



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 06:30 PM GMT

PDB ID : 1B5G
Title : HUMAN THROMBIN COMPLEXED WITH NOVEL SYNTHETIC PEP-
TIDE MIMETIC INHIBITOR AND HIRUGEN
Authors : St Charles, R.; Tulinsky, A.; Kahn, M.
Deposited on : 1998-03-05
Resolution : 2.07 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 ⓘ symbol.

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

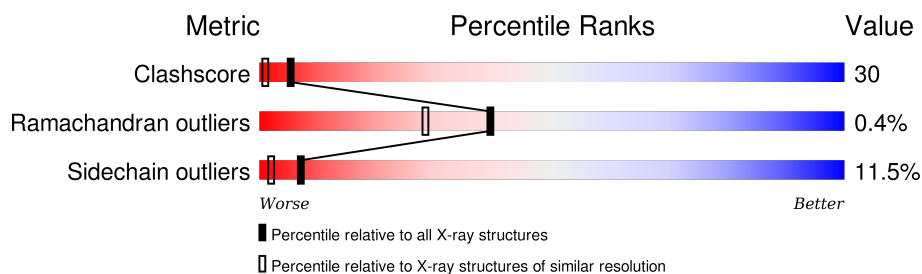
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.07 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	1910 (2.08-2.04)
Ramachandran outliers	100387	1893 (2.08-2.04)
Sidechain outliers	100360	1893 (2.08-2.04)

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 $\leq 5\%$. 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	
3	I	12	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 2526 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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	29	Total	C	N	O	S	0	0	0
			239	149	38	51	1			

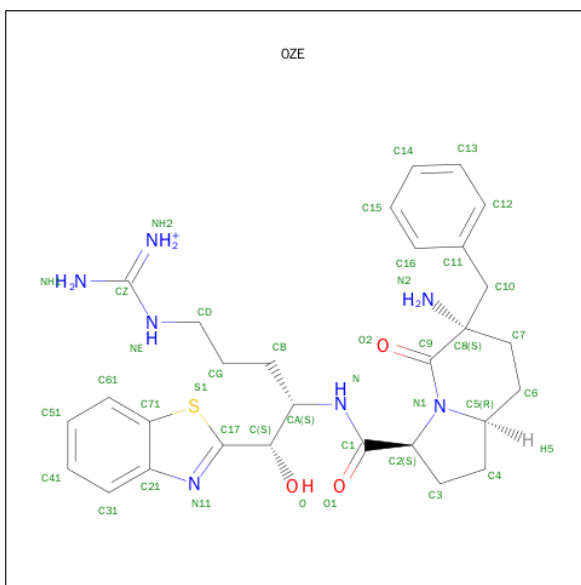
- Molecule 2 is a protein called ALPHA-THROMBIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	250	Total	C	N	O	S	0	0	0
			2022	1290	358	360	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
			90	56	10	23	1			

- Molecule 4 is [[[(4S,5S)-4-[[[(3S,6S,8AR)-6-AZANYL-5-OXO-6-(PHENYLMETHYL)-1,2,3,7,8,8A-HEXAHYDROINDOLIZIN-3-YL]CARBONYLAMINO]-5-(1,3-BENZOTHIAZOL-2-YL)-5-HYDROXY-PENTYL]AMINO]-AZANYL-METHYLIDENE]AZANUM (three-letter code: 0ZE) (formula: C₂₉H₃₈N₇O₃S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	H	1	Total	C	N	O	S	0	0
			40	29	7	3	1		

- Molecule 5 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	H	2	Total	Na	0	0
			2	2		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	L	11	Total	O	0	0
			11	11		
6	H	117	Total	O	0	0
			117	117		
6	I	5	Total	O	0	0
			5	5		

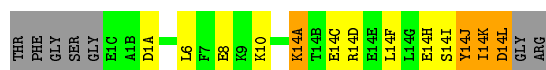
3 Residue-property plots [i](#)

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.

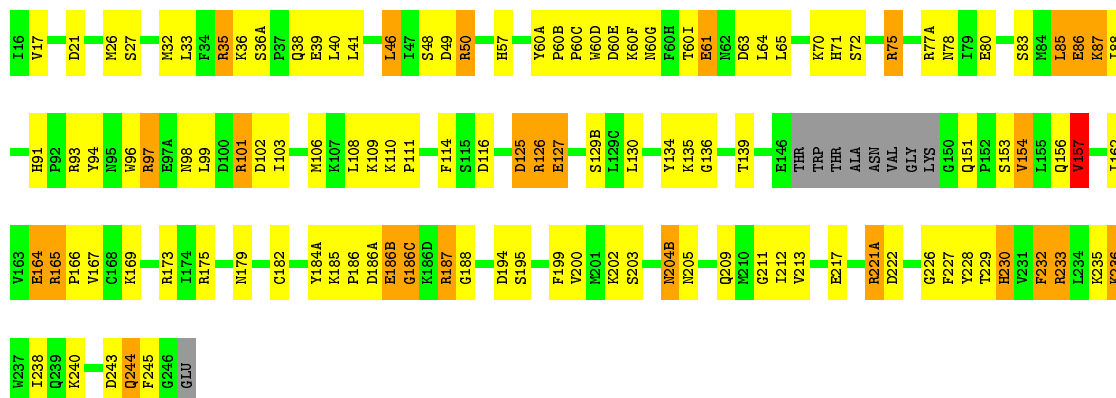
• Molecule 1: ALPHA-THROMBIN

Chain L: 



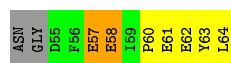
• Molecule 2: ALPHA-THROMBIN

Chain H: 



• Molecule 3: HIRUGEN

Chain I: 



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, α , β , γ	70.84Å 72.08Å 73.00Å 90.00° 100.90° 90.00°	Depositor
Resolution (Å)	7.00 – 2.07	Depositor
% Data completeness (in resolution range)	81.0 (7.00-2.07)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROFFT	Depositor
R, R_{free}	0.164 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2526	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 0ZE, NA, 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	L	1.15	1/241 (0.4%)	2.04	8/321 (2.5%)
2	H	1.04	0/2074	1.94	43/2801 (1.5%)
3	I	1.11	0/74	2.03	2/96 (2.1%)
All	All	1.05	1/2389 (0.0%)	1.95	53/3218 (1.6%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	L	14(C)	GLU	CD-OE2	-5.51	1.19	1.25

All (53) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	233	ARG	NE-CZ-NH1	22.80	131.70	120.30
2	H	233	ARG	NE-CZ-NH2	-14.29	113.15	120.30
2	H	187	ARG	NE-CZ-NH1	-12.40	114.10	120.30
2	H	221(A)	ARG	NE-CZ-NH1	-12.24	114.18	120.30
2	H	101	ARG	NE-CZ-NH2	-12.23	114.19	120.30
2	H	165	ARG	NE-CZ-NH2	11.69	126.15	120.30
2	H	221(A)	ARG	NE-CZ-NH2	11.68	126.14	120.30
2	H	97	ARG	NE-CZ-NH2	-10.19	115.21	120.30
2	H	60(A)	TYR	CB-CG-CD1	-10.10	114.94	121.00
2	H	35	ARG	NE-CZ-NH2	-9.91	115.34	120.30
2	H	175	ARG	NE-CZ-NH1	-9.79	115.41	120.30
2	H	35	ARG	NE-CZ-NH1	9.63	125.12	120.30
1	L	1(A)	ASP	CB-CG-OD1	-9.59	109.67	118.30
2	H	233	ARG	CD-NE-CZ	9.32	136.65	123.60
2	H	175	ARG	CD-NE-CZ	-9.26	110.64	123.60
2	H	173	ARG	NE-CZ-NH2	-9.03	115.78	120.30
2	H	243	ASP	CB-CG-OD2	-9.02	110.18	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	165	ARG	NE-CZ-NH1	-7.82	116.39	120.30
2	H	125	ASP	CB-CG-OD2	7.81	125.33	118.30
2	H	228	TYR	CB-CG-CD2	7.43	125.46	121.00
2	H	97	ARG	NE-CZ-NH1	7.38	123.99	120.30
2	H	243	ASP	CB-CA-C	7.22	124.85	110.40
2	H	101	ARG	NE-CZ-NH1	7.08	123.84	120.30
2	H	93	ARG	NE-CZ-NH1	-7.05	116.78	120.30
2	H	164	GLU	CG-CD-OE2	6.92	132.15	118.30
2	H	186(C)	GLY	C-N-CA	6.64	138.31	121.70
2	H	186	PRO	C-N-CA	6.63	138.26	121.70
1	L	8	GLU	OE1-CD-OE2	-6.43	115.58	123.30
2	H	230	HIS	CA-CB-CG	6.07	123.92	113.60
2	H	126	ARG	NE-CZ-NH1	6.07	123.33	120.30
2	H	173	ARG	NE-CZ-NH1	6.01	123.31	120.30
2	H	194	ASP	CB-CG-OD2	6.01	123.71	118.30
2	H	213	VAL	CA-CB-CG2	5.99	119.88	110.90
2	H	125	ASP	CB-CG-OD1	-5.96	112.93	118.30
2	H	227	PHE	CB-CG-CD1	-5.85	116.70	120.80
2	H	116	ASP	CB-CG-OD1	5.72	123.45	118.30
1	L	14(D)	ARG	NE-CZ-NH1	-5.63	117.48	120.30
2	H	232	PHE	CB-CG-CD1	-5.61	116.87	120.80
2	H	126	ARG	NE-CZ-NH2	-5.61	117.50	120.30
3	I	57	GLU	N-CA-CB	5.58	120.64	110.60
1	L	1(A)	ASP	CB-CG-OD2	5.56	123.31	118.30
1	L	8	GLU	CG-CD-OE2	5.41	129.12	118.30
2	H	179	ASN	CB-CG-OD1	-5.38	110.85	121.60
3	I	58	GLU	CG-CD-OE2	5.35	129.00	118.30
1	L	14(J)	TYR	CA-C-N	-5.29	105.57	117.20
2	H	157	VAL	CA-CB-CG1	5.23	118.74	110.90
2	H	86	GLU	CG-CD-OE1	-5.21	107.88	118.30
1	L	14(L)	ASP	N-CA-CB	5.18	119.93	110.60
2	H	175	ARG	NH1-CZ-NH2	5.11	125.02	119.40
1	L	14(A)	LYS	CA-CB-CG	5.10	124.61	113.40
2	H	80	GLU	CG-CD-OE2	-5.10	108.11	118.30
2	H	222	ASP	CB-CG-OD1	5.09	122.88	118.30
2	H	60(A)	TYR	CB-CG-CD2	5.06	124.04	121.00

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	L	239	0	235	8	1
2	H	2022	0	1996	118	2
3	I	90	0	68	12	2
4	H	40	0	37	16	0
5	H	2	0	0	0	0
6	H	117	0	0	12	0
6	I	5	0	0	0	0
6	L	11	0	0	2	0
All	All	2526	0	2336	141	3

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 30.

All (141) 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:195:SER:OG	4:H:372:0ZE:H37	0.90	1.07
2:H:151:GLN:HG2	6:H:447:HOH:O	1.55	1.06
4:H:372:0ZE:C12	4:H:372:0ZE:H72	1.85	1.03
2:H:75:ARG:HG3	2:H:75:ARG:HH11	1.17	1.02
2:H:165:ARG:HE	2:H:169:LYS:HE3	1.20	1.02
1:L:14(J):TYR:C	1:L:14(K):ILE:HG12	1.84	0.97
2:H:165:ARG:HE	2:H:169:LYS:CE	1.76	0.97
2:H:61:GLU:OE1	2:H:87:LYS:HA	1.66	0.95
2:H:165:ARG:NE	2:H:169:LYS:HZ1	1.63	0.95
2:H:26:MET:HE3	6:H:458:HOH:O	1.67	0.93
1:L:14(H):GLU:HA	1:L:14(L):ASP:HA	1.49	0.91
2:H:195:SER:CB	4:H:372:0ZE:H37	2.01	0.91
1:L:14(I):SER:C	1:L:14(K):ILE:H	1.72	0.91
2:H:195:SER:HG	4:H:372:0ZE:H37	1.18	0.90
2:H:72:SER:OG	2:H:75:ARG:HG2	1.76	0.86
2:H:75:ARG:HG3	2:H:75:ARG:NH1	1.86	0.84
2:H:165:ARG:CZ	2:H:169:LYS:HZ1	1.93	0.82
4:H:372:0ZE:H12	4:H:372:0ZE:H72	1.60	0.81
2:H:60(E):ASP:OD2	6:H:511:HOH:O	1.96	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:165:ARG:NE	2:H:169:LYS:NZ	2.28	0.80
1:L:10:LYS:HE2	6:L:515:HOH:O	1.81	0.79
2:H:165:ARG:CZ	2:H:169:LYS:NZ	2.46	0.79
2:H:85:LEU:HD13	2:H:106:MET:CE	2.14	0.78
2:H:236:LYS:HE2	6:H:540:HOH:O	1.84	0.77
2:H:35:ARG:O	2:H:38:GLN:HA	1.85	0.76
4:H:372:OZE:H4	6:H:497:HOH:O	1.86	0.76
2:H:164:GLU:OE1	2:H:167:VAL:HG21	1.86	0.76
1:L:14(J):TYR:C	1:L:14(K):ILE:CG1	2.54	0.75
2:H:85:LEU:HD13	2:H:106:MET:HE2	1.68	0.75
2:H:130:LEU:CD2	2:H:162:ILE:HD13	2.17	0.75
2:H:26:MET:CE	6:H:458:HOH:O	2.29	0.73
4:H:372:OZE:C41	6:H:497:HOH:O	2.38	0.71
2:H:46:LEU:HD22	2:H:48:SER:O	1.92	0.70
1:L:14(I):SER:C	1:L:14(K):ILE:N	2.45	0.69
2:H:165:ARG:NE	2:H:169:LYS:HE3	2.03	0.69
2:H:130:LEU:HD23	2:H:162:ILE:CD1	2.24	0.68
2:H:195:SER:CB	4:H:372:OZE:C	2.66	0.67
2:H:165:ARG:NH2	2:H:169:LYS:HZ3	1.92	0.67
3:I:58:GLU:H	3:I:58:GLU:CD	1.96	0.67
2:H:50:ARG:HD3	2:H:108:LEU:O	1.96	0.66
2:H:17:VAL:O	2:H:188:GLY:HA2	1.95	0.66
2:H:165:ARG:HE	2:H:169:LYS:NZ	1.93	0.64
2:H:70:LYS:HE3	2:H:72:SER:O	1.97	0.64
2:H:32:MET:HG3	2:H:40:LEU:HD12	1.80	0.63
2:H:186(A):ASP:O	2:H:186(C):GLY:N	2.32	0.63
2:H:60(F):LYS:HG2	6:H:497:HOH:O	1.99	0.62
2:H:135:LYS:HE2	2:H:184(A):TYR:OH	1.98	0.62
3:I:60:PRO:HD2	3:I:63:TYS:HD2	1.81	0.62
1:L:10:LYS:CE	6:L:515:HOH:O	2.44	0.62
2:H:244:GLN:HG2	2:H:245:PHE:CD2	2.35	0.61
2:H:130:LEU:HD22	2:H:162:ILE:HD13	1.84	0.60
2:H:36:LYS:HG2	3:I:64:LEU:HD23	1.83	0.60
2:H:61:GLU:OE1	2:H:87:LYS:CA	2.47	0.60
2:H:204(B):ASN:O	2:H:205:ASN:HB2	2.01	0.59
2:H:236:LYS:HB2	6:H:517:HOH:O	2.02	0.59
2:H:130:LEU:CD2	2:H:162:ILE:CD1	2.79	0.59
2:H:130:LEU:HD23	2:H:162:ILE:HD13	1.83	0.59
2:H:165:ARG:NE	2:H:169:LYS:CE	2.55	0.59
2:H:165:ARG:NH2	2:H:169:LYS:NZ	2.50	0.59
2:H:75:ARG:CG	2:H:75:ARG:NH1	2.60	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:64:LEU:HB2	2:H:85:LEU:HD11	1.84	0.58
3:I:60:PRO:HB2	3:I:62:GLU:OE2	2.03	0.58
2:H:97:ARG:NH2	6:H:498:HOH:O	2.38	0.57
2:H:71:HIS:CD2	2:H:154:VAL:HG22	2.39	0.57
2:H:164:GLU:OE1	2:H:167:VAL:CG2	2.54	0.55
3:I:61:GLU:C	3:I:63:TYS:N	2.59	0.55
2:H:165:ARG:O	2:H:169:LYS:HG3	2.06	0.55
2:H:60(D):TRP:CH2	4:H:372:OZE:H42	2.43	0.54
2:H:185:LYS:N	2:H:186(B):GLU:OE1	2.39	0.54
2:H:60(G):ASN:OD1	2:H:60(G):ASN:O	2.25	0.54
4:H:372:OZE:C7	4:H:372:OZE:C12	2.74	0.54
2:H:204(B):ASN:C	2:H:204(B):ASN:HD22	2.09	0.53
2:H:49:ASP:HB3	2:H:114:PHE:CZ	2.44	0.53
2:H:32:MET:HG3	2:H:40:LEU:CD1	2.39	0.52
2:H:203:SER:HB3	2:H:204(B):ASN:ND2	2.24	0.52
3:I:60:PRO:O	3:I:63:TYS:HB3	2.10	0.52
2:H:91:HIS:CE1	2:H:101:ARG:HD3	2.45	0.52
2:H:230:HIS:CE1	2:H:233:ARG:HG3	2.45	0.51
2:H:98:ASN:OD1	2:H:98:ASN:N	2.42	0.51
2:H:57:HIS:HD2	4:H:372:OZE:H3	1.75	0.51
2:H:85:LEU:HD13	2:H:106:MET:HE1	1.89	0.51
2:H:50:ARG:NH1	2:H:108:LEU:O	2.43	0.51
2:H:126:ARG:HA	2:H:232:PHE:CZ	2.45	0.51
3:I:57:GLU:HG2	3:I:58:GLU:O	2.11	0.50
2:H:186(A):ASP:C	2:H:186(C):GLY:H	2.15	0.50
2:H:244:GLN:OE1	2:H:245:PHE:CE2	2.66	0.49
2:H:186(A):ASP:C	2:H:186(C):GLY:N	2.66	0.49
2:H:94:TYR:CZ	2:H:96:TRP:HB3	2.47	0.49
2:H:139:THR:HG22	2:H:157:VAL:HG13	1.93	0.49
2:H:136:GLY:HA3	2:H:199:PHE:CZ	2.47	0.49
2:H:164:GLU:C	2:H:166:PRO:HD2	2.34	0.48
2:H:36:LYS:O	2:H:38:GLN:HG3	2.12	0.48
2:H:195:SER:CB	4:H:372:OZE:O	2.60	0.48
2:H:50:ARG:CD	2:H:108:LEU:O	2.61	0.48
2:H:75:ARG:HD3	2:H:75:ARG:N	2.26	0.48
2:H:60(I):THR:O	2:H:63:ASP:HB2	2.13	0.48
2:H:21:ASP:OD1	2:H:156:GLN:OE1	2.31	0.47
2:H:217:GLU:HA	4:H:372:OZE:H16	1.95	0.47
1:L:14(K):ILE:HD11	2:H:134:TYR:OH	2.14	0.47
2:H:165:ARG:N	2:H:166:PRO:HD2	2.29	0.47
2:H:97:ARG:HG3	6:H:438:HOH:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:46:LEU:CD2	2:H:48:SER:O	2.62	0.47
2:H:99:LEU:O	2:H:102:ASP:HB2	2.15	0.47
2:H:139:THR:CG2	2:H:157:VAL:HG13	2.45	0.47
2:H:35:ARG:HD3	2:H:39:GLU:OE2	2.15	0.46
3:I:60:PRO:HB2	3:I:62:GLU:CD	2.36	0.46
2:H:195:SER:HG	4:H:372:OZE:C17	2.29	0.46
3:I:61:GLU:O	3:I:63:TYS:N	2.49	0.45
2:H:135:LYS:CE	2:H:184(A):TYR:OH	2.64	0.45
2:H:85:LEU:CD1	2:H:106:MET:HE1	2.47	0.45
2:H:60(G):ASN:O	2:H:60(G):ASN:CG	2.54	0.45
2:H:77(A):ARG:O	2:H:78:ASN:HB2	2.17	0.45
2:H:204(B):ASN:HD22	2:H:205:ASN:N	2.15	0.45
2:H:41:LEU:CD2	2:H:64:LEU:CD2	2.96	0.44
2:H:165:ARG:N	2:H:166:PRO:CD	2.81	0.44
2:H:230:HIS:ND1	2:H:233:ARG:HG3	2.32	0.44
2:H:103:ILE:HD11	2:H:238:ILE:HD11	1.99	0.44
2:H:182:CYS:HA	2:H:226:GLY:O	2.17	0.44
2:H:212:ILE:HD12	2:H:229:THR:HB	2.00	0.44
2:H:211:GLY:HA2	2:H:229:THR:O	2.17	0.43
3:I:57:GLU:CG	3:I:58:GLU:N	2.82	0.43
2:H:202:LYS:HE3	2:H:205:ASN:O	2.18	0.43
3:I:57:GLU:HG2	3:I:58:GLU:N	2.34	0.42
2:H:110:LYS:HE2	6:H:529:HOH:O	2.18	0.42
2:H:195:SER:OG	4:H:372:OZE:C17	2.61	0.42
2:H:204(B):ASN:O	2:H:205:ASN:CB	2.68	0.42
2:H:85:LEU:CD1	2:H:106:MET:CE	2.94	0.42
2:H:127:GLU:CD	2:H:127:GLU:H	2.23	0.42
2:H:88:ILE:HG21	2:H:88:ILE:HD13	1.73	0.42
2:H:60(F):LYS:HE2	4:H:372:OZE:H51	2.02	0.42
2:H:134:TYR:HD1	2:H:134:TYR:N	2.18	0.42
2:H:50:ARG:HG2	2:H:111:PRO:HA	2.01	0.41
2:H:60(B):PRO:N	2:H:60(C):PRO:CD	2.82	0.41
2:H:134:TYR:CD1	2:H:134:TYR:N	2.89	0.41
3:I:62:GLU:OE1	3:I:62:GLU:N	2.33	0.41
2:H:204(B):ASN:ND2	2:H:204(B):ASN:C	2.74	0.41
2:H:200:VAL:HG12	2:H:209:GLN:HA	2.02	0.41
2:H:185:LYS:HG3	2:H:186(B):GLU:OE1	2.20	0.41
2:H:57:HIS:CD2	2:H:57:HIS:C	2.94	0.41
2:H:85:LEU:HD23	2:H:85:LEU:N	2.35	0.41
2:H:125:ASP:OD1	2:H:127:GLU:HB2	2.20	0.41

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry 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]	1.70	0.50
2:H:75:ARG:NH2	3:I:57:GLU:CB[2_555]	2.12	0.08
1:L:14(H):GLU:OE2	1:L:14(L):ASP:OD2[2_556]	2.15	0.05

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%)	22 (82%)	5 (18%)	0	100	100
2	H	246/259 (95%)	236 (96%)	9 (4%)	1 (0%)	39	28
3	I	7/12 (58%)	6 (86%)	1 (14%)	0	100	100
All	All	280/307 (91%)	264 (94%)	15 (5%)	1 (0%)	39	28

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	H	186(B)	GLU

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	27/31 (87%)	23 (85%)	4 (15%)	4	1
2	H	218/225 (97%)	193 (88%)	25 (12%)	7	2

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	I	8/10 (80%)	8 (100%)	0	100	100
All	All	253/266 (95%)	224 (88%)	29 (12%)	7	2

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	6	LEU
1	L	14(A)	LYS
1	L	14(F)	LEU
1	L	14(K)	ILE
2	H	27	SER
2	H	33	LEU
2	H	36(A)	SER
2	H	46	LEU
2	H	50	ARG
2	H	61	GLU
2	H	65	LEU
2	H	75	ARG
2	H	83	SER
2	H	85	LEU
2	H	86	GLU
2	H	87	LYS
2	H	109	LYS
2	H	127	GLU
2	H	129(B)	SER
2	H	153	SER
2	H	154	VAL
2	H	157	VAL
2	H	187	ARG
2	H	204(B)	ASN
2	H	221(A)	ARG
2	H	235	LYS
2	H	236	LYS
2	H	240	LYS
2	H	244	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
2	H	78	ASN
2	H	156	GLN

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Mol	Chain	Res	Type
2	H	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	I	63	3	15,16,17	1.50	2 (13%)	16,22,24	1.78	5 (31%)

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

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	I	63	TYS	OH-CZ	-4.52	1.35	1.42
3	I	63	TYS	CB-CA	2.10	1.58	1.53

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	I	63	TYS	CG-CB-CA	-3.54	106.20	114.21
3	I	63	TYS	O-C-CA	-2.39	119.25	125.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	I	63	TYS	CB-CG-CD1	-2.38	115.91	120.90
3	I	63	TYS	O2-S-O1	2.31	122.52	112.46
3	I	63	TYS	CZ-OH-S	3.19	123.95	118.52

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	I	63	TYS	4	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 2 are monoatomic - leaving 1 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
4	0ZE	H	372	2	35,44,44	1.79	5 (14%)	39,63,63	2.10	10 (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
4	0ZE	H	372	2	-	0/21/57/57	0/5/5/5

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	H	372	0ZE	O-C	-8.02	1.26	1.42
4	H	372	0ZE	C17-S1	2.20	1.80	1.73
4	H	372	0ZE	C41-C31	2.36	1.42	1.36
4	H	372	0ZE	C51-C61	2.38	1.42	1.36
4	H	372	0ZE	CB-CA	2.42	1.57	1.52

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	H	372	0ZE	CG-CB-CA	-4.90	103.75	114.10
4	H	372	0ZE	C8-C10-C11	-3.99	107.45	115.85
4	H	372	0ZE	CA-N-C1	-2.85	117.83	123.13
4	H	372	0ZE	C4-C5-N1	-2.73	99.05	101.93
4	H	372	0ZE	C6-C7-C8	2.16	116.06	112.32
4	H	372	0ZE	C7-C6-C5	2.43	114.59	111.02
4	H	372	0ZE	C71-C21-N11	2.62	114.23	108.16
4	H	372	0ZE	C6-C5-N1	3.32	115.08	110.21
4	H	372	0ZE	O-C-CA	3.46	117.20	107.94
4	H	372	0ZE	C17-C-CA	7.38	118.61	111.14

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	H	372	0ZE	16	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

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

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

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

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

6.4 Ligands [i](#)

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

6.5 Other polymers [i](#)

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