Full wwPDB X-ray Structure Validation Report (i)

Jan 31, 2016 – 10:25 PM GMT

PDB ID	:	1TMB
Title	:	MOLECULAR BASIS FOR THE INHIBITION OF HUMAN ALPHA-THROMBIN BY THE MACROCYCLIC PEPTIDE CYCLOTHEONAMIDE A
Authors	:	Qiu, X.; Padmanabhan, K.P.; Maryanoff, B.E.; Tulinsky, A.
Deposited on	:	1993-05-27
Resolution	:	2.30 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (i)) were used in the production of this report:

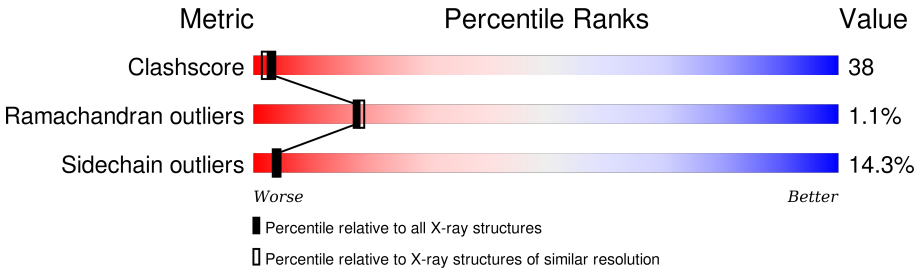
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	22% 47% 8% · 19%
2	H	259	42% 34% 18% · ·
3	I	13	23% 38% 8% 8% 23%
4	T	5	80% 20%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 2644 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	29	Total	C	N	O	S	0	0	0
			235	147	38	49	1			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	252	Total	C	N	O	S	0	0	0
			2032	1295	358	365	14			

- Molecule 3 is a protein called HIRUGEN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	I	10	Total	C	N	O	S	0	0	0
			85	53	10	21	1			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
I	54	GLU	GLY	CONFLICT	UNP P09945

- Molecule 4 is a protein called cyclotheonamide A.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	T	5	Total	C	N	O	0	0	0
			53	36	9	8			

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	L	24	Total	O	0	0
			24	24		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	H	202	Total O 202 202	0	0
5	I	7	Total O 7 7	0	0
5	T	6	Total O 6 6	0	0

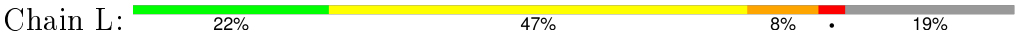
3 Residue-property plots (i)

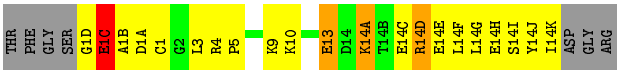
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: ALPHA-THROMBIN (SMALL SUBUNIT)

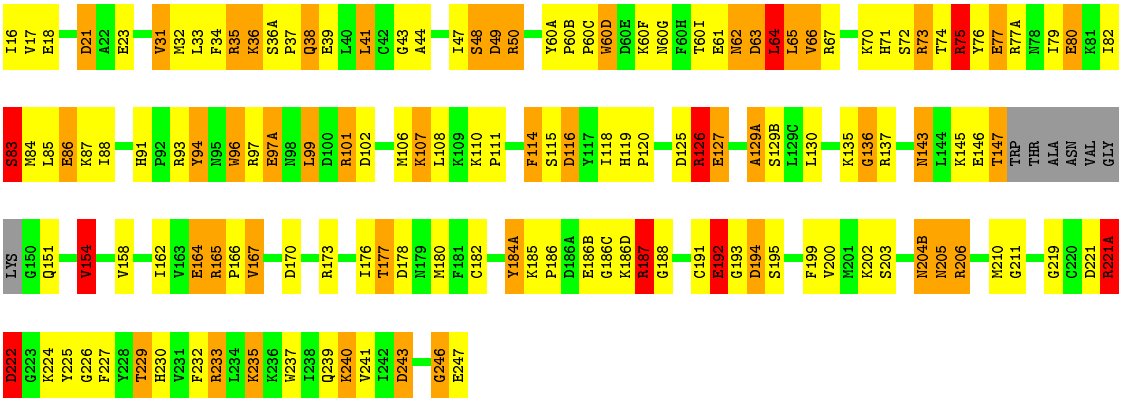
Chain L:





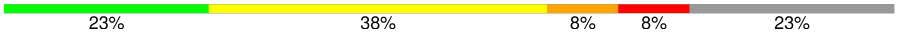
- Molecule 2: ALPHA-THROMBIN (LARGE SUBUNIT)

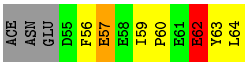
Chain H:



- Molecule 3: HIRUGEN

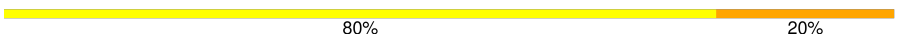
Chain I:





- Molecule 4: cyclotheonamide A

Chain T:



4 Data and refinement statistics (i)

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, α, β, γ	70.61Å 72.38Å 73.35Å 90.00° 101.10° 90.00°	Depositor
Resolution (Å)	(Not available) – 2.30	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-2.30)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROLSQ	Depositor
R, R_{free}	0.138 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2644	wwPDB-VP
Average B, all atoms (Å^2)	29.0	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: DPN, 0MG, 0FL, TYS, VLT

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	1.14	0/237	2.32	11/315 (3.5%)
2	H	1.17	0/2084	2.37	90/2815 (3.2%)
3	I	0.97	0/69	1.89	3/89 (3.4%)
4	T	0.93	0/7	1.80	0/8
All	All	1.16	0/2397	2.35	104/3227 (3.2%)

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
4	T	0	2

There are no bond length outliers.

All (104) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	H	73	ARG	CD-NE-CZ	39.32	178.65	123.60
2	H	206	ARG	NE-CZ-NH1	19.78	130.19	120.30
2	H	75	ARG	CD-NE-CZ	13.79	142.91	123.60
2	H	187	ARG	NE-CZ-NH1	-13.68	113.46	120.30
2	H	35	ARG	NE-CZ-NH2	-12.37	114.11	120.30
2	H	126	ARG	NE-CZ-NH2	-12.19	114.21	120.30
2	H	233	ARG	NE-CZ-NH2	-12.08	114.26	120.30
2	H	73	ARG	NE-CZ-NH1	-11.84	114.38	120.30
2	H	93	ARG	NE-CZ-NH2	-10.87	114.86	120.30
2	H	165	ARG	NE-CZ-NH1	10.61	125.61	120.30
2	H	60(A)	TYR	CB-CG-CD1	-9.99	115.00	121.00
2	H	225	TYR	CB-CG-CD1	-9.95	115.03	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	H	126	ARG	NE-CZ-NH1	9.74	125.17	120.30
2	H	187	ARG	CD-NE-CZ	-9.70	110.02	123.60
2	H	49	ASP	CB-CG-OD2	9.63	126.97	118.30
2	H	75	ARG	NE-CZ-NH1	9.54	125.07	120.30
2	H	116	ASP	CB-CG-OD2	-9.50	109.75	118.30
2	H	206	ARG	NE-CZ-NH2	-9.47	115.57	120.30
2	H	114	PHE	CB-CG-CD1	-9.33	114.27	120.80
2	H	184(A)	TYR	CB-CG-CD1	-9.32	115.41	121.00
2	H	50	ARG	NE-CZ-NH1	9.27	124.93	120.30
1	L	14(D)	ARG	NE-CZ-NH2	-9.22	115.69	120.30
2	H	243	ASP	CB-CG-OD1	9.00	126.40	118.30
2	H	221(A)	ARG	CD-NE-CZ	8.97	136.16	123.60
2	H	137	ARG	NE-CZ-NH1	8.93	124.76	120.30
2	H	164	GLU	OE1-CD-OE2	-8.65	112.92	123.30
2	H	116	ASP	CB-CG-OD1	8.64	126.08	118.30
2	H	125	ASP	CB-CG-OD2	-8.60	110.56	118.30
1	L	13	GLU	OE1-CD-OE2	8.38	133.35	123.30
2	H	233	ARG	NE-CZ-NH1	8.01	124.31	120.30
2	H	21	ASP	CB-CG-OD1	8.01	125.51	118.30
2	H	246	GLY	C-N-CA	7.92	141.51	121.70
2	H	222	ASP	CB-CG-OD1	7.84	125.36	118.30
2	H	192	GLU	OE1-CD-OE2	-7.76	113.98	123.30
2	H	192	GLU	CB-CG-CD	7.72	135.06	114.20
2	H	77(A)	ARG	CA-C-N	7.71	134.16	117.20
2	H	194	ASP	CB-CG-OD2	7.70	125.23	118.30
2	H	173	ARG	NE-CZ-NH1	7.66	124.13	120.30
2	H	73	ARG	NE-CZ-NH2	7.56	124.08	120.30
2	H	194	ASP	CB-CG-OD1	-7.36	111.67	118.30
1	L	1(C)	GLU	OE1-CD-OE2	7.34	132.10	123.30
2	H	200	VAL	CA-CB-CG1	7.14	121.61	110.90
2	H	184(A)	TYR	CB-CG-CD2	7.03	125.22	121.00
2	H	187	ARG	NE-CZ-NH2	7.03	123.82	120.30
2	H	167	VAL	CG1-CB-CG2	7.03	122.14	110.90
2	H	38	GLN	CB-CA-C	-6.96	96.49	110.40
2	H	35	ARG	NE-CZ-NH1	6.95	123.77	120.30
1	L	13	GLU	CB-CA-C	-6.90	96.60	110.40
2	H	125	ASP	CB-CG-OD1	6.67	124.30	118.30
2	H	77(A)	ARG	CA-C-O	-6.66	106.11	120.10
2	H	80	GLU	CG-CD-OE2	6.64	131.57	118.30
2	H	67	ARG	NE-CZ-NH1	6.52	123.56	120.30
1	L	14(C)	GLU	CG-CD-OE1	6.46	131.22	118.30
2	H	67	ARG	CD-NE-CZ	6.38	132.53	123.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	H	38	GLN	CA-CB-CG	6.37	127.42	113.40
2	H	74	THR	OG1-CB-CG2	6.33	124.57	110.00
2	H	115	SER	CB-CA-C	6.29	122.05	110.10
2	H	101	ARG	NE-CZ-NH1	6.29	123.44	120.30
3	I	62	GLU	CG-CD-OE2	6.25	130.80	118.30
2	H	86	GLU	OE1-CD-OE2	6.22	130.76	123.30
2	H	93	ARG	CD-NE-CZ	6.18	132.25	123.60
2	H	154	VAL	N-CA-CB	-6.13	98.02	111.50
2	H	21	ASP	CB-CG-OD2	-6.09	112.82	118.30
2	H	229	THR	OG1-CB-CG2	6.07	123.96	110.00
2	H	102	ASP	CB-CG-OD1	-6.00	112.90	118.30
1	L	4	ARG	CG-CD-NE	5.92	124.24	111.80
1	L	1(A)	ASP	CB-CA-C	5.86	122.11	110.40
2	H	77(A)	ARG	NE-CZ-NH1	-5.85	117.38	120.30
2	H	192	GLU	CG-CD-OE2	5.82	129.93	118.30
2	H	243	ASP	CA-CB-CG	5.80	126.16	113.40
2	H	63	ASP	CB-CG-OD1	-5.79	113.09	118.30
2	H	60(D)	TRP	CA-CB-CG	5.78	124.68	113.70
1	L	13	GLU	O-C-N	5.77	131.94	122.70
2	H	129(A)	ALA	O-C-N	5.77	131.93	122.70
2	H	165	ARG	NE-CZ-NH2	-5.71	117.45	120.30
2	H	170	ASP	CB-CG-OD1	5.70	123.43	118.30
2	H	93	ARG	NE-CZ-NH1	5.65	123.12	120.30
1	L	10	LYS	CD-CE-NZ	-5.62	98.78	111.70
2	H	135	LYS	O-C-N	5.62	132.75	123.20
2	H	136	GLY	N-CA-C	-5.51	99.32	113.10
2	H	80	GLU	CG-CD-OE1	-5.50	107.30	118.30
2	H	77	GLU	CG-CD-OE2	5.50	129.29	118.30
2	H	75	ARG	NE-CZ-NH2	-5.47	117.56	120.30
3	I	57	GLU	CG-CD-OE1	5.46	129.23	118.30
2	H	221(A)	ARG	NE-CZ-NH1	5.41	123.00	120.30
2	H	31	VAL	CA-CB-CG2	5.40	119.00	110.90
2	H	97(A)	GLU	CA-CB-CG	5.38	125.23	113.40
2	H	74	THR	N-CA-CB	-5.37	100.09	110.30
2	H	177	THR	CA-CB-CG2	5.37	119.92	112.40
2	H	73	ARG	C-N-CA	5.35	135.07	121.70
2	H	221(A)	ARG	CA-C-O	-5.35	108.87	120.10
2	H	38	GLN	O-C-N	5.31	131.20	122.70
2	H	64	LEU	CA-C-N	-5.21	105.74	117.20
2	H	74	THR	CA-CB-OG1	-5.18	98.12	109.00
1	L	13	GLU	CG-CD-OE1	-5.16	107.97	118.30
2	H	83	SER	O-C-N	5.13	130.90	122.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	H	158	VAL	CA-CB-CG1	5.12	118.58	110.90
2	H	143	ASN	CA-C-N	-5.12	105.94	117.20
2	H	96	TRP	O-C-N	5.11	130.87	122.70
2	H	48	SER	N-CA-CB	5.10	118.15	110.50
2	H	73	ARG	CG-CD-NE	-5.08	101.14	111.80
2	H	170	ASP	CB-CG-OD2	-5.06	113.75	118.30
1	L	4	ARG	NE-CZ-NH1	5.04	122.82	120.30
3	I	62	GLU	CG-CD-OE1	-5.01	108.28	118.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
4	T	48	0MG	Mainchain,Peptide

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	L	235	0	234	20	0
2	H	2032	0	1995	152	1
3	I	85	0	63	14	1
4	T	53	0	42	12	0
5	H	202	0	0	26	0
5	I	7	0	0	0	0
5	L	24	0	0	7	0
5	T	6	0	0	1	0
All	All	2644	0	2334	177	1

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 (177) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:205:ASN:HB3	5:H:503:HOH:O	1.44	1.14

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:1(C):GLU:HG3	5:L:483:HOH:O	1.50	1.07
2:H:192:GLU:OE1	5:H:405:HOH:O	1.76	1.01
2:H:193:GLY:HA2	4:T:49:DPN:CE2	1.90	1.00
2:H:36:LYS:HE3	5:H:521:HOH:O	1.62	0.99
2:H:85:LEU:HD13	2:H:106:MET:HE2	1.43	0.99
2:H:35:ARG:O	2:H:38:GLN:HA	1.64	0.97
2:H:17:VAL:O	2:H:188:GLY:HA2	1.68	0.92
2:H:147:THR:HG21	5:H:522:HOH:O	1.69	0.92
2:H:64:LEU:HD12	2:H:85:LEU:HD12	1.50	0.92
2:H:60(I):THR:OG1	2:H:62:ASN:ND2	2.01	0.91
2:H:50:ARG:NH1	2:H:108:LEU:O	2.03	0.90
1:L:1(D):GLY:HA2	5:L:621:HOH:O	1.72	0.88
2:H:165:ARG:NH1	5:H:491:HOH:O	2.07	0.87
3:I:59:ILE:HD11	3:I:64:LEU:HD11	1.53	0.87
2:H:85:LEU:HD13	2:H:106:MET:CE	2.06	0.85
3:I:59:ILE:HD11	3:I:64:LEU:CD1	2.06	0.84
2:H:64:LEU:HD12	2:H:85:LEU:CD1	2.07	0.83
1:L:1(C):GLU:CG	5:L:483:HOH:O	2.17	0.82
2:H:187:ARG:NH2	2:H:222:ASP:OD1	2.18	0.77
2:H:37:PRO:O	2:H:39:GLU:HG2	1.86	0.76
2:H:126:ARG:NH1	2:H:127:GLU:OE2	2.19	0.76
3:I:60:PRO:HB2	3:I:62:GLU:HG3	1.70	0.74
3:I:59:ILE:CD1	3:I:64:LEU:HD11	2.16	0.74
2:H:36:LYS:CE	5:H:521:HOH:O	2.25	0.74
2:H:165:ARG:NH2	2:H:180:MET:O	2.22	0.73
2:H:87:LYS:HE3	5:H:556:HOH:O	1.89	0.73
2:H:21:ASP:HB3	2:H:154:VAL:CG1	2.18	0.73
2:H:204(B):ASN:C	2:H:205:ASN:HD22	1.91	0.73
2:H:195:SER:HG	4:T:48:0MG:C2	2.02	0.73
2:H:184(A):TYR:HA	2:H:186(B):GLU:OE1	1.88	0.72
3:I:59:ILE:CG1	3:I:64:LEU:HD11	2.20	0.71
2:H:195:SER:OG	4:T:48:0MG:C	2.38	0.71
2:H:164:GLU:OE1	2:H:185:LYS:NZ	2.21	0.70
2:H:62:ASN:ND2	2:H:63:ASP:OD1	2.24	0.70
2:H:203:SER:O	2:H:205:ASN:HA	1.92	0.69
2:H:219:GLY:HA3	2:H:221(A):ARG:HE	1.57	0.69
2:H:16:ILE:N	2:H:194:ASP:OD2	2.25	0.69
2:H:75:ARG:N	2:H:75:ARG:HD3	2.08	0.68
2:H:162:ILE:HD11	2:H:199:PHE:CZ	2.28	0.68
2:H:129(A):ALA:O	5:H:582:HOH:O	2.11	0.67
2:H:195:SER:OG	4:T:48:0MG:O1	2.11	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:85:LEU:CD1	2:H:106:MET:CE	2.73	0.67
2:H:185:LYS:N	2:H:186(B):GLU:OE1	2.23	0.66
2:H:97:ARG:NE	5:H:496:HOH:O	1.84	0.65
1:L:14(A):LYS:HG2	2:H:23:GLU:OE2	1.96	0.64
1:L:14(G):LEU:HD21	2:H:202:LYS:HD3	1.79	0.64
2:H:205:ASN:OD1	5:H:436:HOH:O	2.15	0.63
2:H:127:GLU:HB2	5:H:586:HOH:O	1.99	0.63
2:H:184(A):TYR:CA	2:H:186(B):GLU:OE1	2.46	0.62
2:H:129(B):SER:O	5:H:511:HOH:O	2.16	0.62
1:L:14(D):ARG:CZ	1:L:14(H):GLU:OE2	2.48	0.62
3:I:60:PRO:HG2	3:I:63:TYS:HE2	1.83	0.61
2:H:178:ASP:O	2:H:233:ARG:HD2	2.02	0.60
2:H:177:THR:HB	5:H:628:HOH:O	2.01	0.60
2:H:147:THR:CG2	5:H:522:HOH:O	2.40	0.60
1:L:3:LEU:HB2	1:L:9:LYS:HE3	1.84	0.60
2:H:70:LYS:HE3	2:H:72:SER:O	2.01	0.59
2:H:64:LEU:CD1	2:H:85:LEU:CD1	2.80	0.59
2:H:35:ARG:O	2:H:38:GLN:CA	2.45	0.59
2:H:146:GLU:OE1	2:H:221(A):ARG:HD3	2.04	0.58
2:H:21:ASP:HB3	2:H:154:VAL:HG11	1.84	0.58
2:H:126:ARG:NH1	2:H:126:ARG:HG2	2.18	0.58
2:H:126:ARG:HB3	2:H:126:ARG:HH11	1.68	0.57
2:H:41:LEU:O	4:T:49:DPN:CE2	2.53	0.57
2:H:85:LEU:CD1	2:H:106:MET:HE1	2.35	0.57
2:H:185:LYS:HB2	2:H:186(B):GLU:HG3	1.86	0.57
1:L:14(G):LEU:HB3	5:L:584:HOH:O	2.05	0.57
2:H:195:SER:OG	4:T:48:0MG:CA	2.48	0.57
2:H:221(A):ARG:HB2	2:H:224:LYS:HG3	1.86	0.56
2:H:86:GLU:HB3	2:H:107:LYS:HG2	1.86	0.56
2:H:37:PRO:CB	5:H:523:HOH:O	2.53	0.56
2:H:41:LEU:O	4:T:49:DPN:CD2	2.54	0.56
2:H:193:GLY:HA2	4:T:49:DPN:CZ	2.37	0.55
2:H:71:HIS:CD2	2:H:154:VAL:HG22	2.41	0.55
2:H:60(C):PRO:HB2	5:H:565:HOH:O	2.05	0.55
1:L:1:CYS:O	2:H:206:ARG:NH1	2.40	0.55
2:H:37:PRO:CG	5:H:523:HOH:O	2.54	0.54
2:H:211:GLY:HA2	2:H:229:THR:O	2.06	0.54
2:H:204(B):ASN:O	2:H:205:ASN:ND2	2.40	0.54
2:H:126:ARG:HH11	2:H:126:ARG:CG	2.21	0.54
2:H:110:LYS:O	5:H:603:HOH:O	2.18	0.54
3:I:64:LEU:CD1	3:I:64:LEU:N	2.70	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:182:CYS:HA	2:H:226:GLY:O	2.08	0.54
2:H:85:LEU:CD1	2:H:106:MET:HE2	2.28	0.53
2:H:87:LYS:HD2	2:H:88:ILE:H	1.73	0.53
2:H:50:ARG:NH1	2:H:107:LYS:HE2	2.24	0.53
2:H:136:GLY:HA3	2:H:199:PHE:CZ	2.44	0.53
3:I:64:LEU:HD12	3:I:64:LEU:N	2.24	0.53
4:T:51:0FL:HA	5:T:610:HOH:O	2.09	0.52
2:H:237:TRP:O	2:H:240:LYS:HB3	2.10	0.52
1:L:1(C):GLU:HB2	2:H:47:ILE:O	2.10	0.52
2:H:66:VAL:CG2	2:H:85:LEU:HD21	2.39	0.52
2:H:187:ARG:NH1	2:H:221:ASP:OD2	2.42	0.52
2:H:186(C):GLY:O	2:H:186(D):LYS:HG3	2.10	0.52
2:H:126:ARG:CB	2:H:126:ARG:HH11	2.23	0.52
2:H:126:ARG:NH1	2:H:126:ARG:CG	2.73	0.51
2:H:204(B):ASN:C	2:H:204(B):ASN:HD22	2.13	0.51
2:H:37:PRO:HB2	5:H:523:HOH:O	2.10	0.51
2:H:195:SER:HG	4:T:48:0MG:C	2.18	0.51
2:H:164:GLU:HB2	2:H:167:VAL:HG23	1.93	0.50
3:I:59:ILE:HD11	3:I:64:LEU:HD12	1.89	0.50
2:H:85:LEU:HD11	2:H:106:MET:HE1	1.93	0.50
2:H:94:TYR:HA	2:H:101:ARG:HB2	1.94	0.50
1:L:14(I):SER:C	1:L:14(K):ILE:H	2.14	0.50
2:H:64:LEU:HD12	2:H:85:LEU:HD11	1.93	0.50
2:H:37:PRO:HG2	5:H:523:HOH:O	2.12	0.49
2:H:230:HIS:ND1	2:H:233:ARG:HG3	2.26	0.49
2:H:127:GLU:O	2:H:129(A):ALA:HB3	2.13	0.49
2:H:60(D):TRP:CH2	4:T:50:VLT:N	2.80	0.49
2:H:165:ARG:N	2:H:166:PRO:HD2	2.28	0.48
2:H:232:PHE:O	2:H:235:LYS:HB2	2.14	0.48
2:H:70:LYS:CE	2:H:72:SER:O	2.61	0.48
2:H:49:ASP:HB3	2:H:114:PHE:CZ	2.49	0.48
1:L:1(D):GLY:H3	2:H:48:SER:HB2	1.77	0.48
2:H:76:TYR:CD2	2:H:76:TYR:O	2.67	0.48
1:L:14(K):ILE:HD12	5:L:469:HOH:O	2.13	0.48
2:H:114:PHE:HA	2:H:118:ILE:HG22	1.96	0.47
2:H:176:ILE:HD12	2:H:227:PHE:CE2	2.48	0.47
2:H:237:TRP:O	2:H:241:VAL:HG13	2.14	0.47
2:H:65:LEU:HG	2:H:82:ILE:CG2	2.45	0.47
2:H:230:HIS:CG	2:H:233:ARG:HG3	2.50	0.46
1:L:9:LYS:NZ	5:L:552:HOH:O	2.48	0.46
2:H:146:GLU:O	2:H:147:THR:C	2.53	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:70:LYS:HB3	2:H:70:LYS:HE3	1.72	0.46
2:H:221(A):ARG:HD3	5:H:465:HOH:O	2.14	0.46
3:I:60:PRO:O	3:I:63:TYS:N	2.49	0.46
2:H:61:GLU:CG	2:H:87:LYS:HA	2.46	0.46
2:H:17:VAL:O	2:H:18:GLU:HB2	2.16	0.46
2:H:49:ASP:O	2:H:111:PRO:HA	2.16	0.46
2:H:205:ASN:HD22	2:H:205:ASN:N	2.08	0.45
2:H:145:LYS:HD2	5:H:544:HOH:O	2.15	0.45
2:H:114:PHE:HD1	2:H:118:ILE:HG22	1.80	0.45
2:H:21:ASP:CB	2:H:154:VAL:CG1	2.91	0.45
2:H:91:HIS:CE1	2:H:101:ARG:HD3	2.51	0.45
1:L:14(D):ARG:NE	1:L:14(H):GLU:OE2	2.49	0.45
2:H:60(B):PRO:HG2	2:H:96:TRP:CD2	2.51	0.45
2:H:60(G):ASN:ND2	5:H:507:HOH:O	2.38	0.45
3:I:59:ILE:CD1	3:I:64:LEU:CD1	2.80	0.45
2:H:65:LEU:HG	2:H:82:ILE:HG22	1.98	0.44
1:L:1(D):GLY:N	2:H:48:SER:HB2	2.32	0.44
2:H:60(B):PRO:N	2:H:60(C):PRO:HD2	2.32	0.44
2:H:143:ASN:HB2	2:H:191:CYS:SG	2.57	0.44
2:H:165:ARG:N	2:H:166:PRO:CD	2.81	0.44
2:H:66:VAL:HG23	2:H:85:LEU:HD21	1.99	0.44
1:L:5:PRO:HB2	2:H:116:ASP:HA	2.00	0.44
3:I:60:PRO:HG2	3:I:63:TYS:CE2	2.47	0.43
2:H:60(F):LYS:CE	5:H:596:HOH:O	2.65	0.43
2:H:205:ASN:CB	5:H:503:HOH:O	2.27	0.43
2:H:65:LEU:HD12	2:H:83:SER:O	2.18	0.43
2:H:167:VAL:H	2:H:167:VAL:HG23	1.49	0.43
2:H:73:ARG:O	3:I:56:PHE:HE2	2.02	0.43
2:H:186:PRO:HB3	2:H:222:ASP:HB3	1.99	0.43
2:H:187:ARG:HH11	2:H:187:ARG:HD3	1.20	0.42
2:H:70:LYS:NZ	2:H:80:GLU:OE2	2.45	0.42
2:H:97:ARG:CD	5:H:496:HOH:O	2.55	0.42
2:H:99:LEU:HD21	4:T:52:PRO:HB3	2.01	0.42
1:L:1(D):GLY:HA3	2:H:114:PHE:HE2	1.83	0.42
2:H:114:PHE:CD1	2:H:114:PHE:N	2.82	0.42
1:L:14(J):TYR:C	1:L:14(K):ILE:HG13	2.40	0.42
2:H:31:VAL:HG12	2:H:32:MET:N	2.33	0.42
2:H:35:ARG:HB2	2:H:41:LEU:HD13	2.00	0.42
2:H:36:LYS:O	2:H:38:GLN:NE2	2.52	0.42
2:H:34:PHE:CE2	2:H:38:GLN:HB3	2.55	0.42
2:H:96:TRP:CH2	2:H:97:ARG:CZ	3.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:61:GLU:OE2	2:H:87:LYS:HD2	2.19	0.42
2:H:119:HIS:HA	2:H:120:PRO:HD3	1.89	0.41
2:H:49:ASP:N	2:H:49:ASP:OD1	2.52	0.41
2:H:130:LEU:HG	2:H:210:MET:HE2	2.01	0.41
2:H:64:LEU:CD1	2:H:85:LEU:HD11	2.49	0.41
2:H:107:LYS:HZ1	2:H:246:GLY:HA3	1.86	0.41
2:H:73:ARG:O	3:I:56:PHE:CE2	2.73	0.41
2:H:203:SER:HB3	2:H:204(B):ASN:ND2	2.35	0.41
2:H:107:LYS:NZ	2:H:246:GLY:HA3	2.35	0.41
2:H:126:ARG:O	2:H:129(A):ALA:HB2	2.20	0.40
2:H:43:GLY:O	2:H:44:ALA:HB2	2.22	0.40
1:L:13:GLU:HA	5:L:406:HOH:O	2.21	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:75:ARG:NH1	3:I:57:GLU:OE1[2_555]	2.18	0.02

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	27/36 (75%)	23 (85%)	2 (7%)	2 (7%)	1 0
2	H	248/259 (96%)	230 (93%)	17 (7%)	1 (0%)	39 48
3	I	7/13 (54%)	5 (71%)	2 (29%)	0	100 100
All	All	282/308 (92%)	258 (92%)	21 (7%)	3 (1%)	17 18

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	1(C)	GLU
2	H	77	GLU
1	L	1(B)	ALA

5.3.2 Protein sidechains (i)

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	26/31 (84%)	23 (88%)	3 (12%)	7 7
2	H	218/225 (97%)	186 (85%)	32 (15%)	4 3
3	I	7/11 (64%)	6 (86%)	1 (14%)	4 4
4	T	1/1 (100%)	1 (100%)	0	100 100
All	All	252/268 (94%)	216 (86%)	36 (14%)	4 4

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	14(A)	LYS
1	L	14(E)	GLU
1	L	14(F)	LEU
2	H	33	LEU
2	H	36	LYS
2	H	36(A)	SER
2	H	41	LEU
2	H	62	ASN
2	H	64	LEU
2	H	65	LEU
2	H	66	VAL
2	H	75	ARG
2	H	79	ILE
2	H	83	SER
2	H	84	MET
2	H	94	TYR
2	H	97(A)	GLU
2	H	99	LEU
2	H	107	LYS

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Mol	Chain	Res	Type
2	H	126	ARG
2	H	127	GLU
2	H	147	THR
2	H	151	GLN
2	H	154	VAL
2	H	187	ARG
2	H	192	GLU
2	H	204(B)	ASN
2	H	205	ASN
2	H	221(A)	ARG
2	H	222	ASP
2	H	235	LYS
2	H	239	GLN
2	H	240	LYS
2	H	243	ASP
2	H	247	GLU
3	I	62	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	H	78	ASN
2	H	156	GLN
2	H	204(B)	ASN
2	H	205	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)

5 non-standard protein/DNA/RNA residues are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
3	TYS	I	63	3	15,16,17	1.57	2 (13%)	16,22,24	2.04	3 (18%)
4	0MG	T	48	2,4	9,12,13	1.52	1 (11%)	8,14,16	2.17	3 (37%)
4	DPN	T	49	4	10,11,12	0.99	0	10,13,15	1.19	1 (10%)
4	VLT	T	50	4	13,14,15	0.84	0	14,17,19	2.97	5 (35%)
4	0FL	T	51	4	5,7,8	1.36	0	3,7,9	5.64	2 (66%)

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	I	63	3	-	0/9/11/13	0/1/1/1
4	0MG	T	48	2,4	-	0/7/13/15	0/0/0/0
4	DPN	T	49	4	-	0/4/6/8	0/1/1/1
4	VLT	T	50	4	-	0/7/8/9	0/1/1/1
4	0FL	T	51	4	-	0/3/7/9	0/0/0/0

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	I	63	TYS	OH-CZ	-4.58	1.35	1.42
3	I	63	TYS	CB-CA	2.64	1.59	1.53
4	T	48	0MG	C-C2	3.88	1.52	1.47

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	T	51	0FL	O1-C1-N	-9.07	111.70	124.76
4	T	50	VLT	C2-CA-N	-6.80	96.83	110.99
3	I	63	TYS	CG-CB-CA	-6.05	100.53	114.21
4	T	50	VLT	O-C-C10	-6.03	110.09	125.51
4	T	50	VLT	C4-C3-CA	-4.71	103.57	114.21
4	T	51	0FL	O-C-CA	-3.51	116.16	125.44
4	T	49	DPN	O-C-CA	-3.13	117.34	125.49
3	I	63	TYS	CB-CG-CD1	-2.66	115.33	120.90
4	T	50	VLT	C3-C4-C5	-2.53	115.60	120.90
4	T	50	VLT	C3-C4-C6	2.05	125.19	120.90
4	T	48	0MG	C3-CA-N	2.60	115.26	109.81
3	I	63	TYS	O3-S-O1	2.73	118.82	108.56
4	T	48	0MG	O-C-C2	3.39	128.81	124.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	T	48	0MG	C2-CA-N	4.17	116.00	109.27

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	I	63	TYS	3	0
4	T	48	0MG	5	0
4	T	49	DPN	4	0
4	T	50	VLT	1	0
4	T	51	0FL	1	0

5.5 Carbohydrates (i)

There are no carbohydrates in this entry.

5.6 Ligand geometry (i)

There are no ligands in this entry.

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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.