



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 07:01 AM GMT

PDB ID : 1GWB
Title : structure of glycoprotein 1b
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Deposited on : 2002-03-14
Resolution : 2.80 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

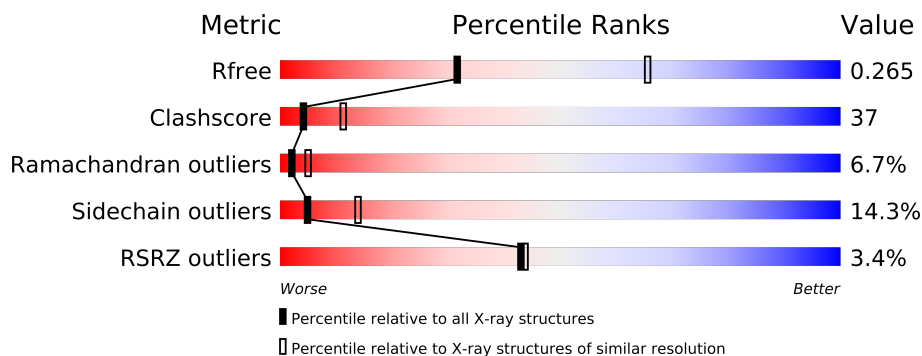
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.80 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}	66092	1799 (2.80-2.80)
Clashscore	79885	2295 (2.80-2.80)
Ramachandran outliers	78287	2252 (2.80-2.80)
Sidechain outliers	78261	2254 (2.80-2.80)
RSRZ outliers	66119	1802 (2.80-2.80)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	281	
1	B	281	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
2	PT	A	404	-	X
3	ACY	A	510	-	X

2 Entry composition i

There are 6 unique types of molecules in this entry. The entry contains 4451 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

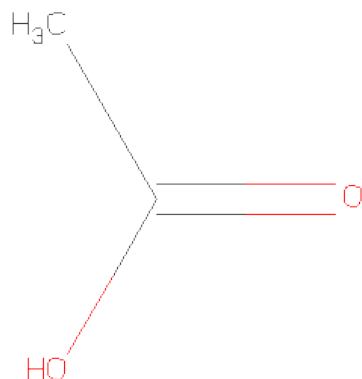
- Molecule 1 is a protein called PLATELET GLYCOPROTEIN IB ALPHA CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	266	Total	C	N	O	S	0	0	1
			2063	1326	343	387	7			
1	B	279	Total	C	N	O	S	0	0	0
			2190	1399	356	425	10			

- Molecule 2 is PLATINUM (II) ION (three-letter code: PT) (formula: Pt).

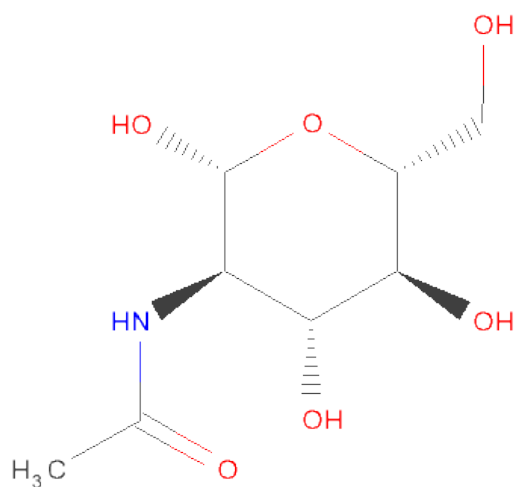
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	4	Total	Pt	0	0
			4	4		

- Molecule 3 is ACETIC ACID (three-letter code: ACY) (formula: C₂H₄O₂).



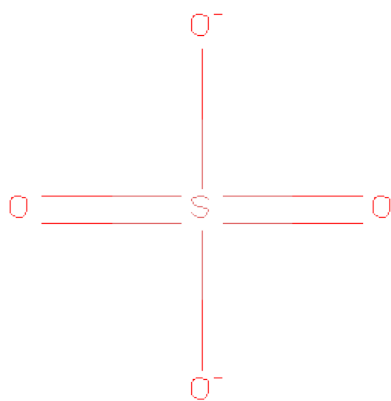
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 4 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



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

- Molecule 5 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 6 is water.

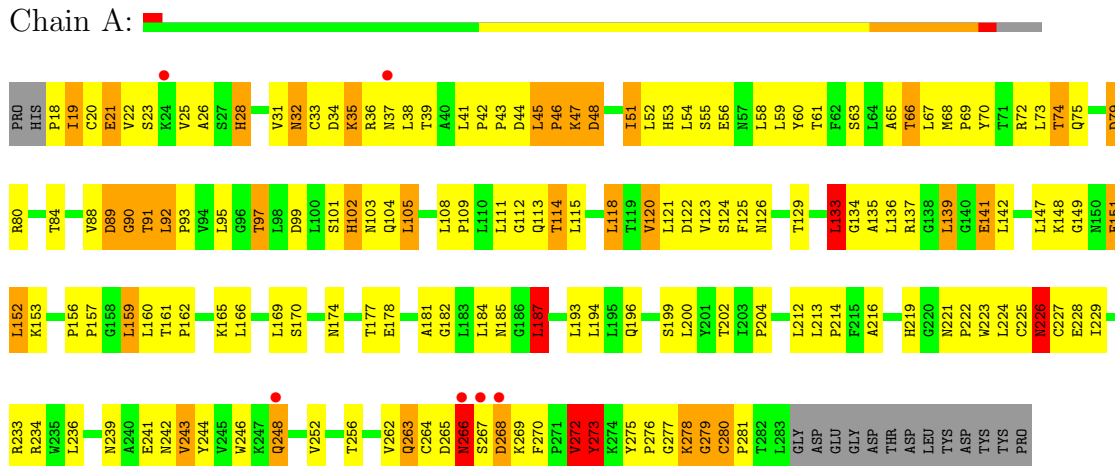
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	57	Total	O	0	0
			57	57		
6	B	86	Total	O	0	0
			86	86		

3 Residue-property plots

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.

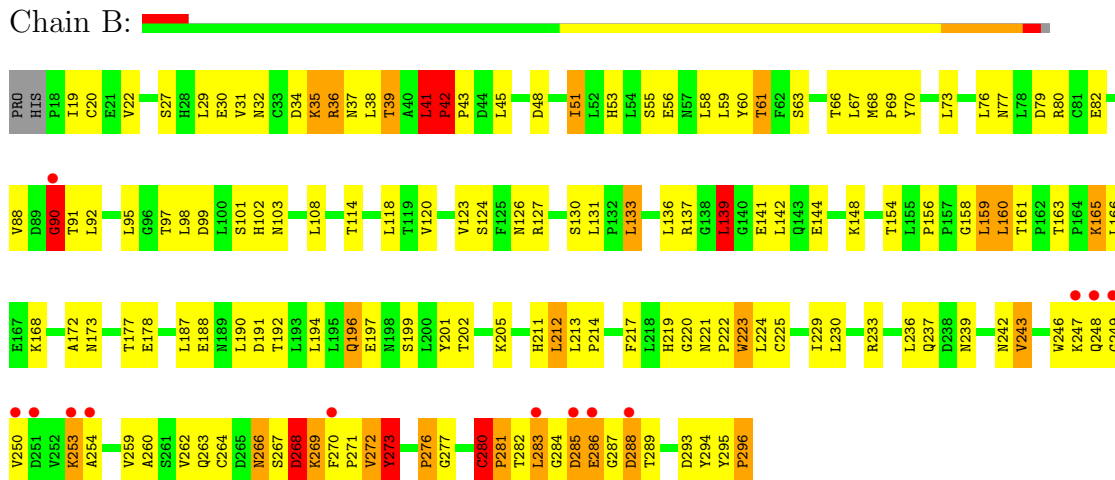
• Molecule 1: PLATELET GLYCOPROTEIN IB ALPHA CHAIN

Chain A:



• Molecule 1: PLATELET GLYCOPROTEIN IB ALPHA CHAIN

Chain B:



4 Data and refinement statistics

Property	Value	Source
Space group	P 64 2 2	Depositor
Cell constants a, b, c, α , β , γ	202.00Å 202.00Å 128.00Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	15.00 – 2.80 39.64 – 2.79	Depositor EDS
% Data completeness (in resolution range)	98.5 (15.00-2.80) 98.0 (39.64-2.79)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.12 (at 2.81Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.245 , 0.274 0.243 , 0.265	Depositor DCC
R_{free} test set	1903 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	50.5	Xtriage
Anisotropy	0.419	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 38.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 38539 reflections	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	4451	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.65% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ACY, TYS, NAG, PT, 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	A	0.38	0/2109	0.68	2/2883 (0.1%)
1	B	0.42	0/2187	0.81	6/2984 (0.2%)
All	All	0.40	0/4296	0.75	8/5867 (0.1%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	41	LEU	C-N-CD	-13.85	90.13	120.60
1	B	90	GLY	N-CA-C	7.46	131.75	113.10
1	B	41	LEU	C-N-CA	6.08	147.55	122.00
1	B	139	LEU	CA-CB-CG	6.07	129.25	115.30
1	A	90	GLY	N-CA-C	5.72	127.40	113.10
1	B	42	PRO	CA-N-CD	-5.41	103.93	111.50
1	A	272	VAL	N-CA-C	-5.38	96.47	111.00
1	B	280	CYS	CA-CB-SG	-5.24	104.56	114.00

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2063	0	2089	172	1
1	B	2190	0	2181	145	1
2	A	4	0	0	0	0
3	A	4	0	3	0	0
4	A	42	0	39	3	0
5	B	5	0	0	0	0
6	A	57	0	0	3	1
6	B	86	0	0	3	1
All	All	4451	0	4312	315	4

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 37.

All (315) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:90:GLY:HA2	1:B:114:THR:HB	1.38	1.04
1:A:105:LEU:HD23	1:A:105:LEU:H	1.19	1.04
1:A:97:THR:HB	1:A:120:VAL:HG13	1.50	0.94
1:B:272:VAL:O	1:B:273:TYR:HB3	1.66	0.92
1:A:151:GLU:HG2	4:A:1175:NAG:H82	1.53	0.90
1:A:153:LYS:HE3	1:A:177:THR:HG21	1.53	0.87
1:A:222:PRO:HA	1:A:265:ASP:HB2	1.56	0.87
1:A:272:VAL:O	1:A:273:TYR:HB3	1.76	0.85
1:A:37:ASN:HA	1:A:58:LEU:HD23	1.55	0.85
1:B:90:GLY:HA2	1:B:114:THR:CB	2.07	0.84
1:B:205:LYS:HD3	1:B:294:TYS:HE1	1.57	0.84
1:A:139:LEU:HD23	1:A:139:LEU:H	1.47	0.79
1:A:225:CYS:HB2	1:A:229:ILE:HG21	1.64	0.79
1:B:91:THR:HG22	6:B:2037:HOH:O	1.81	0.79
1:B:133:LEU:H	1:B:133:LEU:HD22	1.47	0.78
1:A:47:LYS:O	1:A:48:ASP:HB2	1.84	0.78
1:B:41:LEU:HB3	1:B:42:PRO:O	1.85	0.77
1:B:259:VAL:O	1:B:262:VAL:HG12	1.86	0.76
1:B:212:LEU:HG	1:B:239:ASN:HD21	1.49	0.76
1:B:196:GLN:HG3	1:B:217:PHE:HB3	1.69	0.74
1:A:75:GLN:HG2	1:A:97:THR:HG23	1.68	0.74
1:A:41:LEU:HD12	1:A:42:PRO:HD2	1.69	0.74
1:B:233:ARG:HG2	1:B:273:TYR:HA	1.68	0.74
1:B:31:VAL:HG11	1:B:45:LEU:HD12	1.70	0.74
1:A:224:LEU:HA	1:A:266:ASN:HD21	1.53	0.74
1:B:45:LEU:HD22	1:B:45:LEU:H	1.54	0.73
1:B:295:TYS:CG	1:B:296:PRO:HD2	2.19	0.72
1:A:105:LEU:H	1:A:105:LEU:CD2	2.00	0.72

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:56:GLU:HA	1:A:80:ARG:O	1.90	0.71
1:A:151:GLU:CG	4:A:1175:NAG:H82	2.20	0.71
1:A:108:LEU:HD22	1:A:109:PRO:HD2	1.72	0.71
1:A:42:PRO:O	1:A:45:LEU:HD23	1.90	0.71
1:B:99:ASP:OD1	1:B:101:SER:HB3	1.90	0.71
1:B:289:THR:HG23	1:B:293:ASP:OD1	1.91	0.70
1:A:19:ILE:HG21	1:A:44:ASP:OD2	1.91	0.70
1:B:88:VAL:HG12	1:B:114:THR:HG21	1.74	0.70
1:B:97:THR:HG22	1:B:120:VAL:HG13	1.74	0.70
1:B:20:CYS:HA	1:B:36:ARG:HH12	1.56	0.69
1:A:123:VAL:HG22	1:A:147:LEU:HD23	1.75	0.69
1:A:262:VAL:HG12	1:A:272:VAL:CG1	2.23	0.69
1:B:236:LEU:HD22	1:B:262:VAL:HG11	1.74	0.69
1:B:90:GLY:CA	1:B:114:THR:HB	2.21	0.68
1:A:266:ASN:HD22	1:A:266:ASN:N	1.91	0.68
1:A:105:LEU:N	1:A:105:LEU:HD23	2.03	0.68
1:A:153:LYS:HG3	1:A:177:THR:HG23	1.73	0.68
1:B:233:ARG:HD3	1:B:237:GLN:NE2	2.08	0.68
1:A:19:ILE:HG23	1:A:20:CYS:N	2.08	0.68
1:A:219:HIS:CE1	1:A:246:TRP:H	2.12	0.67
1:B:51:ILE:HD11	1:B:53:HIS:NE2	2.09	0.67
1:A:139:LEU:CD2	1:A:139:LEU:H	2.05	0.67
1:A:241:GLU:HG2	6:A:2048:HOH:O	1.94	0.66
1:B:148:LYS:HE3	1:B:173:ASN:HD22	1.59	0.66
1:A:91:THR:O	1:A:92:LEU:HB2	1.96	0.66
1:A:88:VAL:HG12	1:A:114:THR:HG21	1.78	0.65
1:A:97:THR:HB	1:A:120:VAL:CG1	2.25	0.65
1:A:99:ASP:OD1	1:A:101:SER:HB2	1.97	0.65
1:A:88:VAL:HG12	1:A:114:THR:CG2	2.27	0.64
1:B:177:THR:O	1:B:178:GLU:HG2	1.98	0.64
1:A:112:GLY:HA3	1:A:137:ARG:O	1.98	0.63
1:B:148:LYS:HE3	1:B:173:ASN:ND2	2.13	0.63
1:A:48:ASP:OD2	1:A:72:ARG:HD2	1.97	0.63
1:A:272:VAL:O	1:A:273:TYR:CB	2.47	0.63
1:B:172:ALA:HB2	1:B:196:GLN:NE2	2.13	0.63
1:A:233:ARG:HG3	1:A:273:TYR:HA	1.79	0.62
1:B:20:CYS:HA	1:B:36:ARG:NH1	2.13	0.62
1:A:268:ASP:OD2	1:A:270:PHE:HB2	2.00	0.62
1:A:280:CYS:SG	1:A:281:PRO:HD2	2.39	0.62
1:B:68:MET:CB	1:B:91:THR:HG21	2.29	0.62
1:A:92:LEU:H	1:A:93:PRO:HD3	1.64	0.62
1:A:222:PRO:HA	1:A:265:ASP:CB	2.30	0.62

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:205:LYS:HD3	1:B:294:TYS:CE1	2.27	0.61
1:B:41:LEU:HB3	1:B:42:PRO:C	2.21	0.61
1:A:19:ILE:HG23	1:A:20:CYS:H	1.63	0.61
1:B:63:SER:HB3	1:B:66:THR:HG23	1.82	0.61
1:A:221:ASN:HB2	1:A:223:TRP:NE1	2.16	0.61
1:B:268:ASP:C	1:B:268:ASP:OD2	2.38	0.61
4:A:1175:NAG:H3	4:A:1175:NAG:H83	1.81	0.61
1:A:263:GLN:OE1	1:A:269:LYS:HG3	2.00	0.61
1:A:42:PRO:HB2	1:A:45:LEU:CD2	2.30	0.60
1:A:212:LEU:O	1:A:214:PRO:HD3	2.02	0.60
1:B:38:LEU:O	1:B:39:THR:HG22	2.02	0.60
1:B:197:GLU:HG3	1:B:220:GLY:HA3	1.83	0.59
1:A:118:LEU:HD11	1:A:121:LEU:HB2	1.83	0.59
1:A:31:VAL:HG11	1:A:45:LEU:CD1	2.33	0.59
1:A:228:GLU:HB2	6:A:2042:HOH:O	2.03	0.59
1:B:196:GLN:CG	1:B:217:PHE:HB3	2.32	0.59
1:B:22:VAL:HG13	1:B:29:LEU:HD21	1.85	0.59
1:B:282:THR:O	1:B:284:GLY:N	2.36	0.58
1:A:184:LEU:HA	1:A:187:LEU:HD22	1.85	0.58
1:A:32:ASN:HD22	1:A:33:CYS:N	2.01	0.58
1:A:225:CYS:HA	1:A:229:ILE:HB	1.86	0.58
1:B:120:VAL:HB	1:B:144:GLU:HB2	1.85	0.58
1:B:269:LYS:HE3	1:B:269:LYS:HA	1.84	0.58
1:B:34:ASP:O	1:B:36:ARG:HD2	2.03	0.58
1:A:80:ARG:HA	1:A:102:HIS:O	2.03	0.57
1:B:280:CYS:SG	1:B:281:PRO:HD2	2.44	0.57
1:A:234:ARG:HH21	1:A:234:ARG:HG3	1.69	0.57
1:B:225:CYS:HB2	1:B:229:ILE:HG21	1.87	0.57
1:A:125:PHE:CD1	1:A:149:GLY:HA3	2.40	0.57
1:A:265:ASP:O	1:A:267:SER:N	2.38	0.57
1:A:68:MET:N	1:A:69:PRO:HD2	2.20	0.57
1:B:276:PRO:HG2	1:B:277:GLY:H	1.69	0.57
1:A:58:LEU:O	1:A:59:LEU:HD23	2.05	0.56
1:A:124:SER:HB2	1:A:148:LYS:O	2.05	0.56
1:A:67:LEU:C	1:A:69:PRO:HD2	2.26	0.56
1:A:182:GLY:CA	1:A:185:ASN:ND2	2.69	0.56
1:A:125:PHE:CZ	1:B:288:ASP:HB3	2.41	0.55
1:A:44:ASP:O	1:A:45:LEU:C	2.45	0.55
1:B:219:HIS:HE1	1:B:246:TRP:H	1.51	0.55
1:B:263:GLN:HG3	1:B:266:ASN:HA	1.89	0.55
1:B:172:ALA:HB2	1:B:196:GLN:HE22	1.70	0.55
1:A:63:SER:O	1:A:66:THR:HG23	2.06	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:28:HIS:H	1:A:28:HIS:CD2	2.25	0.55
1:B:267:SER:O	1:B:268:ASP:CB	2.55	0.54
1:B:263:GLN:CD	1:B:268:ASP:HA	2.28	0.54
1:B:103:ASN:O	1:B:126:ASN:HA	2.08	0.54
1:A:277:GLY:O	1:A:279:GLY:N	2.41	0.54
1:B:286:GLU:CG	1:B:287:GLY:N	2.68	0.54
1:A:20:CYS:O	1:A:21:GLU:CB	2.56	0.54
1:A:123:VAL:HG23	1:A:123:VAL:O	2.07	0.54
1:B:262:VAL:O	1:B:262:VAL:HG13	2.08	0.54
1:B:35:LYS:HG3	1:B:56:GLU:HB2	1.89	0.54
1:A:92:LEU:H	1:A:93:PRO:CD	2.21	0.53
1:B:42:PRO:HB3	1:B:45:LEU:HD11	1.89	0.53
1:B:247:LYS:HB3	1:B:250:VAL:HB	1.89	0.53
1:A:170:SER:HB2	1:A:194:LEU:HD13	1.90	0.53
1:A:153:LYS:CE	1:A:177:THR:HG21	2.32	0.53
1:B:124:SER:HB2	1:B:148:LYS:O	2.08	0.53
1:A:266:ASN:ND2	1:A:266:ASN:N	2.57	0.52
1:B:39:THR:HA	1:B:58:LEU:O	2.09	0.52
1:A:32:ASN:C	1:A:32:ASN:HD22	2.12	0.52
1:B:191:ASP:HA	1:B:214:PRO:HD2	1.91	0.52
1:B:88:VAL:CG1	1:B:114:THR:HG21	2.39	0.52
1:B:269:LYS:C	1:B:271:PRO:HD3	2.30	0.52
1:A:262:VAL:HG12	1:A:272:VAL:HG11	1.90	0.52
1:A:139:LEU:CD2	1:A:139:LEU:N	2.73	0.52
1:A:136:LEU:HA	1:A:139:LEU:HD21	1.92	0.52
1:B:187:LEU:N	1:B:187:LEU:HD12	2.25	0.52
1:B:267:SER:O	1:B:268:ASP:HB3	2.10	0.51
1:A:151:GLU:O	1:A:174:ASN:HB3	2.10	0.51
1:B:55:SER:OG	1:B:79:ASP:HB3	2.11	0.51
1:A:114:THR:HG23	6:A:2015:HOH:O	2.08	0.51
1:A:153:LYS:HE3	1:A:177:THR:CG2	2.35	0.51
1:A:134:GLY:O	1:A:137:ARG:HB3	2.10	0.51
1:B:19:ILE:HG23	1:B:20:CYS:N	2.26	0.51
1:A:275:TYR:CE2	1:A:277:GLY:HA2	2.46	0.51
1:B:177:THR:C	1:B:178:GLU:HG2	2.31	0.50
1:B:68:MET:N	1:B:69:PRO:HD2	2.26	0.50
1:A:41:LEU:HD12	1:A:42:PRO:CD	2.40	0.50
1:A:202:THR:OG1	1:A:228:GLU:HG3	2.11	0.50
1:B:36:ARG:HB2	1:B:38:LEU:HG	1.93	0.50
1:B:219:HIS:CE1	1:B:246:TRP:H	2.29	0.50
1:B:165:LYS:HE2	6:B:2057:HOH:O	2.10	0.50
1:B:236:LEU:CD2	1:B:262:VAL:HG11	2.40	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:278:LYS:O	1:A:278:LYS:HG3	2.12	0.50
1:B:168:LYS:HG3	1:B:192:THR:HB	1.94	0.50
1:A:18:PRO:O	1:A:19:ILE:HG22	2.12	0.49
1:B:233:ARG:HD3	1:B:237:GLN:HE22	1.76	0.49
1:B:263:GLN:NE2	1:B:268:ASP:CB	2.75	0.49
1:A:123:VAL:HG22	1:A:147:LEU:CD2	2.39	0.49
1:B:286:GLU:HG3	1:B:287:GLY:N	2.27	0.49
1:A:264:CYS:C	1:A:266:ASN:H	2.15	0.49
1:A:268:ASP:O	1:A:269:LYS:HB3	2.13	0.49
1:A:226:ASN:O	1:A:227:CYS:HB2	2.13	0.49
1:A:219:HIS:HE1	1:A:246:TRP:H	1.58	0.49
1:B:187:LEU:HD23	1:B:190:LEU:HD13	1.94	0.49
1:A:58:LEU:HD22	1:A:58:LEU:N	2.27	0.49
1:B:53:HIS:HD2	1:B:77:ASN:HD22	1.61	0.49
1:A:161:THR:N	1:A:162:PRO:CD	2.76	0.49
1:A:19:ILE:CG2	1:A:20:CYS:N	2.75	0.49
1:B:213:LEU:H	1:B:213:LEU:HD23	1.78	0.49
1:A:95:LEU:HD12	1:A:118:LEU:HD23	1.95	0.48
1:B:108:LEU:HD22	1:B:131:LEU:HD22	1.94	0.48
1:A:236:LEU:HD11	1:A:262:VAL:HG11	1.95	0.48
1:A:38:LEU:O	1:A:39:THR:HB	2.13	0.48
1:B:38:LEU:O	1:B:39:THR:CB	2.60	0.48
1:A:199:SER:HA	1:A:222:PRO:CD	2.43	0.48
1:A:59:LEU:C	1:A:61:THR:H	2.15	0.48
1:B:177:THR:C	1:B:178:GLU:CG	2.81	0.48
1:A:19:ILE:CD1	1:A:42:PRO:HB3	2.44	0.48
1:A:125:PHE:N	1:A:148:LYS:O	2.47	0.48
1:B:35:LYS:HA	1:B:56:GLU:O	2.13	0.48
1:B:59:LEU:O	1:B:61:THR:N	2.42	0.48
1:A:37:ASN:HA	1:A:58:LEU:CD2	2.36	0.47
1:B:137:ARG:HG2	6:B:2055:HOH:O	2.14	0.47
1:B:295:TYS:CD1	1:B:296:PRO:HD2	2.43	0.47
1:B:35:LYS:C	1:B:36:ARG:HD2	2.35	0.47
1:B:29:LEU:HD23	1:B:30:GLU:N	2.29	0.47
1:A:69:PRO:HG2	1:A:70:TYR:CE1	2.49	0.47
1:A:226:ASN:HD22	1:A:226:ASN:N	2.13	0.47
1:B:133:LEU:H	1:B:133:LEU:CD2	2.24	0.47
1:A:136:LEU:O	1:A:139:LEU:HD23	2.14	0.47
1:A:103:ASN:HB2	1:A:126:ASN:OD1	2.15	0.47
1:B:60:TYR:HA	1:B:82:GLU:O	2.14	0.47
1:A:22:VAL:HG12	1:A:23:SER:N	2.28	0.47
1:A:276:PRO:HG2	1:A:277:GLY:H	1.80	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:159:LEU:HD13	1:B:160:LEU:HD13	1.97	0.47
1:B:264:CYS:C	1:B:266:ASN:N	2.65	0.47
1:B:194:LEU:HD23	1:B:217:PHE:CG	2.50	0.46
1:B:97:THR:HB	1:B:120:VAL:HG22	1.96	0.46
1:A:91:THR:O	1:A:92:LEU:CB	2.62	0.46
1:B:177:THR:HG22	1:B:199:SER:HB2	1.96	0.46
1:A:51:ILE:O	1:A:51:ILE:HG12	2.15	0.46
1:B:286:GLU:HG3	1:B:287:GLY:H	1.81	0.46
1:B:264:CYS:C	1:B:266:ASN:H	2.19	0.46
1:A:34:ASP:O	1:A:35:LYS:C	2.54	0.46
1:B:53:HIS:HD2	1:B:77:ASN:ND2	2.14	0.46
1:A:234:ARG:HG3	1:A:234:ARG:NH2	2.31	0.46
1:A:46:PRO:O	1:A:48:ASP:N	2.49	0.46
1:B:248:GLN:HG2	1:B:249:GLY:N	2.30	0.46
1:B:136:LEU:O	1:B:139:LEU:HD22	2.16	0.46
1:A:39:THR:HA	1:A:58:LEU:O	2.16	0.45
1:A:248:GLN:HE22	1:B:286:GLU:HA	1.80	0.45
1:B:160:LEU:HG	1:B:166:LEU:HD12	1.98	0.45
1:A:262:VAL:HG12	1:A:272:VAL:HG13	1.98	0.45
1:B:295:TYS:CD2	1:B:296:PRO:HB2	2.47	0.45
1:B:80:ARG:HA	1:B:102:HIS:O	2.16	0.45
1:A:141:GLU:OE1	1:A:141:GLU:HA	2.16	0.45
1:B:213:LEU:HA	1:B:214:PRO:HD3	1.75	0.45
1:A:59:LEU:C	1:A:60:TYR:HD1	2.21	0.45
1:A:177:THR:O	1:A:200:LEU:HA	2.17	0.45
1:B:20:CYS:HB3	1:B:32:ASN:O	2.17	0.45
1:A:239:ASN:HD22	1:A:242:ASN:HD22	1.65	0.45
1:B:196:GLN:HG3	1:B:217:PHE:CB	2.44	0.44
1:A:51:ILE:HD11	1:A:53:HIS:NE2	2.32	0.44
1:A:169:LEU:C	1:A:169:LEU:HD13	2.38	0.44
1:B:259:VAL:HG13	1:B:260:ALA:N	2.32	0.44
1:A:74:THR:HG22	1:A:75:GLN:HG3	1.98	0.44
1:A:36:ARG:O	1:A:37:ASN:C	2.55	0.44
1:A:58:LEU:HB3	1:A:60:TYR:HE1	1.82	0.44
1:B:69:PRO:HG2	1:B:70:TYR:CE1	2.52	0.44
1:A:19:ILE:HD11	1:A:42:PRO:HB3	1.99	0.44
1:B:194:LEU:HD23	1:B:217:PHE:CD1	2.51	0.44
1:A:216:ALA:O	1:A:243:VAL:HA	2.16	0.44
1:B:272:VAL:O	1:B:273:TYR:CB	2.50	0.44
1:A:124:SER:HB3	1:A:148:LYS:H	1.83	0.44
1:A:269:LYS:O	1:A:269:LYS:HG2	2.17	0.44
1:B:139:LEU:O	1:B:163:THR:HG22	2.17	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:91:THR:OG1	1:B:92:LEU:HG	2.18	0.44
1:A:92:LEU:N	1:A:93:PRO:CD	2.78	0.44
1:B:29:LEU:HD23	1:B:29:LEU:C	2.38	0.44
1:A:59:LEU:O	1:A:61:THR:N	2.40	0.44
1:B:92:LEU:HB3	1:B:95:LEU:HB2	2.00	0.44
1:B:38:LEU:O	1:B:39:THR:CG2	2.66	0.43
1:A:121:LEU:HD12	1:A:122:ASP:N	2.32	0.43
1:B:156:PRO:O	1:B:159:LEU:HB2	2.18	0.43
1:A:60:TYR:CD1	1:A:60:TYR:N	2.85	0.43
1:A:42:PRO:HB2	1:A:45:LEU:HD21	2.00	0.43
1:B:34:ASP:OD1	1:B:55:SER:HB2	2.19	0.43
1:B:177:THR:OG1	1:B:178:GLU:HG3	2.17	0.43
1:A:19:ILE:CG2	1:A:20:CYS:H	2.28	0.43
1:A:181:ALA:HA	1:A:204:PRO:CG	2.48	0.43
1:B:269:LYS:CE	1:B:269:LYS:HA	2.45	0.43
1:B:36:ARG:O	1:B:37:ASN:C	2.55	0.43
1:A:32:ASN:ND2	1:A:34:ASP:OD2	2.50	0.43
1:B:270:PHE:N	1:B:271:PRO:HD3	2.33	0.43
1:A:45:LEU:HA	1:A:46:PRO:HD2	1.83	0.43
1:A:125:PHE:CE2	1:B:288:ASP:HB3	2.54	0.43
1:A:124:SER:O	1:A:125:PHE:HB2	2.19	0.43
1:A:129:THR:O	1:A:152:LEU:HA	2.18	0.43
1:A:182:GLY:HA3	1:A:185:ASN:ND2	2.33	0.43
1:A:84:THR:HG22	1:A:104:GLN:HB2	2.01	0.42
1:B:67:LEU:C	1:B:69:PRO:HD2	2.40	0.42
1:A:32:ASN:ND2	1:A:32:ASN:C	2.72	0.42
1:A:226:ASN:ND2	1:A:226:ASN:N	2.66	0.42
1:A:74:THR:CG2	1:A:75:GLN:HG3	2.50	0.42
1:A:244:TYR:N	1:A:244:TYR:CD1	2.87	0.42
1:B:223:TRP:NE1	1:B:262:VAL:HG23	2.35	0.42
1:B:286:GLU:CG	1:B:287:GLY:H	2.33	0.42
1:A:278:LYS:CG	1:A:278:LYS:O	2.68	0.42
1:A:32:ASN:HA	1:A:53:HIS:HB2	2.00	0.42
1:A:161:THR:HB	1:A:162:PRO:HD3	2.00	0.42
1:A:79:ASP:HB2	1:A:99:ASP:OD2	2.19	0.42
1:A:65:ALA:HB2	1:A:89:ASP:HB2	2.02	0.42
1:A:264:CYS:C	1:A:266:ASN:N	2.73	0.42
1:B:268:ASP:OD2	1:B:269:LYS:N	2.53	0.42
1:B:271:PRO:O	1:B:272:VAL:C	2.58	0.42
1:A:39:THR:N	1:A:58:LEU:O	2.52	0.42
1:B:253:LYS:HD2	1:B:254:ALA:N	2.35	0.42
1:B:223:TRP:O	1:B:264:CYS:HA	2.20	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:199:SER:HA	1:A:222:PRO:HD2	2.01	0.41
1:B:248:GLN:HG2	1:B:249:GLY:H	1.85	0.41
1:B:19:ILE:O	1:B:36:ARG:NH1	2.53	0.41
1:B:201:TYR:CE1	1:B:202:THR:HG22	2.56	0.41
1:B:76:LEU:HD23	1:B:98:LEU:CD1	2.51	0.41
1:A:111:LEU:HB3	1:A:115:LEU:HD12	2.02	0.41
1:B:221:ASN:HB2	1:B:223:TRP:CE2	2.55	0.41
1:B:243:VAL:HG11	1:B:259:VAL:HG23	2.03	0.41
1:A:270:PHE:HE2	1:A:275:TYR:HA	1.84	0.41
1:B:253:LYS:HD2	1:B:253:LYS:C	2.40	0.41
1:B:221:ASN:HA	1:B:222:PRO:HD3	1.81	0.41
1:A:25:VAL:O	1:A:26:ALA:HB3	2.20	0.41
1:B:53:HIS:CD2	1:B:77:ASN:ND2	2.88	0.41
1:B:263:GLN:HA	1:B:272:VAL:HG12	2.03	0.41
1:A:156:PRO:HA	1:A:157:PRO:HD3	1.75	0.41
1:A:159:LEU:HB3	1:A:160:LEU:HD22	2.01	0.41
1:A:225:CYS:HA	1:A:229:ILE:CB	2.49	0.41
1:B:194:LEU:HB3	1:B:196:GLN:NE2	2.36	0.41
1:A:102:HIS:HA	1:A:125:PHE:O	2.21	0.41
1:B:34:ASP:O	1:B:36:ARG:CD	2.68	0.41
1:B:214:PRO:O	1:B:242:ASN:HB3	2.20	0.41
1:A:252:VAL:HG12	1:A:252:VAL:O	2.20	0.41
1:A:266:ASN:O	1:A:267:SER:C	2.59	0.41
1:A:225:CYS:CB	1:A:229:ILE:HG21	2.43	0.41
1:A:136:LEU:HD12	1:A:159:LEU:HD22	2.04	0.40
1:B:123:VAL:O	1:B:123:VAL:HG12	2.21	0.40
1:A:121:LEU:HD11	1:A:123:VAL:CG1	2.51	0.40
1:A:34:ASP:OD1	1:A:55:SER:HB2	2.21	0.40
1:A:45:LEU:HA	1:A:45:LEU:HD13	1.93	0.40
1:B:187:LEU:CD1	1:B:187:LEU:N	2.85	0.40
1:A:193:LEU:HB3	1:A:213:LEU:HD21	2.03	0.40
1:A:52:LEU:HD11	1:A:54:LEU:HG	2.03	0.40
1:A:31:VAL:HG11	1:A:45:LEU:HD13	2.02	0.40
1:B:158:GLY:O	1:B:161:THR:HB	2.21	0.40
1:A:133:LEU:HA	1:A:133:LEU:HD22	1.95	0.40

All (4) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
6:A:2028:HOH:O	6:A:2028:HOH:O[12_565]	1.97	0.23
6:B:2057:HOH:O	6:B:2057:HOH:O[11_556]	2.05	0.15

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:39:THR:CG2	1:A:39:THR:CG2[9_555]	2.12	0.08
1:B:188:GLU:OE1	1:B:188:GLU:OE1[11_556]	2.18	0.02

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	264/281 (94%)	200 (76%)	40 (15%)	24 (9%)	1	2
1	B	274/281 (98%)	222 (81%)	40 (15%)	12 (4%)	4	12
All	All	538/562 (96%)	422 (78%)	80 (15%)	36 (7%)	2	5

All (36) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	21	GLU
1	A	28	HIS
1	A	48	ASP
1	A	92	LEU
1	A	248	GLN
1	A	278	LYS
1	B	42	PRO
1	B	90	GLY
1	B	283	LEU
1	B	285	ASP
1	B	286	GLU
1	A	35	LYS
1	A	47	LYS
1	A	91	THR
1	A	133	LEU
1	A	152	LEU
1	A	266	ASN
1	A	273	TYR
1	A	279	GLY
1	B	41	LEU
1	B	268	ASP

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Mol	Chain	Res	Type
1	B	273	TYR
1	A	19	ILE
1	A	43	PRO
1	A	46	PRO
1	A	135	ALA
1	B	223	TRP
1	A	45	LEU
1	A	90	GLY
1	A	102	HIS
1	A	187	LEU
1	A	226	ASN
1	B	35	LYS
1	A	268	ASP
1	B	276	PRO
1	B	281	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	236/249 (95%)	204 (86%)	32 (14%)	5	16
1	B	245/249 (98%)	208 (85%)	37 (15%)	4	12
All	All	481/498 (97%)	412 (86%)	69 (14%)	5	14

All (69) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	ASN
1	A	51	ILE
1	A	66	THR
1	A	73	LEU
1	A	74	THR
1	A	79	ASP
1	A	89	ASP
1	A	97	THR
1	A	105	LEU
1	A	113	GLN

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Mol	Chain	Res	Type
1	A	114	THR
1	A	118	LEU
1	A	120	VAL
1	A	133	LEU
1	A	139	LEU
1	A	141	GLU
1	A	142	LEU
1	A	151	GLU
1	A	159	LEU
1	A	165	LYS
1	A	166	LEU
1	A	178	GLU
1	A	187	LEU
1	A	196	GLN
1	A	226	ASN
1	A	243	VAL
1	A	256	THR
1	A	263	GLN
1	A	266	ASN
1	A	272	VAL
1	A	273	TYR
1	A	280	CYS
1	B	27	SER
1	B	36	ARG
1	B	39	THR
1	B	42	PRO
1	B	43	PRO
1	B	48	ASP
1	B	51	ILE
1	B	61	THR
1	B	73	LEU
1	B	118	LEU
1	B	127	ARG
1	B	130	SER
1	B	133	LEU
1	B	139	LEU
1	B	141	GLU
1	B	142	LEU
1	B	154	THR
1	B	159	LEU
1	B	160	LEU
1	B	165	LYS

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Mol	Chain	Res	Type
1	B	196	GLN
1	B	211	HIS
1	B	212	LEU
1	B	224	LEU
1	B	230	LEU
1	B	243	VAL
1	B	253	LYS
1	B	266	ASN
1	B	268	ASP
1	B	269	LYS
1	B	272	VAL
1	B	273	TYR
1	B	280	CYS
1	B	283	LEU
1	B	285	ASP
1	B	288	ASP
1	B	296	PRO

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	28	HIS
1	A	32	ASN
1	A	75	GLN
1	A	87	GLN
1	A	113	GLN
1	A	143	GLN
1	A	185	ASN
1	A	196	GLN
1	A	211	HIS
1	A	219	HIS
1	A	226	ASN
1	A	242	ASN
1	A	248	GLN
1	A	266	ASN
1	B	53	HIS
1	B	75	GLN
1	B	77	ASN
1	B	87	GLN
1	B	102	HIS
1	B	173	ASN
1	B	189	ASN

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Mol	Chain	Res	Type
1	B	196	GLN
1	B	219	HIS
1	B	237	GLN
1	B	239	ASN
1	B	263	GLN
1	B	266	ASN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

3 non-standard protein/DNA/RNA residues 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	$\# Z > 2$
1	TYS	B	292	1	16,16,17	5.06	4 (25%)	20,22,24	1.16	1 (5%)
1	TYS	B	294	1	16,16,17	5.11	4 (25%)	20,22,24	1.10	1 (5%)
1	TYS	B	295	1	16,16,17	5.14	4 (25%)	20,22,24	1.13	2 (10%)

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
1	TYS	B	292	1	-	0/9/11/13	0/1/1/1
1	TYS	B	294	1	-	0/9/11/13	0/1/1/1
1	TYS	B	295	1	-	0/9/11/13	0/1/1/1

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	295	TYS	O-C	18.78	1.24	1.11
1	B	292	TYS	O-C	17.83	1.23	1.11
1	B	294	TYS	O-C	17.79	1.23	1.11
1	B	294	TYS	OH-S	-8.77	1.47	1.63
1	B	292	TYS	OH-S	-8.41	1.48	1.63
1	B	295	TYS	OH-S	-6.77	1.51	1.63
1	B	294	TYS	OH-CZ	-3.64	1.38	1.42
1	B	292	TYS	OH-CZ	-3.29	1.38	1.42
1	B	295	TYS	CA-C	3.15	1.54	1.48
1	B	294	TYS	CA-C	2.85	1.53	1.48
1	B	292	TYS	CA-C	2.65	1.53	1.48
1	B	295	TYS	CE2-CD2	2.03	1.42	1.38

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	292	TYS	CZ-OH-S	4.78	126.15	118.98
1	B	294	TYS	C-CA-N	-4.14	109.69	113.83
1	B	295	TYS	C-CA-N	-3.86	109.97	113.83
1	B	295	TYS	CZ-OH-S	2.37	122.54	118.98

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 9 ligands modelled in this entry, 4 are monoatomic - leaving 5 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	NAG	A	1137	1	12,14,15	0.45	0	15,19,21	0.65	0
4	NAG	A	1175	1	12,14,15	0.48	0	15,19,21	0.54	0
3	ACY	A	510	-	3,3,3	1.11	0	3,3,3	1.43	0
4	NAG	A	600	1	12,14,15	0.53	0	15,19,21	0.58	0
5	SO4	B	504	-	4,4,4	0.28	0	6,6,6	0.07	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
4	NAG	A	1137	1	1/1/5/7	0/6/23/26	0/1/1/1
4	NAG	A	1175	1	-	1/6/23/26	0/1/1/1
3	ACY	A	510	-	-	0/0/0/0	0/0/0/0
4	NAG	A	600	1	-	0/6/23/26	0/1/1/1
5	SO4	B	504	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
4	A	1137	NAG	C1

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1175	NAG	O7-C7-N2-C2

There are no ring outliers.

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 ⓘ

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	266/281 (94%)	-0.13	6 (2%) 57 58	33, 58, 90, 105	0
1	B	279/281 (99%)	-0.27	13 (4%) 30 30	24, 44, 91, 115	0
All	All	545/562 (96%)	-0.20	19 (3%) 43 42	24, 51, 90, 115	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	250	VAL	6.1
1	B	251	ASP	5.9
1	B	249	GLY	5.0
1	B	247	LYS	3.4
1	B	285	ASP	3.3
1	A	267	SER	3.2
1	A	248	GLN	3.1
1	A	268	ASP	2.8
1	B	254	ALA	2.7
1	B	288	ASP	2.6
1	A	266	ASN	2.6
1	B	286	GLU	2.5
1	B	253	LYS	2.5
1	B	283	LEU	2.3
1	B	248	GLN	2.3
1	B	90	GLY	2.3
1	A	37	ASN	2.2
1	B	270	PHE	2.2
1	A	24	LYS	2.1

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

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
1	TYS	B	295	16/17	0.29	1.38	84,103,114,115	0
1	TYS	B	292	16/17	0.19	0.52	61,63,72,72	0
1	TYS	B	294	16/17	0.12	-1.02	74,77,79,79	0

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	ACY	A	510	4/4	0.36	6.65	91,92,92,92	0
2	PT	A	404	1/1	0.13	5.13	143,143,143,143	0
4	NAG	A	1175	14/15	0.21	1.86	83,87,88,89	0
4	NAG	A	1137	14/15	0.34	0.67	102,103,104,104	0
2	PT	A	402	1/1	0.17	0.17	117,117,117,117	0
4	NAG	A	600	14/15	0.13	-0.12	51,63,74,75	0
2	PT	A	403	1/1	0.15	-0.30	98,98,98,98	0
5	SO4	B	504	5/5	0.12	-0.33	70,70,70,71	0
2	PT	A	400	1/1	0.12	-1.43	112,112,112,112	0

6.5 Other polymers ⓘ

There are no such residues in this entry.