Full wwPDB X-ray Structure Validation Report (i)

Jan 31, 2016 – 07:22 PM GMT

PDB ID	:	1FAK
Title	:	HUMAN TISSUE FACTOR COMPLEXED WITH COAGULATION FACTOR VIIA INHIBITED WITH A BPTI-MUTANT
Authors	:	Zhang, E.; St Charles, R.; Tulinsky, A.
Deposited on	:	1998-12-28
Resolution	:	2.10 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:

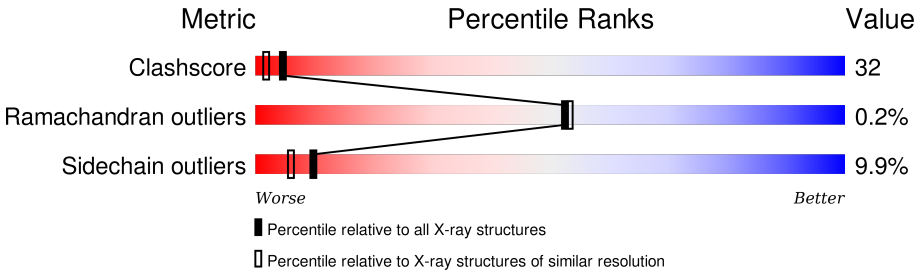
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
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 2.10 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	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)

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	152	36% 26% 7% • 29%
2	H	254	50% 37% 10% ••
3	T	206	35% 42% 11% 11%
4	I	55	20% 58% 20% •

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
5	GLC	L	600	X	-	-	-

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 5079 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 PROTEIN (BLOOD COAGULATION FACTOR VIIA).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	108	Total	C	N	O	S	0	0	0
			827	504	143	167	13			

- Molecule 2 is a protein called PROTEIN (BLOOD COAGULATION FACTOR VIIA).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	251	Total	C	N	O	S	0	0	0
			1955	1242	348	352	13			

- Molecule 3 is a protein called PROTEIN (SOLUBLE TISSUE FACTOR).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	T	183	Total	C	N	O	S	0	0	0
			1495	957	240	293	5			

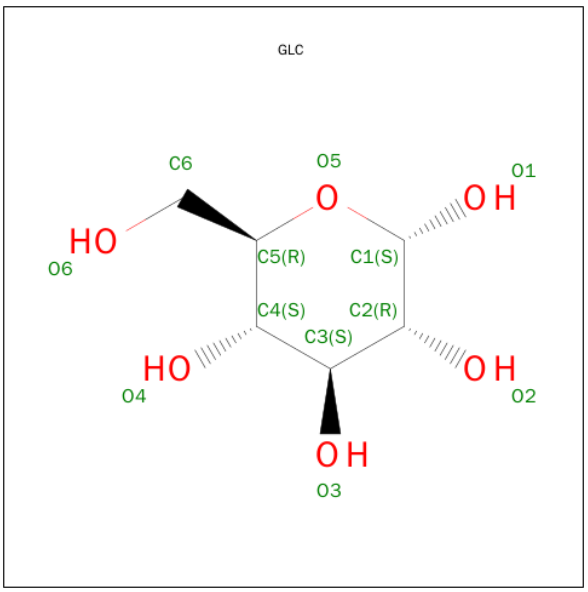
- Molecule 4 is a protein called PROTEIN (5L15).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	I	55	Total	C	N	O	S	0	0	0
			439	277	75	80	7			

There are 8 discrepancies between the modelled and reference sequences:

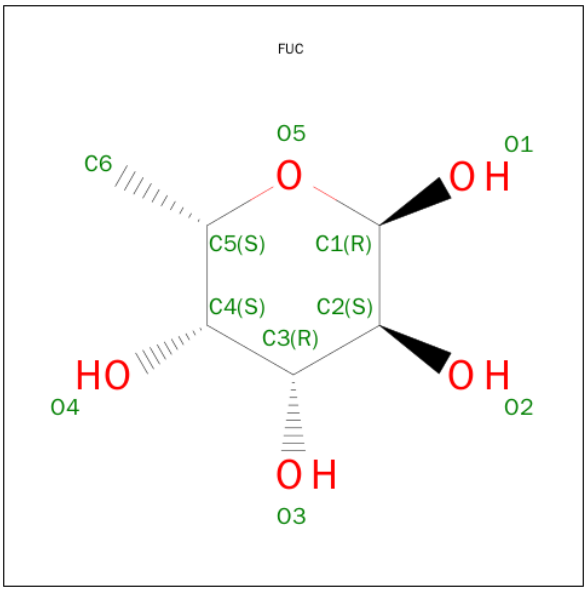
Chain	Residue	Modelled	Actual	Comment	Reference
I	11	ASP	THR	CONFLICT	UNP P00974
I	15	ARG	LYS	CONFLICT	UNP P00974
I	17	LEU	ARG	CONFLICT	UNP P00974
I	18	HIS	ILE	CONFLICT	UNP P00974
I	19	LEU	ILE	CONFLICT	UNP P00974
I	34	TYR	VAL	CONFLICT	UNP P00974
I	39	LEU	ARG	CONFLICT	UNP P00974
I	46	GLU	LYS	CONFLICT	UNP P00974

- Molecule 5 is SUGAR (GLUCOSE) (three-letter code: GLC) (formula: $C_6H_{12}O_6$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	L	1	Total C O 11 6 5	0	0

- Molecule 6 is SUGAR (FUCOSE) (three-letter code: FUC) (formula: $C_6H_{12}O_5$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	L	1	Total C O 10 6 4	0	0

- Molecule 7 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	H	1	Total Ca 1 1	0	0
7	L	1	Total Ca 1 1	0	0

- Molecule 8 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	H	164	Total O 164 164	0	0
8	I	23	Total O 23 23	0	0
8	L	78	Total O 78 78	0	0
8	T	75	Total O 75 75	0	0

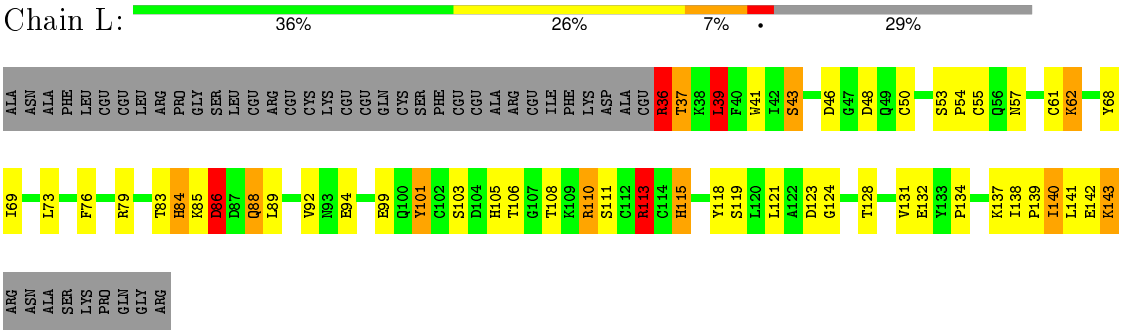
3 Residue-property plots (i)

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.

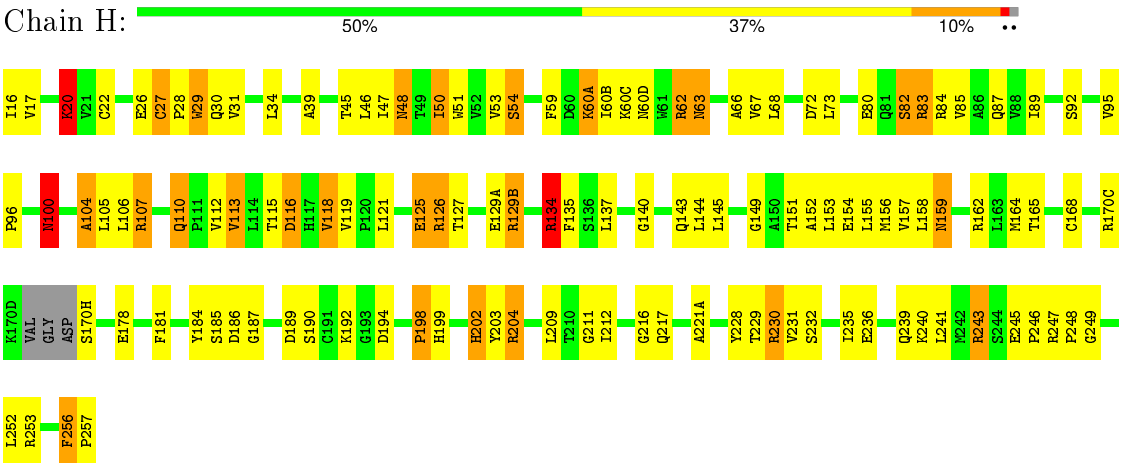
- Molecule 1: PROTEIN (BLOOD COAGULATION FACTOR VIIA)

Chain L:



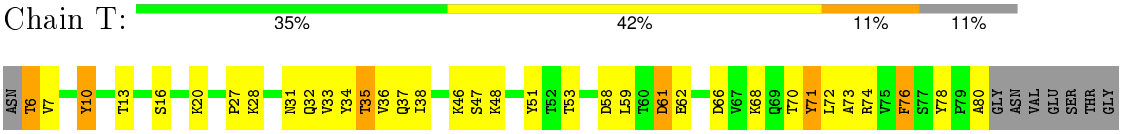
- Molecule 2: PROTEIN (BLOOD COAGULATION FACTOR VIIA)

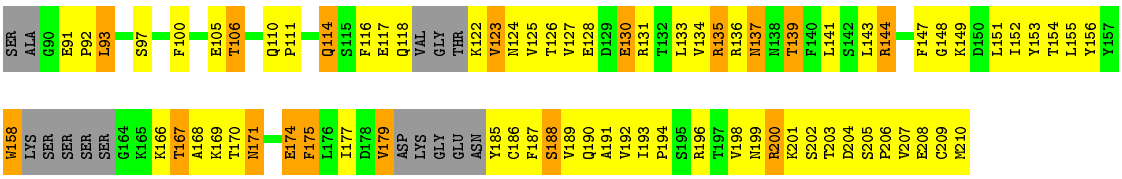
Chain H:



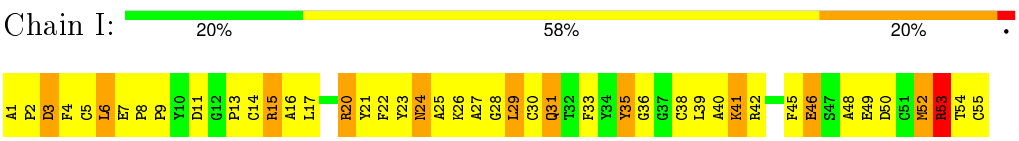
- Molecule 3: PROTEIN (SOLUBLE TISSUE FACTOR)

Chain T:





- Molecule 4: PROTEIN (5L15)



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 2 2 21	Depositor
Cell constants a, b, c, α, β, γ	63.49Å 190.00Å 175.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	9.00 – 2.10	Depositor
% Data completeness (in resolution range)	80.0 (9.00-2.10)	Depositor
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROLSQ	Depositor
R, R_{free}	0.237 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5079	wwPDB-VP
Average B, all atoms (Å^2)	27.0	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: CA, GLC, FUC

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	1.16	0/844	2.03	23/1139 (2.0%)
2	H	1.37	4/2004 (0.2%)	2.17	74/2726 (2.7%)
3	T	1.11	0/1527	1.85	30/2075 (1.4%)
4	I	1.14	0/452	1.98	12/609 (2.0%)
All	All	1.24	4/4827 (0.1%)	2.03	139/6549 (2.1%)

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
4	I	0	1
All	All	0	3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	92	SER	CB-OG	7.46	1.51	1.42
2	H	168	CYS	CB-SG	5.26	1.91	1.82
2	H	228	TYR	CA-CB	5.19	1.65	1.53
2	H	152	ALA	CA-CB	5.11	1.63	1.52

All (139) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	H	253	ARG	NE-CZ-NH1	20.24	130.42	120.30
2	H	107	ARG	CD-NE-CZ	18.48	149.47	123.60
1	L	36	ARG	NE-CZ-NH1	-15.34	112.63	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	H	126	ARG	NE-CZ-NH1	15.32	127.96	120.30
2	H	247	ARG	NE-CZ-NH1	15.17	127.89	120.30
2	H	107	ARG	NE-CZ-NH1	14.56	127.58	120.30
2	H	243	ARG	NE-CZ-NH1	13.59	127.09	120.30
3	T	135	ARG	NE-CZ-NH1	13.53	127.06	120.30
1	L	79	ARG	NE-CZ-NH2	-12.74	113.93	120.30
2	H	134	ARG	NE-CZ-NH1	-12.08	114.26	120.30
2	H	125	GLU	OE1-CD-OE2	12.06	137.77	123.30
2	H	170(C)	ARG	NE-CZ-NH1	11.46	126.03	120.30
1	L	101	TYR	CB-CG-CD2	-10.80	114.52	121.00
1	L	101	TYR	CB-CG-CD1	10.76	127.45	121.00
1	L	123	ASP	CB-CG-OD1	10.72	127.94	118.30
2	H	62	ARG	NE-CZ-NH2	-10.50	115.05	120.30
2	H	170(C)	ARG	CD-NE-CZ	10.49	138.29	123.60
3	T	131	ARG	CD-NE-CZ	10.49	138.29	123.60
2	H	116	ASP	CB-CG-OD1	9.94	127.25	118.30
4	I	42	ARG	NE-CZ-NH1	9.91	125.26	120.30
2	H	126	ARG	CD-NE-CZ	9.38	136.73	123.60
2	H	62	ARG	NE-CZ-NH1	9.31	124.96	120.30
1	L	36	ARG	NE-CZ-NH2	9.15	124.88	120.30
2	H	83	ARG	NE-CZ-NH2	-9.13	115.73	120.30
2	H	247	ARG	CD-NE-CZ	9.13	136.38	123.60
1	L	79	ARG	NE-CZ-NH1	9.12	124.86	120.30
3	T	58	ASP	CB-CG-OD2	-8.87	110.32	118.30
3	T	204	ASP	CB-CG-OD1	8.84	126.25	118.30
2	H	83	ARG	NE-CZ-NH1	8.64	124.62	120.30
2	H	107	ARG	NE-CZ-NH2	-8.63	115.98	120.30
3	T	66	ASP	CB-CG-OD1	8.39	125.85	118.30
4	I	20	ARG	NE-CZ-NH1	8.29	124.44	120.30
1	L	113	ARG	CA-CB-CG	8.25	131.55	113.40
4	I	50	ASP	CB-CG-OD2	8.10	125.59	118.30
4	I	20	ARG	CD-NE-CZ	8.08	134.91	123.60
3	T	200	ARG	NE-CZ-NH1	8.05	124.32	120.30
2	H	204	ARG	NE-CZ-NH2	-8.01	116.29	120.30
3	T	144	ARG	CD-NE-CZ	-7.93	112.50	123.60
1	L	36	ARG	CD-NE-CZ	-7.83	112.64	123.60
2	H	116	ASP	CB-CG-OD2	-7.81	111.27	118.30
2	H	253	ARG	CD-NE-CZ	7.74	134.43	123.60
2	H	186	ASP	CB-CG-OD1	7.49	125.04	118.30
3	T	144	ARG	NE-CZ-NH2	7.42	124.01	120.30
3	T	66	ASP	CB-CG-OD2	-7.10	111.91	118.30
3	T	70	THR	CA-CB-CG2	7.05	122.28	112.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	H	85	VAL	CA-CB-CG1	7.04	121.46	110.90
2	H	248	PRO	N-CA-C	6.99	130.28	112.10
3	T	135	ARG	NE-CZ-NH2	-6.93	116.84	120.30
1	L	99	GLU	CG-CD-OE2	-6.88	104.55	118.30
2	H	253	ARG	NH1-CZ-NH2	-6.87	111.84	119.40
2	H	243	ARG	CD-NE-CZ	6.86	133.21	123.60
2	H	243	ARG	NE-CZ-NH2	-6.78	116.91	120.30
3	T	61	ASP	CB-CG-OD1	6.74	124.37	118.30
1	L	99	GLU	CG-CD-OE1	6.62	131.53	118.30
3	T	179	VAL	CB-CA-C	6.58	123.90	111.40
4	I	14	CYS	CB-CA-C	6.41	123.22	110.40
2	H	217	GLN	CB-CA-C	6.41	123.22	110.40
2	H	151	THR	CA-CB-CG2	6.37	121.32	112.40
2	H	184	TYR	CB-CG-CD2	6.36	124.82	121.00
1	L	46	ASP	CB-CG-OD2	-6.35	112.59	118.30
2	H	80	GLU	CA-CB-CG	6.33	127.33	113.40
2	H	185	SER	CA-CB-OG	6.31	128.24	111.20
3	T	158	TRP	CA-CB-CG	6.31	125.68	113.70
2	H	39	ALA	N-CA-CB	6.26	118.87	110.10
2	H	181	PHE	N-CA-CB	6.25	121.86	110.60
2	H	217	GLN	CA-CB-CG	6.21	127.07	113.40
2	H	100	ASN	CB-CA-C	6.21	122.81	110.40
2	H	39	ALA	O-C-N	6.14	132.53	122.70
2	H	54	SER	O-C-N	6.13	132.51	122.70
1	L	94	GLU	CG-CD-OE2	6.08	130.47	118.30
2	H	185	SER	N-CA-CB	6.03	119.55	110.50
1	L	84	HIS	CA-CB-CG	6.03	123.85	113.60
2	H	129(B)	ARG	NE-CZ-NH2	-6.00	117.30	120.30
2	H	190	SER	N-CA-CB	5.95	119.43	110.50
3	T	179	VAL	CA-CB-CG2	5.95	119.83	110.90
2	H	199	HIS	N-CA-C	-5.89	95.09	111.00
3	T	131	ARG	NE-CZ-NH1	5.88	123.24	120.30
1	L	57	ASN	N-CA-CB	-5.87	100.03	110.60
2	H	17	VAL	C-N-CA	5.82	134.52	122.30
2	H	198	PRO	C-N-CA	5.77	136.14	121.70
3	T	74	ARG	NE-CZ-NH2	5.77	123.19	120.30
2	H	84	ARG	NE-CZ-NH1	5.75	123.17	120.30
2	H	50	ILE	CA-CB-CG2	5.75	122.39	110.90
2	H	118	VAL	CA-CB-CG2	5.73	119.49	110.90
1	L	123	ASP	CB-CG-OD2	-5.70	113.17	118.30
3	T	117	GLU	CA-CB-CG	5.68	125.90	113.40
3	T	35	THR	CA-CB-OG1	-5.63	97.18	109.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	H	125	GLU	CB-CA-C	-5.62	99.15	110.40
2	H	247	ARG	NE-CZ-NH2	-5.62	117.49	120.30
3	T	106	THR	CA-CB-CG2	5.60	120.24	112.40
3	T	131	ARG	NE-CZ-NH2	-5.59	117.51	120.30
3	T	200	ARG	NE-CZ-NH2	-5.58	117.51	120.30
1	L	142	GLU	C-N-CA	5.57	135.63	121.70
2	H	104	ALA	N-CA-CB	-5.55	102.33	110.10
2	H	125	GLU	CG-CD-OE2	-5.54	107.21	118.30
2	H	247	ARG	CG-CD-NE	5.52	123.39	111.80
1	L	137	LYS	N-CA-CB	5.50	120.50	110.60
2	H	67	VAL	O-C-N	5.50	131.50	122.70
4	I	49	GLU	CA-CB-CG	5.50	125.50	113.40
1	L	94	GLU	CG-CD-OE1	-5.49	107.31	118.30
1	L	86	ASP	CB-CG-OD2	5.48	123.23	118.30
4	I	53	ARG	NE-CZ-NH1	5.47	123.03	120.30
3	T	144	ARG	NE-CZ-NH1	-5.45	117.57	120.30
1	L	99	GLU	O-C-N	5.43	131.39	122.70
2	H	113	VAL	CB-CA-C	5.43	121.71	111.40
4	I	42	ARG	NE-CZ-NH2	-5.43	117.59	120.30
2	H	82	SER	N-CA-CB	-5.42	102.37	110.50
2	H	184	TYR	CB-CG-CD1	-5.38	117.77	121.00
2	H	27	CYS	N-CA-CB	-5.37	100.94	110.60
2	H	256	PHE	CB-CG-CD1	-5.36	117.05	120.80
2	H	248	PRO	CA-C-N	5.35	126.90	116.20
3	T	137	ASN	CB-CA-C	5.29	120.98	110.40
2	H	202	HIS	CA-CB-CG	-5.29	104.61	113.60
2	H	189	ASP	N-CA-CB	-5.28	101.09	110.60
3	T	76	PHE	CA-CB-CG	5.27	126.56	113.90
2	H	178	GLU	CG-CD-OE1	5.27	128.84	118.30
3	T	10	TYR	CB-CG-CD1	-5.25	117.85	121.00
2	H	253	ARG	NE-CZ-NH2	-5.23	117.68	120.30
2	H	154	GLU	OE1-CD-OE2	5.23	129.58	123.30
2	H	155	LEU	O-C-N	5.22	131.04	122.70
2	H	162	ARG	CA-CB-CG	5.21	124.87	113.40
4	I	42	ARG	CD-NE-CZ	5.19	130.87	123.60
4	I	35	TYR	CB-CG-CD2	-5.18	117.89	121.00
2	H	186	ASP	CB-CG-OD2	-5.16	113.66	118.30
3	T	143	LEU	CB-CA-C	5.13	119.95	110.20
2	H	60(A)	LYS	N-CA-CB	5.12	119.82	110.60
4	I	11	ASP	CB-CA-C	5.10	120.60	110.40
2	H	72	ASP	CB-CG-OD1	5.09	122.88	118.30
3	T	53	THR	N-CA-CB	5.09	119.97	110.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	H	230	ARG	C-N-CA	5.08	134.41	121.70
3	T	130	GLU	CA-CB-CG	5.08	124.58	113.40
2	H	126	ARG	NE-CZ-NH2	-5.07	117.77	120.30
1	L	115	HIS	CA-CB-CG	5.07	122.22	113.60
2	H	48	ASN	N-CA-CB	5.07	119.72	110.60
2	H	20	LYS	CB-CA-C	-5.05	100.30	110.40
3	T	71	TYR	C-N-CA	5.04	134.31	121.70
1	L	39	LEU	CB-CA-C	5.03	119.76	110.20
2	H	162	ARG	NE-CZ-NH2	-5.03	117.78	120.30
4	I	15	ARG	NE-CZ-NH1	5.01	122.81	120.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	H	134	ARG	Sidechain
2	H	230	ARG	Sidechain
4	I	53	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	827	0	755	44	0
2	H	1955	0	1933	92	0
3	T	1495	0	1455	134	0
4	I	439	0	401	54	0
5	L	11	0	10	1	0
6	L	10	0	10	0	0
7	H	1	0	0	0	0
7	L	1	0	0	0	0
8	H	164	0	0	9	0
8	I	23	0	0	2	0
8	L	78	0	0	0	0
8	T	75	0	0	6	0
All	All	5079	0	4564	298	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 32.

All (298) 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)
3:T:158:TRP:CH2	3:T:207:VAL:HG11	1.77	1.20
3:T:155:LEU:HB2	3:T:175:PHE:CE2	1.82	1.15
3:T:185:TYR:O	3:T:209:CYS:HA	1.44	1.14
3:T:186:CYS:HA	3:T:208:GLU:O	1.58	1.04
2:H:46:LEU:HD13	2:H:68:LEU:HD11	1.39	1.03
3:T:7:VAL:HG23	3:T:32:GLN:HE22	1.22	1.03
3:T:158:TRP:HH2	3:T:207:VAL:CG1	1.74	1.01
2:H:68:LEU:HD12	2:H:112:VAL:HG11	1.38	1.01
3:T:158:TRP:HH2	3:T:207:VAL:HG11	1.21	0.99
3:T:155:LEU:CB	3:T:175:PHE:CE2	2.44	0.99
3:T:158:TRP:CH2	3:T:207:VAL:CG1	2.45	0.99
1:L:140:ILE:CD1	2:H:26:GLU:HG2	1.93	0.98
3:T:185:TYR:HB2	3:T:210:MET:SD	2.07	0.95
3:T:155:LEU:HB2	3:T:175:PHE:CD2	2.03	0.94
4:I:7:GLU:OE2	4:I:41:LYS:NZ	2.01	0.93
3:T:136:ARG:O	3:T:139:THR:HG23	1.70	0.91
1:L:140:ILE:HD11	2:H:26:GLU:HG2	1.52	0.90
3:T:155:LEU:CB	3:T:175:PHE:CD2	2.56	0.89
3:T:20:LYS:HB2	3:T:133:LEU:HD13	1.54	0.89
3:T:7:VAL:HG23	3:T:32:GLN:NE2	1.88	0.89
3:T:7:VAL:CG2	3:T:32:GLN:NE2	2.38	0.87
3:T:155:LEU:HD11	3:T:187:PHE:HB3	1.57	0.86
2:H:60(D):ASN:HB3	2:H:63:ASN:HD21	1.38	0.86
3:T:185:TYR:O	3:T:209:CYS:CA	2.23	0.85
2:H:187:GLY:HA2	2:H:221(A):ALA:O	1.77	0.84
3:T:192:VAL:HG13	3:T:201:LYS:HG2	1.61	0.82
2:H:164:MET:CE	3:T:91:GLU:HB3	2.09	0.82
4:I:27:ALA:HB1	4:I:29:LEU:CD1	2.10	0.82
2:H:60(C):LYS:NZ	4:I:46:GLU:HG2	1.95	0.81
2:H:60(D):ASN:CB	2:H:63:ASN:HD21	1.92	0.81
1:L:138:ILE:HG22	1:L:140:ILE:HG13	1.62	0.81
4:I:27:ALA:CB	4:I:29:LEU:HD12	2.12	0.80
2:H:45:THR:OG1	2:H:198:PRO:HB3	1.82	0.79
2:H:164:MET:HE3	3:T:91:GLU:HG2	1.63	0.79
3:T:137:ASN:O	3:T:139:THR:HG22	1.82	0.79
4:I:27:ALA:HB3	4:I:29:LEU:HD12	1.64	0.78
4:I:27:ALA:HB1	4:I:29:LEU:HD11	1.65	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:T:196:ARG:O	3:T:200:ARG:HD2	1.83	0.78
4:I:27:ALA:CB	4:I:29:LEU:CD1	2.61	0.77
1:L:143:LYS:HG3	1:L:143:LYS:O	1.82	0.77
2:H:60(D):ASN:HB3	2:H:63:ASN:ND2	2.00	0.76
2:H:60(D):ASN:CB	2:H:63:ASN:ND2	2.50	0.75
3:T:136:ARG:HB3	3:T:141:LEU:HD21	1.69	0.74
4:I:24:ASN:OD1	4:I:31:GLN:NE2	2.21	0.74
2:H:165:THR:HG21	8:H:364:HOH:O	1.86	0.74
3:T:155:LEU:HB3	3:T:175:PHE:CD2	2.23	0.73
4:I:35:TYR:CZ	4:I:40:ALA:HB2	2.24	0.73
3:T:155:LEU:N	3:T:175:PHE:HE2	1.87	0.72
2:H:83:ARG:HG2	8:H:383:HOH:O	1.89	0.72
3:T:32:GLN:HA	3:T:78:TYR:O	1.91	0.71
1:L:140:ILE:HG12	8:H:317:HOH:O	1.91	0.71
3:T:196:ARG:O	3:T:200:ARG:CD	2.39	0.71
3:T:185:TYR:O	3:T:210:MET:N	2.25	0.70
1:L:121:LEU:HD21	1:L:128:THR:CG2	2.23	0.69
4:I:24:ASN:O	4:I:28:GLY:N	2.26	0.69
3:T:35:THR:HG21	8:T:230:HOH:O	1.93	0.68
3:T:158:TRP:CZ2	3:T:207:VAL:HG11	2.29	0.68
4:I:29:LEU:HA	4:I:52:MET:HE1	1.75	0.68
3:T:186:CYS:CA	3:T:208:GLU:O	2.39	0.67
4:I:25:ALA:O	4:I:28:GLY:N	2.28	0.67
4:I:6:LEU:N	4:I:6:LEU:HD23	2.09	0.66
3:T:6:THR:HG21	3:T:78:TYR:CA	2.26	0.66
3:T:196:ARG:HB2	3:T:200:ARG:HG2	1.76	0.66
3:T:179:VAL:HB	3:T:185:TYR:CE2	2.31	0.65
2:H:164:MET:HE3	3:T:91:GLU:CG	2.27	0.65
4:I:23:TYR:OH	4:I:28:GLY:HA2	1.95	0.65
2:H:50:ILE:HD12	2:H:107:ARG:HG3	1.80	0.63
3:T:37:GLN:O	3:T:73:ALA:HA	1.98	0.63
1:L:92:VAL:HG12	2:H:129(B):ARG:HG2	1.81	0.63
3:T:155:LEU:N	3:T:175:PHE:CE2	2.66	0.63
2:H:89:ILE:HG23	2:H:252:LEU:HB3	1.81	0.62
3:T:123:VAL:CG1	3:T:177:ILE:HD11	2.29	0.62
1:L:101:TYR:CE2	2:H:125:GLU:HG3	2.34	0.62
1:L:110:ARG:HD2	1:L:110:ARG:C	2.20	0.61
3:T:156:TYR:HA	3:T:166:LYS:O	1.99	0.61
3:T:156:TYR:O	3:T:187:PHE:HA	2.00	0.61
2:H:236:GLU:OE2	2:H:240:LYS:HE3	2.00	0.61
3:T:155:LEU:O	3:T:167:THR:HA	2.01	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:T:114:GLN:HG3	3:T:128:GLU:HB2	1.83	0.60
4:I:39:LEU:HD12	4:I:39:LEU:N	2.16	0.60
3:T:179:VAL:HB	3:T:185:TYR:HE2	1.66	0.60
3:T:6:THR:HG21	3:T:78:TYR:N	2.16	0.60
3:T:193:ILE:HG21	8:T:241:HOH:O	2.02	0.60
4:I:27:ALA:CB	4:I:29:LEU:HD11	2.29	0.59
2:H:73:LEU:HD12	8:H:321:HOH:O	2.03	0.59
2:H:60(C):LYS:HZ3	4:I:46:GLU:HG2	1.66	0.59
4:I:38:CYS:SG	4:I:39:LEU:HD13	2.42	0.59
3:T:158:TRP:CH2	3:T:207:VAL:HG13	2.37	0.58
3:T:202:SER:OG	3:T:203:THR:O	2.21	0.58
3:T:203:THR:HG22	8:T:220:HOH:O	2.03	0.58
3:T:48:LYS:HG3	3:T:59:LEU:CD2	2.33	0.58
3:T:192:VAL:O	3:T:194:PRO:HD3	2.03	0.58
1:L:140:ILE:HD12	2:H:26:GLU:HG2	1.85	0.58
3:T:124:ASN:ND2	8:T:263:HOH:O	2.37	0.58
2:H:241:LEU:HG	8:H:382:HOH:O	2.02	0.58
2:H:59:PHE:HA	2:H:60(B):ILE:HG12	1.86	0.58
2:H:140:GLY:HA3	2:H:194:ASP:OD1	2.04	0.57
3:T:185:TYR:HB3	3:T:187:PHE:CZ	2.39	0.57
3:T:36:VAL:O	3:T:47:SER:HA	2.03	0.57
3:T:207:VAL:HG12	3:T:208:GLU:N	2.20	0.57
3:T:7:VAL:HG22	3:T:32:GLN:NE2	2.20	0.57
3:T:114:GLN:HB2	3:T:126:THR:HG22	1.87	0.56
3:T:72:LEU:HD11	3:T:97:SER:O	2.05	0.56
1:L:36:ARG:NH2	3:T:190:GLN:NE2	2.53	0.56
1:L:140:ILE:HD11	2:H:26:GLU:HA	1.88	0.56
2:H:54:SER:O	2:H:212:ILE:HD13	2.05	0.56
1:L:121:LEU:HD21	1:L:128:THR:HG21	1.87	0.56
1:L:105:HIS:NE2	1:L:111:SER:OG	2.25	0.56
1:L:140:ILE:CD1	2:H:26:GLU:CG	2.79	0.56
3:T:196:ARG:HB2	3:T:200:ARG:CG	2.35	0.56
2:H:143:GLN:HG2	2:H:149:GLY:O	2.05	0.56
4:I:27:ALA:HB3	4:I:29:LEU:CD1	2.31	0.56
2:H:216:GLY:HA3	4:I:15:ARG:NH2	2.20	0.56
3:T:187:PHE:O	3:T:207:VAL:HG13	2.06	0.55
3:T:6:THR:HG22	3:T:32:GLN:OE1	2.06	0.55
3:T:155:LEU:CD1	3:T:187:PHE:HB3	2.34	0.55
2:H:164:MET:CE	3:T:91:GLU:CG	2.85	0.55
2:H:126:ARG:O	2:H:129(A):GLU:HG3	2.06	0.55
4:I:7:GLU:HG3	8:I:112:HOH:O	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:127:THR:O	2:H:129(B):ARG:HB2	2.06	0.54
4:I:38:CYS:SG	4:I:39:LEU:CD1	2.95	0.54
3:T:144:ARG:CZ	3:T:151:LEU:HD23	2.37	0.54
4:I:4:PHE:HZ	4:I:54:THR:HG21	1.71	0.54
3:T:16:SER:HA	3:T:20:LYS:O	2.07	0.54
3:T:76:PHE:HD2	3:T:92:PRO:HB2	1.72	0.54
2:H:27:CYS:N	2:H:28:PRO:CD	2.70	0.54
3:T:76:PHE:CD2	3:T:92:PRO:HB2	2.43	0.54
1:L:110:ARG:HD2	1:L:110:ARG:O	2.08	0.54
3:T:153:TYR:O	3:T:169:LYS:HA	2.07	0.54
2:H:164:MET:CE	3:T:91:GLU:CB	2.85	0.54
2:H:245:GLU:CD	2:H:246:PRO:HD2	2.28	0.53
3:T:34:TYR:HA	3:T:76:PHE:O	2.08	0.53
2:H:46:LEU:HD11	2:H:48:ASN:O	2.08	0.53
4:I:23:TYR:CZ	4:I:28:GLY:HA2	2.42	0.53
2:H:22:CYS:HB2	2:H:157:VAL:HB	1.90	0.53
3:T:37:GLN:HA	3:T:46:LYS:O	2.09	0.53
2:H:60(D):ASN:HB2	2:H:63:ASN:ND2	2.22	0.52
3:T:48:LYS:HG3	3:T:59:LEU:HD23	1.90	0.52
4:I:1:ALA:HB2	4:I:55:CYS:HA	1.90	0.52
3:T:33:VAL:CG2	3:T:78:TYR:HB2	2.39	0.52
2:H:60(D):ASN:HB2	2:H:63:ASN:HD21	1.74	0.52
2:H:50:ILE:HG13	2:H:51:TRP:CD1	2.44	0.52
1:L:141:LEU:HD23	1:L:141:LEU:N	2.24	0.52
1:L:131:VAL:O	1:L:134:PRO:HD3	2.10	0.51
2:H:83:ARG:HD2	2:H:110:GLN:HG3	1.92	0.51
1:L:101:TYR:CD2	2:H:125:GLU:HG3	2.45	0.51
2:H:20:LYS:O	2:H:156:MET:HA	2.10	0.51
4:I:1:ALA:CB	4:I:55:CYS:HA	2.40	0.51
2:H:135:PHE:HB3	2:H:159:ASN:ND2	2.25	0.51
3:T:71:TYR:O	3:T:100:PHE:N	2.41	0.51
2:H:30:GLN:HG3	2:H:31:VAL:N	2.25	0.51
4:I:48:ALA:O	4:I:52:MET:N	2.36	0.51
4:I:7:GLU:CG	8:I:112:HOH:O	2.60	0.50
3:T:144:ARG:NH1	3:T:151:LEU:HB3	2.26	0.50
4:I:16:ALA:HB3	4:I:36:GLY:O	2.12	0.50
3:T:72:LEU:CD1	3:T:97:SER:O	2.60	0.50
2:H:235:ILE:O	2:H:239:GLN:HG3	2.11	0.50
3:T:6:THR:CG2	3:T:78:TYR:C	2.80	0.50
3:T:114:GLN:HG3	3:T:128:GLU:CB	2.42	0.50
1:L:41:TRP:CE3	1:L:41:TRP:HA	2.47	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:46:LEU:CD1	2:H:48:ASN:O	2.60	0.49
3:T:33:VAL:HG23	3:T:78:TYR:HD2	1.75	0.49
3:T:207:VAL:CG1	3:T:208:GLU:N	2.75	0.49
1:L:85:LYS:HE2	3:T:61:ASP:OD2	2.12	0.49
3:T:155:LEU:HB3	3:T:175:PHE:CE2	2.40	0.49
2:H:60(C):LYS:NZ	4:I:46:GLU:CG	2.70	0.49
2:H:159:ASN:ND2	8:H:326:HOH:O	2.44	0.49
2:H:16:ILE:N	2:H:194:ASP:OD2	2.46	0.48
3:T:147:PHE:O	3:T:148:GLY:C	2.49	0.48
2:H:126:ARG:HG3	2:H:232:SER:O	2.13	0.48
3:T:31:ASN:HA	3:T:80:ALA:HB3	1.95	0.48
3:T:118:GLN:NE2	3:T:210:MET:HE1	2.29	0.48
3:T:155:LEU:HB2	3:T:175:PHE:CZ	2.44	0.48
3:T:127:VAL:CG2	3:T:189:VAL:HG11	2.44	0.48
1:L:115:HIS:O	1:L:118:TYR:HB2	2.13	0.48
3:T:116:PHE:HA	3:T:124:ASN:O	2.14	0.48
1:L:39:LEU:O	1:L:43:SER:HB2	2.13	0.48
3:T:170:THR:HG21	3:T:174:GLU:O	2.14	0.48
3:T:155:LEU:CA	3:T:175:PHE:CE2	2.96	0.47
2:H:45:THR:HG21	2:H:121:LEU:HD23	1.96	0.47
1:L:85:LYS:CE	3:T:61:ASP:OD2	2.62	0.47
2:H:209:LEU:HG	2:H:231:VAL:HG11	1.96	0.47
2:H:137:LEU:HA	2:H:158:LEU:O	2.14	0.47
4:I:38:CYS:C	4:I:39:LEU:HD12	2.35	0.47
4:I:23:TYR:HA	4:I:30:CYS:HA	1.96	0.47
3:T:170:THR:OG1	3:T:171:ASN:N	2.47	0.47
2:H:60(A):LYS:HE2	8:H:421:HOH:O	2.13	0.47
3:T:106:THR:O	3:T:196:ARG:NH1	2.43	0.47
3:T:33:VAL:HG22	3:T:78:TYR:HB2	1.97	0.47
4:I:25:ALA:C	4:I:28:GLY:H	2.18	0.47
2:H:143:GLN:HB3	2:H:145:LEU:O	2.15	0.47
4:I:4:PHE:O	4:I:7:GLU:HB3	2.14	0.47
3:T:38:ILE:HA	3:T:73:ALA:HA	1.96	0.47
2:H:135:PHE:HB3	2:H:159:ASN:HD21	1.80	0.47
3:T:118:GLN:NE2	3:T:210:MET:CE	2.78	0.47
4:I:24:ASN:HD21	4:I:26:LYS:HB2	1.79	0.46
4:I:24:ASN:ND2	4:I:26:LYS:HB2	2.30	0.46
8:H:390:HOH:O	4:I:46:GLU:HG3	2.16	0.46
2:H:87:GLN:NE2	2:H:252:LEU:CD1	2.78	0.46
3:T:114:GLN:HE21	3:T:114:GLN:HB3	1.42	0.46
1:L:36:ARG:NH2	3:T:190:GLN:HE22	2.13	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:192:LYS:HE3	8:H:419:HOH:O	2.15	0.46
2:H:68:LEU:HD13	2:H:118:VAL:HG13	1.96	0.46
3:T:122:LYS:HA	3:T:177:ILE:O	2.15	0.46
5:L:600:GLC:O6	5:L:600:GLC:O4	2.21	0.46
2:H:68:LEU:HD12	2:H:112:VAL:CG1	2.28	0.46
3:T:205:SER:HB2	3:T:206:PRO:CD	2.45	0.46
2:H:256:PHE:CD1	2:H:257:PRO:HA	2.51	0.46
2:H:192:LYS:HD2	4:I:17:LEU:HD23	1.98	0.46
2:H:47:ILE:HD13	2:H:53:VAL:CG2	2.46	0.46
1:L:110:ARG:C	1:L:110:ARG:CD	2.85	0.45
1:L:61:CYS:SG	1:L:68:TYR:HB2	2.56	0.45
1:L:140:ILE:HD11	2:H:26:GLU:CG	2.35	0.45
1:L:48:ASP:OD1	1:L:50:CYS:HB2	2.16	0.45
1:L:138:ILE:CG2	1:L:140:ILE:HG13	2.39	0.45
4:I:24:ASN:HB3	4:I:27:ALA:HB3	1.97	0.45
3:T:93:LEU:HA	3:T:93:LEU:HD23	1.60	0.45
3:T:154:THR:O	3:T:189:VAL:HA	2.17	0.45
1:L:39:LEU:HA	1:L:39:LEU:HD13	1.78	0.45
2:H:164:MET:HE2	3:T:91:GLU:HB3	1.95	0.44
2:H:60(C):LYS:NZ	4:I:46:GLU:OE2	2.50	0.44
1:L:132:GLU:O	1:L:134:PRO:HD2	2.18	0.44
2:H:113:VAL:O	2:H:115:THR:HG23	2.17	0.44
2:H:239:GLN:O	2:H:243:ARG:HG2	2.18	0.44
1:L:55:CYS:SG	1:L:61:CYS:HB2	2.58	0.44
3:T:123:VAL:HB	3:T:177:ILE:CD1	2.48	0.44
1:L:84:HIS:C	1:L:86:ASP:H	2.20	0.44
2:H:100:ASN:HD22	2:H:100:ASN:HA	1.52	0.44
3:T:135:ARG:HA	3:T:139:THR:O	2.16	0.44
1:L:88:GLN:HE21	1:L:88:GLN:HB3	1.65	0.44
4:I:20:ARG:HB2	4:I:33:PHE:CE1	2.52	0.44
3:T:34:TYR:O	3:T:51:TYR:HA	2.17	0.44
3:T:125:VAL:O	3:T:174:GLU:HA	2.18	0.44
3:T:196:ARG:HD3	3:T:198:VAL:O	2.17	0.43
3:T:110:GLN:HA	3:T:111:PRO:HD3	1.84	0.43
3:T:134:VAL:HG12	3:T:141:LEU:HD12	1.99	0.43
2:H:164:MET:HE1	3:T:91:GLU:HB3	1.97	0.43
2:H:51:TRP:CE3	2:H:105:LEU:HG	2.53	0.43
3:T:91:GLU:HA	3:T:92:PRO:HD3	1.88	0.43
1:L:62:LYS:HB3	1:L:69:ILE:HG13	2.01	0.43
3:T:194:PRO:HA	3:T:200:ARG:NH2	2.33	0.43
3:T:114:GLN:NE2	8:T:252:HOH:O	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:T:59:LEU:O	3:T:62:GLU:HB2	2.19	0.43
3:T:10:TYR:HB2	8:T:257:HOH:O	2.17	0.43
2:H:87:GLN:CD	2:H:252:LEU:HD13	2.39	0.43
3:T:111:PRO:HB2	3:T:189:VAL:HG23	2.00	0.43
3:T:192:VAL:C	3:T:194:PRO:HD3	2.38	0.42
2:H:87:GLN:O	2:H:106:LEU:HA	2.19	0.42
2:H:95:VAL:HA	2:H:96:PRO:HD2	1.80	0.42
2:H:153:LEU:HD23	2:H:153:LEU:HA	1.71	0.42
2:H:29:TRP:HA	2:H:119:VAL:O	2.19	0.42
2:H:212:ILE:HB	2:H:229:THR:HB	2.00	0.42
1:L:105:HIS:HB2	1:L:108:THR:HG23	2.02	0.42
3:T:37:GLN:HG3	3:T:76:PHE:HE1	1.83	0.42
2:H:47:ILE:HD13	2:H:53:VAL:HG23	2.00	0.42
4:I:45:PHE:N	4:I:45:PHE:CD1	2.88	0.42
4:I:3:ASP:O	4:I:6:LEU:HG	2.20	0.42
2:H:236:GLU:OE2	2:H:240:LYS:CE	2.68	0.42
3:T:191:ALA:HB3	3:T:202:SER:HB3	2.02	0.42
2:H:203:TYR:O	2:H:204:ARG:C	2.55	0.42
2:H:144:LEU:HA	2:H:144:LEU:HD23	1.87	0.42
4:I:6:LEU:N	4:I:6:LEU:CD2	2.80	0.42
2:H:187:GLY:CA	2:H:221(A):ALA:O	2.58	0.41
4:I:7:GLU:HG3	4:I:8:PRO:O	2.20	0.41
2:H:68:LEU:HD13	2:H:118:VAL:CG1	2.51	0.41
3:T:168:ALA:HB3	3:T:175:PHE:HD2	1.85	0.41
4:I:4:PHE:CZ	4:I:54:THR:HG21	2.54	0.41
2:H:211:GLY:HA2	2:H:229:THR:O	2.20	0.41
3:T:205:SER:HB2	3:T:206:PRO:HD2	2.03	0.41
1:L:103:SER:OG	1:L:113:ARG:NE	2.53	0.41
3:T:151:LEU:HD21	3:T:153:TYR:OH	2.21	0.41
4:I:2:PRO:HD2	4:I:5:CYS:SG	2.60	0.41
3:T:187:PHE:N	3:T:208:GLU:O	2.48	0.41
1:L:121:LEU:O	1:L:124:GLY:N	2.45	0.41
1:L:73:LEU:HD23	1:L:73:LEU:HA	1.61	0.41
4:I:9:PRO:HG3	4:I:22:PHE:CE2	2.56	0.41
3:T:105:GLU:O	3:T:199:ASN:ND2	2.50	0.41
3:T:179:VAL:CB	3:T:185:TYR:CE2	3.03	0.41
2:H:89:ILE:O	2:H:104:ALA:HA	2.20	0.41
4:I:9:PRO:HG3	4:I:22:PHE:CZ	2.55	0.41
3:T:114:GLN:N	3:T:126:THR:O	2.38	0.40
3:T:147:PHE:CD1	3:T:193:ILE:HD13	2.55	0.40
3:T:37:GLN:O	3:T:73:ALA:CA	2.67	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:I:38:CYS:CB	4:I:39:LEU:HD12	2.50	0.40
4:I:21:TYR:N	4:I:21:TYR:CD1	2.89	0.40
3:T:158:TRP:CE2	3:T:186:CYS:HB2	2.57	0.40
3:T:186:CYS:HA	3:T:209:CYS:HA	2.03	0.40
2:H:68:LEU:CD1	2:H:112:VAL:HG21	2.52	0.40
4:I:23:TYR:CE1	4:I:29:LEU:N	2.89	0.40
1:L:37:THR:HG23	1:L:41:TRP:CD1	2.57	0.40
1:L:53:SER:N	1:L:54:PRO:HD3	2.37	0.40
1:L:139:PRO:HG2	2:H:116:ASP:HA	2.04	0.40
4:I:23:TYR:HH	4:I:28:GLY:HA2	1.85	0.40
3:T:123:VAL:HB	3:T:177:ILE:HD11	2.04	0.40
4:I:39:LEU:N	4:I:39:LEU:CD1	2.85	0.40
3:T:155:LEU:HA	3:T:188:SER:O	2.21	0.40
2:H:66:ALA:O	2:H:82:SER:HA	2.22	0.40
1:L:76:PHE:HA	1:L:83:THR:O	2.21	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	106/152 (70%)	100 (94%)	6 (6%)	0	100	100
2	H	247/254 (97%)	235 (95%)	11 (4%)	1 (0%)	39	37
3	T	173/206 (84%)	163 (94%)	10 (6%)	0	100	100
4	I	53/55 (96%)	50 (94%)	3 (6%)	0	100	100
All	All	579/667 (87%)	548 (95%)	30 (5%)	1 (0%)	52	53

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	H	249	GLY

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	95/122 (78%)	81 (85%)	14 (15%)	4 2
2	H	214/216 (99%)	203 (95%)	11 (5%)	29 26
3	T	172/190 (90%)	155 (90%)	17 (10%)	10 6
4	I	45/45 (100%)	35 (78%)	10 (22%)	1 0
All	All	526/573 (92%)	474 (90%)	52 (10%)	10 6

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	36	ARG
1	L	37	THR
1	L	39	LEU
1	L	43	SER
1	L	62	LYS
1	L	86	ASP
1	L	88	GLN
1	L	89	LEU
1	L	106	THR
1	L	110	ARG
1	L	113	ARG
1	L	119	SER
1	L	140	ILE
1	L	143	LYS
2	H	20	LYS
2	H	29	TRP
2	H	34	LEU
2	H	62	ARG
2	H	63	ASN
2	H	100	ASN
2	H	110	GLN

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Mol	Chain	Res	Type
2	H	134	ARG
2	H	159	ASN
2	H	170(H)	SER
2	H	202	HIS
3	T	6	THR
3	T	13	THR
3	T	27	PRO
3	T	28	LYS
3	T	68	LYS
3	T	93	LEU
3	T	114	GLN
3	T	123	VAL
3	T	130	GLU
3	T	139	THR
3	T	149	LYS
3	T	152	ILE
3	T	167	THR
3	T	171	ASN
3	T	174	GLU
3	T	175	PHE
3	T	188	SER
4	I	3	ASP
4	I	6	LEU
4	I	13	PRO
4	I	24	ASN
4	I	29	LEU
4	I	31	GLN
4	I	41	LYS
4	I	46	GLU
4	I	52	MET
4	I	53	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	66	GLN
1	L	80	ASN
1	L	88	GLN
1	L	115	HIS
2	H	60(D)	ASN
2	H	63	ASN
2	H	100	ASN

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Mol	Chain	Res	Type
2	H	110	GLN
2	H	159	ASN
3	T	114	GLN
3	T	118	GLN
3	T	137	ASN
3	T	171	ASN
4	I	24	ASN
4	I	31	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

There are no carbohydrates in this entry.

5.6 Ligand geometry

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

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	$\#\|Z\| > 2$
5	GLC	L	600	1	11,11,12	1.28	1 (9%)	14,15,17	2.67	4 (28%)
6	FUC	L	601	1	10,10,11	1.44	1 (10%)	14,14,16	1.96	3 (21%)

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
5	GLC	L	600	1	1/1/4/5	0/2/19/22	0/1/1/1
6	FUC	L	601	1	-	0/0/17/20	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	L	600	GLC	C4-C5	2.10	1.57	1.53
6	L	601	FUC	C4-C5	2.37	1.57	1.52

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
6	L	601	FUC	C2-C3-C4	-2.33	107.09	111.04
5	L	600	GLC	O3-C3-C2	2.04	113.69	110.00
6	L	601	FUC	O5-C1-C2	2.69	115.22	110.86
5	L	600	GLC	O5-C5-C6	2.70	113.19	107.35
5	L	600	GLC	O5-C1-C2	4.99	118.95	110.86
6	L	601	FUC	C1-O5-C5	5.49	120.86	112.38
5	L	600	GLC	C1-O5-C5	7.05	121.19	112.25

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
5	L	600	GLC	C1

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	L	600	GLC	1	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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.