N	O	S	0	0	0
			240	150	39	50	1			

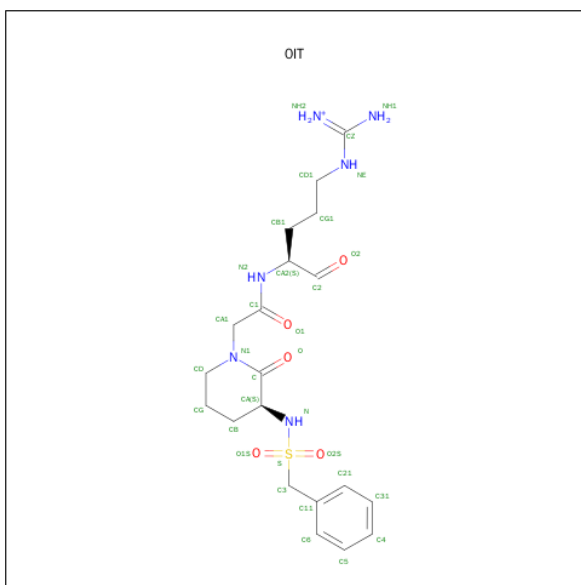
- Molecule 2 is a protein called THROMBIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	252	Total	C	N	O	S	0	0	0
			2039	1299	360	366	14			

- Molecule 3 is a protein called HIRUGEN.

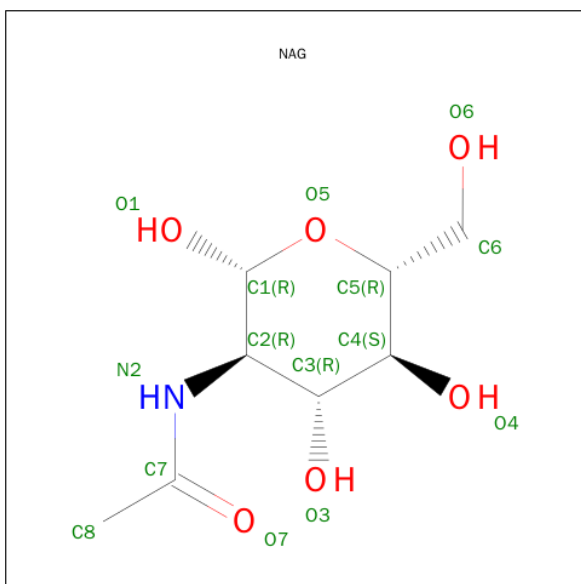
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	8	Total	C	N	O	S	0	0	0
			68	43	8	16	1			

- Molecule 4 is AMINO({(4S)-4-[({(3S)-3-[(BENZYL SULFONYL)AMINO]-2-OXOPIPERIDIN-1-YL}ACETYL)AMINO]-5-OXOPENTYL}AMINO)METHANIMINIUM (three-letter code: 0IT) (formula: C<sub>20</sub>H<sub>31</sub>N<sub>6</sub>O<sub>5</sub>S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	B	1	Total	C	N	O	S	0	0
			32	20	6	5	1		

- Molecule 5 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula:  $\text{C}_8\text{H}_{15}\text{NO}_6$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	16	Total 16	O 16	0	0
6	B	140	Total 140	O 140	0	0
6	C	2	Total 2	O 2	0	0

### 3 Residue-property plots [i](#)

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

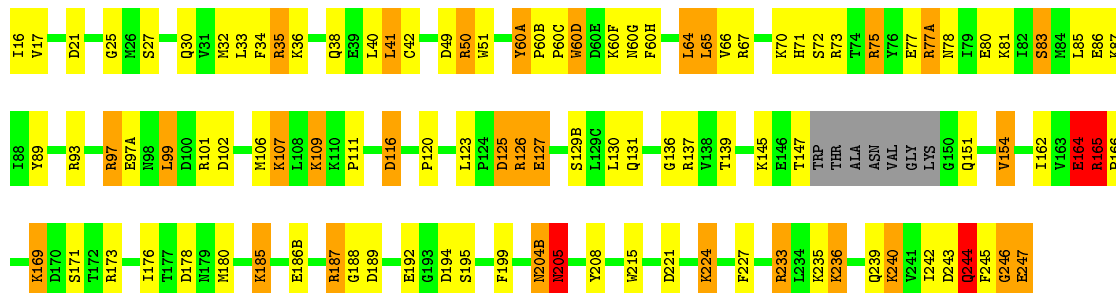
#### • Molecule 1: THROMBIN

Chain A: 

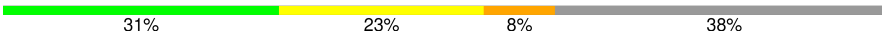


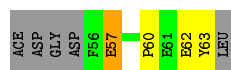
#### • Molecule 2: THROMBIN

Chain B: 



#### • Molecule 3: HIRUGEN

Chain C: 



## 4 Data and refinement statistics

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

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	71.26 Å   72.29 Å   72.99 Å 90.00°   100.90°   90.00°	Depositor
Resolution (Å)	7.00 – 1.80	Depositor
% Data completeness (in resolution range)	75.0 (7.00-1.80)	Depositor
$R_{merge}$	0.04	Depositor
$R_{sym}$	0.12	Depositor
Refinement program	PROLSQ, X-PLOR	Depositor
R, $R_{free}$	0.152 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2551	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	26.0	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: 0IT, NAG, TYS

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	1.36	3/242 (1.2%)	2.08	10/322 (3.1%)
2	B	1.08	0/2091	2.06	64/2823 (2.3%)
3	C	0.93	0/53	1.80	1/70 (1.4%)
All	All	1.11	3/2386 (0.1%)	2.05	75/3215 (2.3%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	14(K)	ILE	CB-CG2	7.73	1.76	1.52
1	A	14(K)	ILE	CB-CG1	6.64	1.72	1.54
1	A	8	GLU	CD-OE2	-5.26	1.19	1.25

All (75) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	233	ARG	NE-CZ-NH2	-16.44	112.08	120.30
2	B	101	ARG	NE-CZ-NH1	15.90	128.25	120.30
2	B	233	ARG	NE-CZ-NH1	14.20	127.40	120.30
2	B	77(A)	ARG	NE-CZ-NH2	-13.78	113.41	120.30
2	B	173	ARG	NE-CZ-NH1	13.36	126.98	120.30
2	B	93	ARG	NE-CZ-NH2	-13.18	113.71	120.30
2	B	67	ARG	NE-CZ-NH1	13.17	126.88	120.30
2	B	187	ARG	NE-CZ-NH1	12.24	126.42	120.30
2	B	173	ARG	NE-CZ-NH2	-11.60	114.50	120.30
2	B	35	ARG	NE-CZ-NH2	-11.20	114.70	120.30
2	B	187	ARG	NE-CZ-NH2	-11.16	114.72	120.30
2	B	97	ARG	NE-CZ-NH2	-11.07	114.76	120.30
1	A	1(A)	ASP	CB-CG-OD1	-9.47	109.78	118.30
2	B	50	ARG	NE-CZ-NH1	9.31	124.96	120.30
2	B	75	ARG	NE-CZ-NH1	9.17	124.88	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	233	ARG	CD-NE-CZ	9.15	136.41	123.60
2	B	101	ARG	NE-CZ-NH2	-9.00	115.80	120.30
2	B	75	ARG	NE-CZ-NH2	-8.92	115.84	120.30
2	B	93	ARG	NE-CZ-NH1	8.73	124.66	120.30
2	B	83	SER	N-CA-CB	8.11	122.67	110.50
2	B	21	ASP	CB-CG-OD2	-7.91	111.18	118.30
2	B	60(A)	TYR	CB-CG-CD1	-7.91	116.25	121.00
2	B	50	ARG	CD-NE-CZ	7.86	134.60	123.60
2	B	67	ARG	NE-CZ-NH2	-7.56	116.52	120.30
2	B	137	ARG	NE-CZ-NH2	-7.46	116.57	120.30
2	B	176	ILE	CA-CB-CG2	7.34	125.58	110.90
2	B	178	ASP	CB-CG-OD2	-7.28	111.75	118.30
2	B	21	ASP	CB-CG-OD1	7.23	124.81	118.30
2	B	246	GLY	O-C-N	7.21	134.24	122.70
2	B	125	ASP	CB-CG-OD2	7.16	124.74	118.30
2	B	243	ASP	CB-CA-C	7.00	124.41	110.40
2	B	60(A)	TYR	CB-CG-CD2	6.96	125.18	121.00
2	B	73	ARG	CD-NE-CZ	6.81	133.13	123.60
1	A	1(E)	SER	CA-C-N	6.61	129.42	116.20
2	B	189	ASP	CB-CG-OD2	6.57	124.21	118.30
2	B	165	ARG	CD-NE-CZ	-6.56	114.42	123.60
2	B	125	ASP	CB-CG-OD1	-6.48	112.47	118.30
2	B	205	ASN	CB-CA-C	6.31	123.02	110.40
1	A	14(D)	ARG	NE-CZ-NH2	-6.29	117.16	120.30
2	B	131	GLN	CA-CB-CG	-6.25	99.64	113.40
2	B	245	PHE	C-N-CA	-6.19	109.31	122.30
2	B	77(A)	ARG	NH1-CZ-NH2	6.18	126.20	119.40
2	B	97(A)	GLU	CA-CB-CG	6.18	127.00	113.40
2	B	116	ASP	CB-CG-OD1	6.08	123.77	118.30
2	B	73	ARG	NE-CZ-NH1	6.02	123.31	120.30
1	A	14(K)	ILE	CB-CG1-CD1	-6.02	97.04	113.90
1	A	4	ARG	NE-CZ-NH2	5.97	123.28	120.30
2	B	165	ARG	NE-CZ-NH1	-5.89	117.36	120.30
2	B	27	SER	CB-CA-C	5.88	121.27	110.10
2	B	180	MET	O-C-N	5.84	132.04	122.70
2	B	126	ARG	NE-CZ-NH2	-5.74	117.43	120.30
2	B	35	ARG	NE-CZ-NH1	5.73	123.16	120.30
2	B	42	CYS	N-CA-CB	-5.70	100.34	110.60
2	B	77	GLU	C-N-CA	5.64	135.81	121.70
1	A	14(J)	TYR	CB-CG-CD2	5.62	124.37	121.00
2	B	35	ARG	CD-NE-CZ	-5.53	115.86	123.60
2	B	102	ASP	CB-CG-OD2	-5.51	113.34	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	154	VAL	CG1-CB-CG2	5.49	119.68	110.90
1	A	1(E)	SER	CA-C-O	-5.46	108.64	120.10
2	B	127	GLU	CB-CG-CD	5.43	128.85	114.20
1	A	14(G)	LEU	CB-CA-C	5.37	120.41	110.20
2	B	78	ASN	CA-CB-CG	5.37	125.22	113.40
2	B	97(A)	GLU	CG-CD-OE2	-5.32	107.65	118.30
2	B	127	GLU	CG-CD-OE1	5.27	128.84	118.30
2	B	173	ARG	N-CA-CB	5.26	120.07	110.60
2	B	77	GLU	CA-C-O	5.23	131.09	120.10
3	C	57	GLU	CG-CD-OE2	-5.20	107.90	118.30
2	B	164	GLU	OE1-CD-OE2	-5.18	117.09	123.30
1	A	14	ASP	CB-CA-C	5.17	120.73	110.40
2	B	60(D)	TRP	N-CA-CB	5.15	119.87	110.60
2	B	154	VAL	N-CA-CB	-5.12	100.23	111.50
2	B	244	GLN	CB-CA-C	-5.11	100.18	110.40
2	B	227	PHE	CB-CG-CD1	-5.09	117.24	120.80
1	A	14(C)	GLU	CG-CD-OE1	-5.05	108.20	118.30
2	B	165	ARG	NE-CZ-NH2	5.02	122.81	120.30

There are no chirality outliers.

There are no planarity outliers.

## 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	A	240	0	239	35	0
2	B	2039	0	2007	114	0
3	C	68	0	49	6	0
4	B	32	0	31	8	0
5	B	14	0	13	1	0
6	A	16	0	0	1	0
6	B	140	0	0	6	0
6	C	2	0	0	0	0
All	All	2551	0	2339	138	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 29.

All (138) 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:A:14(K):ILE:CB	1:A:14(K):ILE:CG2	1.76	1.56
1:A:14(K):ILE:CG1	1:A:14(K):ILE:CD1	1.80	1.52
2:B:195:SER:OG	4:B:1:OIT:C2	1.67	1.40
2:B:86:GLU:HG3	2:B:109:LYS:NZ	1.39	1.33
2:B:75:ARG:NH2	3:C:57:GLU:HB2	1.58	1.17
1:A:1(C):GLU:HB3	1:A:1:CYS:HB3	1.21	1.16
2:B:86:GLU:CG	2:B:109:LYS:NZ	2.15	1.10
2:B:240:LYS:HB3	2:B:240:LYS:NZ	1.68	1.07
1:A:1(C):GLU:HG3	2:B:120:PRO:HG2	1.40	1.04
1:A:1(C):GLU:CB	1:A:1:CYS:HB3	1.90	1.01
1:A:1(C):GLU:HG3	2:B:120:PRO:CG	1.94	0.98
2:B:86:GLU:CB	2:B:109:LYS:HZ2	1.78	0.96
2:B:86:GLU:CG	2:B:109:LYS:HZ1	1.77	0.95
2:B:85:LEU:HD13	2:B:106:MET:HE2	1.46	0.94
2:B:50:ARG:NH1	2:B:86:GLU:OE1	2.02	0.92
1:A:1(D):GLY:H	2:B:123:LEU:H	1.18	0.91
2:B:86:GLU:CG	2:B:109:LYS:HZ2	1.80	0.89
2:B:72:SER:OG	2:B:75:ARG:HG2	1.74	0.87
2:B:86:GLU:HG3	2:B:109:LYS:HZ1	0.93	0.86
2:B:75:ARG:HH22	3:C:57:GLU:HB2	1.41	0.85
2:B:85:LEU:CD1	2:B:106:MET:HE2	2.10	0.81
2:B:240:LYS:HZ2	2:B:240:LYS:HB3	1.45	0.79
1:A:14(D):ARG:O	1:A:14(H):GLU:HG3	1.82	0.79
2:B:240:LYS:HZ3	2:B:240:LYS:HB3	1.43	0.79
2:B:85:LEU:CD1	2:B:106:MET:CE	2.61	0.78
2:B:75:ARG:NH2	3:C:57:GLU:CB	2.46	0.77
2:B:85:LEU:HD11	2:B:106:MET:HE1	1.68	0.76
1:A:14(K):ILE:CB	1:A:14(K):ILE:CD1	2.64	0.76
1:A:14(K):ILE:HD12	1:A:14(K):ILE:HG23	1.68	0.75
2:B:32:MET:HG3	2:B:40:LEU:HD13	1.69	0.75
2:B:86:GLU:HB3	2:B:107:LYS:HG3	1.66	0.75
2:B:64:LEU:HD12	2:B:85:LEU:HD12	1.68	0.74
2:B:233:ARG:O	2:B:236:LYS:HE2	1.88	0.74
2:B:130:LEU:HD23	2:B:162:ILE:CD1	2.18	0.73
2:B:86:GLU:HG3	2:B:109:LYS:HZ2	1.38	0.73
2:B:235:LYS:HE2	2:B:239:GLN:OE1	1.89	0.72
2:B:195:SER:OG	4:B:1:OIT:O2	2.05	0.72
2:B:246:GLY:C	2:B:247:GLU:HG3	2.10	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:195:SER:OG	4:B:1:0IT:HXT	1.84	0.71
1:A:1(D):GLY:N	2:B:123:LEU:H	1.88	0.71
2:B:85:LEU:HD11	2:B:106:MET:CE	2.20	0.69
2:B:192:GLU:HB2	6:B:517:HOH:O	1.89	0.69
1:A:1(C):GLU:HB3	1:A:1:CYS:CB	2.13	0.69
2:B:195:SER:CB	4:B:1:0IT:C2	2.71	0.69
3:C:60:PRO:HG2	3:C:63:TYS:HE2	1.75	0.68
1:A:14(K):ILE:HD12	1:A:14(K):ILE:CG2	2.24	0.68
2:B:80:GLU:O	2:B:81:LYS:HD3	1.93	0.68
2:B:35:ARG:HB2	2:B:41:LEU:HD13	1.75	0.67
2:B:246:GLY:O	2:B:247:GLU:O	2.12	0.67
2:B:30:GLN:NE2	2:B:139:THR:OG1	2.28	0.66
1:A:14(K):ILE:CD1	1:A:14(K):ILE:CG2	2.75	0.66
2:B:64:LEU:CD1	2:B:85:LEU:HD12	2.26	0.65
2:B:165:ARG:O	2:B:169:LYS:HD2	1.96	0.64
1:A:1(E):SER:HB2	2:B:208:TYR:HE1	1.63	0.64
2:B:195:SER:OG	4:B:1:0IT:CA2	2.41	0.64
1:A:14(K):ILE:CG1	1:A:14(K):ILE:CG2	2.76	0.63
2:B:86:GLU:HB2	2:B:109:LYS:HZ2	1.60	0.63
2:B:130:LEU:HD23	2:B:162:ILE:HD11	1.79	0.63
2:B:36:LYS:O	2:B:38:GLN:HG2	1.98	0.62
2:B:240:LYS:O	2:B:244:GLN:HB2	2.00	0.62
2:B:205:ASN:ND2	6:B:541:HOH:O	2.33	0.62
1:A:14(D):ARG:NE	1:A:14(H):GLU:OE2	2.33	0.61
2:B:204(B):ASN:HD22	2:B:204(B):ASN:C	2.03	0.61
1:A:14(K):ILE:CD1	1:A:14(K):ILE:HG23	2.30	0.61
2:B:246:GLY:O	2:B:247:GLU:HG3	2.01	0.61
2:B:130:LEU:HD23	2:B:162:ILE:HD13	1.83	0.60
2:B:71:HIS:HD2	6:B:437:HOH:O	1.84	0.60
1:A:1(D):GLY:O	1:A:1(C):GLU:HB2	2.00	0.60
2:B:165:ARG:NH1	6:B:467:HOH:O	2.33	0.60
2:B:60(G):ASN:OD1	2:B:60(G):ASN:C	2.41	0.59
2:B:236:LYS:O	2:B:240:LYS:HG3	2.02	0.58
2:B:240:LYS:CB	2:B:240:LYS:NZ	2.56	0.57
2:B:236:LYS:HD2	2:B:236:LYS:H	1.70	0.56
2:B:246:GLY:C	2:B:247:GLU:O	2.44	0.56
1:A:6:LEU:HD12	2:B:25:GLY:HA3	1.88	0.55
2:B:85:LEU:CD1	2:B:106:MET:HE1	2.29	0.55
2:B:36:LYS:HG3	2:B:65:LEU:HD22	1.88	0.55
2:B:16:ILE:N	2:B:194:ASP:OD2	2.38	0.55
2:B:72:SER:HG	2:B:75:ARG:HG2	1.71	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1(D):GLY:H	2:B:123:LEU:N	1.96	0.54
2:B:32:MET:HG3	2:B:40:LEU:CD1	2.36	0.54
1:A:1(C):GLU:CG	2:B:120:PRO:CG	2.79	0.53
2:B:17:VAL:O	2:B:188:GLY:HA2	2.08	0.53
2:B:35:ARG:O	2:B:38:GLN:HA	2.09	0.52
1:A:1(E):SER:CB	2:B:123:LEU:O	2.58	0.52
2:B:242:ILE:HG23	2:B:247:GLU:HG2	1.92	0.51
2:B:187:ARG:HD3	2:B:221:ASP:OD2	2.11	0.51
5:B:400:NAG:O7	5:B:400:NAG:C1	2.58	0.51
2:B:70:LYS:HE3	2:B:72:SER:O	2.11	0.51
1:A:1(C):GLU:CB	1:A:1:CYS:CB	2.79	0.50
2:B:50:ARG:HD2	2:B:107:LYS:HE2	1.94	0.50
2:B:185:LYS:HB2	2:B:186(B):GLU:OE1	2.12	0.50
1:A:1(C):GLU:HG3	2:B:120:PRO:CB	2.41	0.49
2:B:60(G):ASN:OD1	2:B:60(G):ASN:O	2.30	0.49
2:B:86:GLU:CA	2:B:109:LYS:HZ2	2.26	0.48
2:B:165:ARG:N	2:B:166:PRO:HD2	2.28	0.48
2:B:49:ASP:O	2:B:111:PRO:HA	2.13	0.48
2:B:215:TRP:CE3	4:B:1:OIT:H21	2.48	0.48
1:A:1(D):GLY:HA2	6:A:462:HOH:O	2.13	0.48
2:B:136:GLY:HA3	2:B:199:PHE:CZ	2.48	0.48
2:B:64:LEU:CD1	2:B:85:LEU:CD1	2.91	0.47
2:B:165:ARG:O	2:B:169:LYS:CD	2.63	0.47
1:A:1(E):SER:HB2	2:B:123:LEU:O	2.13	0.47
2:B:60(F):LYS:HD2	2:B:60(H):PHE:HE2	1.80	0.47
2:B:86:GLU:HB3	2:B:107:LYS:CG	2.42	0.47
2:B:85:LEU:HD13	2:B:106:MET:CE	2.23	0.46
1:A:9:LYS:HG3	1:A:9:LYS:HZ3	1.44	0.46
1:A:1(D):GLY:CA	2:B:123:LEU:H	2.28	0.46
2:B:130:LEU:CD2	2:B:162:ILE:CD1	2.92	0.46
2:B:60(B):PRO:N	2:B:60(C):PRO:CD	2.77	0.46
2:B:151:GLN:HG2	6:B:419:HOH:O	2.15	0.46
2:B:60(A):TYR:C	2:B:60(C):PRO:HD2	2.36	0.45
2:B:60(D):TRP:CH2	4:B:1:OIT:HG2	2.51	0.45
2:B:87:LYS:HB3	2:B:89:TYR:CE1	2.52	0.45
1:A:1(E):SER:HB3	2:B:123:LEU:O	2.17	0.45
2:B:130:LEU:CD2	2:B:162:ILE:HD13	2.47	0.45
2:B:60(B):PRO:N	2:B:60(C):PRO:HD2	2.32	0.45
2:B:107:LYS:HG3	2:B:107:LYS:O	2.16	0.45
1:A:14(K):ILE:CA	1:A:14(K):ILE:CG2	2.83	0.44
1:A:1(B):ALA:H	1:A:1:CYS:H	1.66	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:14(D):ARG:HG2	1:A:14(H):GLU:HG3	2.00	0.43
2:B:60(A):TYR:CZ	2:B:60(C):PRO:HG2	2.53	0.43
2:B:126:ARG:NH1	2:B:126:ARG:HB3	2.34	0.43
2:B:164:GLU:OE1	2:B:164:GLU:N	2.29	0.43
1:A:14(D):ARG:CZ	1:A:14(H):GLU:OE2	2.67	0.43
2:B:51:TRP:CZ3	2:B:107:LYS:HB3	2.54	0.42
2:B:81:LYS:HD2	3:C:63:TYS:O2	2.19	0.42
2:B:204(B):ASN:HD22	2:B:205:ASN:N	2.17	0.42
2:B:171:SER:O	2:B:224:LYS:CE	2.67	0.42
2:B:99:LEU:HA	2:B:99:LEU:HD12	1.90	0.42
2:B:33:LEU:HD23	2:B:33:LEU:HA	1.85	0.42
2:B:165:ARG:N	2:B:166:PRO:CD	2.83	0.42
2:B:165:ARG:NH1	2:B:169:LYS:HE3	2.35	0.41
3:C:62:GLU:H	3:C:62:GLU:HG2	1.52	0.41
2:B:97:ARG:HG3	6:B:415:HOH:O	2.21	0.40
1:A:5:PRO:HB2	2:B:116:ASP:HA	2.03	0.40
2:B:34:PHE:CZ	2:B:38:GLN:HB3	2.57	0.40
2:B:60(D):TRP:CZ2	4:B:10IT:HG2	2.57	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	A	28/36 (78%)	24 (86%)	2 (7%)	2 (7%)	1	0
2	B	248/259 (96%)	241 (97%)	6 (2%)	1 (0%)	39	23
3	C	6/13 (46%)	6 (100%)	0	0	100	100
All	All	282/308 (92%)	271 (96%)	8 (3%)	3 (1%)	17	5

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1(C)	GLU
1	A	1(B)	ALA
2	B	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	A	27/31 (87%)	25 (93%)	2 (7%)	17	5
2	B	220/225 (98%)	195 (89%)	25 (11%)	7	1
3	C	5/10 (50%)	5 (100%)	0	100	100
All	All	252/266 (95%)	225 (89%)	27 (11%)	8	2

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1(C)	GLU
1	A	14(F)	LEU
2	B	41	LEU
2	B	64	LEU
2	B	65	LEU
2	B	66	VAL
2	B	83	SER
2	B	99	LEU
2	B	107	LYS
2	B	109	LYS
2	B	125	ASP
2	B	127	GLU
2	B	129(B)	SER
2	B	145	LYS
2	B	147	THR
2	B	154	VAL
2	B	164	GLU
2	B	165	ARG
2	B	169	LYS
2	B	185	LYS

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Mol	Chain	Res	Type
2	B	204(B)	ASN
2	B	205	ASN
2	B	224	LYS
2	B	236	LYS
2	B	240	LYS
2	B	244	GLN
2	B	247	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	30	GLN
2	B	71	HIS
2	B	156	GLN
2	B	204(B)	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue 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	TYS	C	63	3	15,16,17	0.89	0	16,22,24	1.80	4 (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	TYS	C	63	3	-	0/9/11/13	0/1/1/1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	63	TYS	O-C-CA	-4.80	113.00	125.49
3	C	63	TYS	CG-CB-CA	-2.81	107.87	114.21
3	C	63	TYS	O2-S-O1	2.06	121.43	112.46
3	C	63	TYS	CE2-CZ-CE1	2.23	123.82	120.20

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	63	TYS	2	0

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

2 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	0IT	B	1	-	29,33,33	1.53	5 (17%)	30,44,44	2.26	11 (36%)
5	NAG	B	400	2	14,14,15	0.77	0	15,19,21	2.25	4 (26%)

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	0IT	B	1	-	-	0/23/41/41	0/2/2/2
5	NAG	B	400	2	-	0/6/23/26	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	1	0IT	CA-N	-3.02	1.41	1.46
4	B	1	0IT	CB-CA	-2.63	1.48	1.53
4	B	1	0IT	CA2-N2	2.11	1.49	1.46
4	B	1	0IT	CA-C	3.25	1.57	1.52
4	B	1	0IT	S-N	3.51	1.67	1.62

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	400	NAG	C2-N2-C7	-6.13	115.17	123.04
4	B	1	0IT	CA1-N1-C	-5.10	114.98	119.73
4	B	1	0IT	O2S-S-N	-4.03	101.58	107.03
4	B	1	0IT	C3-C11-C21	-2.72	116.81	120.52
4	B	1	0IT	O-C-CA	-2.12	115.94	120.48
4	B	1	0IT	O2-C2-CA2	-2.12	119.84	125.44
4	B	1	0IT	C4-C31-C21	2.02	123.14	120.19
4	B	1	0IT	C3-C11-C6	2.41	123.81	120.52
5	B	400	NAG	C1-O5-C5	2.86	115.88	112.25
4	B	1	0IT	CG-CB-CA	2.87	115.50	111.04
5	B	400	NAG	O3-C3-C4	3.25	117.66	110.34
5	B	400	NAG	C3-C4-C5	3.49	116.28	110.20
4	B	1	0IT	O2S-S-O1S	3.78	125.87	119.34
4	B	1	0IT	O-C-N1	4.65	128.08	122.38
4	B	1	0IT	CB-CG-CD	4.79	117.37	110.97

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	1	0IT	8	0
5	B	400	NAG	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

### 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.