



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 26, 2017 – 12:01 PM EDT

PDB ID : 1IQE
Title : Human coagulation factor Xa in complex with M55590
Authors : Shiromizu, I.; Matsusue, T.
Deposited on : unknown
Resolution : 2.90 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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20030345

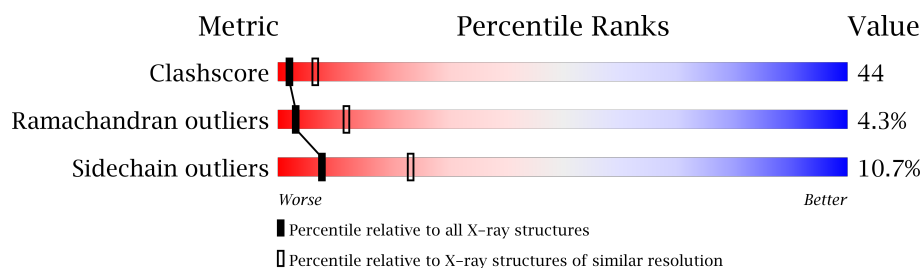
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	112137	1807 (2.90-2.90)
Ramachandran outliers	110173	1768 (2.90-2.90)
Sidechain outliers	110143	1770 (2.90-2.90)

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	A	235	
2	L	96	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 2281 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 Xa.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	235	Total	C	N	O	S	0	0	0
			1864	1172	327	351	14			

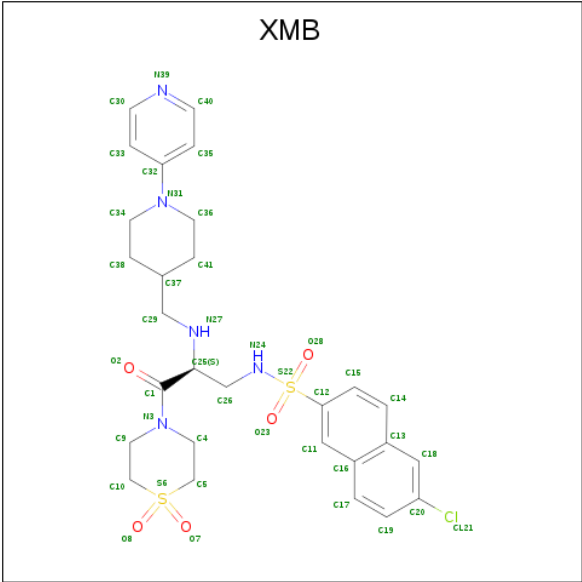
- Molecule 2 is a protein called coagulation Factor Xa.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	51	Total	C	N	O	S	0	0	0
			375	226	65	77	7			

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Ca	0	0
			1	1		

- Molecule 4 is 4-[(2R)-3-[[[(6-CHLORO-2-NAPHTHALENYL)SULFONYL]AMINO]-1-OXO-2-[[[1-(4-PYRIDINYL)-4-PIPERIDINYL]METHYL]AMINO]PROPYL]-THIOMORPHOLINE-1,1-DIOXIDE (three-letter code: XMB) (formula: C₂₈H₃₄ClN₅O₅S₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
4	A	1	Total	C	Cl	N	O	S	0	0
			41	28	1	5	5	2		

4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	72.49 Å 78.65 Å 55.75 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.90	Depositor
% Data completeness (in resolution range)	81.9 (8.00-2.90)	Depositor
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNX 2000	Depositor
R, R_{free}	0.213 , 0.297	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2281	wwPDB-VP
Average B, all atoms (Å ²)	22.0	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, XMB

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	0.41	0/1902	0.66	0/2560
2	L	0.37	0/381	0.68	0/515
All	All	0.40	0/2283	0.66	0/3075

There are no bond length outliers.

There are no bond angle outliers.

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	1864	0	1821	183	0
2	L	375	0	343	23	0
3	A	1	0	0	0	0
4	A	41	0	34	9	0
All	All	2281	0	2198	199	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 44.

All (199) 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:96:LYS:H	1:A:96:LYS:HD3	1.22	1.05
2:L:113:ARG:NH1	2:L:113:ARG:H	1.59	1.00
1:A:232:THR:HA	1:A:235:LEU:HD22	1.49	0.90
1:A:240:ARG:HD3	1:A:245:ARG:HD2	1.53	0.90
1:A:215:TRP:CE3	4:A:401:XMB:H362	2.07	0.89
1:A:63:ARG:HH12	1:A:65:LYS:HE3	1.37	0.89
1:A:158:LEU:HD11	1:A:188:GLU:HB3	1.53	0.88
2:L:113:ARG:HH11	2:L:113:ARG:H	1.16	0.87
2:L:87:LYS:O	2:L:88:LEU:HB2	1.73	0.85
1:A:57:HIS:NE2	1:A:195:SER:OG	2.10	0.85
1:A:96:LYS:N	1:A:96:LYS:HD3	1.90	0.85
1:A:32:LEU:HD13	1:A:141:PHE:HD2	1.39	0.84
1:A:211:GLY:HA2	1:A:231:VAL:HG23	1.60	0.83
1:A:217:GLU:OE1	1:A:224:LYS:HD3	1.76	0.83
1:A:50:PHE:CE1	1:A:111:PRO:HG3	2.15	0.81
1:A:72:ASN:ND2	1:A:154:ARG:HE	1.79	0.79
1:A:168:CYS:SG	1:A:176:ILE:HD13	2.22	0.78
1:A:74:GLU:HB2	1:A:75:GLN:OE1	1.86	0.76
2:L:113:ARG:HH11	2:L:113:ARG:N	1.85	0.74
1:A:32:LEU:HD13	1:A:141:PHE:CD2	2.23	0.73
1:A:189:ASP:OD1	4:A:401:XMB:H19	1.88	0.73
1:A:244:THR:O	1:A:245:ARG:HB2	1.88	0.72
1:A:143:ARG:HD2	1:A:192:GLN:OE1	1.90	0.72
1:A:72:ASN:HD21	1:A:75:GLN:NE2	1.87	0.72
1:A:158:LEU:HD22	1:A:159:GLU:N	2.06	0.71
1:A:87:VAL:HG12	1:A:107:ARG:HB3	1.72	0.70
1:A:50:PHE:N	1:A:50:PHE:CD2	2.60	0.70
1:A:217:GLU:HB2	1:A:224:LYS:HE2	1.74	0.69
1:A:56:ALA:HB2	1:A:103:ILE:H	1.59	0.68
1:A:77:GLU:HG2	1:A:80:GLU:OE1	1.94	0.67
2:L:87:LYS:HG3	2:L:102:GLU:HG2	1.76	0.67
1:A:96:LYS:H	1:A:96:LYS:CD	1.90	0.66
1:A:94:PHE:O	1:A:95:THR:HG23	1.96	0.65
1:A:54:THR:HG23	1:A:55:ALA:O	1.97	0.64
1:A:87:VAL:CG1	1:A:107:ARG:HB3	2.28	0.63
1:A:214:SER:HB3	1:A:227:ILE:O	1.97	0.63
1:A:50:PHE:N	1:A:50:PHE:HD2	1.97	0.62
1:A:107:ARG:HH11	1:A:107:ARG:HG3	1.64	0.62
1:A:131(B):MET:HA	1:A:162:TYR:CZ	2.35	0.61
1:A:35:ASN:O	1:A:38:ASN:N	2.29	0.61
1:A:50:PHE:CZ	1:A:111:PRO:HG3	2.36	0.61
2:L:97:ASP:CG	2:L:123:ALA:HA	2.21	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:190:ALA:HB3	4:A:401:XMB:C19	2.31	0.60
1:A:209:VAL:HG12	1:A:231:VAL:HG21	1.83	0.60
1:A:158:LEU:HD22	1:A:159:GLU:H	1.66	0.59
1:A:65:LYS:HE2	1:A:84:GLU:HG2	1.84	0.59
1:A:131(A):LEU:O	1:A:133:GLN:HB2	2.03	0.59
1:A:231:VAL:HG12	1:A:235:LEU:HD13	1.83	0.59
1:A:45:THR:OG1	1:A:198:PRO:HB3	2.03	0.59
1:A:201:THR:OG1	1:A:210:THR:HG21	2.02	0.59
1:A:56:ALA:HB2	1:A:103:ILE:N	2.17	0.59
1:A:63:ARG:HH11	1:A:63:ARG:HG2	1.67	0.59
1:A:50:PHE:O	1:A:107:ARG:HA	2.03	0.59
1:A:74:GLU:HG3	1:A:153:THR:HG23	1.85	0.59
1:A:41:PHE:O	1:A:193:GLY:HA2	2.03	0.58
1:A:143:ARG:NH1	1:A:148:LYS:HA	2.20	0.57
1:A:218:GLY:HA3	1:A:222:ARG:NE	2.19	0.57
2:L:88:LEU:HD23	2:L:91:LEU:HB2	1.86	0.56
2:L:88:LEU:HD23	2:L:91:LEU:HD23	1.86	0.56
1:A:240:ARG:HG2	1:A:245:ARG:HB3	1.86	0.56
1:A:215:TRP:HA	4:A:401:XMB:C14	2.35	0.56
1:A:27:CYS:SG	1:A:157:MET:CE	2.95	0.55
2:L:89:CYS:O	2:L:93:ASN:HA	2.06	0.55
1:A:27:CYS:SG	1:A:157:MET:HE2	2.46	0.55
1:A:71:ARG:HD3	1:A:117:ASN:HD21	1.71	0.55
1:A:165:ARG:HH22	1:A:178:GLN:N	2.05	0.55
1:A:69:GLY:CA	1:A:118:VAL:HG22	2.37	0.55
1:A:19:GLY:HA2	1:A:158:LEU:HD23	1.89	0.54
1:A:22:CYS:HA	1:A:157:MET:HE2	1.89	0.54
1:A:217:GLU:HB2	1:A:224:LYS:HG3	1.90	0.54
1:A:72:ASN:ND2	1:A:154:ARG:NE	2.53	0.54
1:A:206:THR:HG23	2:L:133:GLY:HA3	1.88	0.53
1:A:217:GLU:O	1:A:222:ARG:HD2	2.08	0.53
1:A:25:GLY:O	1:A:28:PRO:HD3	2.08	0.53
1:A:63:ARG:NH1	1:A:65:LYS:HE3	2.14	0.53
2:L:87:LYS:HG2	2:L:88:LEU:N	2.23	0.53
1:A:211:GLY:CA	1:A:231:VAL:HG23	2.36	0.53
1:A:71:ARG:HD3	1:A:117:ASN:ND2	2.24	0.53
1:A:55:ALA:HB2	1:A:196:GLY:HA2	1.90	0.53
1:A:21:GLU:OE1	1:A:23:LYS:NZ	2.33	0.52
1:A:52:ILE:HD11	1:A:112:ILE:HD11	1.92	0.52
1:A:160:VAL:HG23	1:A:160:VAL:O	2.09	0.52
1:A:202:ARG:HB2	1:A:207:TYR:CE2	2.45	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:74:GLU:OE1	1:A:153:THR:HG21	2.10	0.52
1:A:195:SER:HA	1:A:213:VAL:CG1	2.40	0.52
1:A:127:TRP:HZ2	1:A:203:PHE:CD2	2.27	0.51
1:A:98:THR:O	4:A:401:XMB:H30	2.10	0.51
1:A:44:GLY:N	1:A:197:GLY:HA2	2.25	0.51
1:A:186:LYS:HE2	1:A:188:GLU:OE1	2.11	0.51
1:A:73:THR:OG1	1:A:153:THR:HA	2.11	0.50
1:A:22:CYS:SG	1:A:157:MET:HE2	2.50	0.50
1:A:51:TYR:CZ	1:A:242:MET:HA	2.46	0.50
1:A:215:TRP:NE1	1:A:227:ILE:HG21	2.26	0.50
1:A:22:CYS:O	1:A:71:ARG:NH1	2.44	0.50
1:A:163:VAL:HG21	1:A:225:TYR:CD1	2.46	0.50
1:A:190:ALA:O	1:A:191:CYS:HB2	2.12	0.50
1:A:183:ALA:HB3	1:A:228:TYR:CE1	2.46	0.50
1:A:124:PRO:HA	2:L:99:PHE:CE1	2.47	0.50
1:A:28:PRO:HG2	1:A:29:TRP:CE3	2.47	0.50
1:A:71:ARG:O	1:A:154:ARG:HA	2.12	0.50
1:A:189:ASP:OD1	1:A:190:ALA:N	2.40	0.49
1:A:122:CYS:SG	2:L:132:CYS:C	2.91	0.49
1:A:16:ILE:N	1:A:194:ASP:OD1	2.45	0.49
1:A:65:LYS:CD	1:A:84:GLU:HG2	2.42	0.49
2:L:91:LEU:O	2:L:91:LEU:HD13	2.12	0.49
1:A:122:CYS:SG	2:L:132:CYS:O	2.71	0.49
1:A:51:TYR:HA	1:A:106:LEU:O	2.12	0.49
1:A:72:ASN:HD22	1:A:154:ARG:HE	1.59	0.49
1:A:166:ASN:C	1:A:168:CYS:H	2.16	0.49
1:A:22:CYS:HA	1:A:157:MET:CE	2.43	0.49
1:A:189:ASP:OD1	4:A:401:XMB:C19	2.61	0.49
1:A:22:CYS:SG	1:A:155:LEU:HD23	2.53	0.48
1:A:98:THR:O	1:A:99:TYR:HB2	2.11	0.48
1:A:35:ASN:OD1	1:A:39:GLU:HB2	2.13	0.48
1:A:51:TYR:OH	1:A:107:ARG:HD2	2.13	0.48
1:A:191:CYS:HA	4:A:401:XMB:C11	2.43	0.48
2:L:87:LYS:O	2:L:88:LEU:CB	2.54	0.48
1:A:125:ARG:NH2	1:A:126:ASP:OD2	2.47	0.48
1:A:99:TYR:OH	4:A:401:XMB:H92	2.14	0.48
1:A:51:TYR:CE1	1:A:107:ARG:HB2	2.49	0.48
1:A:69:GLY:HA3	1:A:118:VAL:HG22	1.96	0.47
1:A:135:THR:HB	1:A:159:GLU:HG2	1.97	0.47
1:A:195:SER:HA	1:A:213:VAL:HG12	1.95	0.47
1:A:165:ARG:HH22	1:A:177:THR:C	2.18	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:26:GLU:O	1:A:28:PRO:HD2	2.15	0.47
1:A:141:PHE:CE2	1:A:155:LEU:HD13	2.50	0.46
1:A:127:TRP:HZ2	1:A:203:PHE:CE2	2.33	0.46
1:A:55:ALA:H	1:A:196:GLY:HA2	1.80	0.46
1:A:227:ILE:N	1:A:227:ILE:HD13	2.31	0.46
1:A:192:GLN:N	4:A:401:XMB:O28	2.49	0.46
1:A:76:GLU:HG3	1:A:76:GLU:O	2.16	0.46
1:A:116:MET:SD	2:L:136:THR:O	2.74	0.46
1:A:16:ILE:HB	1:A:156:LYS:HB2	1.98	0.46
1:A:63:ARG:NH1	1:A:63:ARG:HG2	2.29	0.46
1:A:151:GLN:HE22	1:A:193:GLY:HA3	1.81	0.46
1:A:50:PHE:H	1:A:50:PHE:HD2	1.64	0.45
1:A:60:TYR:OH	1:A:90:LYS:HE2	2.17	0.45
1:A:77:GLU:HG2	1:A:80:GLU:CD	2.37	0.45
1:A:34:ILE:HG21	1:A:67:ARG:HH21	1.81	0.45
1:A:209:VAL:HG12	1:A:209:VAL:O	2.18	0.44
1:A:20:GLN:HA	1:A:20:GLN:OE1	2.17	0.44
1:A:242:MET:C	1:A:244:THR:H	2.20	0.44
1:A:110:THR:HA	1:A:111:PRO:HD3	1.80	0.44
1:A:45:THR:HB	1:A:53:LEU:HB3	2.00	0.44
1:A:27:CYS:SG	1:A:157:MET:HE1	2.57	0.44
1:A:206:THR:CG2	2:L:133:GLY:HA3	2.47	0.44
1:A:100:ASP:O	1:A:101:PHE:HB2	2.18	0.43
1:A:24:ASP:OD2	1:A:117:ASN:OD1	2.36	0.43
1:A:123:LEU:HD21	1:A:238:ILE:HG21	2.00	0.43
1:A:206:THR:HA	2:L:133:GLY:O	2.18	0.43
1:A:133:GLN:O	1:A:162:TYR:HD2	2.02	0.43
1:A:67:ARG:NH1	1:A:70:ASP:OD2	2.46	0.43
1:A:245:ARG:O	1:A:245:ARG:HD3	2.17	0.43
1:A:151:GLN:HG3	1:A:152:SER:H	1.83	0.43
1:A:57:HIS:CD2	1:A:58:CYS:N	2.87	0.43
1:A:185(A):ASP:O	1:A:223:LYS:HA	2.18	0.43
1:A:72:ASN:HD21	1:A:75:GLN:CD	2.22	0.43
1:A:116:MET:O	1:A:117:ASN:HB2	2.18	0.43
1:A:158:LEU:CD1	1:A:188:GLU:HB3	2.36	0.43
1:A:125:ARG:CA	1:A:235:LEU:HD23	2.49	0.43
1:A:56:ALA:HB3	1:A:102:ASP:HA	2.00	0.42
1:A:141:PHE:CZ	1:A:155:LEU:HB2	2.54	0.42
1:A:182:CYS:O	1:A:183:ALA:HB2	2.19	0.42
1:A:30:GLN:HG2	1:A:155:LEU:CD1	2.49	0.42
1:A:212:ILE:HB	1:A:229:THR:HB	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:88:VAL:O	1:A:88:VAL:HG13	2.19	0.42
1:A:139:SER:HB2	1:A:156:LYS:O	2.19	0.42
1:A:150:ARG:HG2	1:A:150:ARG:H	1.64	0.42
1:A:93:ARG:NE	1:A:101:PHE:CZ	2.87	0.42
2:L:91:LEU:C	2:L:91:LEU:HD13	2.40	0.42
1:A:92:ASN:C	1:A:94:PHE:H	2.22	0.42
1:A:123:LEU:HD12	1:A:123:LEU:HA	1.73	0.42
1:A:163:VAL:HB	1:A:182:CYS:HB3	2.02	0.42
1:A:234:PHE:O	1:A:237:TRP:HB3	2.19	0.42
1:A:65:LYS:CE	1:A:84:GLU:HG2	2.48	0.42
1:A:16:ILE:HG23	1:A:158:LEU:HB2	2.01	0.42
1:A:217:GLU:CD	1:A:224:LYS:HD3	2.38	0.42
1:A:136:GLY:O	1:A:159:GLU:HA	2.20	0.41
1:A:157:MET:HG3	1:A:157:MET:O	2.21	0.41
1:A:100:ASP:OD2	1:A:177:THR:HG21	2.20	0.41
1:A:151:GLN:HG3	1:A:152:SER:N	2.36	0.41
1:A:127:TRP:CZ2	1:A:203:PHE:CD2	3.07	0.41
1:A:20:GLN:CA	1:A:20:GLN:OE1	2.69	0.41
2:L:88:LEU:HB3	2:L:91:LEU:HB2	2.02	0.41
2:L:93:ASN:O	2:L:100:CYS:SG	2.78	0.41
2:L:103:GLU:O	2:L:105:ASN:N	2.53	0.41
1:A:123:LEU:HD12	1:A:209:VAL:HB	2.03	0.41
1:A:72:ASN:OD1	1:A:75:GLN:OE1	2.38	0.41
1:A:72:ASN:ND2	1:A:75:GLN:NE2	2.63	0.41
1:A:141:PHE:O	1:A:151:GLN:HG3	2.19	0.41
1:A:215:TRP:CE2	1:A:227:ILE:HG21	2.55	0.41
1:A:177:THR:C	1:A:179:ASN:H	2.24	0.41
1:A:164:ASP:HB3	1:A:167:SER:OG	2.20	0.41
1:A:143:ARG:HA	1:A:150:ARG:O	2.21	0.41
1:A:57:HIS:CE1	1:A:195:SER:OG	2.70	0.40
1:A:241:SER:O	1:A:244:THR:HA	2.21	0.40
1:A:55:ALA:N	1:A:196:GLY:HA2	2.36	0.40
1:A:16:ILE:HD11	1:A:139:SER:HA	2.04	0.40
1:A:86:GLU:HB2	1:A:109:LYS:HA	2.02	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	233/235 (99%)	186 (80%)	39 (17%)	8 (3%)	4	18
2	L	49/96 (51%)	37 (76%)	8 (16%)	4 (8%)	1	2
All	All	282/331 (85%)	223 (79%)	47 (17%)	12 (4%)	3	12

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	L	88	LEU
1	A	244	THR
1	A	183	ALA
2	L	92	ASP
2	L	105	ASN
1	A	19	GLY
1	A	77	GLU
1	A	93	ARG
2	L	118	ALA
1	A	78	GLY
1	A	118	VAL
1	A	120	PRO

5.3.2 Protein sidechains

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	200/200 (100%)	181 (90%)	19 (10%)	10	29

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	L	43/82 (52%)	36 (84%)	7 (16%)	3	8
All	All	243/282 (86%)	217 (89%)	26 (11%)	8	23

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	50	PHE
1	A	61	GLN
1	A	66	VAL
1	A	73	THR
1	A	96	LYS
1	A	123	LEU
1	A	131	THR
1	A	131(A)	LEU
1	A	131(B)	MET
1	A	134	LYS
1	A	144	THR
1	A	145	HIS
1	A	158	LEU
1	A	178	GLN
1	A	198	PRO
1	A	205	ASP
1	A	210	THR
1	A	235	LEU
1	A	236	LYS
2	L	97	ASP
2	L	101	HIS
2	L	104	GLN
2	L	105	ASN
2	L	107	VAL
2	L	113	ARG
2	L	137	LEU

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

Mol	Chain	Res	Type
1	A	61	GLN
1	A	72	ASN
1	A	117	ASN
1	A	133	GLN
1	A	151	GLN

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Mol	Chain	Res	Type
2	L	104	GLN
2	L	135	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 2 ligands modelled in this entry, 1 is 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	XMB	A	401	-	44,45,45	3.94	15 (34%)	56,65,65	2.34	10 (17%)

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	XMB	A	401	-	-	0/29/51/51	0/4/5/5

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	401	XMB	C25-N27	2.06	1.50	1.46
4	A	401	XMB	C9-N3	2.15	1.50	1.47
4	A	401	XMB	C18-C20	2.20	1.40	1.36
4	A	401	XMB	C1-N3	2.39	1.38	1.34
4	A	401	XMB	C4-N3	2.56	1.51	1.47
4	A	401	XMB	C36-N31	3.02	1.51	1.46
4	A	401	XMB	C34-N31	3.23	1.51	1.46
4	A	401	XMB	C10-S6	3.47	1.79	1.76
4	A	401	XMB	C32-N31	3.97	1.49	1.38
4	A	401	XMB	C5-S6	6.39	1.81	1.76
4	A	401	XMB	S22-N24	6.45	1.70	1.61
4	A	401	XMB	O7-S6	6.49	1.57	1.44
4	A	401	XMB	O8-S6	6.76	1.58	1.44
4	A	401	XMB	O23-S22	13.86	1.58	1.43
4	A	401	XMB	O28-S22	14.50	1.59	1.43

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	401	XMB	O8-S6-O7	-8.92	109.08	117.83
4	A	401	XMB	O28-S22-O23	-8.26	108.98	119.55
4	A	401	XMB	O2-C1-C25	-3.24	113.73	119.65
4	A	401	XMB	O28-S22-C12	2.50	111.04	107.95
4	A	401	XMB	O23-S22-N24	2.57	111.18	107.03
4	A	401	XMB	O23-S22-C12	2.77	111.38	107.95
4	A	401	XMB	O8-S6-C10	3.88	113.72	109.24
4	A	401	XMB	O7-S6-C5	4.31	114.21	109.24
4	A	401	XMB	C38-C34-N31	5.04	121.05	111.22
4	A	401	XMB	C41-C36-N31	5.09	121.16	111.22

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	401	XMB	9	0

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.