



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:12 PM UTC

PDB ID : 5I25 / pdb_00005i25
Title : human recombinant coagulation FXI in complex with a peptide derived from human high molecular weight kininogen (HKP)
Authors : Hall, G.A.F.; Wong, S.S.; Emsley, J.
Deposited on : 2016-02-08
Resolution : 2.85 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 ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

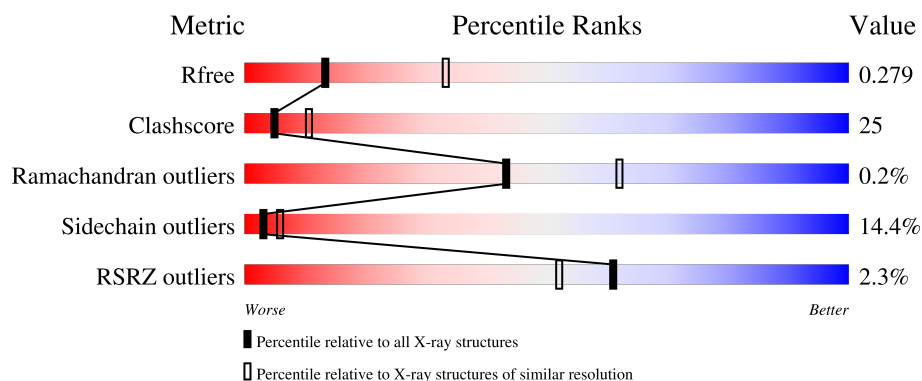
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.85 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}	180053	1407 (2.88-2.84)
Clashscore	190562	1446 (2.88-2.84)
Ramachandran outliers	187476	1406 (2.88-2.84)
Sidechain outliers	187428	1407 (2.88-2.84)
RSRZ outliers	180081	1408 (2.88-2.84)

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 $\leq 5\%$. 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.

Mol	Chain	Length	Quality of chain
1	A	607	2% 52% 37% 9% .
2	B	8	50% 25% 25%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4842 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 XI.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	596	Total	C	N	O	S	0	0	0
			4684	2950	815	878	41			

- Molecule 2 is a protein called ASN-PRO-ILE-SER-ASP-PHE-PRO-ASP.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	8	Total	C	N	O	0	0	0
			63	40	9	14			

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



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

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

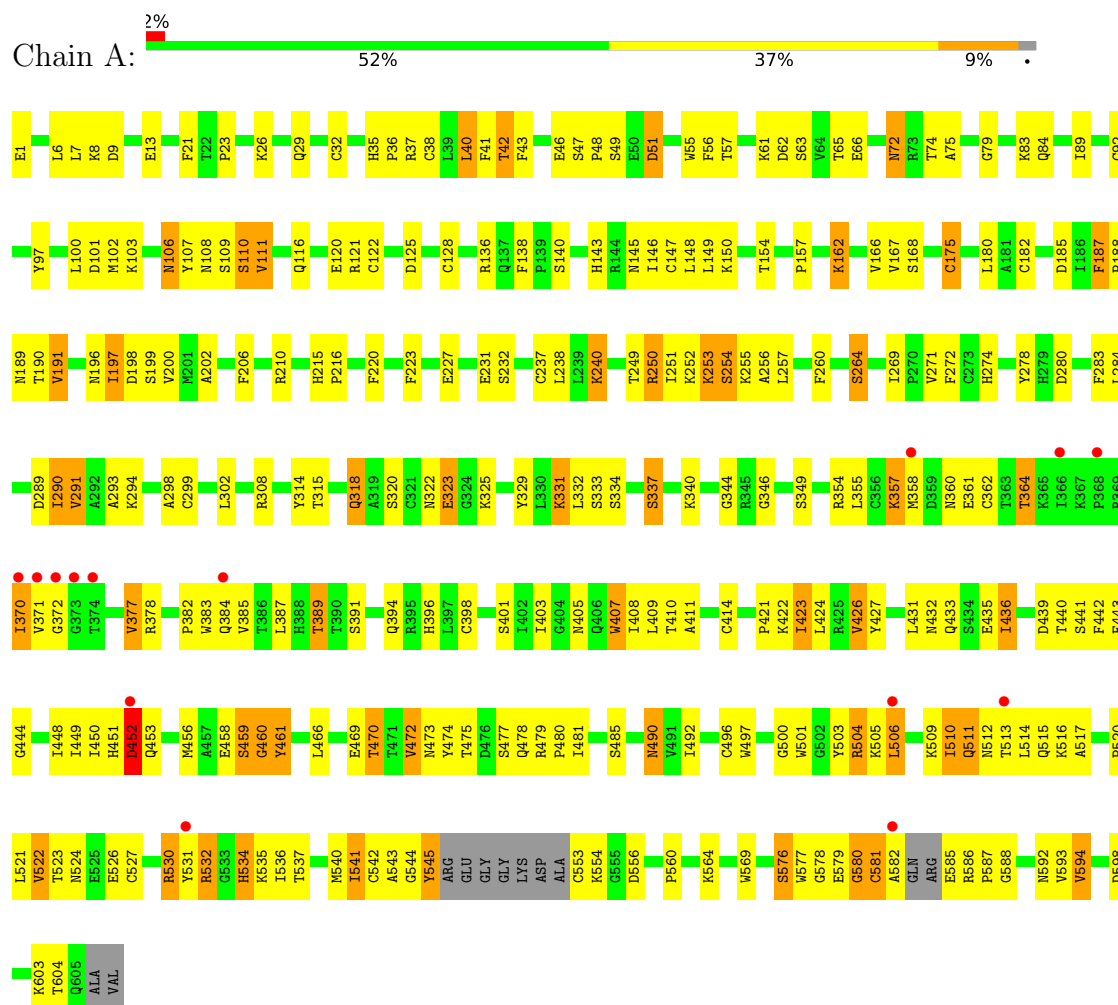
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	51	Total	O	0	0
			51	51		
4	B	2	Total	O	0	0
			2	2		

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: Coagulation factor XI



• Molecule 2: ASN-PRO-ILE-SER-ASP-PHE-PRO-ASP



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	80.59Å 80.59Å 251.81Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.79 – 2.85 39.79 – 2.85	Depositor EDS
% Data completeness (in resolution range)	99.6 (39.79-2.85) 99.6 (39.79-2.85)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.48 (at 2.85Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.226 , 0.288 0.219 , 0.279	Depositor DCC
R_{free} test set	2057 reflections (10.15%)	wwPDB-VP
Wilson B-factor (Å ²)	73.1	Xtriage
Anisotropy	0.031	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 55.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4842	wwPDB-VP
Average B, all atoms (Å ²)	74.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.37% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ 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:
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	0.69	1/4793 (0.0%)	1.06	12/6490 (0.2%)
2	B	0.67	0/65	1.51	2/89 (2.2%)
All	All	0.69	1/4858 (0.0%)	1.07	14/6579 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	452	ASP	CG-OD2	5.01	1.34	1.25

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	460	GLY	N-CA-C	10.83	126.64	112.77
1	A	459	SER	N-CA-C	-8.40	99.21	110.55
1	A	337	SER	CA-C-N	7.79	127.84	119.90
1	A	337	SER	C-N-CA	7.79	127.84	119.90
2	B	2	ASN	CA-C-N	7.67	127.30	119.56
2	B	2	ASN	C-N-CA	7.67	127.30	119.56
1	A	202	ALA	CA-C-N	6.63	126.58	119.28
1	A	202	ALA	C-N-CA	6.63	126.58	119.28
1	A	407	TRP	N-CA-C	6.58	119.63	108.90
1	A	581	CYS	N-CA-C	-6.50	100.68	110.24
1	A	421	PRO	N-CA-C	-5.56	106.61	113.84
1	A	580	GLY	N-CA-C	-5.15	98.36	113.30
1	A	398	CYS	N-CA-C	5.09	116.41	109.18
1	A	291	VAL	N-CA-C	5.06	117.11	108.86

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	4684	0	4571	236	0
2	B	63	0	52	7	0
3	A	42	0	39	3	0
4	A	51	0	0	2	0
4	B	2	0	0	2	0
All	All	4842	0	4662	237	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 25.

All (237) 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:452:ASP:OD1	1:A:453:GLN:HG2	1.53	1.09
1:A:541:ILE:HD11	1:A:592:ASN:CG	1.78	1.08
1:A:579:GLU:HG3	1:A:579:GLU:O	1.42	1.05
1:A:9:ASP:OD1	1:A:74:THR:HA	1.60	1.01
1:A:189:ASN:HD22	1:A:254:SER:HA	1.37	0.90
1:A:377:VAL:HG22	1:A:378:ARG:H	1.45	0.81
1:A:582:ALA:C	1:A:585:GLU:HA	2.07	0.80
1:A:189:ASN:ND2	1:A:254:SER:HA	1.97	0.80
1:A:407:TRP:CE2	1:A:604:THR:HG22	2.17	0.79
1:A:433:GLN:O	1:A:436:ILE:HG13	1.82	0.77
1:A:83:LYS:HB3	1:A:354:ARG:HD3	1.68	0.76
1:A:474:TYR:CZ	1:A:480:PRO:HG3	2.21	0.76
1:A:187:PHE:N	1:A:187:PHE:HD2	1.83	0.76
1:A:384:GLN:HE22	1:A:560:PRO:HD3	1.51	0.76
1:A:472:VAL:HG12	1:A:478:GLN:NE2	2.02	0.75
1:A:187:PHE:N	1:A:187:PHE:CD2	2.50	0.75
1:A:405:ASN:O	1:A:472:VAL:HG23	1.86	0.75
1:A:511:GLN:HE21	1:A:512:ASN:H	1.36	0.74
1:A:354:ARG:NH2	4:A:1001:HOH:O	2.20	0.74
1:A:452:ASP:CG	1:A:453:GLN:HG2	2.12	0.74
1:A:505:LYS:HD2	1:A:506:LEU:H	1.50	0.74
1:A:522:VAL:HG13	1:A:543:ALA:HA	1.70	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:283:PHE:O	1:A:331:LYS:HE3	1.87	0.73
1:A:461:TYR:CD1	1:A:461:TYR:N	2.52	0.72
1:A:72:ASN:ND2	3:A:901:NAG:O5	2.23	0.72
1:A:580:GLY:O	1:A:581:CYS:C	2.32	0.72
1:A:40:LEU:HD12	1:A:61:LYS:HB2	1.70	0.71
1:A:432:ASN:ND2	3:A:903:NAG:O5	2.24	0.71
1:A:92:CYS:SG	1:A:175:CYS:N	2.64	0.70
1:A:293:ALA:HB3	1:A:314:TYR:HE1	1.57	0.70
1:A:473:ASN:O	1:A:478:GLN:NE2	2.23	0.70
1:A:196:ASN:OD1	1:A:238:LEU:HD13	1.91	0.69
1:A:384:GLN:HE21	1:A:384:GLN:HA	1.57	0.69
1:A:541:ILE:HD11	1:A:592:ASN:CB	2.22	0.68
1:A:371:VAL:O	1:A:371:VAL:HG23	1.93	0.68
1:A:384:GLN:NE2	1:A:560:PRO:HG3	2.09	0.68
1:A:227:GLU:O	1:A:255:LYS:HD3	1.93	0.67
1:A:387:LEU:HD13	1:A:426:VAL:HG13	1.76	0.67
1:A:511:GLN:HE21	1:A:512:ASN:N	1.91	0.67
1:A:540:MET:C	1:A:541:ILE:HD13	2.18	0.67
1:A:389:THR:O	1:A:394:GLN:HA	1.93	0.67
1:A:469:GLU:HB2	1:A:470:THR:HG22	1.77	0.66
1:A:250:ARG:O	1:A:250:ARG:HG2	1.95	0.66
1:A:197:ILE:O	1:A:198:ASP:HB2	1.95	0.65
1:A:122:CYS:O	1:A:128:CYS:HB3	1.96	0.65
1:A:503:TYR:HE1	1:A:506:LEU:HA	1.61	0.65
1:A:51:ASP:OD2	1:A:51:ASP:N	2.20	0.64
1:A:322:ASN:OD1	1:A:323:GLU:HG3	1.98	0.64
1:A:510:ILE:HG23	1:A:511:GLN:N	2.12	0.64
1:A:370:ILE:HD12	1:A:371:VAL:H	1.62	0.64
1:A:360:ASN:ND2	1:A:362:CYS:SG	2.71	0.64
1:A:401:SER:OG	1:A:560:PRO:HB3	1.99	0.63
1:A:531:TYR:OH	1:A:587:PRO:O	2.17	0.63
1:A:461:TYR:H	1:A:461:TYR:HD1	1.45	0.62
1:A:280:ASP:OD1	1:A:344:GLY:HA2	1.99	0.62
1:A:424:LEU:CD1	1:A:448:ILE:HD11	2.29	0.62
1:A:514:LEU:HD12	1:A:515:GLN:N	2.14	0.62
1:A:541:ILE:HD11	1:A:592:ASN:ND2	2.15	0.62
1:A:456:MET:O	1:A:459:SER:HB2	2.00	0.62
1:A:198:ASP:OD1	1:A:199:SER:N	2.32	0.61
1:A:384:GLN:NE2	1:A:560:PRO:HD3	2.15	0.61
1:A:443:PHE:CZ	1:A:472:VAL:HG13	2.35	0.61
1:A:385:VAL:HG21	1:A:408:ILE:HD12	1.81	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:564:LYS:HB2	1:A:569:TRP:CZ3	2.35	0.61
1:A:424:LEU:HD11	1:A:448:ILE:HD11	1.82	0.61
1:A:436:ILE:HG22	1:A:440:THR:HG21	1.83	0.60
1:A:384:GLN:HA	1:A:384:GLN:NE2	2.15	0.60
1:A:188:PRO:HA	1:A:257:LEU:HG	1.83	0.60
1:A:187:PHE:HD2	1:A:187:PHE:H	1.50	0.60
1:A:576:SER:OG	1:A:577:TRP:CD1	2.54	0.60
1:A:405:ASN:O	1:A:472:VAL:CG2	2.49	0.60
1:A:182:CYS:SG	1:A:264:SER:HB3	2.42	0.59
1:A:372:GLY:O	1:A:516:LYS:CE	2.50	0.59
1:A:299:CYS:SG	1:A:314:TYR:HB2	2.43	0.59
1:A:576:SER:OG	1:A:577:TRP:HD1	1.85	0.59
1:A:532:ARG:HH22	1:A:586:ARG:NH2	2.00	0.58
1:A:23:PRO:HB2	1:A:97:TYR:CE2	2.38	0.58
1:A:116:GLN:O	1:A:120:GLU:HG3	2.03	0.58
1:A:407:TRP:CD2	1:A:604:THR:HG22	2.38	0.58
1:A:377:VAL:HA	1:A:431:LEU:HD13	1.85	0.58
1:A:444:GLY:HA3	1:A:469:GLU:HG3	1.85	0.58
1:A:579:GLU:O	1:A:579:GLU:CG	2.30	0.58
1:A:206:PHE:O	1:A:210:ARG:HG3	2.05	0.57
1:A:492:ILE:N	1:A:492:ILE:HD12	2.19	0.57
1:A:196:ASN:OD1	1:A:240:LYS:HD2	2.04	0.57
1:A:102:MET:HE2	1:A:168:SER:HB3	1.87	0.57
1:A:564:LYS:HB2	1:A:569:TRP:CH2	2.40	0.57
1:A:511:GLN:NE2	1:A:512:ASN:H	2.00	0.57
1:A:461:TYR:N	1:A:461:TYR:HD1	1.97	0.56
1:A:522:VAL:HG22	1:A:542:CYS:HB2	1.86	0.56
1:A:318:GLN:HA	1:A:318:GLN:NE2	2.19	0.56
1:A:6:LEU:C	1:A:7:LEU:HD23	2.30	0.56
1:A:272:PHE:HB3	1:A:355:LEU:HB3	1.87	0.56
1:A:427:TYR:CE2	1:A:442:PHE:HB3	2.40	0.56
1:A:503:TYR:CE1	1:A:506:LEU:HA	2.40	0.55
1:A:511:GLN:HE21	1:A:511:GLN:CA	2.20	0.55
1:A:517:ALA:HB2	1:A:554:LYS:HE2	1.89	0.55
1:A:534:HIS:N	1:A:534:HIS:ND1	2.53	0.55
1:A:294:LYS:HE2	1:A:325:LYS:HE2	1.88	0.55
1:A:432:ASN:O	1:A:435:GLU:HB3	2.06	0.55
1:A:43:PHE:CE2	1:A:56:PHE:HA	2.42	0.55
1:A:293:ALA:HB3	1:A:314:TYR:CE1	2.41	0.54
1:A:40:LEU:CD1	1:A:61:LYS:HB2	2.37	0.54
1:A:535:LYS:HG3	1:A:536:ILE:N	2.22	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:531:TYR:HD2	1:A:536:ILE:HD11	1.72	0.53
1:A:108:ASN:HD21	3:A:902:NAG:C5	2.21	0.53
1:A:505:LYS:CD	1:A:506:LEU:H	2.20	0.53
1:A:540:MET:O	1:A:541:ILE:HD13	2.08	0.53
1:A:377:VAL:HG22	1:A:378:ARG:N	2.18	0.53
1:A:106:ASN:HB2	2:B:4:ILE:HB	1.90	0.52
1:A:426:VAL:HG23	1:A:443:PHE:O	2.09	0.52
1:A:148:LEU:HG	1:A:150:LYS:HE2	1.92	0.52
1:A:143:HIS:CD2	2:B:8:PRO:CD	2.93	0.52
1:A:396:HIS:HB3	1:A:510:ILE:HD11	1.91	0.52
1:A:46:GLU:OE1	1:A:162:LYS:HE3	2.09	0.52
1:A:298:ALA:O	1:A:302:LEU:HD23	2.09	0.52
1:A:110:SER:HB2	1:A:121:ARG:NH1	2.24	0.52
1:A:191:VAL:HG22	1:A:251:ILE:HB	1.92	0.51
1:A:426:VAL:HG23	1:A:443:PHE:HB2	1.92	0.51
1:A:42:THR:HG22	1:A:42:THR:O	2.09	0.51
1:A:223:PHE:O	1:A:256:ALA:HA	2.10	0.51
1:A:252:LYS:O	1:A:253:LYS:HG2	2.10	0.51
1:A:501:TRP:O	1:A:511:GLN:HB2	2.10	0.51
1:A:514:LEU:HD12	1:A:515:GLN:H	1.74	0.51
1:A:21:PHE:O	1:A:157:PRO:HD2	2.11	0.51
1:A:106:ASN:ND2	4:B:101:HOH:O	2.41	0.51
1:A:532:ARG:NH2	1:A:586:ARG:NH2	2.58	0.51
1:A:541:ILE:HD11	1:A:592:ASN:HB2	1.93	0.51
1:A:149:LEU:C	1:A:150:LYS:HG2	2.36	0.51
1:A:490:ASN:N	1:A:490:ASN:OD1	2.43	0.50
1:A:284:LEU:HD12	1:A:340:LYS:HB3	1.93	0.50
1:A:382:PRO:HB2	1:A:479:ARG:H	1.77	0.50
1:A:504:ARG:CG	1:A:504:ARG:HH11	2.25	0.50
1:A:46:GLU:HG3	1:A:55:TRP:CZ3	2.47	0.50
1:A:46:GLU:HA	1:A:55:TRP:CD1	2.47	0.49
1:A:29:GLN:HG3	1:A:41:PHE:CE2	2.47	0.49
1:A:382:PRO:HB2	1:A:479:ARG:N	2.27	0.49
1:A:537:THR:OG1	1:A:540:MET:HE3	2.13	0.49
1:A:101:ASP:OD1	1:A:103:LYS:HE2	2.13	0.49
1:A:143:HIS:NE2	2:B:8:PRO:HD3	2.28	0.49
1:A:136:ARG:C	1:A:138:PHE:N	2.69	0.49
1:A:372:GLY:O	1:A:516:LYS:NZ	2.46	0.49
1:A:577:TRP:O	1:A:577:TRP:CD2	2.66	0.49
2:B:6:ASP:OD2	4:B:101:HOH:O	2.20	0.48
1:A:200:VAL:HG22	1:A:237:CYS:HB3	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:449:ILE:HG21	1:A:603:LYS:HD2	1.94	0.48
1:A:451:HIS:HD2	1:A:453:GLN:H	1.60	0.48
1:A:35:HIS:CG	1:A:36:PRO:HD2	2.48	0.48
1:A:384:GLN:NE2	1:A:560:PRO:CD	2.76	0.48
1:A:47:SER:HB2	1:A:48:PRO:HD2	1.96	0.48
1:A:432:ASN:OD1	1:A:513:THR:HA	2.14	0.48
1:A:451:HIS:CD2	1:A:453:GLN:H	2.32	0.47
1:A:423:ILE:O	1:A:423:ILE:HG13	2.13	0.47
1:A:475:THR:O	1:A:478:GLN:N	2.47	0.47
1:A:136:ARG:O	1:A:138:PHE:N	2.47	0.47
1:A:384:GLN:NE2	1:A:560:PRO:CG	2.77	0.47
1:A:107:TYR:OH	1:A:125:ASP:OD1	2.29	0.47
1:A:143:HIS:CD2	2:B:8:PRO:HD3	2.50	0.47
1:A:332:LEU:C	1:A:332:LEU:HD12	2.39	0.47
1:A:475:THR:C	1:A:477:SER:N	2.71	0.47
1:A:408:ILE:HB	1:A:466:LEU:HB2	1.96	0.47
1:A:355:LEU:HD23	1:A:358:MET:HE3	1.96	0.47
1:A:410:THR:OG1	1:A:411:ALA:N	2.48	0.47
1:A:530:ARG:NE	1:A:545:TYR:OH	2.48	0.47
1:A:106:ASN:OD1	1:A:106:ASN:N	2.48	0.46
1:A:9:ASP:OD1	1:A:74:THR:CA	2.49	0.46
1:A:231:GLU:CD	1:A:231:GLU:C	2.84	0.46
1:A:269:ILE:HG13	1:A:271:VAL:HG23	1.96	0.46
1:A:148:LEU:H	1:A:148:LEU:HD23	1.78	0.46
1:A:357:LYS:HA	1:A:357:LYS:HD3	1.38	0.46
1:A:154:THR:O	1:A:216:PRO:HG2	2.16	0.46
1:A:280:ASP:CG	1:A:344:GLY:HA2	2.41	0.46
1:A:387:LEU:CD1	1:A:426:VAL:HG13	2.42	0.46
1:A:62:ASP:CG	1:A:308:ARG:HH21	2.24	0.46
1:A:109:SER:HB2	2:B:4:ILE:CG2	2.45	0.46
1:A:8:LYS:O	1:A:9:ASP:C	2.59	0.46
1:A:46:GLU:HA	1:A:55:TRP:CG	2.51	0.46
1:A:65:THR:C	1:A:66:GLU:HG3	2.40	0.46
1:A:29:GLN:HG3	1:A:41:PHE:HE2	1.81	0.45
1:A:200:VAL:CG2	1:A:237:CYS:HB3	2.45	0.45
1:A:460:GLY:HA2	1:A:540:MET:HE2	1.97	0.45
1:A:1:GLU:OE2	1:A:1:GLU:HA	2.16	0.45
1:A:377:VAL:HA	1:A:431:LEU:CD1	2.46	0.45
1:A:210:ARG:NH2	4:A:1003:HOH:O	2.49	0.45
1:A:41:PHE:CE2	1:A:79:GLY:HA3	2.51	0.45
1:A:289:ASP:OD1	1:A:290:ILE:N	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:110:SER:O	1:A:146:ILE:HA	2.17	0.44
1:A:100:LEU:HD12	1:A:100:LEU:HA	1.80	0.44
1:A:290:ILE:HD13	1:A:329:TYR:CE1	2.53	0.44
1:A:383:TRP:CE2	1:A:481:ILE:HD12	2.52	0.44
1:A:496:CYS:C	1:A:497:TRP:CD1	2.96	0.44
1:A:422:LYS:HA	1:A:422:LYS:HD3	1.79	0.44
1:A:396:HIS:HB3	1:A:510:ILE:CD1	2.48	0.43
1:A:403:ILE:HD13	1:A:409:LEU:HB2	1.99	0.43
1:A:111:VAL:HA	1:A:145:ASN:O	2.18	0.43
1:A:278:TYR:N	1:A:349:SER:O	2.48	0.43
1:A:577:TRP:CD1	1:A:577:TRP:N	2.86	0.43
1:A:361:GLU:O	1:A:364:THR:HG23	2.19	0.43
1:A:6:LEU:HD23	1:A:6:LEU:N	2.33	0.43
1:A:577:TRP:O	1:A:577:TRP:CG	2.70	0.43
1:A:32:CYS:O	1:A:38:CYS:HB3	2.18	0.43
1:A:215:HIS:HA	1:A:216:PRO:HD3	1.84	0.43
1:A:436:ILE:O	1:A:436:ILE:HD12	2.19	0.43
1:A:504:ARG:CG	1:A:504:ARG:NH1	2.82	0.43
1:A:505:LYS:HD2	1:A:506:LEU:N	2.26	0.42
1:A:520:PRO:HD2	1:A:544:GLY:O	2.19	0.42
1:A:206:PHE:N	1:A:206:PHE:CD1	2.86	0.42
1:A:46:GLU:HG3	1:A:55:TRP:CE3	2.54	0.42
1:A:74:THR:O	1:A:75:ALA:HB3	2.19	0.42
1:A:136:ARG:C	1:A:138:PHE:H	2.26	0.42
1:A:109:SER:HB2	2:B:4:ILE:HG23	2.01	0.42
1:A:500:GLY:H	1:A:514:LEU:HD11	1.84	0.42
1:A:475:THR:O	1:A:477:SER:N	2.52	0.42
1:A:424:LEU:HD12	1:A:448:ILE:HD11	2.01	0.42
1:A:544:GLY:O	1:A:545:TYR:HB2	2.20	0.42
1:A:21:PHE:HA	1:A:56:PHE:O	2.19	0.42
1:A:444:GLY:O	1:A:469:GLU:HG2	2.20	0.42
1:A:56:PHE:CZ	1:A:100:LEU:HD11	2.54	0.41
1:A:511:GLN:NE2	1:A:511:GLN:CA	2.83	0.41
1:A:578:GLY:HA3	1:A:588:GLY:HA2	2.01	0.41
1:A:191:VAL:CG1	1:A:253:LYS:HB2	2.50	0.41
1:A:220:PHE:HA	1:A:260:PHE:HA	2.02	0.41
1:A:280:ASP:OD2	1:A:346:GLY:HA2	2.21	0.41
1:A:231:GLU:OE2	1:A:232:SER:HA	2.20	0.41
1:A:474:TYR:C	1:A:475:THR:CG2	2.93	0.41
1:A:377:VAL:CG2	1:A:378:ARG:H	2.23	0.41
1:A:411:ALA:HB3	1:A:414:CYS:SG	2.61	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:458:GLU:HG3	1:A:577:TRP:CD2	2.56	0.41
1:A:524:ASN:O	1:A:527:CYS:N	2.54	0.41
1:A:517:ALA:HB2	1:A:554:LYS:CE	2.49	0.41
1:A:371:VAL:O	1:A:371:VAL:CG2	2.65	0.40
1:A:185:ASP:O	1:A:187:PHE:HE2	2.04	0.40
1:A:72:ASN:OD1	1:A:72:ASN:N	2.49	0.40
1:A:210:ARG:HD2	1:A:274:HIS:CE1	2.56	0.40
1:A:333:SER:OG	1:A:334:SER:N	2.53	0.40
1:A:509:LYS:HE2	1:A:509:LYS:HB2	1.83	0.40
1:A:593:VAL:O	1:A:594:VAL:C	2.65	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	590/607 (97%)	567 (96%)	22 (4%)	1 (0%)	43	62
2	B	6/8 (75%)	6 (100%)	0	0	100	100
All	All	596/615 (97%)	573 (96%)	22 (4%)	1 (0%)	43	62

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	377	VAL

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	528/535 (99%)	453 (86%)	75 (14%)	3	6
2	B	8/8 (100%)	6 (75%)	2 (25%)	0	0
All	All	536/543 (99%)	459 (86%)	77 (14%)	3	6

All (77) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	GLU
1	A	26	LYS
1	A	37	ARG
1	A	40	LEU
1	A	42	THR
1	A	49	SER
1	A	51	ASP
1	A	57	THR
1	A	63	SER
1	A	72	ASN
1	A	84	GLN
1	A	89	ILE
1	A	106	ASN
1	A	110	SER
1	A	111	VAL
1	A	140	SER
1	A	147	CYS
1	A	162	LYS
1	A	166	VAL
1	A	167	VAL
1	A	175	CYS
1	A	180	LEU
1	A	187	PHE
1	A	190	THR
1	A	191	VAL
1	A	197	ILE
1	A	240	LYS
1	A	249	THR
1	A	250	ARG
1	A	253	LYS
1	A	254	SER
1	A	264	SER
1	A	290	ILE

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Mol	Chain	Res	Type
1	A	291	VAL
1	A	315	THR
1	A	318	GLN
1	A	320	SER
1	A	323	GLU
1	A	331	LYS
1	A	337	SER
1	A	357	LYS
1	A	364	THR
1	A	370	ILE
1	A	389	THR
1	A	391	SER
1	A	423	ILE
1	A	426	VAL
1	A	436	ILE
1	A	439	ASP
1	A	441	SER
1	A	450	ILE
1	A	452	ASP
1	A	461	TYR
1	A	470	THR
1	A	472	VAL
1	A	485	SER
1	A	490	ASN
1	A	504	ARG
1	A	506	LEU
1	A	510	ILE
1	A	511	GLN
1	A	521	LEU
1	A	522	VAL
1	A	523	THR
1	A	526	GLU
1	A	530	ARG
1	A	532	ARG
1	A	534	HIS
1	A	541	ILE
1	A	545	TYR
1	A	553	CYS
1	A	556	ASP
1	A	576	SER
1	A	594	VAL
1	A	598	ASP

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Mol	Chain	Res	Type
2	B	2	ASN
2	B	6	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	143	HIS
1	A	189	ASN
1	A	235	ASN
1	A	274	HIS
1	A	318	GLN
1	A	384	GLN
1	A	451	HIS
1	A	511	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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

3 ligands 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
3	NAG	A	902	-	14,14,15	0.81	0	17,19,21	1.87	5 (29%)
3	NAG	A	903	-	14,14,15	0.51	0	17,19,21	0.61	0
3	NAG	A	901	-	14,14,15	0.56	0	17,19,21	0.96	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	A	902	-	-	5/6/23/26	0/1/1/1
3	NAG	A	903	-	-	3/6/23/26	0/1/1/1
3	NAG	A	901	-	-	1/6/23/26	0/1/1/1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	902	NAG	C2-N2-C7	3.27	127.28	122.90
3	A	902	NAG	C1-O5-C5	3.17	116.44	112.19
3	A	902	NAG	O5-C1-C2	2.75	115.54	111.29
3	A	902	NAG	O7-C7-C8	-2.29	117.97	122.05
3	A	902	NAG	O3-C3-C2	2.26	114.09	109.40

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	902	NAG	C1-C2-N2-C7
3	A	902	NAG	C8-C7-N2-C2
3	A	902	NAG	O7-C7-N2-C2
3	A	903	NAG	O7-C7-N2-C2
3	A	903	NAG	C8-C7-N2-C2
3	A	902	NAG	O5-C5-C6-O6
3	A	902	NAG	C4-C5-C6-O6
3	A	901	NAG	O5-C5-C6-O6
3	A	903	NAG	O5-C5-C6-O6

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	902	NAG	1	0
3	A	903	NAG	1	0
3	A	901	NAG	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 [i](#)

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

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, 95th 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	596/607 (98%)	0.06	14 (2%) 61 52	38, 70, 112, 145	0
2	B	8/8 (100%)	0.60	0 100 100	89, 100, 119, 127	0
All	All	604/615 (98%)	0.07	14 (2%) 61 52	38, 70, 112, 145	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	371	VAL	4.4
1	A	366	ILE	4.2
1	A	452	ASP	3.8
1	A	374	THR	3.8
1	A	582	ALA	3.6
1	A	370	ILE	3.3
1	A	384	GLN	3.3
1	A	368	PRO	2.9
1	A	513	THR	2.8
1	A	506	LEU	2.7
1	A	373	GLY	2.2
1	A	531	TYR	2.2
1	A	372	GLY	2.1
1	A	358	MET	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 oligosaccharides 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, 95th 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
3	NAG	A	903	14/15	0.18	0.14	105,124,127,128	0
3	NAG	A	901	14/15	0.62	0.14	103,119,128,128	0
3	NAG	A	902	14/15	0.68	0.16	55,76,84,87	0

6.5 Other polymers [i](#)

There are no such residues in this entry.