Full wwPDB X-ray Structure Validation Report

Jan 31, 2016 – 11:14 PM GMT

PDB ID	:	1WU1
Title	:	Factor Xa in complex with the inhibitor 4-[(5-chloroindol-2-yl)sulfonyl]-2-(2-methylpropyl)-1-[[5-(pyridin-4-yl) pyrimidin-2-yl]carbonyl]piperazine
Authors	:	Suzuki, M.
Deposited on	:	2004-11-29
Resolution	:	2.30 Å(reported)

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

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

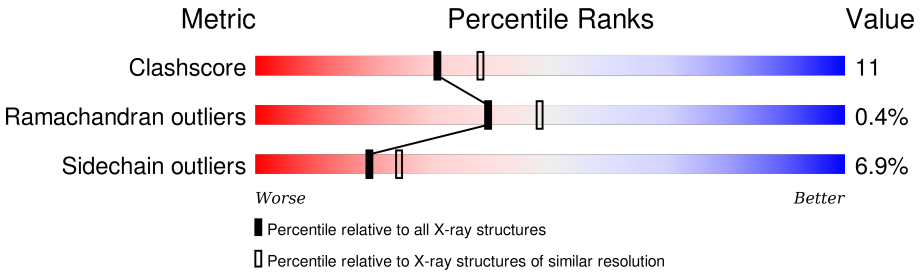
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)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

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	233	77% 21%
2	B	95	42% 14% 43%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 2266 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 Coagulation factor X, heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	233	Total	C	N	O	S	0	0	0
			1777	1131	300	332	14			

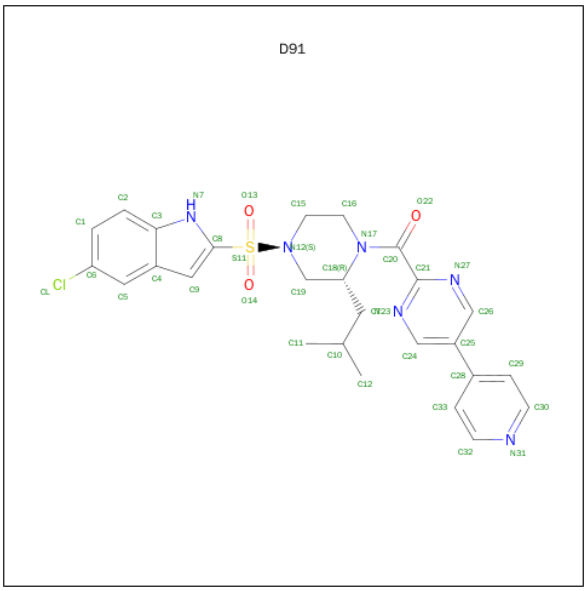
- Molecule 2 is a protein called Coagulation factor X, light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	54	Total	C	N	O	S	0	0	0
			378	228	63	80	7			

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Ca	0	0
			2	2		

- Molecule 4 is 5-CHLORO-2-({3-ISOBUTYL-4-[(5-PYRIDIN-4-YLPYRIMIDIN-2-YL)CARBONYL]PIPERAZIN-1-YL}SULFONYL)-1H-INDOLE (three-letter code: D91) (formula: $C_{26}H_{27}ClN_6O_3S$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
4	A	1	Total	C	Cl	N	O	S	0	0
			33	22	1	6	3	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	59	Total	O	0	0
			59	59		
5	B	17	Total	O	0	0
			17	17		

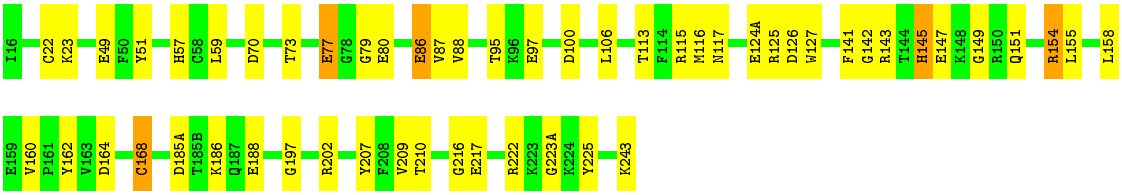
3 Residue-property plots (i)

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.

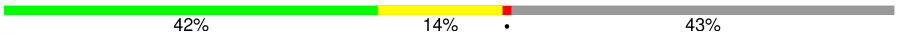
- Molecule 1: Coagulation factor X, heavy chain

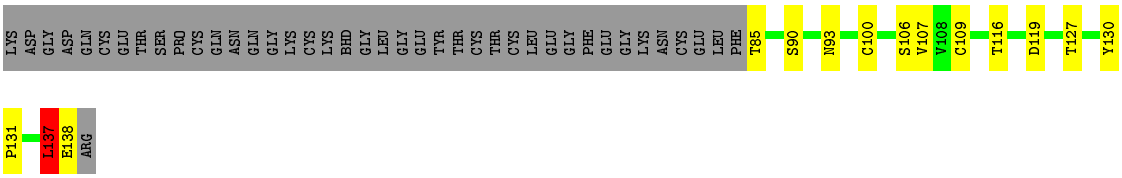
Chain A:



- Molecule 2: Coagulation factor X, light chain

Chain B:





4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	56.57Å 72.49Å 79.02Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.00 – 2.30	Depositor
% Data completeness (in resolution range)	100.0 (25.00-2.30)	Depositor
R_{merge}	0.06	Depositor
R_{sym}	0.34	Depositor
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.189 , 0.235	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2266	wwPDB-VP
Average B, all atoms (Å^2)	47.0	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, D91

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	1.03	4/1815 (0.2%)	1.00	3/2453 (0.1%)
2	B	1.13	1/384 (0.3%)	1.04	0/523
All	All	1.05	5/2199 (0.2%)	1.01	3/2976 (0.1%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	109	CYS	CB-SG	-7.95	1.68	1.82
1	A	162	TYR	CD2-CE2	5.86	1.48	1.39
1	A	168	CYS	CB-SG	-5.62	1.72	1.81
1	A	160	VAL	CB-CG1	-5.18	1.42	1.52
1	A	22	CYS	CB-SG	-5.02	1.73	1.81

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	154	ARG	NE-CZ-NH2	7.09	123.85	120.30
1	A	100	ASP	CB-CG-OD1	5.99	123.69	118.30
1	A	154	ARG	NE-CZ-NH1	-5.30	117.65	120.30

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	1777	0	1678	38	0
2	B	378	0	326	10	0
3	A	2	0	0	0	0
4	A	33	0	17	2	0
5	A	59	0	0	5	0
5	B	17	0	0	1	0
All	All	2266	0	2021	46	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 11.

All (46) 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:A:86:GLU:HG2	1:A:87:VAL:HG23	1.38	1.02
1:A:49:GLU:HG2	5:A:344:HOH:O	1.68	0.91
1:A:124(A):GLU:HB3	5:A:324:HOH:O	1.73	0.86
1:A:145:HIS:CD2	1:A:149:GLY:HA3	2.14	0.81
2:B:137:LEU:HD12	2:B:138:GLU:N	1.99	0.77
1:A:77:GLU:HG2	1:A:80:GLU:OE1	1.83	0.77
1:A:217:GLU:O	1:A:222:ARG:HD2	1.87	0.74
1:A:185(A):ASP:OD1	1:A:223(A):GLY:HA2	1.89	0.73
2:B:116:THR:HG23	2:B:127:THR:CG2	2.20	0.70
1:A:158:LEU:HD11	1:A:188:GLU:HB3	1.77	0.66
1:A:145:HIS:CD2	1:A:149:GLY:CA	2.81	0.63
1:A:86:GLU:CG	1:A:87:VAL:HG23	2.21	0.61
1:A:127:TRP:HE1	2:B:93:ASN:HD21	1.49	0.60
2:B:116:THR:HG23	2:B:127:THR:HG23	1.84	0.59
1:A:86:GLU:HG2	1:A:87:VAL:CG2	2.23	0.58
1:A:59:LEU:HD13	1:A:88:VAL:HG11	1.86	0.58
1:A:79:GLY:O	1:A:115:ARG:NH2	2.37	0.57
1:A:127:TRP:HE1	2:B:93:ASN:ND2	2.03	0.56
1:A:147:GLU:OE2	1:A:222:ARG:NH1	2.39	0.55
1:A:197:GLY:HA3	5:A:303:HOH:O	2.07	0.55
2:B:100:CYS:SG	2:B:107:VAL:HG13	2.47	0.54
1:A:116:MET:O	1:A:117:ASN:HB2	2.07	0.53
1:A:97:GLU:O	4:A:700:D91:H32	2.08	0.53
1:A:113:THR:O	1:A:113:THR:HG22	2.08	0.53
1:A:145:HIS:HD2	1:A:149:GLY:CA	2.20	0.52
1:A:145:HIS:HD2	1:A:149:GLY:N	2.07	0.52
1:A:142:GLY:HA2	1:A:151:GLN:HE21	1.76	0.51
2:B:130:TYR:N	2:B:131:PRO:CD	2.75	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:137:LEU:HD12	2:B:137:LEU:C	2.32	0.49
1:A:158:LEU:CD1	1:A:188:GLU:HB3	2.42	0.49
2:B:106:SER:HB2	5:B:320:HOH:O	2.12	0.49
1:A:186:LYS:HE2	1:A:188:GLU:OE1	2.13	0.48
2:B:116:THR:CG2	2:B:127:THR:CG2	2.91	0.46
1:A:125:ARG:NH2	1:A:126:ASP:OD1	2.36	0.46
1:A:77:GLU:HG2	1:A:80:GLU:CD	2.37	0.45
1:A:141:PHE:CZ	1:A:155:LEU:HB2	2.52	0.45
1:A:79:GLY:O	1:A:115:ARG:NE	2.50	0.44
1:A:154:ARG:HA	1:A:154:ARG:HD2	1.91	0.43
1:A:51:TYR:HA	1:A:106:LEU:O	2.19	0.42
1:A:216:GLY:O	4:A:700:D91:H191	2.20	0.42
1:A:70:ASP:OD2	1:A:73:THR:HG22	2.21	0.41
1:A:23:LYS:HD3	1:A:23:LYS:HA	1.70	0.41
1:A:124(A):GLU:CB	5:A:324:HOH:O	2.50	0.40
1:A:225:TYR:C	5:A:349:HOH:O	2.60	0.40
1:A:142:GLY:HA2	1:A:151:GLN:NE2	2.35	0.40
1:A:202:ARG:HD3	1:A:207:TYR:CZ	2.57	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	231/233 (99%)	219 (95%)	12 (5%)	0	100 100
2	B	52/95 (55%)	45 (86%)	6 (12%)	1 (2%)	10 8
All	All	283/328 (86%)	264 (93%)	18 (6%)	1 (0%)	39 48

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	137	LEU

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	177/198 (89%)	166 (94%)	11 (6%)	23 30
2	B	41/80 (51%)	37 (90%)	4 (10%)	10 11
All	All	218/278 (78%)	203 (93%)	15 (7%)	19 24

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	57	HIS
1	A	77	GLU
1	A	86	GLU
1	A	95	THR
1	A	143	ARG
1	A	145	HIS
1	A	164	ASP
1	A	168	CYS
1	A	209	VAL
1	A	210	THR
1	A	243	LYS
2	B	85	THR
2	B	90	SER
2	B	119	ASP
2	B	137	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	145	HIS
1	A	151	GLN
1	A	166	ASN

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Mol	Chain	Res	Type
2	B	93	ASN
2	B	135	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

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	D91	A	700	-	33,37,41	2.96	7 (21%)	45,54,60	2.41	14 (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
4	D91	A	700	-	-	0/14/34/41	0/5/5/5

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	700	D91	C21-C20	-13.53	1.46	1.53
4	A	700	D91	C5-C4	-3.53	1.34	1.42
4	A	700	D91	C2-C3	-2.56	1.37	1.41
4	A	700	D91	C3-N7	-2.25	1.31	1.39
4	A	700	D91	C9-C4	2.69	1.51	1.41
4	A	700	D91	C20-N17	2.70	1.40	1.34
4	A	700	D91	S11-N12	7.30	1.74	1.63

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	A	700	D91	C8-C9-C4	-6.95	97.88	106.55
4	A	700	D91	C19-C18-N17	-5.53	97.91	110.49
4	A	700	D91	C5-C6-CL	-4.51	113.31	119.74
4	A	700	D91	O22-C20-N17	-4.25	115.66	122.42
4	A	700	D91	C15-C16-N17	-3.80	101.83	110.49
4	A	700	D91	O14-S11-C8	-3.22	103.05	107.75
4	A	700	D91	C2-C1-C6	-2.43	116.25	119.28
4	A	700	D91	C29-C28-C25	-2.22	117.42	121.39
4	A	700	D91	C1-C6-C5	2.22	124.78	121.87
4	A	700	D91	C16-N17-C18	2.50	117.19	112.56
4	A	700	D91	C15-N12-S11	3.00	122.89	117.12
4	A	700	D91	O13-S11-N12	3.10	109.75	106.69
4	A	700	D91	C21-C20-N17	3.13	127.13	119.61
4	A	700	D91	C15-N12-C19	3.91	116.94	112.20

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	700	D91	2	0

5.7 Other polymers (i)

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.