



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 04:39 PM GMT

PDB ID : 1A7L
Title : DOMINANT B-CELL EPITOPE FROM THE PRES2 REGION OF HEPATITIS B VIRUS IN THE FORM OF AN INSERTED PEPTIDE SEGMENT IN MALTODEXTRIN-BINDING PROTEIN
Authors : Saul, F.A.; Vulliez-Lenormand, B.; Lema, F.; Bentley, G.A.
Deposited on : 1998-03-16
Resolution : 2.90 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>

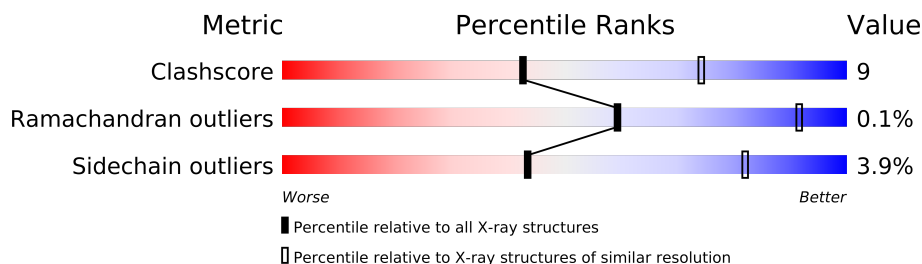
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 21963
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 2.90 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	79885	1326 (2.90-2.90)
Ramachandran outliers	78287	1290 (2.90-2.90)
Sidechain outliers	78261	1292 (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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	389	
1	B	389	
1	C	389	

2 Entry composition i

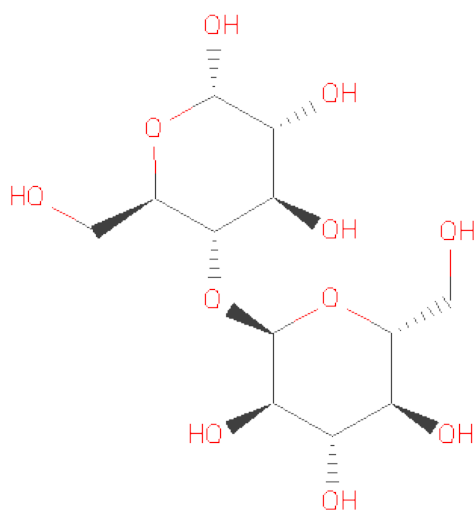
There are 3 unique types of molecules in this entry. The entry contains 8750 atoms, of which 0 are hydrogen and 0 are deuterium.

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 MALE-B363.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	380	Total	C	N	O	S	0	0	0
			2936	1888	480	560	8			
1	B	372	Total	C	N	O	S	0	0	0
			2893	1864	472	551	6			
1	C	362	Total	C	N	O	S	0	0	0
			2805	1807	455	537	6			

- Molecule 2 is SUGAR (MALTOSE) (three-letter code: MAL) (formula: C₁₂H₂₂O₁₁).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	C	1	Total	C	O	0	0
			22	11	11		
2	A	1	Total	C	O	0	0
			23	12	11		
2	B	1	Total	C	O	0	0
			23	12	11		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	20	Total 20	O 20	0	0
3	B	17	Total 17	O 17	0	0
3	C	11	Total 11	O 11	0	0

V343	V347	L361	S3651
			GLN
			ASP
			PRO
			ARG
			VAL
			ARG
			GLY
			LEU
			TYR
			PHE
			PRO
			ALA
			GLY
			GLY
			SER
			GLU
			CYS
			CYS
			GLU
			MET
			PRO
			ASP
			ALA
			ALA

4 Data and refinement statistics

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

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	64.60Å 71.30Å 123.20Å 90.00° 94.90° 90.00°	Depositor
Resolution (Å)	10.00 – 2.90	Depositor
% Data completeness (in resolution range)	98.0 (10.00-2.90)	Depositor
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.8	Depositor
R, R_{free}	0.192 , 0.286	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	8750	wwPDB-VP
Average B, all atoms (Å ²)	26.0	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: MAL

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	0.39	1/3009 (0.0%)	0.57	0/4085
1	B	0.36	0/2965	0.58	0/4025
1	C	0.34	0/2874	0.55	0/3902
All	All	0.37	1/8848 (0.0%)	0.57	0/12012

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	5	GLY	N-CA	-8.37	1.33	1.46

There are no bond angle outliers.

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	2936	0	2891	44	0
1	B	2893	0	2856	43	0
1	C	2805	0	2770	63	0
2	A	23	0	22	0	0
2	B	23	0	22	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	22	0	18	4	0
3	A	20	0	0	0	0
3	B	17	0	0	0	0
3	C	11	0	0	1	0
All	All	8750	0	8579	153	0

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 9.

All (153) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:79:ILE:HG22	1:A:81:PRO:HD3	1.44	0.96
1:B:79:ILE:HG22	1:B:81:PRO:HD3	1.45	0.95
1:B:321:MET:O	1:B:325:GLN:HG2	1.78	0.82
1:C:331:PRO:HG2	1:C:336:MET:SD	2.21	0.81
2:C:400:MAL:C5	2:C:400:MAL:C3	2.58	0.81
1:C:332:ASN:HD22	1:C:332:ASN:H	1.32	0.77
1:C:246:VAL:HA	1:C:323:ASN:HD21	1.51	0.75
1:C:178:ILE:HD12	1:C:178:ILE:H	1.51	0.74
1:B:192:LEU:HD23	1:B:357:VAL:HG13	1.69	0.73
1:A:137(E):ARG:HG3	1:A:137(E):ARG:HH11	1.53	0.73
1:C:64:HIS:HE1	1:C:260:GLY:HA2	1.56	0.70
1:C:258:PHE:HB3	1:C:330:MET:HE2	1.73	0.70
1:A:195:LEU:HA	1:A:198:LEU:HD12	1.74	0.70
1:B:34:LYS:HD2	1:B:34:LYS:H	1.57	0.69
1:A:178:ILE:H	1:A:178:ILE:HD13	1.58	0.69
1:A:192:LEU:HD23	1:A:357:VAL:HG13	1.75	0.69
1:C:5:GLY:H	1:C:272:ASN:HD21	1.41	0.67
1:A:249:THR:HG22	1:A:254:PRO:HA	1.77	0.66
1:C:122:LEU:HD21	1:C:126:PRO:HD3	1.78	0.66
1:C:6:LYS:HA	1:C:33:ILE:HG23	1.78	0.66
1:A:5:GLY:H	1:A:272:ASN:HD21	1.43	0.65
2:C:400:MAL:O4	2:C:400:MAL:C3	2.45	0.65
1:A:331:PRO:HG2	1:A:336:MET:SD	2.37	0.64
1:A:89:LEU:HD13	1:A:107:PRO:HG2	1.79	0.64
1:C:64:HIS:CE1	1:C:260:GLY:HA2	2.32	0.63
2:C:400:MAL:O4	2:C:400:MAL:C5	2.47	0.62
1:A:44:GLU:HG2	1:A:45:GLU:N	2.16	0.61
1:B:229:PRO:HA	1:B:232:TRP:CE2	2.36	0.60
1:A:192:LEU:O	1:A:196:VAL:HG23	2.02	0.60
1:C:31:THR:HB	1:C:33:ILE:HD13	1.84	0.59
1:A:218:ASN:HD21	1:A:235:ILE:HG12	1.67	0.59

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:258:PHE:HB3	1:C:330:MET:CE	2.32	0.59
1:C:139:LEU:HD23	1:C:142:LYS:HD3	1.83	0.59
1:C:192:LEU:O	1:C:196:VAL:HG23	2.03	0.58
1:B:133(E):ASP:HB3	1:B:136(E):VAL:HG23	1.86	0.58
1:C:170:LYS:O	1:C:176:TYR:HA	2.04	0.58
1:C:6:LYS:HB2	1:C:34:LYS:O	2.05	0.57
1:A:218:ASN:ND2	1:A:235:ILE:HA	2.20	0.57
1:B:11:ILE:HG21	1:B:17:TYR:HB3	1.87	0.57
1:C:79:ILE:HG22	1:C:81:PRO:HD3	1.85	0.56
1:B:178:ILE:HD12	1:B:179:LYS:NZ	2.20	0.56
1:C:246:VAL:HA	1:C:323:ASN:ND2	2.19	0.56
1:B:183:VAL:O	1:B:361:LEU:HD13	2.05	0.56
1:A:272:ASN:HD22	1:A:275:LEU:HD12	1.71	0.56
1:C:218:ASN:HB3	1:C:238:SER:OG	2.05	0.56
1:B:356:THR:HG22	1:B:358:ASP:H	1.72	0.55
1:B:11:ILE:CG2	1:B:17:TYR:HB3	2.37	0.54
1:A:137(E):ARG:HG3	1:A:137(E):ARG:NH1	2.23	0.54
1:C:174:GLY:O	1:C:175:LYS:HD2	2.07	0.54
1:A:50:VAL:HG23	1:A:51:ALA:N	2.23	0.54
1:C:249:THR:HG22	1:C:254:PRO:HA	1.89	0.54
1:C:229:PRO:HA	1:C:232:TRP:CE2	2.43	0.53
1:B:57:PRO:O	1:B:267:ASN:HB2	2.09	0.53
1:B:122:LEU:HD21	1:B:126:PRO:HD3	1.91	0.53
1:B:4:GLU:HA	1:B:272:ASN:HD21	1.74	0.52
1:A:133(E):ASP:O	1:A:137(E):ARG:HG2	2.10	0.52
1:C:48:PRO:HA	1:C:75:LEU:HD13	1.91	0.52
1:B:68:GLY:HA3	1:B:332:ASN:O	2.08	0.52
1:C:340:TRP:CE3	2:C:400:MAL:H61	2.45	0.52
1:C:332:ASN:HD22	1:C:332:ASN:N	2.03	0.52
1:A:6:LYS:HA	1:A:33:ILE:HG23	1.91	0.51
1:B:151:LEU:HD21	1:B:195:LEU:HD11	1.93	0.51
1:B:331:PRO:HG2	1:B:336:MET:SD	2.51	0.51
1:C:334:PRO:HB2	1:C:335:GLN:NE2	2.26	0.51
1:B:6:LYS:HA	1:B:33:ILE:HG23	1.92	0.51
1:C:217:PHE:HA	1:C:222:THR:HG22	1.93	0.50
1:B:44:GLU:HG2	1:B:45:GLU:HG3	1.93	0.50
1:C:176:TYR:CZ	1:C:331:PRO:HB3	2.47	0.50
1:A:219:LYS:HE2	1:A:219:LYS:HA	1.93	0.50
1:C:47:PHE:HA	1:C:50:VAL:HG22	1.93	0.50
1:B:179:LYS:HE3	1:B:179:LYS:HA	1.94	0.50
1:C:158:TRP:N	1:C:159:PRO:HD2	2.27	0.49
1:C:62:TRP:HB3	1:C:67:PHE:HE1	1.77	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:218:ASN:HD22	1:B:218:ASN:N	2.09	0.49
1:A:218:ASN:HD21	1:A:235:ILE:HA	1.77	0.49
1:C:11:ILE:CG2	1:C:17:TYR:HB3	2.43	0.49
1:A:143:GLY:O	1:A:144:LYS:HD2	2.12	0.49
1:B:27:PHE:HD1	1:B:283:TYR:CD2	2.30	0.49
1:B:11:ILE:O	1:B:39:HIS:HB3	2.13	0.49
1:A:278:GLU:O	1:A:282:ASN:HB2	2.13	0.48
1:A:129:TRP:NE1	1:A:248:PRO:HG2	2.28	0.48
1:C:85:PHE:CD1	1:C:88:LYS:HD3	2.49	0.48
1:C:68:GLY:HA3	1:C:332:ASN:O	2.13	0.47
1:C:81:PRO:HB3	1:C:85:PHE:HD2	1.79	0.47
1:A:108:ILE:HD13	1:A:285:LEU:HD21	1.96	0.47
1:A:7:LEU:HB2	1:A:35:VAL:HG22	1.95	0.47
1:C:53:THR:HG22	1:C:53:THR:O	2.15	0.47
1:C:5:GLY:N	1:C:272:ASN:HD21	2.11	0.47
1:A:132:ILE:N	1:A:133:PRO:HD2	2.30	0.47
1:C:126:PRO:HD2	1:C:224:MET:SD	2.55	0.47
1:C:51:ALA:HA	1:C:55:ASP:O	2.15	0.46
1:A:25:LYS:HA	1:A:25:LYS:HE3	1.96	0.46
1:A:89:LEU:CD1	1:A:107:PRO:HG2	2.45	0.46
1:B:258:PHE:CG	1:B:330:MET:HG2	2.50	0.46
1:C:85:PHE:HD1	1:C:88:LYS:HD3	1.80	0.46
1:B:116:ILE:HB	1:B:225:THR:CG2	2.45	0.46
1:B:136:ASP:O	1:B:140:LYS:HB2	2.15	0.46
1:A:68:GLY:HA3	1:A:332:ASN:O	2.15	0.46
1:A:272:ASN:ND2	1:A:275:LEU:HD12	2.30	0.46
1:C:256:LYS:HG2	1:C:326:LYS:O	2.16	0.46
1:C:291:GLU:HG2	1:C:295:LYS:HD3	1.97	0.45
1:A:67:PHE:HB3	1:A:104:ILE:HD12	1.98	0.45
1:A:143:GLY:C	1:A:144:LYS:HD2	2.36	0.45
1:A:122:LEU:HD21	1:A:126:PRO:HD3	1.98	0.45
1:C:40:PRO:HG2	1:C:43:LEU:HB3	1.99	0.45
1:C:20:LEU:HD23	1:C:37:VAL:HG11	1.99	0.45
1:B:179:LYS:HA	1:B:179:LYS:CE	2.47	0.45
1:C:215:ALA:O	1:C:219:LYS:HG3	2.16	0.45
1:A:381(I):GLU:O	1:A:382(I):CYS:HB2	2.17	0.45
1:C:113:LEU:HD11	1:C:156:PHE:HA	1.99	0.44
1:B:115:LEU:HB2	1:B:247:LEU:HD23	1.99	0.44
1:C:93:THR:HB	1:C:107:PRO:HB3	2.00	0.44
1:A:33:ILE:HD13	1:A:275:LEU:HD22	1.99	0.44
1:C:184:ASP:O	1:C:189:LYS:HE3	2.17	0.44
1:C:61:PHE:HA	1:C:263:SER:O	2.17	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:62:TRP:CD1	1:C:66:ARG:HG3	2.53	0.44
1:B:67:PHE:HB3	1:B:104:ILE:HD12	1.99	0.44
1:B:167:TYR:CE1	1:B:182:GLY:HA3	2.53	0.44
1:B:90:TYR:HA	1:B:91:PRO:HD3	1.87	0.43
1:C:270:SER:O	1:C:273:LYS:HG3	2.18	0.43
1:A:184:ASP:O	1:A:189:LYS:HE3	2.18	0.43
1:A:178:ILE:H	1:A:178:ILE:CD1	2.28	0.43
1:B:27:PHE:HD1	1:B:283:TYR:CE2	2.36	0.43
1:B:178:ILE:HD12	1:B:179:LYS:HZ2	1.82	0.43
1:B:219:LYS:HA	1:B:219:LYS:HD3	1.81	0.43
1:C:220:GLY:HA2	3:C:402:HOH:O	2.17	0.43
1:C:331:PRO:O	1:C:336:MET:HG3	2.19	0.43
1:B:61:PHE:HA	1:B:263:SER:O	2.19	0.43
1:B:167:TYR:CZ	1:B:170:LYS:HE3	2.54	0.43
1:A:183:VAL:O	1:A:361:LEU:HD13	2.18	0.43
1:A:246:VAL:HG12	1:A:247:LEU:O	2.19	0.42
1:C:48:PRO:HA	1:C:75:LEU:CD1	2.50	0.42
1:B:279:PHE:O	1:B:283:TYR:HB2	2.20	0.42
1:C:148:MET:HG2	1:C:213:ALA:HA	2.01	0.42
1:A:137(E):ARG:HH11	1:A:137(E):ARG:CG	2.29	0.41
1:B:14:ASP:OD1	1:B:15:LYS:HD2	2.20	0.41
1:B:168:ALA:O	1:B:181:VAL:HA	2.21	0.41
1:B:270:SER:O	1:B:273:LYS:HG3	2.20	0.41
1:B:184:ASP:O	1:B:189:LYS:HE3	2.21	0.41
1:C:61:PHE:O	1:C:62:TRP:HB2	2.20	0.41
1:C:198:LEU:O	1:C:203:HIS:HB2	2.20	0.41
1:A:21:ALA:O	1:A:25:LYS:HD2	2.21	0.41
1:C:183:VAL:O	1:C:361:LEU:HD13	2.21	0.41
1:B:128:THR:HG22	1:B:249:THR:OG1	2.20	0.41
1:A:198:LEU:O	1:A:203:HIS:HB2	2.20	0.41
1:C:106:TYR:CD2	1:C:280:LEU:HD13	2.55	0.41
1:C:195:LEU:HD12	1:C:198:LEU:HD12	2.04	0.40
1:C:12:ASN:HA	1:C:43:LEU:HD21	2.03	0.40
1:C:318:ALA:O	1:C:321:MET:HB2	2.21	0.40
1:C:343:VAL:O	1:C:347:VAL:HG23	2.22	0.40
1:B:34:LYS:CD	1:B:34:LYS:H	2.23	0.40
1:A:125:PRO:HA	1:A:126:PRO:HD3	1.90	0.40
1:A:307:TYR:O	1:A:311:LEU:HD13	2.22	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	378/389 (97%)	367 (97%)	11 (3%)	0	100	100
1	B	370/389 (95%)	358 (97%)	12 (3%)	0	100	100
1	C	360/389 (92%)	346 (96%)	13 (4%)	1 (0%)	50	85
All	All	1108/1167 (95%)	1071 (97%)	36 (3%)	1 (0%)	59	91

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	172	GLU

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	302/310 (97%)	291 (96%)	11 (4%)	47	85
1	B	298/310 (96%)	287 (96%)	11 (4%)	45	84
1	C	289/310 (93%)	276 (96%)	13 (4%)	38	77
All	All	889/930 (96%)	854 (96%)	35 (4%)	43	82

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	LYS
1	A	82	ASP
1	A	91	PRO
1	A	100	ASN
1	A	115	LEU

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Mol	Chain	Res	Type
1	A	178	ILE
1	A	202	LYS
1	A	205	ASN
1	A	258	PHE
1	A	309	GLU
1	A	329	ILE
1	B	4	GLU
1	B	6	LYS
1	B	34	LYS
1	B	160	LEU
1	B	179	LYS
1	B	185	ASN
1	B	218	ASN
1	B	241	ASN
1	B	258	PHE
1	B	309	GLU
1	B	321	MET
1	C	15	LYS
1	C	25	LYS
1	C	29	LYS
1	C	41	ASP
1	C	82	ASP
1	C	102	LYS
1	C	115	LEU
1	C	160	LEU
1	C	180	ASP
1	C	258	PHE
1	C	309	GLU
1	C	311	LEU
1	C	332	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	18	ASN
1	A	100	ASN
1	A	218	ASN
1	A	272	ASN
1	A	325	GLN
1	B	18	ASN
1	B	72	GLN
1	B	86	GLN

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Mol	Chain	Res	Type
1	B	100	ASN
1	B	152	GLN
1	B	185	ASN
1	B	218	ASN
1	B	241	ASN
1	B	272	ASN
1	B	325	GLN
1	C	272	ASN
1	C	332	ASN

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

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	$\# Z > 2$
2	MAL	A	400	-	24,24,24	0.75	0	35,35,35	0.66	0
2	MAL	B	400	-	24,24,24	0.86	1 (4%)	35,35,35	0.62	0
2	MAL	C	400	-	18,21,24	3.97	1 (5%)	22,28,35	0.61	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
2	MAL	A	400	-	1/1/10/10	0/8/48/48	0/2/2/2
2	MAL	B	400	-	1/1/10/10	0/8/48/48	0/2/2/2
2	MAL	C	400	-	1/1/7/10	0/12/36/48	0/1/1/2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	400	MAL	O3-C3	16.58	1.43	1.25
2	B	400	MAL	C4-C5	2.05	1.57	1.53

There are no bond angle outliers.

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	B	400	MAL	C1'
2	A	400	MAL	C1'
2	C	400	MAL	C1'

There are no torsion outliers.

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 ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.