Full wwPDB X-ray Structure Validation Report (i)

Aug 7, 2023 – 02:47 AM EDT

PDB ID	:	1JOU
Title	:	Crystal Structure of Native S195A Thrombin with an Unoccupied Active Site
Authors	:	Huntington, J.A.; Esmon, C.T.
Deposited on	:	2001-07-31
Resolution	:	1.80 Å(reported)

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

We welcome your comments at *validation@mail.wwpdb.org*
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 (1)) were used in the production of this report:

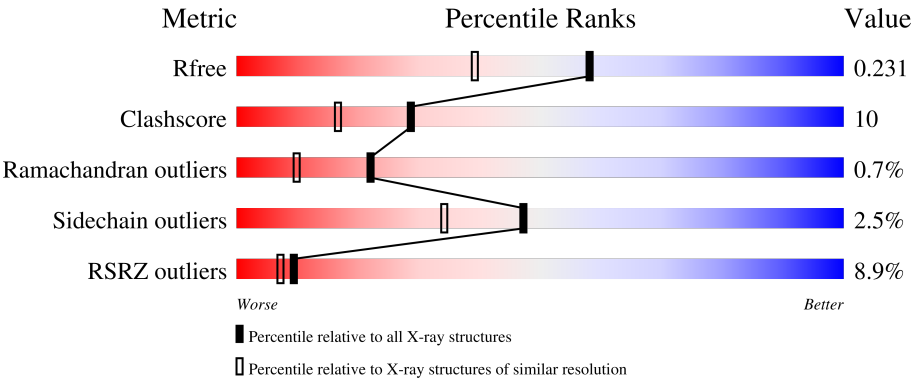
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

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 1.80 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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% 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.

Mol	Chain	Length	Quality of chain
1	A	49	8% 80% 10% • 8%
1	C	49	12% 63% 33% •
1	E	49	14% 65% 27% • 6%
2	B	259	3% 86% 13% •
2	D	259	13% 79% 19% ••

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Mol	Chain	Length	Quality of chain
2	F	259	8% 80% 18%

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
3	NAG	B	1	X	-	-	X
3	NAG	F	2	X	-	-	-

2 Entry composition (i)

There are 7 unique types of molecules in this entry. The entry contains 7838 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, light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	45	Total	C	N	O	S	0	0	0
			366	230	59	76	1			
1	C	47	Total	C	N	O	S	6	0	0
			385	240	64	80	1			
1	E	46	Total	C	N	O	S	0	0	0
			378	236	63	78	1			

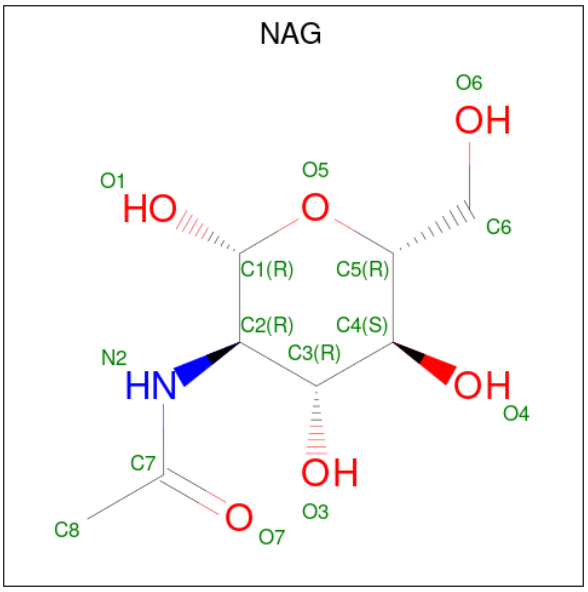
- Molecule 2 is a protein called Thrombin, heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	259	Total	C	N	O	S	8	0	0
			2092	1334	370	374	14			
2	D	257	Total	C	N	O	S	15	0	0
			2078	1327	368	369	14			
2	F	259	Total	C	N	O	S	13	0	0
			2092	1334	370	374	14			

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	195	ALA	SER	engineered mutation	UNP P00734
D	195	ALA	SER	engineered mutation	UNP P00734
F	195	ALA	SER	engineered mutation	UNP P00734

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).

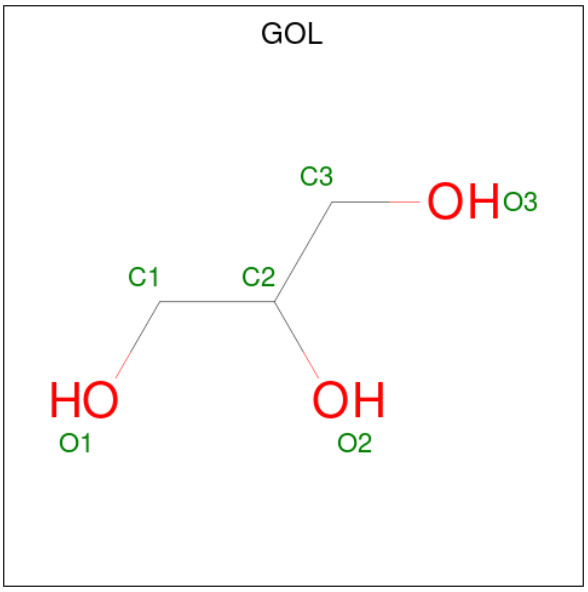


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total C N O 14 8 1 5	0	0
3	F	1	Total C N O 14 8 1 5	0	0

- Molecule 4 is SODIUM ION (three-letter code: NA) (formula: Na).

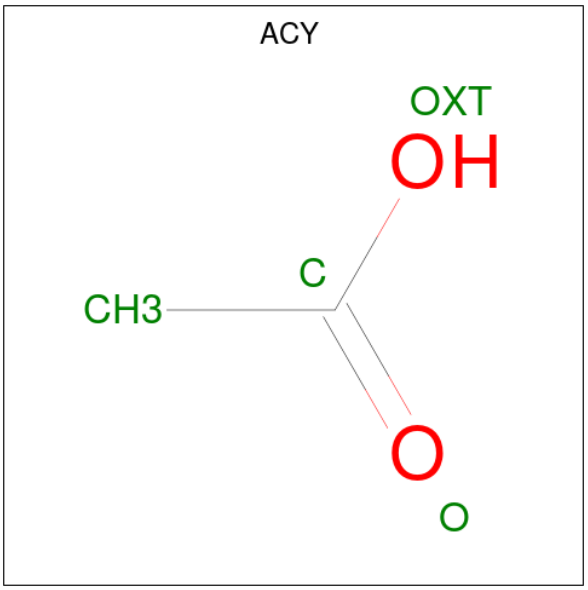
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total Na 1 1	0	0
4	D	1	Total Na 1 1	0	0
4	F	1	Total Na 1 1	0	0

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	1	Total C O 6 3 3	0	0
5	B	1	Total C O 6 3 3	0	0
5	B	1	Total C O 6 3 3	0	0
5	B	1	Total C O 6 3 3	0	0

- Molecule 6 is ACETIC ACID (three-letter code: ACY) (formula: $C_2H_4O_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	B	1	Total C O 4 2 2	0	0
6	B	1	Total C O 4 2 2	0	0
6	D	1	Total C O 4 2 2	0	0
6	F	1	Total C O 4 2 2	0	0
6	F	1	Total C O 4 2 2	0	0

- Molecule 7 is water.

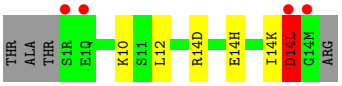
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	28	Total O 28 28	0	0
7	B	142	Total O 142 142	0	0
7	C	13	Total O 13 13	0	0
7	D	72	Total O 72 72	0	0
7	E	17	Total O 17 17	0	0
7	F	100	Total O 100 100	0	0

3 Residue-property plots

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 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.

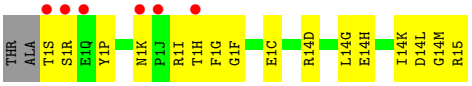
• Molecule 1: Thrombin, light chain

Chain A:



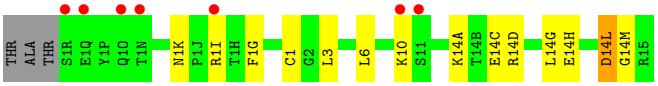
• Molecule 1: Thrombin, light chain

Chain C:



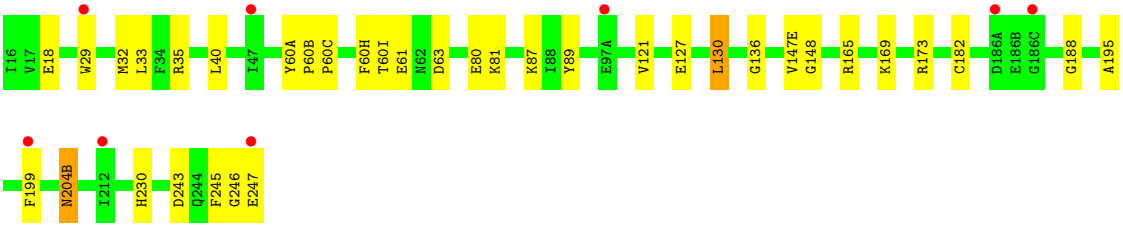
• Molecule 1: Thrombin, light chain

Chain E:



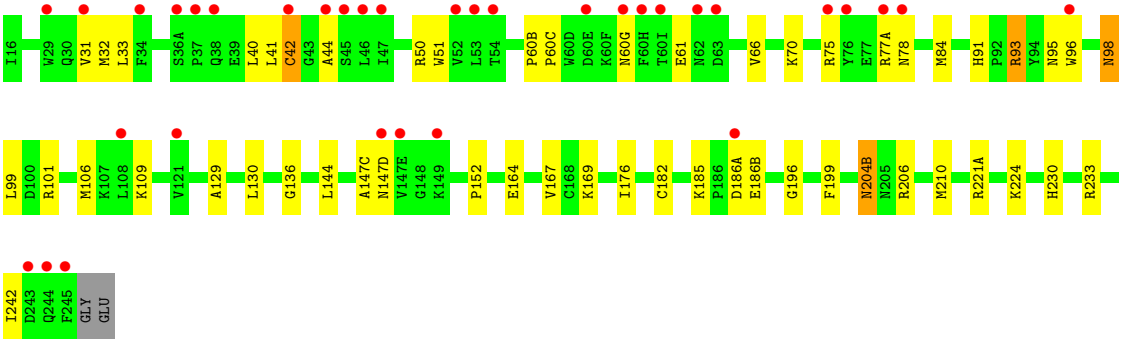
• Molecule 2: Thrombin, heavy chain

Chain B:

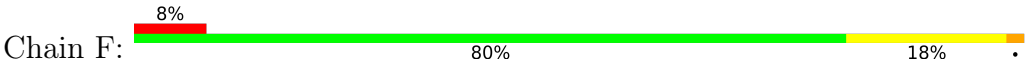


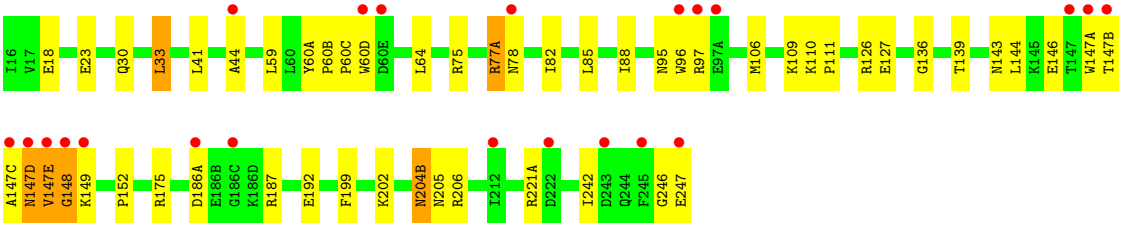
• Molecule 2: Thrombin, heavy chain

Chain D:



• Molecule 2: Thrombin, heavy chain





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	71.23Å 76.58Å 97.32Å 90.00° 106.64° 90.00°	Depositor
Resolution (Å)	24.69 – 1.80 24.69 – 1.80	Depositor EDS
% Data completeness (in resolution range)	(Not available) (24.69-1.80) 98.8 (24.69-1.80)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$< I/\sigma(I) >$ [1]	1.28 (at 1.80Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.222 , 0.245 0.212 , 0.231	Depositor DCC
R_{free} test set	1000 reflections (1.09%)	wwPDB-VP
Wilson B-factor (Å^2)	31.5	Xtriage
Anisotropy	0.277	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.38 , 54.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	7838	wwPDB-VP
Average B, all atoms (Å^2)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.33% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ACY, GOL, NA, NAG

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\#\|Z\| > 5$	RMSZ	$\#\|Z\| > 5$
1	A	0.39	0/373	0.61	0/499
1	C	0.36	0/392	0.65	0/523
1	E	0.36	0/385	0.59	0/513
2	B	0.36	0/2147	0.64	0/2902
2	D	0.31	0/2133	0.58	0/2885
2	F	0.34	0/2147	0.62	1/2902 (0.0%)
All	All	0.34	0/7577	0.61	1/10224 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	F	147(C)	ALA	N-CA-C	-6.25	94.12	111.00

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	A	366	0	344	6	0
1	C	385	0	364	16	0
1	E	378	0	357	14	0
2	B	2092	0	2062	27	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	2078	0	2055	38	0
2	F	2092	0	2062	52	0
3	B	14	0	13	2	0
3	F	14	0	13	0	0
4	B	1	0	0	0	0
4	D	1	0	0	0	0
4	F	1	0	0	0	0
5	B	24	0	32	1	0
6	B	8	0	6	1	0
6	D	4	0	3	0	0
6	F	8	0	6	2	0
7	A	28	0	0	0	0
7	B	142	0	0	3	0
7	C	13	0	0	0	0
7	D	72	0	0	0	0
7	E	17	0	0	0	0
7	F	100	0	0	4	0
All	All	7838	0	7317	142	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 10.

All (142) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:14(L):ASP:CG	1:E:14(M):GLY:H	1.60	1.04
2:D:93:ARG:HB3	2:D:101:ARG:HD2	1.45	0.94
2:B:246:GLY:O	2:B:247:GLU:HB2	1.66	0.94
1:C:14(L):ASP:CG	1:C:14(M):GLY:H	1.69	0.90
1:E:14(L):ASP:CG	1:E:14(M):GLY:N	2.30	0.84
2:F:110:LYS:HD2	2:F:111:PRO:HD2	1.62	0.82
2:B:60(I):THR:HG22	2:B:63:ASP:OD2	1.84	0.78
2:D:32:MET:HE3	2:D:70:LYS:HD3	1.66	0.78
1:C:14(L):ASP:CG	1:C:14(M):GLY:N	2.39	0.75
2:F:147(E):VAL:HG12	2:F:148:GLY:N	2.03	0.73
1:E:1(K):ASN:ND2	1:E:1(I):ARG:HB3	2.09	0.68
1:A:10:LYS:NZ	1:A:12:LEU:HD12	2.08	0.67
1:A:14(K):ILE:HG13	1:A:14(L):ASP:H	1.60	0.67
2:F:18:GLU:HG3	2:F:187:ARG:HG3	1.78	0.66
2:F:33:LEU:HD21	2:F:106:MET:CE	2.26	0.65
2:D:75:ARG:HB3	2:D:75:ARG:NH1	2.12	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:60(C):PRO:HD3	2:D:96:TRP:CZ3	2.32	0.64
2:D:50:ARG:HG3	2:D:51:TRP:CD1	2.33	0.64
2:D:167:VAL:HG11	2:D:185:LYS:HE2	1.79	0.64
2:F:33:LEU:HD21	2:F:106:MET:HE2	1.80	0.63
2:D:91:HIS:CE1	2:D:93:ARG:HB2	2.33	0.63
2:B:230:HIS:NE2	5:B:1003:GOL:H11	2.15	0.61
2:B:61:GLU:HG2	2:B:87:LYS:HA	1.83	0.60
2:F:204(B):ASN:C	2:F:204(B):ASN:HD22	2.04	0.60
2:D:98:ASN:O	2:D:99:LEU:HB2	2.01	0.60
2:F:147(E):VAL:CG1	2:F:148:GLY:N	2.65	0.60
2:F:110:LYS:HD2	2:F:111:PRO:CD	2.31	0.59
2:F:75:ARG:HH21	2:F:75:ARG:HB3	1.67	0.59
2:B:60(C):PRO:HA	2:D:164:GLU:HG2	1.84	0.58
2:B:81:LYS:HG2	7:B:3135:HOH:O	2.03	0.58
2:B:80:GLU:O	2:B:81:LYS:HD2	2.04	0.57
2:F:202:LYS:HE2	2:F:205:ASN:OD1	2.05	0.57
1:E:1(G):PHE:HD1	2:F:242:ILE:HD13	1.69	0.56
2:F:246:GLY:O	2:F:247:GLU:HG2	2.04	0.56
1:E:6:LEU:HA	1:E:10:LYS:HE2	1.88	0.56
1:E:14(L):ASP:OD1	1:E:14(M):GLY:N	2.19	0.56
1:C:1(G):PHE:HE2	1:C:1(C):GLU:HB2	1.70	0.56
1:C:1(K):ASN:ND2	1:C:1(I):ARG:HB3	2.20	0.56
2:D:75:ARG:HB3	2:D:75:ARG:HH11	1.71	0.56
2:F:204(B):ASN:ND2	2:F:206:ARG:H	2.04	0.56
2:D:169:LYS:HA	2:D:176:ILE:HD12	1.89	0.55
2:F:136:GLY:HA3	2:F:199:PHE:CZ	2.41	0.55
1:E:1(K):ASN:HD21	1:E:1(I):ARG:HB3	1.71	0.54
2:F:147(B):THR:CB	2:F:147(E):VAL:HA	2.37	0.54
1:C:1(G):PHE:HD1	2:D:242:ILE:HD13	1.71	0.54
2:B:165:ARG:HG2	2:B:169:LYS:HE3	1.88	0.54
2:F:60(A):TYR:CE2	2:F:60(C):PRO:HG2	2.43	0.53
2:F:82:ILE:HD12	2:F:82:ILE:N	2.23	0.53
2:F:147(B):THR:HA	2:F:147(E):VAL:N	2.23	0.53
2:D:204(B):ASN:HD22	2:D:204(B):ASN:C	2.12	0.53
1:A:14(D):ARG:O	1:A:14(H):GLU:HG3	2.08	0.53
1:C:14(D):ARG:O	1:C:14(H):GLU:HG3	2.09	0.52
2:F:147(B):THR:HA	2:F:147(E):VAL:HA	1.91	0.52
2:D:93:ARG:HB3	2:D:101:ARG:CD	2.30	0.52
2:B:35:ARG:HH21	2:B:60(H):PHE:HZ	1.56	0.52
2:D:91:HIS:HE1	2:D:93:ARG:HB2	1.75	0.51
2:F:41:LEU:HD12	2:F:64:LEU:HD22	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:221(A):ARG:HB2	2:D:224:LYS:HB2	1.93	0.51
2:D:32:MET:CE	2:D:70:LYS:HD3	2.38	0.50
2:D:185:LYS:HB2	2:D:186(B):GLU:HG3	1.94	0.50
1:A:14(K):ILE:O	1:A:14(L):ASP:CB	2.60	0.49
2:D:61:GLU:CD	2:D:61:GLU:H	2.15	0.49
1:E:1:CYS:O	2:F:206:ARG:HD3	2.13	0.49
2:F:30:GLN:NE2	2:F:139:THR:OG1	2.45	0.49
2:B:60(B):PRO:HB2	2:B:60(C):PRO:HD3	1.95	0.49
2:F:143:ASN:ND2	2:F:192:GLU:HB3	2.28	0.49
2:D:144:LEU:HD21	2:D:152:PRO:HB3	1.95	0.48
3:B:1:NAG:O7	3:B:1:NAG:H3	2.13	0.48
2:F:147(E):VAL:CG1	2:F:148:GLY:H	2.25	0.48
2:B:61:GLU:CG	2:B:87:LYS:HA	2.44	0.48
2:F:147(E):VAL:HG12	2:F:148:GLY:H	1.78	0.48
1:E:1(G):PHE:CD1	2:F:242:ILE:HD13	2.48	0.48
2:F:60(D):TRP:CE3	6:F:2004:ACY:H3	2.49	0.48
1:E:3:LEU:HD21	2:F:206:ARG:HG2	1.95	0.48
2:F:144:LEU:HD21	2:F:152:PRO:HB3	1.95	0.48
2:B:204(B):ASN:C	2:B:204(B):ASN:HD22	2.16	0.47
1:C:1(H):THR:C	1:C:1(F):GLY:H	2.17	0.47
2:D:41:LEU:O	2:D:42:CYS:HB3	2.14	0.47
1:A:14(K):ILE:HG13	1:A:14(L):ASP:N	2.27	0.47
2:D:31:VAL:CG1	2:D:66:VAL:HG13	2.44	0.47
1:E:14(D):ARG:O	1:E:14(H):GLU:HG3	2.15	0.47
1:C:1(H):THR:O	1:C:1(G):PHE:HB3	2.15	0.47
6:F:2003:ACY:H2	7:F:3030:HOH:O	2.15	0.47
2:B:195:ALA:HB2	1:C:15:ARG:C	2.35	0.46
2:F:59:LEU:HD13	2:F:88:ILE:HG21	1.97	0.46
2:B:89:TYR:OH	2:B:245:PHE:HB3	2.15	0.46
1:E:14(A):LYS:HG2	2:F:23:GLU:OE2	2.15	0.46
2:D:32:MET:HG3	2:D:40:LEU:HD13	1.97	0.46
1:C:1(P):TYR:HE2	1:C:1(C):GLU:HG2	1.81	0.46
2:D:136:GLY:HA3	2:D:199:PHE:CZ	2.51	0.46
2:D:130:LEU:HD23	2:D:210:MET:HE2	1.98	0.45
2:F:204(B):ASN:HD22	2:F:206:ARG:H	1.64	0.45
2:F:175:ARG:N	2:F:175:ARG:HD2	2.31	0.45
1:C:14(G):LEU:HD22	1:C:14(G):LEU:N	2.32	0.45
1:E:14(C):GLU:O	1:E:14(G):LEU:HD23	2.16	0.45
2:F:41:LEU:CD1	2:F:64:LEU:HD22	2.47	0.45
2:F:147(B):THR:HB	2:F:147(E):VAL:HA	2.00	0.44
2:D:77(A):ARG:O	2:D:78:ASN:HB2	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:33:LEU:HD23	2:F:44:ALA:HB2	1.98	0.44
2:B:136:GLY:HA3	2:B:199:PHE:CZ	2.53	0.44
2:D:147(C):ALA:O	2:D:147(D):ASN:HB2	2.18	0.44
2:F:146:GLU:OE2	2:F:221(A):ARG:NH2	2.51	0.44
2:D:44:ALA:HA	2:D:196:GLY:O	2.17	0.44
2:B:81:LYS:HE3	7:B:3066:HOH:O	2.17	0.43
1:C:1(G):PHE:HA	2:D:51:TRP:CD1	2.53	0.43
2:F:33:LEU:HD21	2:F:106:MET:HE1	1.98	0.43
2:D:130:LEU:HD21	2:D:210:MET:HE3	2.01	0.43
2:F:75:ARG:HB3	2:F:75:ARG:NH2	2.33	0.43
2:D:84:MET:HB2	2:D:109:LYS:HD2	1.99	0.43
2:F:96:TRP:NE1	7:F:3102:HOH:O	2.32	0.43
2:F:85:LEU:HD22	2:F:106:MET:HB3	2.00	0.43
2:B:80:GLU:C	2:B:81:LYS:HD2	2.39	0.43
2:F:246:GLY:O	2:F:247:GLU:CB	2.66	0.43
2:F:147(B):THR:C	2:F:147(D):ASN:N	2.63	0.43
2:F:147(A):TRP:HB2	2:F:147(B):THR:H	1.63	0.43
2:D:60(B):PRO:HD2	2:D:96:TRP:CE3	2.54	0.43
2:D:93:ARG:CB	2:D:101:ARG:HD2	2.33	0.43
2:F:147(B):THR:HA	2:F:147(E):VAL:CA	2.49	0.42
2:D:33:LEU:HD11	2:D:106:MET:CE	2.49	0.42
2:F:77(A):ARG:O	2:F:78:ASN:HB2	2.19	0.42
2:F:97:ARG:NH2	7:F:3102:HOH:O	2.52	0.42
1:A:14(K):ILE:O	1:A:14(L):ASP:HB2	2.18	0.42
2:B:32:MET:HG2	2:B:40:LEU:CD1	2.49	0.42
1:E:14(A):LYS:HG2	2:F:23:GLU:CD	2.40	0.42
2:F:204(B):ASN:C	2:F:204(B):ASN:ND2	2.73	0.42
2:D:93:ARG:HG2	2:D:101:ARG:CZ	2.50	0.42
2:F:60(B):PRO:N	2:F:60(C):PRO:HD2	2.35	0.41
1:C:1(R):SER:OG	2:D:206:ARG:NE	2.53	0.41
2:F:126:ARG:HG3	2:F:127:GLU:OE2	2.21	0.41
2:B:147(E):VAL:HG13	2:B:148:GLY:N	2.35	0.41
2:B:18:GLU:HB2	2:B:188:GLY:HA2	2.02	0.41
2:B:60(A):TYR:OH	1:C:14(K):ILE:HD11	2.21	0.41
2:D:230:HIS:CE1	2:D:233:ARG:HG3	2.56	0.41
2:B:173:ARG:HH11	1:C:14(D):ARG:NH2	2.18	0.41
2:D:129:ALA:O	2:D:130:LEU:HB2	2.21	0.41
2:B:127:GLU:HG3	7:B:3114:HOH:O	2.21	0.41
2:B:29:TRP:CG	2:B:121:VAL:HB	2.56	0.40
2:B:148:GLY:HA3	6:B:2005:ACY:OXT	2.21	0.40
2:B:60(B):PRO:HB3	3:B:1:NAG:H4	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:109:LYS:NZ	7:F:3055:HOH:O	2.54	0.40
2:B:130:LEU:HD12	2:B:130:LEU:HA	1.92	0.40
1:C:1(H):THR:O	1:C:1(F):GLY:N	2.53	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	43/49 (88%)	39 (91%)	3 (7%)	1 (2%)	6 1
1	C	45/49 (92%)	42 (93%)	3 (7%)	0	100 100
1	E	44/49 (90%)	41 (93%)	3 (7%)	0	100 100
2	B	257/259 (99%)	250 (97%)	7 (3%)	0	100 100
2	D	255/259 (98%)	234 (92%)	20 (8%)	1 (0%)	34 21
2	F	257/259 (99%)	239 (93%)	14 (5%)	4 (2%)	9 2
All	All	901/924 (98%)	845 (94%)	50 (6%)	6 (1%)	22 10

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	14(L)	ASP
2	F	147(E)	VAL
2	F	147(D)	ASN
2	F	149	LYS
2	F	148	GLY
2	D	98	ASN

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	A	40/43 (93%)	39 (98%)	1 (2%)	47	34
1	C	42/43 (98%)	41 (98%)	1 (2%)	49	36
1	E	41/43 (95%)	40 (98%)	1 (2%)	49	36
2	B	224/224 (100%)	219 (98%)	5 (2%)	52	39
2	D	223/224 (100%)	216 (97%)	7 (3%)	40	25
2	F	224/224 (100%)	219 (98%)	5 (2%)	52	39
All	All	794/801 (99%)	774 (98%)	20 (2%)	47	34

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14(L)	ASP
2	B	33	LEU
2	B	130	LEU
2	B	182	CYS
2	B	204(B)	ASN
2	B	243	ASP
1	C	1(S)	THR
2	D	42	CYS
2	D	60(G)	ASN
2	D	93	ARG
2	D	95	ASN
2	D	182	CYS
2	D	186(A)	ASP
2	D	204(B)	ASN
1	E	14(L)	ASP
2	F	33	LEU
2	F	77(A)	ARG
2	F	95	ASN
2	F	186(A)	ASP
2	F	204(B)	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14)

such sidechains are listed below:

Mol	Chain	Res	Type
2	B	204(B)	ASN
2	B	239	GLN
2	D	60(G)	ASN
2	D	95	ASN
2	D	156	GLN
2	D	204(B)	ASN
2	D	239	GLN
2	D	244	GLN
1	E	1(O)	GLN
2	F	30	GLN
2	F	78	ASN
2	F	95	ASN
2	F	156	GLN
2	F	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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates

There are no monosaccharides in this entry.

5.6 Ligand geometry

Of 14 ligands modelled in this entry, 3 are monoatomic - leaving 11 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	$\#\|Z\|>2$
6	ACY	B	2005	-	3,3,3	1.29	0	3,3,3	1.59	1 (33%)
3	NAG	F	2	2	14,14,15	0.52	0	17,19,21	0.66	0
3	NAG	B	1	2	14,14,15	0.59	0	17,19,21	0.60	0
6	ACY	D	2002	-	3,3,3	1.22	0	3,3,3	1.66	1 (33%)
6	ACY	F	2003	-	3,3,3	1.18	0	3,3,3	1.79	1 (33%)
5	GOL	B	1002	-	5,5,5	0.74	0	5,5,5	0.44	0
6	ACY	B	2001	-	3,3,3	1.24	0	3,3,3	1.67	1 (33%)
5	GOL	B	1004	-	5,5,5	0.84	0	5,5,5	0.67	0
5	GOL	B	1001	-	5,5,5	0.75	0	5,5,5	0.42	0
6	ACY	F	2004	-	3,3,3	1.18	0	3,3,3	1.68	1 (33%)
5	GOL	B	1003	-	5,5,5	0.69	0	5,5,5	0.46	0

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	NAG	F	2	2	1/1/5/7	4/6/23/26	0/1/1/1
3	NAG	B	1	2	1/1/5/7	5/6/23/26	0/1/1/1
5	GOL	B	1002	-	-	2/4/4/4	-
5	GOL	B	1004	-	-	2/4/4/4	-
5	GOL	B	1001	-	-	2/4/4/4	-
5	GOL	B	1003	-	-	4/4/4/4	-

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
6	F	2003	ACY	O-C-CH3	-2.45	112.78	122.33
6	F	2004	ACY	O-C-CH3	-2.31	113.35	122.33
6	B	2001	ACY	O-C-CH3	-2.30	113.36	122.33
6	D	2002	ACY	O-C-CH3	-2.28	113.45	122.33
6	B	2005	ACY	O-C-CH3	-2.19	113.81	122.33

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	B	1	NAG	C1

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Mol	Chain	Res	Type	Atom
3	F	2	NAG	C1

All (19) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	1	NAG	C3-C2-N2-C7
3	B	1	NAG	C8-C7-N2-C2
3	B	1	NAG	O7-C7-N2-C2
3	F	2	NAG	C8-C7-N2-C2
3	F	2	NAG	O7-C7-N2-C2
5	B	1004	GOL	C1-C2-C3-O3
3	F	2	NAG	O5-C5-C6-O6
3	F	2	NAG	C4-C5-C6-O6
5	B	1004	GOL	O2-C2-C3-O3
3	B	1	NAG	O5-C5-C6-O6
5	B	1001	GOL	C1-C2-C3-O3
5	B	1002	GOL	C1-C2-C3-O3
5	B	1003	GOL	C1-C2-C3-O3
5	B	1002	GOL	O2-C2-C3-O3
5	B	1001	GOL	O2-C2-C3-O3
5	B	1003	GOL	O1-C1-C2-O2
5	B	1003	GOL	O2-C2-C3-O3
3	B	1	NAG	C4-C5-C6-O6
5	B	1003	GOL	O1-C1-C2-C3

There are no ring outliers.

5 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	B	2005	ACY	1	0
3	B	1	NAG	2	0
6	F	2003	ACY	1	0
6	F	2004	ACY	1	0
5	B	1003	GOL	1	0

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95^{th} percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å^2)	Q<0.9
1	A	45/49 (91%)	0.07	4 (8%) 9 7	21, 32, 56, 63	0
1	C	47/49 (95%)	0.77	6 (12%) 3 2	24, 45, 66, 72	1 (2%)
1	E	46/49 (93%)	0.67	7 (15%) 2 1	33, 43, 56, 61	0
2	B	259/259 (100%)	0.25	8 (3%) 49 43	18, 28, 50, 68	2 (0%)
2	D	257/259 (99%)	0.66	34 (13%) 3 2	24, 40, 63, 74	5 (1%)
2	F	259/259 (100%)	0.52	22 (8%) 10 8	22, 35, 58, 85	3 (1%)
All	All	913/924 (98%)	0.48	81 (8%) 9 7	18, 35, 60, 85	11 (1%)

All (81) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	147(B)	THR	10.5
2	F	147(C)	ALA	9.7
2	F	147(A)	TRP	7.1
2	F	147(D)	ASN	6.9
1	C	1(S)	THR	6.3
2	F	247	GLU	6.2
2	B	247	GLU	6.1
2	F	147(E)	VAL	6.0
2	F	186(C)	GLY	5.7
2	D	31	VAL	4.9
2	D	76	TYR	4.8
2	D	96	TRP	4.8
2	F	148	GLY	4.5
2	F	60(E)	ASP	4.3
1	A	1(R)	SER	4.1
2	D	60(E)	ASP	4.1
1	C	1(R)	SER	4.0
2	D	46	LEU	3.9
2	D	60(I)	THR	3.8

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Mol	Chain	Res	Type	RSRZ
2	D	243	ASP	3.6
2	D	245	PHE	3.6
2	D	77(A)	ARG	3.6
1	C	1(K)	ASN	3.6
2	F	147	THR	3.5
2	F	186(A)	ASP	3.4
2	D	52	VAL	3.4
2	D	37	PRO	3.2
1	A	1(Q)	GLU	3.2
2	D	38	GLN	3.2
2	D	244	GLN	3.1
2	D	78	ASN	3.1
1	E	1(R)	SER	3.1
2	D	62	ASN	3.1
1	E	1(I)	ARG	3.1
2	F	78	ASN	3.0
2	D	53	LEU	2.9
2	F	97(A)	GLU	2.9
2	D	29	TRP	2.9
2	D	47	ILE	2.9
1	A	14(M)	GLY	2.8
2	D	44	ALA	2.8
2	D	45	SER	2.8
2	F	149	LYS	2.8
2	D	34	PHE	2.7
2	B	186(C)	GLY	2.7
2	D	42	CYS	2.7
1	C	1(Q)	GLU	2.7
1	A	14(L)	ASP	2.6
1	C	1(H)	THR	2.6
2	B	97(A)	GLU	2.6
2	D	75	ARG	2.6
2	D	149	LYS	2.6
2	F	243	ASP	2.6
2	F	97	ARG	2.5
1	E	11	SER	2.5
1	E	1(N)	THR	2.4
2	D	186(A)	ASP	2.4
2	F	44	ALA	2.4
1	E	10	LYS	2.4
2	D	60(H)	PHE	2.4
2	F	245	PHE	2.4

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Mol	Chain	Res	Type	RSRZ
2	D	108	LEU	2.4
2	F	212	ILE	2.4
2	D	36(A)	SER	2.3
2	D	60(G)	ASN	2.3
2	F	96	TRP	2.3
2	D	121	VAL	2.3
2	B	47	ILE	2.3
2	D	54	THR	2.3
2	B	199	PHE	2.3
1	E	1(Q)	GLU	2.2
2	F	60(D)	TRP	2.2
2	B	186(A)	ASP	2.2
2	D	147(E)	VAL	2.2
2	D	63	ASP	2.1
2	F	222	ASP	2.1
2	B	29	TRP	2.1
2	B	212	ILE	2.1
2	D	147(D)	ASN	2.1
1	E	1(O)	GLN	2.0
1	C	1(J)	PRO	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
5	GOL	B	1004	6/6	0.50	0.23	66,68,69,69	0
5	GOL	B	1003	6/6	0.63	0.35	66,67,69,69	0
3	NAG	B	1	14/15	0.68	0.51	68,71,72,73	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	NAG	F	2	14/15	0.68	0.35	68,69,72,72	0
5	GOL	B	1002	6/6	0.79	0.15	66,69,69,70	0
6	ACY	B	2001	4/4	0.84	0.25	49,50,50,51	0
6	ACY	F	2003	4/4	0.85	0.13	36,37,38,39	0
5	GOL	B	1001	6/6	0.87	0.13	37,44,44,47	0
6	ACY	B	2005	4/4	0.91	0.09	35,37,37,40	0
6	ACY	D	2002	4/4	0.92	0.13	56,56,56,57	0
6	ACY	F	2004	4/4	0.92	0.13	47,48,50,50	0
4	NA	D	3002	1/1	0.94	0.14	40,40,40,40	0
4	NA	F	3003	1/1	0.99	0.09	38,38,38,38	0
4	NA	B	3001	1/1	0.99	0.07	27,27,27,27	0

6.5 Other polymers (i)

There are no such residues in this entry.