



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:33 AM GMT

PDB ID : 3B9F
Title : 1.6 Å structure of the PCI-thrombin-heparin complex
Authors : Li, W.; Adams, T.E.; Huntington, J.A.
Deposited on : 2007-11-05
Resolution : 1.60 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) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

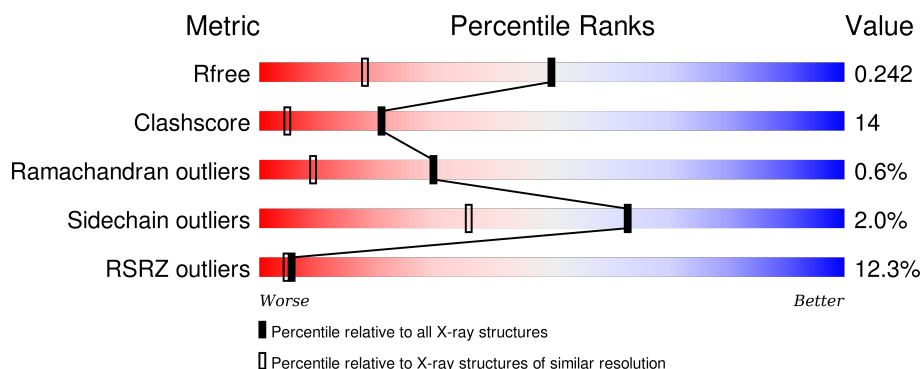
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 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}	91344	2475 (1.60-1.60)
Clashscore	102246	2732 (1.60-1.60)
Ramachandran outliers	100387	2654 (1.60-1.60)
Sidechain outliers	100360	2653 (1.60-1.60)
RSRZ outliers	91569	2479 (1.60-1.60)

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.

Mol	Chain	Length	Quality of chain
1	L	49	<div> <div>6%</div> <div>82%</div> <div>8%</div> <div>10%</div> </div>
2	H	259	<div> <div>5%</div> <div>71%</div> <div>26%</div> <div>••</div> </div>
3	I	395	<div> <div>16%</div> <div>68%</div> <div>22%</div> <div>• 10%</div> </div>

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
5	SGN	H	1	-	-	-	X
5	IDS	H	2	-	-	-	X
7	GOL	H	3	-	-	-	X
7	GOL	H	4	-	-	-	X
7	GOL	I	390	-	-	X	-

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 6092 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 Prothrombin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	44	Total	C	N	O	S	0	2	0
			377	236	61	79	1			

- Molecule 2 is a protein called Prothrombin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	253	Total	C	N	O	S	0	34	0
			2294	1456	412	410	16			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
H	195	ALA	SER	ENGINEERED	UNP P00734

- Molecule 3 is a protein called protein C inhibitor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	I	356	Total	C	N	O	S	5	9	0
			2798	1796	461	524	17			

- Molecule 4 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	H	2	Total	C	N	O	0	0
			24	14	1	9		

- Molecule 5 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	H	2	Total	C	N	O	S	0	0
			31	12	1	16	2		

- Molecule 6 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	H	1	Total	O	S	0	0
			5	4	1		
6	I	1	Total	O	S	0	0
			5	4	1		
6	I	1	Total	O	S	0	0
			5	4	1		

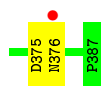
- Molecule 7 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	H	1	Total C O 6 3 3	0	0
7	H	1	Total C O 6 3 3	0	0
7	H	1	Total C O 6 3 3	0	0
7	I	1	Total C O 6 3 3	0	0
7	I	1	Total C O 6 3 3	0	0

- Molecule 8 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	L	83	Total O 83 83	0	0
8	H	272	Total O 272 272	0	0
8	I	168	Total O 168 168	0	0



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	44.05Å 48.83Å 97.85Å 78.72° 81.52° 77.69°	Depositor
Resolution (Å)	30.95 – 1.60 30.95 – 1.60	Depositor EDS
% Data completeness (in resolution range)	93.4 (30.95-1.60) 88.6 (30.95-1.60)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.94 (at 1.60Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.208 , 0.234 0.218 , 0.242	Depositor DCC
R_{free} test set	2928 reflections (3.05%)	DCC
Wilson B-factor (Å ²)	19.0	Xtriage
Anisotropy	0.068	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 61.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 95905 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6092	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.18% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

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: GOL, NAG, IDS, FUC, SGN, SO4

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	L	0.35	0/384	0.63	0/514
2	H	0.31	0/2347	0.70	0/3169
3	I	0.29	0/2851	0.56	0/3860
All	All	0.30	0/5582	0.63	0/7543

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 [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	L	377	0	352	3	0
2	H	2294	0	2222	84	0
3	I	2798	0	2718	76	0
4	H	24	0	22	1	0
5	H	31	0	16	9	0
6	H	5	0	0	0	0
6	I	10	0	0	0	0
7	H	18	0	24	1	0
7	I	12	0	16	7	0
8	H	272	0	0	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	I	168	0	0	6	0
8	L	83	0	0	1	0
All	All	6092	0	5370	158	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 14.

All (158) 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:H:185[A]:LYS:HB2	2:H:186[A]:PRO:C	1.77	1.05
2:H:185[A]:LYS:HB2	2:H:186(A)[A]:ASP:N	1.73	1.02
2:H:139:THR:HG22	2:H:157[A]:VAL:HG22	1.54	0.89
2:H:67[B]:ARG:HG2	2:H:82[B]:ILE:HD12	1.59	0.84
3:I:179:ILE:HD12	3:I:327:GLU:HB2	1.64	0.78
3:I:59:MET:HG2	3:I:130[A]:MET:SD	2.25	0.77
3:I:234:ARG:NH2	3:I:361:GLN:HE21	1.83	0.76
2:H:185[A]:LYS:HG3	2:H:186[A]:PRO:HA	1.70	0.73
2:H:185[A]:LYS:HB2	2:H:186[A]:PRO:CA	2.20	0.72
2:H:61:GLU:HG2	2:H:87:LYS:HA	1.71	0.71
3:I:353:PHE:H	7:I:390:GOL:C1	2.06	0.68
2:H:237:TRP:HB2	5:H:1:SGN:H61	1.76	0.68
2:H:185[A]:LYS:HB3	2:H:186(B)[A]:GLU:HB2	1.76	0.67
2:H:35:ARG:HB3	2:H:41[A]:LEU:HD11	1.76	0.67
2:H:59:LEU:HD13	2:H:88[B]:ILE:HG21	1.77	0.66
2:H:61:GLU:CG	2:H:87:LYS:HA	2.25	0.66
2:H:68:ILE:HD12	2:H:112[A]:VAL:HG21	1.77	0.65
2:H:108:LEU:HD13	2:H:112[A]:VAL:HG23	1.79	0.65
2:H:174[B]:ILE:CD1	3:I:351:PHE:HB3	2.28	0.64
3:I:55[B]:MET:HG2	3:I:92:PHE:CD1	2.33	0.64
3:I:157:VAL:HG21	3:I:168:LEU:HD22	1.80	0.63
2:H:34:PHE:CE1	2:H:67[A]:ARG:HD3	2.33	0.63
2:H:179:ASN:ND2	5:H:1:SGN:O3S	2.29	0.62
2:H:64:LEU:HD12	2:H:88[B]:ILE:HD11	1.82	0.62
3:I:161:THR:HG21	3:I:165:ILE:HB	1.80	0.62
3:I:353:PHE:H	7:I:390:GOL:H12	1.64	0.61
3:I:202:GLN:HB2	3:I:216[B]:MET:HG3	1.83	0.61
2:H:32[A]:MET:SD	2:H:70:LYS:HD3	2.40	0.60
1:L:1(Q):GLU:HG2	8:L:55:HOH:O	2.01	0.60
2:H:92:PRO:HG2	5:H:2:IDS:O61	2.01	0.60
3:I:234:ARG:HH21	3:I:361:GLN:HE21	1.48	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:I:93:GLN:O	3:I:97:GLN:HG3	2.02	0.59
2:H:237:TRP:HD1	5:H:2:IDS:H3	1.69	0.58
3:I:178:VAL:HG21	3:I:323:ILE:HD11	1.85	0.58
2:H:93[A]:ARG:HG3	5:H:2:IDS:O61	2.03	0.58
2:H:174[A]:ILE:HD11	3:I:350:ILE:HB	1.85	0.58
2:H:185[A]:LYS:HE2	2:H:186(D)[A]:LYS:O	2.04	0.58
2:H:93[A]:ARG:NH2	5:H:1:SGN:HN	2.03	0.57
2:H:173:ARG:NH2	3:I:350:ILE:HG23	2.20	0.57
3:I:116:THR:HG22	3:I:178:VAL:HG22	1.86	0.56
2:H:237:TRP:HB2	5:H:1:SGN:C6	2.35	0.56
2:H:61:GLU:CD	2:H:61:GLU:H	2.09	0.56
3:I:111:GLY:HA3	3:I:183:TYR:CE1	2.40	0.56
2:H:185[A]:LYS:HB3	2:H:186(B)[A]:GLU:N	2.21	0.55
2:H:60(D):TRP:CH2	7:I:390:GOL:H2	2.41	0.55
3:I:344:ALA:HB3	8:I:528:HOH:O	2.07	0.55
2:H:34:PHE:HB2	2:H:65:LEU:HD22	1.87	0.54
3:I:180:MET:HE2	3:I:328:MET:HG3	1.89	0.54
2:H:185[A]:LYS:CB	2:H:186(A)[A]:ASP:N	2.61	0.54
3:I:352:THR:HA	7:I:390:GOL:H11	1.90	0.54
2:H:185[A]:LYS:CB	2:H:186[A]:PRO:CA	2.86	0.53
3:I:118:LEU:HD11	3:I:142:ASN:HB2	1.90	0.53
3:I:67:SER:O	3:I:71:GLN:HG3	2.08	0.53
4:H:778:NAG:H61	8:H:1014:HOH:O	2.07	0.53
3:I:103:ARG:O	3:I:105:GLY:N	2.41	0.53
3:I:109:SER:O	3:I:184:ILE:HA	2.07	0.53
2:H:75[B]:ARG:HG2	2:H:75[B]:ARG:HH11	1.74	0.52
2:H:99:LEU:HD12	2:H:215:TRP:HB3	1.89	0.52
3:I:66:SER:O	3:I:70:MET:HG3	2.10	0.52
5:H:1:SGN:H1	5:H:1:SGN:HOS3	1.75	0.52
2:H:240[B]:LYS:CG	2:H:244[B]:GLN:HE21	2.23	0.52
1:L:14(C):GLU:O	1:L:14(G):LEU:HD23	2.10	0.51
3:I:161:THR:HG21	3:I:165:ILE:HD12	1.91	0.51
3:I:217:MET:SD	3:I:338:GLU:HG3	2.50	0.51
2:H:34:PHE:HB2	2:H:65:LEU:CD2	2.40	0.51
3:I:114:LEU:HD13	3:I:130[A]:MET:HE1	1.93	0.51
2:H:22:ALA:HB2	2:H:157[A]:VAL:HG23	1.91	0.51
3:I:108:LEU:C	3:I:108:LEU:HD23	2.31	0.51
3:I:353:PHE:H	7:I:390:GOL:H11	1.74	0.51
2:H:174[B]:ILE:HD12	2:H:215:TRP:CZ3	2.46	0.51
3:I:165:ILE:HG23	3:I:331:LYS:HD3	1.93	0.51
2:H:111:PRO:HD3	7:H:4:GOL:H32	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:I:63:GLY:HA3	3:I:317:ILE:HG13	1.93	0.50
3:I:344:ALA:O	3:I:345:ALA:HB2	2.12	0.50
2:H:85[A]:LEU:HD22	2:H:85[A]:LEU:N	2.26	0.50
3:I:125[A]:THR:HG23	8:I:490:HOH:O	2.10	0.50
3:I:253:GLU:HG2	8:I:520:HOH:O	2.12	0.50
3:I:29:PHE:HB3	7:I:391:GOL:O1	2.12	0.50
2:H:50[B]:ARG:HG3	2:H:51:TRP:CD1	2.47	0.49
2:H:35:ARG:NH1	2:H:37:PRO:HD2	2.27	0.49
2:H:64:LEU:HD12	2:H:88[B]:ILE:CD1	2.41	0.49
3:I:192:THR:OG1	3:I:344:ALA:CB	2.60	0.49
2:H:40[A]:LEU:C	2:H:40[A]:LEU:HD23	2.33	0.49
2:H:17:VAL:HG12	2:H:18:GLU:HG2	1.94	0.49
2:H:32[A]:MET:HG2	2:H:141:TRP:CZ3	2.47	0.49
3:I:179:ILE:CD1	3:I:327:GLU:HB2	2.37	0.49
3:I:125[A]:THR:HG22	8:I:447:HOH:O	2.12	0.49
2:H:73[A]:ARG:HD3	2:H:73[A]:ARG:O	2.13	0.49
3:I:319:ASN:C	3:I:319:ASN:HD22	2.16	0.48
3:I:249:ILE:N	3:I:249:ILE:HD12	2.28	0.48
3:I:59:MET:CG	3:I:130[A]:MET:SD	2.99	0.48
2:H:65:LEU:C	2:H:65:LEU:HD23	2.32	0.48
2:H:61:GLU:HB3	2:H:85[A]:LEU:O	2.14	0.48
2:H:103:ILE:HG21	2:H:234[A]:LEU:HD13	1.95	0.48
2:H:185[A]:LYS:CG	2:H:186[A]:PRO:HA	2.42	0.48
2:H:34:PHE:CE1	2:H:67[B]:ARG:HD2	2.48	0.47
3:I:257:GLN:NE2	3:I:261:ASN:ND2	2.61	0.47
2:H:67[B]:ARG:NH1	2:H:82[B]:ILE:CD1	2.77	0.47
2:H:172:THR:HG21	2:H:176:ILE:HD11	1.97	0.47
3:I:175:ASN:N	3:I:175:ASN:HD22	2.13	0.47
2:H:185[A]:LYS:HB3	2:H:186(B)[A]:GLU:CB	2.44	0.47
2:H:67[A]:ARG:HG2	2:H:82[A]:ILE:HG12	1.96	0.47
3:I:107:GLN:O	3:I:186:PHE:HA	2.15	0.47
2:H:60(B):PRO:HB2	2:H:60(C):PRO:HD3	1.97	0.47
2:H:35:ARG:O	2:H:38:GLN:HA	2.15	0.47
3:I:178:VAL:CG2	3:I:323:ILE:HD11	2.45	0.47
2:H:173:ARG:HH21	3:I:350:ILE:HG23	1.80	0.46
2:H:174[B]:ILE:HD11	3:I:351:PHE:HB3	1.97	0.46
3:I:95:LEU:HD23	3:I:95:LEU:C	2.36	0.46
2:H:38:GLN:O	2:H:38:GLN:HG3	2.16	0.46
2:H:68:ILE:CD1	2:H:112[A]:VAL:HG21	2.44	0.46
3:I:186:PHE:HB3	3:I:334:VAL:HG22	1.96	0.45
3:I:298:VAL:HG12	8:I:559:HOH:O	2.15	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:I:60:LEU:O	3:I:60:LEU:HD23	2.15	0.45
3:I:290:GLU:HB2	3:I:333:VAL:HG22	1.98	0.45
2:H:29:TRP:CG	2:H:121:VAL:HB	2.52	0.44
3:I:234:ARG:CZ	3:I:363:LEU:HD13	2.48	0.44
2:H:35:ARG:HH11	2:H:37:PRO:HD2	1.82	0.44
2:H:32[B]:MET:SD	2:H:34:PHE:CZ	3.11	0.44
2:H:35:ARG:HD3	2:H:37:PRO:O	2.18	0.44
3:I:226[A]:LEU:C	3:I:226[A]:LEU:HD13	2.38	0.44
2:H:75[A]:ARG:NH2	8:H:1016:HOH:O	2.50	0.43
3:I:144:ARG:HD3	3:I:144:ARG:HA	1.81	0.43
2:H:67[B]:ARG:HH11	2:H:82[B]:ILE:CD1	2.31	0.43
2:H:88[A]:ILE:HD12	2:H:88[A]:ILE:C	2.38	0.43
3:I:117:ASP:HA	3:I:141:THR:O	2.18	0.43
2:H:109:LYS:HB3	8:H:1018:HOH:O	2.19	0.43
1:L:1(G):PHE:HD1	2:H:242:ILE:HD13	1.83	0.43
3:I:194:PHE:O	3:I:338:GLU:HB3	2.19	0.43
3:I:319:ASN:HD22	3:I:320:HIS:N	2.17	0.43
2:H:129:ALA:HA	2:H:210:MET:HE1	2.01	0.42
3:I:309:THR:C	3:I:324:GLN:NE2	2.73	0.42
3:I:165:ILE:HD13	3:I:183:TYR:HB2	2.01	0.42
2:H:66:VAL:HG22	2:H:83[B]:SER:O	2.19	0.42
3:I:192:THR:OG1	3:I:344:ALA:HB1	2.20	0.42
3:I:202:GLN:HG3	8:I:455:HOH:O	2.21	0.41
2:H:59:LEU:HD22	2:H:64:LEU:HD11	2.02	0.41
3:I:192:THR:OG1	3:I:344:ALA:HB2	2.20	0.41
3:I:375:ASP:HB3	3:I:376:ASN:H	1.56	0.41
3:I:299:LEU:N	3:I:300:PRO:CD	2.84	0.41
2:H:178:ASP:O	2:H:233:ARG:HD2	2.20	0.41
2:H:32[A]:MET:CG	2:H:141:TRP:CZ3	3.02	0.41
2:H:235:LYS:NZ	8:H:951:HOH:O	2.54	0.41
2:H:75[B]:ARG:HG2	2:H:75[B]:ARG:NH1	2.34	0.41
2:H:186(D)[B]:LYS:NZ	8:H:1029:HOH:O	2.54	0.41
3:I:51:VAL:O	3:I:55[B]:MET:HG3	2.21	0.41
3:I:257:GLN:HE22	3:I:261:ASN:HD21	1.69	0.41
3:I:290:GLU:CB	3:I:333:VAL:HG22	2.51	0.41
3:I:114:LEU:HD13	3:I:130[A]:MET:CE	2.51	0.41
2:H:101:ARG:HE	5:H:1:SGN:H2	1.86	0.41
2:H:240[B]:LYS:O	2:H:244[B]:GLN:HG3	2.21	0.41
3:I:95:LEU:HD23	3:I:95:LEU:O	2.20	0.40
3:I:57:LEU:HB2	3:I:76:LEU:HD21	2.03	0.40
2:H:211:GLY:HA2	2:H:229:THR:O	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:I:353:PHE:CD1	7:I:390:GOL:H12	2.57	0.40
2:H:174[B]:ILE:HD13	3:I:351:PHE:HB3	2.00	0.40
3:I:192:THR:OG1	3:I:219:ARG:NH1	2.54	0.40
3:I:257:GLN:NE2	3:I:261:ASN:HD21	2.19	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	L	44/49 (90%)	43 (98%)	1 (2%)	0	100	100
2	H	282/259 (109%)	275 (98%)	7 (2%)	0	100	100
3	I	361/395 (91%)	337 (93%)	20 (6%)	4 (1%)	17	3
All	All	687/703 (98%)	655 (95%)	28 (4%)	4 (1%)	30	9

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	I	104	ASP
3	I	346	ALA
3	I	105	GLY
3	I	102	PRO

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	L	42/43 (98%)	42 (100%)	0	100	100
2	H	241/224 (108%)	237 (98%)	4 (2%)	68	44
3	I	296/347 (85%)	289 (98%)	7 (2%)	57	27
All	All	579/614 (94%)	568 (98%)	11 (2%)	63	39

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	H	35	ARG
2	H	65	LEU
2	H	94	TYR
2	H	182	CYS
3	I	78	LEU
3	I	192	THR
3	I	195	ASN
3	I	196	HIS
3	I	203	ASP
3	I	275	PHE
3	I	319	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
2	H	38	GLN
2	H	179	ASN
3	I	97	GLN
3	I	152	GLN
3	I	175	ASN
3	I	195	ASN
3	I	196	HIS
3	I	257	GLN
3	I	261	ASN
3	I	294	GLN
3	I	319	ASN
3	I	324	GLN
3	I	361	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 ⓘ

4 carbohydrates 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$
5	SGN	H	1	5	12,15,20	4.65	2 (16%)	11,22,31	1.17	1 (9%)
5	IDS	H	2	5	12,16,17	1.60	3 (25%)	14,24,26	0.69	0
4	NAG	H	778	2,4	14,14,15	0.51	0	15,19,21	0.75	1 (6%)
4	FUC	H	780	4	10,10,11	0.46	0	14,14,16	0.52	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
5	SGN	H	1	5	-	0/7/23/31	0/1/1/1
5	IDS	H	2	5	-	0/5/25/29	1/1/1/1
4	NAG	H	778	2,4	-	0/6/23/26	0/1/1/1
4	FUC	H	780	4	-	0/0/17/20	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	H	2	IDS	O2-S	-4.10	1.44	1.57
5	H	2	IDS	O2S-S	2.25	1.53	1.45
5	H	2	IDS	O1S-S	2.26	1.53	1.45
5	H	1	SGN	O1S-S1	11.28	1.53	1.42
5	H	1	SGN	O2S-S1	11.39	1.53	1.42

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	H	778	NAG	C2-N2-C7	-2.22	120.19	123.04
5	H	1	SGN	O5-C5-C6	2.25	109.38	106.62

There are no chirality outliers.

There are no torsion outliers.

All (1) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	H	2	IDS	C1-C2-C3-C4-C5-O5

3 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	H	1	SGN	6	0
5	H	2	IDS	3	0
4	H	778	NAG	1	0

5.6 Ligand geometry

8 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	$\# Z > 2$
7	GOL	H	3	-	5,5,5	0.25	0	5,5,5	0.24	0
7	GOL	H	4	-	5,5,5	0.26	0	5,5,5	0.21	0
6	SO4	H	781	-	4,4,4	0.20	0	6,6,6	0.10	0
7	GOL	H	782	-	5,5,5	0.27	0	5,5,5	0.13	0
6	SO4	I	388	-	4,4,4	0.21	0	6,6,6	0.08	0
6	SO4	I	389	-	4,4,4	0.20	0	6,6,6	0.08	0
7	GOL	I	390	-	5,5,5	0.25	0	5,5,5	0.24	0
7	GOL	I	391	-	5,5,5	0.26	0	5,5,5	0.24	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
7	GOL	H	3	-	-	0/4/4/4	0/0/0/0
7	GOL	H	4	-	-	0/4/4/4	0/0/0/0
6	SO4	H	781	-	-	0/0/0/0	0/0/0/0
7	GOL	H	782	-	-	0/4/4/4	0/0/0/0
6	SO4	I	388	-	-	0/0/0/0	0/0/0/0
6	SO4	I	389	-	-	0/0/0/0	0/0/0/0
7	GOL	I	390	-	-	0/4/4/4	0/0/0/0
7	GOL	I	391	-	-	0/4/4/4	0/0/0/0

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.

3 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	H	4	GOL	1	0
7	I	390	GOL	6	0
7	I	391	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, 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	L	44/49 (89%)	-0.24	3 (6%) 20 19	12, 18, 28, 45	2 (4%)
2	H	253/259 (97%)	0.17	14 (5%) 29 25	9, 17, 35, 53	23 (9%)
3	I	355/395 (89%)	1.11	63 (17%) 2 1	14, 37, 59, 69	4 (1%)
All	All	652/703 (92%)	0.65	80 (12%) 5 4	9, 28, 54, 69	29 (4%)

All (80) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	I	343	ALA	12.9
3	I	345	ALA	11.0
3	I	102	PRO	7.3
3	I	103	ARG	7.1
3	I	106	PHE	6.8
3	I	104	ASP	6.5
3	I	83	SER	6.2
3	I	196	HIS	5.8
3	I	344	ALA	5.7
3	I	100	ASN	5.3
3	I	84	SER	5.3
3	I	347	THR	5.2
2	H	36(A)	SER	5.1
3	I	341	THR	5.0
3	I	78	LEU	4.5
3	I	101	GLN	4.5
3	I	105	GLY	4.3
3	I	195	ASN	4.1
3	I	77	GLY	3.8
1	L	14(L)	ASP	3.7
3	I	171	ASN	3.7
3	I	348	GLY	3.7
2	H	45[A]	SER	3.7

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Mol	Chain	Res	Type	RSRZ
3	I	107	GLN	3.7
3	I	371[A]	MET	3.6
3	I	242	GLY	3.5
3	I	200	GLN	3.4
3	I	376	ASN	3.4
3	I	39	SER	3.3
3	I	342	ARG	3.2
3	I	322	ASN	3.2
2	H	147	THR	3.1
3	I	42	PRO	3.1
3	I	257	GLN	3.1
2	H	212	ILE	3.0
3	I	172	LEU	3.0
3	I	88	LEU	2.8
2	H	121	VAL	2.8
3	I	230	ASN	2.8
3	I	205	TYR	2.8
3	I	40	ALA	2.8
2	H	200	VAL	2.7
3	I	28	ASP	2.7
2	H	149[A]	LYS	2.7
2	H	44	ALA	2.6
3	I	311	HIS	2.6
2	H	199	PHE	2.6
3	I	99	LEU	2.6
3	I	170	LYS	2.5
3	I	147	ALA	2.5
3	I	70	MET	2.5
3	I	211	VAL	2.5
3	I	197	LYS	2.5
3	I	241	GLN	2.5
3	I	199	THR	2.4
3	I	44	GLN	2.4
3	I	167	ASP	2.4
3	I	189	LYS	2.4
3	I	94	GLN	2.4
3	I	210	THR	2.3
3	I	274	MET	2.3
3	I	173	ASP	2.3
1	L	1(R)	SER	2.3
3	I	73	LEU	2.3
3	I	166	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
3	I	273	LYS	2.2
2	H	47	ILE	2.2
3	I	358	LEU	2.2
3	I	114	LEU	2.1
3	I	86	LYS	2.1
3	I	67	SER	2.1
1	L	1(Q)	GLU	2.1
2	H	46	LEU	2.1
2	H	228	TYR	2.1
3	I	91	GLY	2.1
3	I	66	SER	2.0
3	I	97	GLN	2.0
2	H	41[A]	LEU	2.0
3	I	133	LEU	2.0
2	H	36	LYS	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](#)

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
5	SGN	H	1	15/20	0.30	0.63	15.66	65,72,76,76	0
5	IDS	H	2	16/17	0.54	0.43	4.12	71,72,76,76	0
4	NAG	H	778	14/15	0.67	0.23	-	50,54,59,63	0
4	FUC	H	780	10/11	0.84	0.48	-	66,68,68,69	0

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
7	GOL	H	4	6/6	0.77	0.21	2.60	55,55,55,55	2
7	GOL	H	3	6/6	0.85	0.19	2.49	48,49,49,50	3
7	GOL	H	782	6/6	0.91	0.12	1.24	20,24,25,27	3
7	GOL	I	391	6/6	0.66	0.17	0.24	60,60,61,61	0
7	GOL	I	390	6/6	0.87	0.12	-0.11	44,45,45,46	2
6	SO4	I	388	5/5	0.96	0.10	-0.99	70,70,70,71	3
6	SO4	I	389	5/5	0.82	0.27	-	76,76,76,76	2
6	SO4	H	781	5/5	0.86	0.14	-	77,77,77,77	3

6.5 Other polymers [i](#)

There are no such residues in this entry.