Full wwPDB X-ray Structure Validation Report (i)

Sep 14, 2023 – 12:48 AM EDT

PDB ID	:	8KME
Title	:	CRYSTAL STRUCTURE OF HUMAN ALPHA-THROMBIN INHIBITED WITH SEL2770.
Authors	:	Mochalkin, I.; Tulinsky, A.
Deposited on	:	1999-03-21
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
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

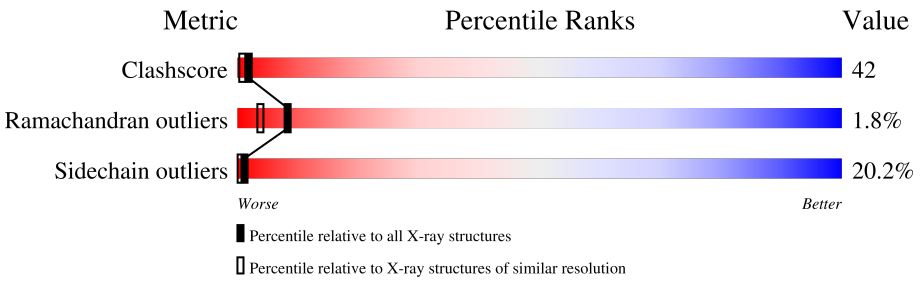
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

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	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (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 of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. 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%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	1	36	33% 28% 8% • 28%
2	2	259	33% 44% 15% • •
3	3	10	40% 50% 10%
4	4	7	14% 29% 29% 29%

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
4	0BN	4	380	-	X	-	-

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 2457 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
			Total	C	N	O	S			
1	1	26	214	134	35	44	1	0	0	0

- Molecule 2 is a protein called THROMBIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	2	250	2005	1282	351	358	14	0	0	0

- Molecule 3 is a protein called N-ACETYLHIRUDIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	3	10	79	50	10	18	1	0	0	0

- Molecule 4 is a protein called SEL2770.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
4	4	7	54	39	9	6	0	0	1

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

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

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
6	1	13	13	13	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
6	2	80	80	80	0	0
			Total	O		
6	3	8	8	8	0	0
			Total	O		
6	4	2	2	2	0	0

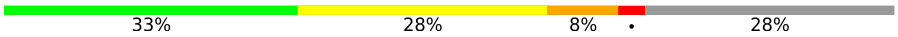
3 Residue-property plots (i)

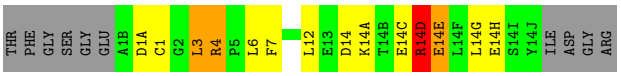
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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

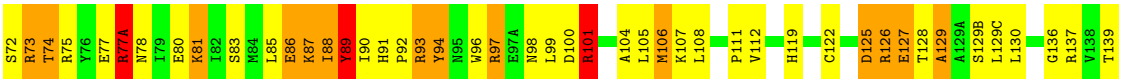
Chain 1:

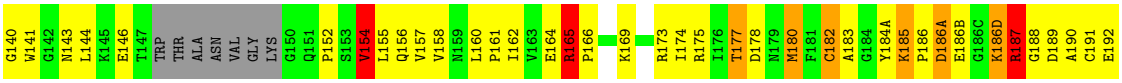


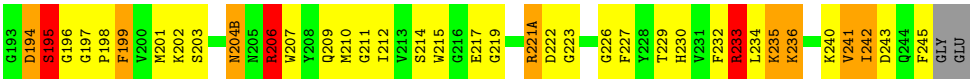


• Molecule 2: THROMBIN

Chain 2:

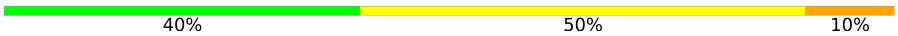


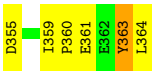




• Molecule 3: N-ACETYLHIRUDIN

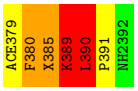
Chain 3:





• Molecule 4: SEL2770

Chain 4:



4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	71.66Å 72.00Å 73.45Å 90.00° 101.20° 90.00°	Depositor
Resolution (Å)	9.00 – 2.10	Depositor
% Data completeness (in resolution range)	60.0 (9.00-2.10)	Depositor
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROFFT	Depositor
R, R_{free}	0.167 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2457	wwPDB-VP
Average B, all atoms (Å^2)	34.0	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: NH2, NA, ACE, MLY, TYS, 0BN, CHG

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	1	1.07	0/216	1.63	2/287 (0.7%)
2	2	0.95	0/2057	1.76	36/2781 (1.3%)
3	3	1.53	0/63	1.62	0/82
4	4	1.03	0/16	2.21	1/23 (4.3%)
All	All	0.98	0/2352	1.75	39/3173 (1.2%)

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
1	1	0	1
2	2	0	12
4	4	0	2
All	All	0	15

There are no bond length outliers.

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	2	126	ARG	NE-CZ-NH1	17.25	128.93	120.30
2	2	187	ARG	NE-CZ-NH1	-17.13	111.73	120.30
2	2	221(A)	ARG	NE-CZ-NH1	14.76	127.68	120.30
2	2	187	ARG	NE-CZ-NH2	14.00	127.30	120.30
2	2	73	ARG	NE-CZ-NH1	13.68	127.14	120.30
2	2	126	ARG	NE-CZ-NH2	-11.49	114.55	120.30
2	2	233	ARG	NE-CZ-NH2	10.96	125.78	120.30
2	2	233	ARG	NE-CZ-NH1	-10.37	115.12	120.30
2	2	206	ARG	NE-CZ-NH1	10.12	125.36	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	2	126	ARG	CD-NE-CZ	8.72	135.81	123.60
2	2	194	ASP	CB-CG-OD2	7.85	125.37	118.30
2	2	187	ARG	CD-NE-CZ	7.44	134.02	123.60
2	2	101	ARG	NE-CZ-NH2	-7.39	116.60	120.30
2	2	125	ASP	CB-CG-OD1	7.37	124.94	118.30
2	2	175	ARG	NE-CZ-NH2	7.31	123.95	120.30
1	1	4	ARG	NE-CZ-NH2	6.91	123.75	120.30
2	2	221(A)	ARG	CD-NE-CZ	6.67	132.93	123.60
2	2	187	ARG	CG-CD-NE	-6.66	97.82	111.80
2	2	180	MET	O-C-N	6.40	132.94	122.70
2	2	77(A)	ARG	NE-CZ-NH2	6.39	123.50	120.30
2	2	73	ARG	CD-NE-CZ	6.37	132.52	123.60
2	2	194	ASP	CB-CG-OD1	-6.26	112.66	118.30
2	2	73	ARG	NE-CZ-NH2	-6.17	117.21	120.30
4	4	389	MLY	C-N-CA	6.00	136.70	121.70
2	2	77	GLU	O-C-N	5.94	132.21	122.70
2	2	42	CYS	CB-CA-C	-5.85	98.70	110.40
1	1	14(D)	ARG	NE-CZ-NH2	5.67	123.14	120.30
2	2	178	ASP	CB-CG-OD2	-5.57	113.29	118.30
2	2	154	VAL	N-CA-CB	-5.50	99.39	111.50
2	2	173	ARG	NE-CZ-NH2	5.41	123.01	120.30
2	2	60(H)	PHE	CB-CG-CD1	-5.31	117.08	120.80
2	2	21	ASP	CB-CG-OD2	-5.26	113.56	118.30
2	2	97	ARG	NE-CZ-NH2	5.26	122.93	120.30
2	2	207	TRP	O-C-N	5.18	130.99	122.70
2	2	146	GLU	C-N-CA	5.16	134.60	121.70
2	2	89	TYR	N-CA-CB	5.14	119.85	110.60
2	2	137	ARG	NE-CZ-NH1	5.07	122.83	120.30
2	2	210	MET	CA-CB-CG	-5.04	104.73	113.30
2	2	49	ASP	CB-CG-OD2	5.01	122.81	118.30

There are no chirality outliers.

All (15) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	1	3	LEU	Mainchain
2	2	100	ASP	Mainchain
2	2	106	MET	Mainchain
2	2	129	ALA	Mainchain
2	2	158	VAL	Mainchain
2	2	160	LEU	Mainchain
2	2	165	ARG	Sidechain

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Mol	Chain	Res	Type	Group
2	2	187	ARG	Sidechain
2	2	199	PHE	Mainchain
2	2	206	ARG	Sidechain
2	2	66	VAL	Mainchain
2	2	88	ILE	Mainchain
2	2	93	ARG	Sidechain
4	4	389	MLY	Peptide
4	4	390	LEU	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	1	214	0	214	17	0
2	2	2005	0	1970	178	0
3	3	79	0	59	7	0
4	4	54	0	58	12	0
5	2	2	0	0	0	0
6	1	13	0	0	1	0
6	2	80	0	0	5	0
6	3	8	0	0	1	0
6	4	2	0	0	0	0
All	All	2457	0	2301	196	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 42.

All (196) 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)
2:2:236:LYS:HE2	2:2:236:LYS:H	1.13	1.14
2:2:21:ASP:HB3	2:2:154:VAL:HG12	1.50	0.92
2:2:174:ILE:HD11	4:4:391:PRO:HD3	1.49	0.92
2:2:91:HIS:CG	2:2:92:PRO:HD2	2.10	0.87
2:2:140:GLY:HA3	2:2:194:ASP:OD1	1.75	0.86
2:2:59:LEU:HD13	2:2:88:ILE:CG2	2.05	0.86

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:2:236:LYS:HE2	2:2:236:LYS:N	1.89	0.86
2:2:91:HIS:ND1	2:2:92:PRO:HD2	1.91	0.86
2:2:139:THR:HG22	2:2:157:VAL:HG22	1.58	0.83
2:2:165:ARG:HD3	2:2:169:LYS:HZ1	1.43	0.82
2:2:36:LYS:HD2	3:3:364:LEU:HD23	1.60	0.82
2:2:232:PHE:O	2:2:235:LYS:HB2	1.81	0.81
2:2:52:VAL:HG23	2:2:108:LEU:HD21	1.61	0.79
2:2:86:GLU:O	2:2:87:LYS:HG2	1.82	0.79
2:2:236:LYS:H	2:2:236:LYS:CE	1.94	0.78
1:1:3:LEU:HD11	2:2:206:ARG:HH11	1.47	0.78
2:2:235:LYS:HD3	2:2:235:LYS:C	2.03	0.77
2:2:60(I):THR:HG23	2:2:61:GLU:HG2	1.68	0.76
2:2:59:LEU:HD22	2:2:88:ILE:HD13	1.67	0.75
2:2:59:LEU:HD13	2:2:88:ILE:HG21	1.66	0.75
2:2:24:ILE:HD12	2:2:24:ILE:H	1.50	0.75
2:2:17:VAL:O	2:2:188:GLY:HA2	1.87	0.74
2:2:32:MET:HE3	2:2:70:LYS:HD3	1.68	0.73
2:2:21:ASP:HB3	2:2:154:VAL:CG1	2.19	0.73
2:2:143:ASN:HD21	2:2:192:GLU:HB3	1.54	0.73
2:2:60(B):PRO:HG2	2:2:96:TRP:CE2	2.22	0.73
2:2:80:GLU:C	2:2:81:LYS:HE2	2.08	0.72
1:1:3:LEU:HD11	2:2:206:ARG:NH1	2.03	0.72
1:1:4:ARG:HG2	2:2:28:PRO:CG	2.20	0.72
2:2:53:LEU:HD23	2:2:209:GLN:OE1	1.91	0.71
2:2:143:ASN:ND2	2:2:192:GLU:HB3	2.06	0.71
2:2:24:ILE:HD12	2:2:24:ILE:N	2.06	0.70
2:2:36:LYS:CD	3:3:364:LEU:HD23	2.21	0.70
1:1:3:LEU:CD1	2:2:206:ARG:HH11	2.04	0.70
2:2:52:VAL:CG2	2:2:108:LEU:HD21	2.21	0.70
2:2:165:ARG:HD3	2:2:169:LYS:NZ	2.07	0.70
2:2:186:PRO:HD3	6:2:440:HOH:O	1.91	0.70
2:2:94:TYR:HE1	2:2:99:LEU:CD2	2.04	0.69
2:2:49:ASP:O	2:2:111:PRO:HA	1.92	0.69
2:2:235:LYS:HD3	2:2:235:LYS:O	1.92	0.69
2:2:186:PRO:HG3	2:2:223:GLY:N	2.06	0.69
1:1:14(G):LEU:HD21	2:2:202:LYS:HG2	1.75	0.68
2:2:60(H):PHE:HB3	2:2:64:LEU:HD21	1.74	0.68
2:2:98:ASN:HD21	2:2:177:THR:CG2	2.07	0.68
2:2:185:LYS:HG2	2:2:186(B):GLU:OE1	1.95	0.67
1:1:4:ARG:HG2	2:2:28:PRO:HG2	1.77	0.66
3:3:360:PRO:HA	6:3:467:HOH:O	1.95	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:1:1:CYS:O	2:2:206:ARG:HD3	1.95	0.66
4:4:389:MLY:N	4:4:389:MLY:HD3	2.12	0.65
2:2:217:GLU:HA	4:4:389:MLY:O	1.96	0.64
2:2:203:SER:HB3	2:2:204(B):ASN:ND2	2.13	0.64
6:2:709:HOH:O	4:4:385:CHG:H4	1.97	0.64
2:2:81:LYS:HE2	2:2:81:LYS:N	2.13	0.63
2:2:186(B):GLU:O	2:2:186(D):LYS:HE2	2.00	0.62
4:4:385:CHG:H6	4:4:385:CHG:O	1.99	0.61
2:2:94:TYR:CE1	2:2:99:LEU:CD2	2.83	0.61
2:2:51:TRP:CZ2	2:2:107:LYS:HD2	2.37	0.60
4:4:379:ACE:O	4:4:389:MLY:HH13	2.01	0.60
2:2:122:CYS:SG	2:2:206:ARG:HD2	2.41	0.60
2:2:129(C):LEU:HB3	2:2:201:MET:SD	2.42	0.60
2:2:108:LEU:CD1	2:2:112:VAL:CG1	2.79	0.60
1:1:7:PHE:HB3	1:1:12:LEU:O	2.00	0.60
2:2:204(B):ASN:C	2:2:204(B):ASN:HD22	2.05	0.59
2:2:36:LYS:HE3	6:2:715:HOH:O	2.03	0.59
2:2:94:TYR:HE1	2:2:99:LEU:HD22	1.66	0.59
2:2:165:ARG:HD3	2:2:169:LYS:CE	2.33	0.59
2:2:101:ARG:HG2	2:2:234:LEU:HD21	1.84	0.59
1:1:14:ASP:OD1	1:1:14(C):GLU:HB3	2.02	0.59
2:2:85:LEU:HD22	2:2:106:MET:HB3	1.84	0.58
2:2:59:LEU:CD2	2:2:88:ILE:HD13	2.31	0.58
2:2:22:ALA:CB	2:2:155:LEU:HD23	2.33	0.58
2:2:60(B):PRO:HG2	2:2:96:TRP:CD2	2.39	0.58
2:2:174:ILE:HD11	4:4:391:PRO:CD	2.29	0.58
1:1:3:LEU:CD1	2:2:206:ARG:NH1	2.64	0.58
2:2:203:SER:HB3	2:2:204(B):ASN:HD21	1.68	0.58
1:1:14(E):GLU:HG3	6:1:702:HOH:O	2.04	0.57
2:2:30:GLN:HE22	2:2:198:PRO:CD	2.17	0.57
4:4:389:MLY:HH22	4:4:389:MLY:HA	1.86	0.57
2:2:28:PRO:HB2	2:2:119:HIS:HB3	1.87	0.56
2:2:36:LYS:HD2	3:3:364:LEU:CD2	2.34	0.56
2:2:89:TYR:O	2:2:104:ALA:HA	2.05	0.56
2:2:57:HIS:CD2	2:2:57:HIS:O	2.59	0.56
2:2:98:ASN:HD22	2:2:180:MET:HE1	1.71	0.56
2:2:204(B):ASN:HD22	2:2:206:ARG:H	1.52	0.56
4:4:380:0BN:C8	4:4:380:0BN:C	2.83	0.56
1:1:14:ASP:HB2	2:2:23:GLU:OE2	2.06	0.56
2:2:31:VAL:HB	2:2:44:ALA:HB3	1.86	0.56
2:2:91:HIS:CE1	2:2:92:PRO:HD2	2.40	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:2:32:MET:HG2	2:2:67:ARG:HB2	1.87	0.55
2:2:60(H):PHE:CB	2:2:64:LEU:HD21	2.36	0.55
2:2:126:ARG:HG2	2:2:126:ARG:O	2.04	0.55
2:2:108:LEU:HD12	2:2:112:VAL:CG1	2.36	0.55
2:2:50:ARG:HH21	2:2:107:LYS:HE3	1.72	0.55
2:2:60(I):THR:CG2	2:2:61:GLU:N	2.70	0.55
2:2:86:GLU:O	2:2:87:LYS:CG	2.53	0.55
2:2:81:LYS:HE2	2:2:81:LYS:CA	2.37	0.54
2:2:165:ARG:HG2	2:2:166:PRO:HD3	1.89	0.54
2:2:217:GLU:O	2:2:221(A):ARG:HD2	2.08	0.54
2:2:32:MET:HE1	2:2:70:LYS:HE2	1.90	0.53
2:2:60(B):PRO:N	2:2:60(C):PRO:HD2	2.23	0.53
2:2:60(I):THR:HG23	2:2:61:GLU:H	1.73	0.53
2:2:215:TRP:HA	4:4:380:0BN:C4	2.38	0.53
2:2:32:MET:HG2	2:2:67:ARG:HD3	1.90	0.53
2:2:98:ASN:O	2:2:99:LEU:HB2	2.08	0.53
2:2:235:LYS:C	2:2:235:LYS:CD	2.77	0.53
2:2:64:LEU:CD1	2:2:88:ILE:HD11	2.39	0.53
2:2:22:ALA:HB2	2:2:157:VAL:HG23	1.90	0.52
2:2:188:GLY:O	2:2:189:ASP:HB2	2.10	0.52
2:2:98:ASN:ND2	2:2:177:THR:CG2	2.72	0.52
2:2:94:TYR:CE1	2:2:99:LEU:HD23	2.45	0.52
2:2:36:LYS:O	2:2:38:GLN:HG2	2.09	0.52
2:2:164:GLU:H	2:2:164:GLU:CD	2.12	0.52
2:2:32:MET:CE	2:2:70:LYS:HE2	2.40	0.51
2:2:91:HIS:ND1	2:2:92:PRO:CD	2.70	0.51
2:2:55:ALA:H	2:2:196:GLY:HA2	1.76	0.51
2:2:41:LEU:HD21	2:2:60(H):PHE:CD2	2.45	0.51
2:2:86:GLU:O	2:2:87:LYS:CB	2.59	0.51
2:2:105:LEU:HD13	2:2:242:ILE:CD1	2.40	0.51
2:2:177:THR:HG23	2:2:180:MET:HE3	1.93	0.51
2:2:33:LEU:O	2:2:40:LEU:HD12	2.11	0.50
2:2:87:LYS:HB3	2:2:89:TYR:CE1	2.46	0.50
2:2:91:HIS:CG	2:2:92:PRO:CD	2.89	0.50
2:2:98:ASN:ND2	2:2:177:THR:HG23	2.25	0.50
2:2:32:MET:HE3	2:2:70:LYS:CD	2.39	0.50
2:2:204(B):ASN:ND2	2:2:206:ARG:H	2.10	0.50
2:2:41:LEU:HD22	2:2:64:LEU:HD22	1.93	0.50
2:2:94:TYR:CE2	2:2:96:TRP:HB3	2.46	0.50
2:2:60(I):THR:CG2	2:2:61:GLU:H	2.25	0.49
2:2:59:LEU:HD22	2:2:88:ILE:CD1	2.41	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:2:80:GLU:O	2:2:81:LYS:HE2	2.11	0.49
2:2:71:HIS:CE1	2:2:154:VAL:HG13	2.48	0.48
2:2:31:VAL:HG12	2:2:32:MET:N	2.28	0.48
2:2:81:LYS:CA	2:2:81:LYS:CE	2.87	0.48
2:2:144:LEU:HD21	2:2:156:GLN:NE2	2.28	0.48
2:2:74:THR:HG22	2:2:75:ARG:N	2.28	0.48
2:2:161:PRO:O	2:2:183:ALA:HB1	2.15	0.47
2:2:127:GLU:H	2:2:127:GLU:CD	2.17	0.47
2:2:189:ASP:HB2	6:2:444:HOH:O	2.14	0.47
2:2:16:ILE:HD13	2:2:190:ALA:HA	1.97	0.47
2:2:64:LEU:HD12	2:2:88:ILE:HD11	1.96	0.46
2:2:165:ARG:CG	2:2:166:PRO:HD3	2.45	0.46
2:2:214:SER:HB3	2:2:227:PHE:O	2.15	0.46
2:2:98:ASN:HD21	2:2:177:THR:HG21	1.81	0.46
2:2:98:ASN:HD22	2:2:180:MET:CE	2.29	0.46
2:2:60(B):PRO:N	2:2:60(C):PRO:CD	2.78	0.45
1:1:14(D):ARG:HD3	1:1:14(H):GLU:OE2	2.15	0.45
2:2:184(A):TYR:O	6:2:570:HOH:O	2.21	0.45
2:2:186:PRO:HB3	2:2:222:ASP:HB3	1.98	0.45
2:2:219:GLY:O	4:4:380:0BN:H8	2.15	0.45
2:2:24:ILE:H	2:2:24:ILE:CD1	2.25	0.45
2:2:73:ARG:CG	2:2:141:TRP:HB3	2.46	0.45
2:2:161:PRO:HG3	2:2:184(A):TYR:CE1	2.51	0.45
1:1:7:PHE:CB	1:1:12:LEU:O	2.63	0.45
2:2:182:CYS:HA	2:2:226:GLY:O	2.17	0.45
2:2:215:TRP:HB2	4:4:389:MLY:HD2	1.98	0.45
2:2:211:GLY:HA2	2:2:229:THR:O	2.16	0.44
2:2:31:VAL:CG1	2:2:32:MET:N	2.80	0.44
2:2:93:ARG:HH11	2:2:93:ARG:HG2	1.83	0.44
2:2:240:LYS:O	2:2:240:LYS:HG3	2.17	0.44
2:2:61:GLU:OE1	2:2:87:LYS:HA	2.18	0.43
2:2:165:ARG:N	2:2:166:PRO:HD2	2.33	0.43
2:2:68:ILE:CD1	2:2:112:VAL:HG11	2.48	0.43
2:2:136:GLY:HA3	2:2:199:PHE:CZ	2.53	0.43
2:2:165:ARG:HD3	2:2:169:LYS:HE2	2.01	0.43
2:2:204(B):ASN:ND2	2:2:206:ARG:HB2	2.33	0.43
2:2:105:LEU:HD12	2:2:241:VAL:HG11	2.01	0.43
2:2:185:LYS:O	2:2:186(A):ASP:N	2.51	0.43
3:3:361:GLU:C	3:3:363:TYS:N	2.70	0.43
2:2:32:MET:HE3	2:2:70:LYS:CE	2.49	0.42
2:2:70:LYS:HB3	2:2:70:LYS:HE3	1.45	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:2:53:LEU:HD11	2:2:212:ILE:HD11	2.00	0.42
2:2:30:GLN:HE22	2:2:198:PRO:HD2	1.84	0.42
2:2:60(A):TYR:CE1	2:2:60(C):PRO:HG2	2.54	0.42
1:1:1(A):ASP:OD1	1:1:1(A):ASP:O	2.37	0.42
2:2:61:GLU:HG2	2:2:61:GLU:H	1.45	0.42
1:1:7:PHE:CA	1:1:12:LEU:O	2.68	0.42
2:2:243:ASP:C	2:2:245:PHE:N	2.73	0.42
2:2:52:VAL:HG23	2:2:108:LEU:CD2	2.41	0.42
2:2:60(B):PRO:HG2	2:2:96:TRP:CZ2	2.55	0.41
2:2:125:ASP:OD1	2:2:128:THR:HB	2.20	0.41
2:2:164:GLU:C	2:2:166:PRO:HD2	2.40	0.41
2:2:230:HIS:CG	2:2:233:ARG:HG3	2.55	0.41
3:3:363:TYS:HD1	3:3:363:TYS:HA	1.69	0.41
1:1:14(A):LYS:H	1:1:14(A):LYS:HG2	1.69	0.41
2:2:46:LEU:HD23	2:2:46:LEU:HA	1.92	0.41
2:2:59:LEU:HD13	2:2:88:ILE:HG23	1.99	0.41
2:2:94:TYR:CZ	2:2:96:TRP:HB3	2.56	0.41
2:2:105:LEU:HD12	2:2:241:VAL:CG1	2.51	0.41
2:2:129:ALA:O	2:2:130:LEU:HB2	2.21	0.41
2:2:195:SER:C	2:2:197:GLY:H	2.25	0.41
2:2:203:SER:CB	2:2:204(B):ASN:ND2	2.83	0.40
3:3:359:ILE:HD12	3:3:363:TYS:HB3	2.03	0.40
2:2:165:ARG:N	2:2:166:PRO:CD	2.85	0.40
2:2:32:MET:CE	2:2:70:LYS:CE	2.99	0.40
2:2:81:LYS:N	2:2:81:LYS:CE	2.82	0.40
2:2:191:CYS:O	2:2:194:ASP:HB2	2.20	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	1	24/36 (67%)	21 (88%)	3 (12%)	0	100	100
2	2	246/259 (95%)	219 (89%)	23 (9%)	4 (2%)	9	5
3	3	7/10 (70%)	7 (100%)	0	0	100	100
4	4	2/7 (29%)	1 (50%)	0	1 (50%)	0	0
All	All	279/312 (89%)	248 (89%)	26 (9%)	5 (2%)	8	4

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	2	87	LYS
2	2	77(A)	ARG
4	4	390	LEU
2	2	101	ARG
2	2	195	SER

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	1	24/31 (77%)	21 (88%)	3 (12%)	4	2
2	2	215/225 (96%)	170 (79%)	45 (21%)	1	0
3	3	6/9 (67%)	5 (83%)	1 (17%)	2	1
4	4	2/2 (100%)	1 (50%)	1 (50%)	0	0
All	All	247/267 (92%)	197 (80%)	50 (20%)	1	0

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	1	6	LEU
1	1	14(D)	ARG
1	1	14(E)	GLU
2	2	21	ASP
2	2	32	MET
2	2	33	LEU

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Mol	Chain	Res	Type
2	2	36(A)	SER
2	2	41	LEU
2	2	46	LEU
2	2	47	ILE
2	2	60(E)	ASP
2	2	61	GLU
2	2	62	ASN
2	2	64	LEU
2	2	65	LEU
2	2	66	VAL
2	2	68	ILE
2	2	70	LYS
2	2	72	SER
2	2	74	THR
2	2	77(A)	ARG
2	2	78	ASN
2	2	81	LYS
2	2	83	SER
2	2	86	GLU
2	2	89	TYR
2	2	90	ILE
2	2	94	TYR
2	2	97	ARG
2	2	127	GLU
2	2	129(B)	SER
2	2	152	PRO
2	2	154	VAL
2	2	162	ILE
2	2	165	ARG
2	2	177	THR
2	2	182	CYS
2	2	185	LYS
2	2	186(A)	ASP
2	2	186(D)	LYS
2	2	187	ARG
2	2	195	SER
2	2	204(B)	ASN
2	2	233	ARG
2	2	235	LYS
2	2	236	LYS
2	2	241	VAL
2	2	242	ILE

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Mol	Chain	Res	Type
3	3	355	ASP
4	4	390	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
2	2	30	GLN
2	2	71	HIS
2	2	78	ASN
2	2	98	ASN
2	2	143	ASN
2	2	156	GLN
2	2	204(B)	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)

4 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	$\#\|Z\| > 2$
4	0BN	4	380	4	13,14,15	3.07	10 (76%)	11,18,20	1.92	5 (45%)
4	CHG	4	385	4	9,10,11	3.32	4 (44%)	7,12,14	1.32	0
3	TYS	3	363	3	15,16,17	1.94	3 (20%)	18,22,24	1.92	4 (22%)
4	MLY	4	389	4	9,10,11	3.91	4 (44%)	6,11,13	1.50	1 (16%)

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	0BN	4	380	4	-	4/9/10/12	0/1/1/1
4	CHG	4	385	4	-	0/5/14/16	0/1/1/1
3	TYS	3	363	3	-	2/10/11/13	0/1/1/1
4	MLY	4	389	4	-	6/8/9/11	-

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	4	389	MLY	CB-CA	9.98	1.66	1.53
4	4	385	CHG	C1-CA	-7.55	1.41	1.53
3	3	363	TYS	OH-S	5.65	1.66	1.58
4	4	380	0BN	C9-C3	-5.56	1.37	1.51
4	4	380	0BN	C8-C3	4.43	1.48	1.38
4	4	385	CHG	C5-C6	-3.82	1.43	1.53
4	4	380	0BN	C6-C10	-3.79	1.40	1.47
4	4	389	MLY	CH2-NZ	-3.65	1.34	1.46
4	4	385	CHG	C6-C1	-3.59	1.44	1.53
4	4	389	MLY	CH1-NZ	-3.50	1.35	1.46
3	3	363	TYS	OH-CZ	-3.44	1.37	1.42
4	4	380	0BN	CA-N	-3.23	1.38	1.48
4	4	380	0BN	C10-N2	-2.94	1.26	1.33
4	4	380	0BN	C9-CA	2.91	1.60	1.53
4	4	385	CHG	C3-C2	-2.70	1.46	1.53
4	4	380	0BN	C5-C4	-2.61	1.34	1.38
4	4	380	0BN	C8-C7	-2.60	1.34	1.38
4	4	380	0BN	O-C	2.60	1.30	1.19
4	4	380	0BN	C5-C6	2.47	1.43	1.39
4	4	389	MLY	CG-CB	-2.17	1.42	1.52
3	3	363	TYS	O-C	2.03	1.28	1.19

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	3	363	TYS	CG-CB-CA	-5.59	102.78	114.10
3	3	363	TYS	OH-S-O2	-2.74	99.72	107.71
4	4	380	0BN	C9-C3-C8	-2.70	115.55	120.91
3	3	363	TYS	OH-S-O1	2.69	115.54	107.71
4	4	380	0BN	C7-C8-C3	-2.66	117.37	121.03
4	4	380	0BN	C8-C3-C4	2.63	122.31	118.17
4	4	389	MLY	CH2-NZ-CH1	2.53	116.28	109.73
4	4	380	0BN	C5-C4-C3	-2.51	117.58	121.03

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	3	363	TYS	CB-CG-CD1	-2.43	116.09	120.91
4	4	380	0BN	C7-C6-C5	2.37	121.97	118.59

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	4	389	MLY	CD-CE-NZ-CH1
4	4	389	MLY	CD-CE-NZ-CH2
4	4	389	MLY	CG-CD-CE-NZ
4	4	389	MLY	CE-CD-CG-CB
3	3	363	TYS	CA-CB-CG-CD1
4	4	380	0BN	N2-C10-C6-C5
4	4	380	0BN	N2-C10-C6-C7
3	3	363	TYS	C-CA-CB-CG
4	4	389	MLY	CA-CB-CG-CD
4	4	389	MLY	C-CA-CB-CG
4	4	380	0BN	N3-C10-C6-C5
4	4	380	0BN	N3-C10-C6-C7

There are no ring outliers.

4 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	4	380	0BN	3	0
4	4	385	CHG	2	0
3	3	363	TYS	3	0
4	4	389	MLY	5	0

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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 is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.