



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 08:58 PM GMT

PDB ID : 1MZ6  
Title : Trypanosoma rangeli sialidase in complex with the inhibitor DANA  
Authors : Buschiazzo, A.; Tavares, G.A.; Campetella, O.; Spinelli, S.; Cremona, M.L.;  
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Deposited on : 2002-10-05  
Resolution : 2.90 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto: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.

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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) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

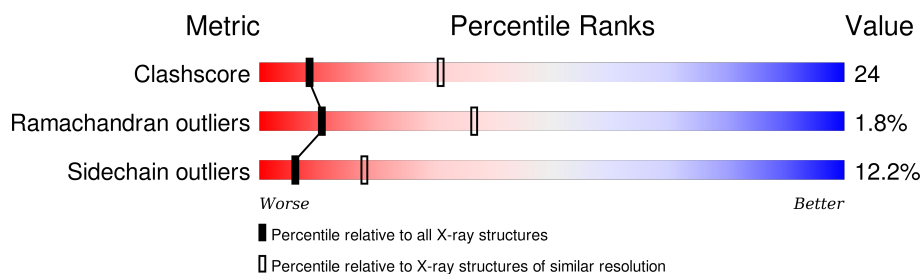
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 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	102246	1668 (2.90-2.90)
Ramachandran outliers	100387	1630 (2.90-2.90)
Sidechain outliers	100360	1632 (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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	638	

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 4889 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 sialidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	620	4759	3007	835	901	16	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	177	VAL	ILE	CONFLICT	UNP O44049

- Molecule 2 is 2,3-didehydro-2-deoxy-N-acetylneuraminic acid (DANA) (three-letter code: NAG) (formula:  $C_8H_{15}NO_6$ ).



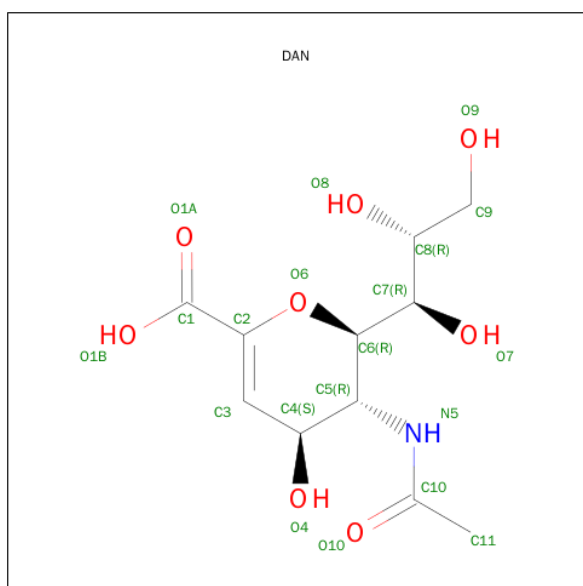
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	A	1	14	8	1	5	0	0
2	A	1	14	8	1	5	0	0

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

- Molecule 3 is 2-DEOXY-2,3-DEHYDRO-N-ACETYL-NEURAMINIC ACID (three-letter code: DAN) (formula:  $C_{11}H_{17}NO_8$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			20	11	1	8		

- Molecule 4 is water.

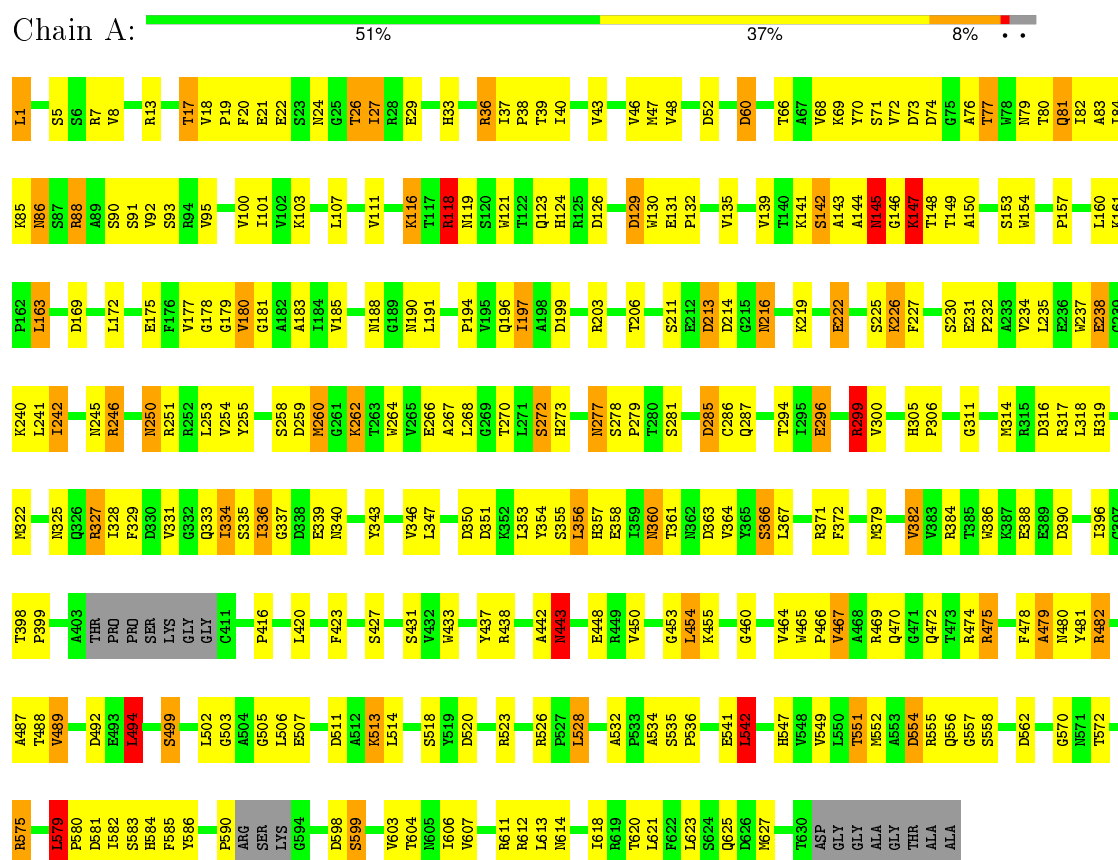
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	40	Total	O	0	0
			40	40		

### 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.

Note EDS was not executed.

- Molecule 1: sialidase



## 4 Data and refinement statistics

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

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	76.20 Å 93.80 Å 105.30 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 2.90	Depositor
% Data completeness (in resolution range)	99.3 (15.00-2.90)	Depositor
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.14	Depositor
Refinement program	REFMAC 5	Depositor
R, $R_{free}$	0.205 , 0.287	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4889	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	22.0	wwPDB-VP

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: DAN, 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	0.81	0/4865	1.25	38/6616 (0.6%)

There are no bond length outliers.

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	60	ASP	CB-CG-OD2	9.55	126.90	118.30
1	A	129	ASP	CB-CG-OD2	9.46	126.81	118.30
1	A	494	LEU	CA-CB-CG	8.90	135.77	115.30
1	A	351	ASP	CB-CG-OD2	8.26	125.73	118.30
1	A	7	ARG	NE-CZ-NH1	7.78	124.19	120.30
1	A	246	ARG	NE-CZ-NH2	-7.43	116.58	120.30
1	A	390	ASP	CB-CG-OD2	7.26	124.83	118.30
1	A	542	LEU	CA-CB-CG	7.26	132.00	115.30
1	A	213	ASP	CB-CG-OD2	7.16	124.75	118.30
1	A	7	ARG	NE-CZ-NH2	-7.14	116.73	120.30
1	A	350	ASP	CB-CG-OD2	7.02	124.62	118.30
1	A	492	ASP	CB-CG-OD2	6.96	124.56	118.30
1	A	554	ASP	CB-CG-OD2	6.82	124.44	118.30
1	A	299	ARG	NE-CZ-NH1	-6.53	117.04	120.30
1	A	36	ARG	NE-CZ-NH2	6.35	123.47	120.30
1	A	316	ASP	CB-CG-OD2	6.23	123.90	118.30
1	A	542	LEU	CB-CG-CD1	-6.17	100.50	111.00
1	A	285	ASP	CB-CG-OD2	6.07	123.76	118.30
1	A	129	ASP	CB-CG-OD1	-6.00	112.90	118.30
1	A	100	VAL	CB-CA-C	-5.95	100.10	111.40
1	A	242	ILE	CB-CA-C	5.74	123.09	111.60
1	A	454	LEU	CA-CB-CG	5.72	128.46	115.30
1	A	52	ASP	CB-CG-OD1	5.68	123.42	118.30
1	A	469	ARG	NE-CZ-NH1	-5.66	117.47	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	163	LEU	CA-CB-CG	5.64	128.28	115.30
1	A	579	LEU	CA-CB-CG	5.57	128.10	115.30
1	A	88	ARG	NE-CZ-NH1	-5.54	117.53	120.30
1	A	197	ILE	CB-CA-C	-5.50	100.60	111.60
1	A	575	ARG	NE-CZ-NH2	-5.47	117.56	120.30
1	A	489	VAL	CB-CA-C	-5.40	101.14	111.40
1	A	118	ARG	NE-CZ-NH1	5.39	123.00	120.30
1	A	443	ASN	CB-CA-C	5.25	120.90	110.40
1	A	118	ARG	NE-CZ-NH2	-5.24	117.68	120.30
1	A	259	ASP	CB-CG-OD2	5.21	122.99	118.30
1	A	562	ASP	CB-CG-OD2	5.18	122.96	118.30
1	A	598	ASP	CB-CG-OD2	5.12	122.91	118.30
1	A	475	ARG	NE-CZ-NH2	-5.03	117.79	120.30
1	A	185	VAL	CB-CA-C	-5.00	101.90	111.40

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	4759	0	4653	228	0
2	A	70	0	65	7	0
3	A	20	0	16	2	0
4	A	40	0	0	7	0
All	All	4889	0	4734	229	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 24.

All (229) 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:487:ALA:HB2	1:A:606:ILE:HD13	1.25	1.17
1:A:226:LYS:HD3	1:A:255:TYR:OH	1.52	1.09

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:196:GLN:NE2	1:A:231:GLU:H	1.58	1.02
1:A:317:ARG:HE	1:A:333:GLN:HE22	1.03	0.97
1:A:547:HIS:HD2	4:A:735:HOH:O	1.47	0.96
1:A:487:ALA:HB2	1:A:606:ILE:CD1	1.99	0.93
1:A:46:VAL:HG22	1:A:72:VAL:HG12	1.55	0.87
1:A:482:ARG:HG3	1:A:611:ARG:HA	1.57	0.86
1:A:206:THR:HG22	1:A:264:TRP:CH2	2.12	0.85
1:A:1:LEU:HD11	1:A:5:SER:HB3	1.59	0.85
1:A:360:ASN:HD21	1:A:363:ASP:H	1.18	0.84
1:A:360:ASN:HD22	1:A:360:ASN:C	1.80	0.84
1:A:360:ASN:HD22	1:A:361:THR:N	1.76	0.83
1:A:194:PRO:HG3	1:A:234:VAL:HG12	1.59	0.82
1:A:305:HIS:CD2	1:A:306:PRO:HD2	2.15	0.81
1:A:279:PRO:HA	1:A:472:GLN:HG3	1.65	0.79
1:A:396:ILE:HG22	1:A:627:MET:HE1	1.66	0.77
1:A:26:THR:HG22	2:A:655:NAG:C1	2.16	0.76
1:A:487:ALA:CB	1:A:606:ILE:HD13	2.09	0.76
1:A:489:VAL:HG12	1:A:603:VAL:HG22	1.66	0.76
1:A:250:ASN:HD22	1:A:251:ARG:H	1.31	0.75
1:A:206:THR:HG22	1:A:264:TRP:HH2	1.48	0.75
1:A:360:ASN:HD21	1:A:363:ASP:N	1.86	0.73
1:A:443:ASN:C	1:A:443:ASN:HD22	1.91	0.73
1:A:277:ASN:ND2	1:A:278:SER:OG	2.22	0.72
1:A:396:ILE:CG2	1:A:627:MET:HE1	2.19	0.72
1:A:196:GLN:NE2	1:A:231:GLU:N	2.37	0.72
1:A:17:THR:CG2	1:A:29:GLU:OE2	2.38	0.72
1:A:13:ARG:HG2	1:A:366:SER:OG	1.89	0.71
1:A:277:ASN:HD22	1:A:277:ASN:C	1.93	0.71
1:A:355:SER:HB2	1:A:372:PHE:CE1	2.26	0.70
1:A:296:GLU:OE1	1:A:384:ARG:HD3	1.91	0.70
1:A:433:TRP:HB3	1:A:442:ALA:HB3	1.73	0.70
1:A:146:GLY:O	1:A:147:LYS:HD3	1.92	0.70
1:A:335:SER:HB2	1:A:339:GLU:OE1	1.91	0.70
1:A:68:VAL:HG22	1:A:81:GLN:HG3	1.73	0.69
1:A:322:MET:HE3	1:A:379:MET:HB3	1.75	0.69
1:A:300:VAL:HG21	1:A:322:MET:HG3	1.72	0.69
1:A:237:TRP:CD1	1:A:238:GLU:HG3	2.29	0.68
1:A:360:ASN:ND2	1:A:360:ASN:C	2.47	0.68
1:A:336:ILE:N	1:A:339:GLU:OE1	2.26	0.68
1:A:254:VAL:HG23	1:A:272:SER:HA	1.76	0.68
1:A:199:ASP:OD2	1:A:203:ARG:HB2	1.93	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:547:HIS:CD2	4:A:735:HOH:O	2.32	0.67
1:A:314:MET:O	1:A:340:ASN:HB3	1.95	0.67
1:A:319:HIS:NE2	1:A:333:GLN:HG2	2.10	0.67
1:A:130:TRP:CZ3	1:A:161:LYS:HD2	2.30	0.66
1:A:26:THR:HG21	2:A:655:NAG:H3	1.76	0.66
1:A:230:SER:O	1:A:246:ARG:HB3	1.96	0.66
1:A:277:ASN:HD22	1:A:278:SER:N	1.95	0.65
1:A:464:VAL:HG13	1:A:584:HIS:CD2	2.30	0.65
1:A:551:THR:HG22	1:A:558:SER:HB2	1.80	0.64
1:A:334:ILE:HB	1:A:372:PHE:CE2	2.33	0.64
1:A:532:ALA:HB2	4:A:708:HOH:O	1.96	0.63
1:A:73:ASP:OD2	1:A:77:THR:HB	1.98	0.63
1:A:260:MET:HA	1:A:260:MET:HE2	1.79	0.63
1:A:396:ILE:CG2	1:A:627:MET:CE	2.77	0.63
1:A:26:THR:CG2	2:A:655:NAG:C1	2.76	0.63
1:A:250:ASN:HD22	1:A:251:ARG:N	1.98	0.62
1:A:505:GLY:HA2	1:A:513:LYS:HA	1.82	0.61
1:A:73:ASP:CG	1:A:76:ALA:HB3	2.21	0.61
1:A:38:PRO:HB2	1:A:356:LEU:HD13	1.82	0.61
1:A:206:THR:HB	1:A:225:SER:OG	2.01	0.61
1:A:612:ARG:HD2	4:A:714:HOH:O	2.00	0.61
1:A:66:THR:CB	1:A:84:ILE:HD12	2.32	0.60
1:A:82:ILE:O	1:A:82:ILE:HG22	2.00	0.60
1:A:194:PRO:HG3	1:A:234:VAL:CG1	2.31	0.60
1:A:196:GLN:HE22	1:A:231:GLU:HG3	1.66	0.60
1:A:481:TYR:CD2	1:A:482:ARG:HD3	2.37	0.59
1:A:95:VAL:HB	1:A:111:VAL:HG21	1.83	0.59
1:A:38:PRO:CB	1:A:356:LEU:HD13	2.32	0.59
1:A:480:ASN:OD1	1:A:582:ILE:HG13	2.02	0.59
1:A:520:ASP:OD2	1:A:526:ARG:NE	2.35	0.58
1:A:141:LYS:HA	1:A:150:ALA:HB2	1.86	0.58
1:A:611:ARG:NH1	1:A:613:LEU:HD23	2.20	0.57
1:A:172:LEU:HD23	1:A:172:LEU:N	2.18	0.57
1:A:277:ASN:ND2	1:A:277:ASN:C	2.58	0.57
1:A:251:ARG:NH1	1:A:474:ARG:HG3	2.19	0.57
1:A:487:ALA:CB	1:A:606:ILE:CD1	2.78	0.56
1:A:552:MET:HG2	1:A:556:GLN:O	2.04	0.56
1:A:18:VAL:HB	1:A:82:ILE:HD11	1.87	0.56
1:A:131:GLU:HG3	2:A:653:NAG:H83	1.87	0.56
1:A:69:LYS:HA	1:A:79:ASN:O	2.06	0.56
1:A:478:PHE:O	1:A:480:ASN:N	2.38	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:423:PHE:CD2	1:A:607:VAL:HG22	2.40	0.56
1:A:86:ASN:C	1:A:86:ASN:HD22	2.08	0.56
1:A:222:GLU:HB3	1:A:262:LYS:O	2.07	0.55
1:A:357:HIS:HD2	1:A:358:GLU:O	1.90	0.55
1:A:466:PRO:HB2	1:A:470:GLN:HG3	1.88	0.55
1:A:277:ASN:ND2	1:A:278:SER:N	2.54	0.55
1:A:327:ARG:NH2	1:A:437:TYR:CE2	2.75	0.55
1:A:226:LYS:CD	1:A:255:TYR:OH	2.40	0.55
1:A:317:ARG:HE	1:A:333:GLN:NE2	1.88	0.54
1:A:467:VAL:O	1:A:470:GLN:HB2	2.06	0.54
1:A:135:VAL:HG22	1:A:157:PRO:HA	1.88	0.54
1:A:86:ASN:HD21	1:A:88:ARG:HD3	1.72	0.54
1:A:478:PHE:CD2	1:A:479:ALA:N	2.76	0.54
1:A:541:GLU:O	1:A:542:LEU:C	2.46	0.53
1:A:69:LYS:HG2	1:A:80:THR:HG23	1.91	0.53
1:A:488:THR:HG23	1:A:625:GLN:NE2	2.24	0.53
1:A:287:GLN:HB3	1:A:343:TYR:CG	2.43	0.53
1:A:503:GLY:O	1:A:585:PHE:HA	2.10	0.51
1:A:206:THR:CG2	1:A:264:TRP:HH2	2.20	0.51
1:A:382:VAL:HG12	1:A:438:ARG:CZ	2.40	0.51
3:A:700:DAN:H92	4:A:725:HOH:O	2.10	0.51
1:A:478:PHE:O	1:A:479:ALA:C	2.50	0.51
1:A:179:GLY:O	1:A:180:VAL:HB	2.11	0.51
1:A:203:ARG:HD2	1:A:227:PHE:CD2	2.47	0.50
1:A:287:GLN:HB3	1:A:343:TYR:CD1	2.46	0.50
1:A:40:ILE:HG23	1:A:40:ILE:O	2.12	0.50
1:A:354:TYR:HE2	1:A:371:ARG:NH1	2.10	0.49
1:A:306:PRO:HA	1:A:318:LEU:HA	1.94	0.49
1:A:40:ILE:HD13	1:A:347:LEU:HD23	1.92	0.49
1:A:194:PRO:HB2	1:A:232:PRO:HG2	1.95	0.49
1:A:614:ASN:O	1:A:618:ILE:HG13	2.13	0.49
1:A:317:ARG:NE	1:A:333:GLN:HE22	1.89	0.48
1:A:142:SER:O	1:A:149:THR:O	2.30	0.48
1:A:19:PRO:HB2	1:A:27:ILE:HD12	1.94	0.48
1:A:396:ILE:HG21	1:A:627:MET:HE3	1.95	0.48
1:A:467:VAL:HG23	1:A:584:HIS:HA	1.96	0.48
1:A:506:LEU:HD23	1:A:514:LEU:N	2.28	0.48
1:A:130:TRP:CG	1:A:131:GLU:N	2.81	0.48
1:A:24:ASN:C	1:A:24:ASN:OD1	2.50	0.48
1:A:268:LEU:HD13	1:A:579:LEU:HD21	1.93	0.48
1:A:103:LYS:NZ	1:A:214:ASP:OD2	2.44	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:334:ILE:HB	1:A:372:PHE:CZ	2.48	0.48
1:A:241:LEU:HD11	1:A:260:MET:HE2	1.96	0.48
1:A:130:TRP:CZ2	1:A:132:PRO:HD3	2.49	0.48
1:A:353:LEU:HG	1:A:372:PHE:HD1	1.78	0.48
1:A:47:MET:N	1:A:71:SER:O	2.45	0.48
1:A:305:HIS:HD2	1:A:306:PRO:O	1.97	0.48
1:A:191:LEU:HB2	1:A:211:SER:HB3	1.96	0.48
1:A:528:LEU:HD23	1:A:534:ALA:HB2	1.95	0.48
1:A:213:ASP:CG	1:A:216:ASN:HB3	2.35	0.47
1:A:499:SER:HB3	1:A:599:SER:OG	2.14	0.47
1:A:494:LEU:HD11	1:A:523:ARG:HD3	1.96	0.47
1:A:47:MET:O	1:A:70:TYR:HA	2.14	0.47
1:A:277:ASN:HD22	1:A:278:SER:CB	2.27	0.47
1:A:355:SER:HB2	1:A:372:PHE:HE1	1.73	0.47
1:A:294:THR:OG1	1:A:299:ARG:NH1	2.45	0.47
1:A:506:LEU:HD21	1:A:514:LEU:HB2	1.96	0.47
1:A:360:ASN:ND2	1:A:363:ASP:H	1.97	0.47
1:A:82:ILE:HA	1:A:82:ILE:HD13	1.68	0.47
1:A:549:VAL:HG21	1:A:618:ILE:HG23	1.95	0.47
1:A:143:ALA:HA	1:A:148:THR:HA	1.96	0.47
1:A:416:PRO:HD3	1:A:620:THR:HG21	1.97	0.47
1:A:416:PRO:HD2	1:A:621:LEU:HD21	1.97	0.47
1:A:357:HIS:CD2	1:A:358:GLU:O	2.68	0.47
1:A:107:LEU:HD13	1:A:154:TRP:CH2	2.50	0.47
1:A:336:ILE:O	1:A:339:GLU:HG2	2.16	0.46
1:A:26:THR:HG21	2:A:655:NAG:H5	1.97	0.46
1:A:43:VAL:HG11	1:A:139:VAL:HB	1.97	0.46
1:A:251:ARG:CZ	1:A:474:ARG:HD2	2.45	0.46
1:A:39:THR:OG1	1:A:181:GLY:HA2	2.16	0.46
1:A:579:LEU:HB3	1:A:580:PRO:HD2	1.98	0.46
1:A:33:HIS:CE1	1:A:364:VAL:HG13	2.50	0.46
1:A:322:MET:HE3	1:A:379:MET:CE	2.45	0.46
1:A:237:TRP:O	1:A:240:LYS:N	2.33	0.46
1:A:266:GLU:HG2	1:A:268:LEU:HD23	1.97	0.46
1:A:260:MET:HG3	1:A:260:MET:O	2.16	0.46
1:A:40:ILE:HA	1:A:48:VAL:O	2.16	0.46
1:A:507:GLU:HG3	1:A:581:ASP:OD2	2.16	0.46
1:A:557:GLY:O	1:A:570:GLY:N	2.39	0.46
1:A:206:THR:HG21	1:A:245:ASN:OD1	2.16	0.45
1:A:305:HIS:CD2	1:A:306:PRO:CD	2.93	0.45
1:A:431:SER:OG	2:A:652:NAG:O7	2.30	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1:LEU:CD1	1:A:5:SER:HB3	2.40	0.45
1:A:234:VAL:HG22	1:A:235:LEU:N	2.31	0.45
1:A:199:ASP:OD1	1:A:199:ASP:C	2.55	0.45
1:A:131:GLU:CG	2:A:653:NAG:H83	2.46	0.45
1:A:141:LYS:HA	1:A:150:ALA:CB	2.45	0.45
1:A:143:ALA:CB	1:A:148:THR:HA	2.46	0.45
1:A:453:GLY:HA3	1:A:604:THR:HG22	1.98	0.45
1:A:443:ASN:C	1:A:443:ASN:ND2	2.64	0.44
1:A:38:PRO:HB3	1:A:367:LEU:HD13	1.99	0.44
1:A:526:ARG:CZ	1:A:536:PRO:HG3	2.47	0.44
1:A:253:LEU:HD23	1:A:273:HIS:CD2	2.51	0.44
1:A:583:SER:OG	1:A:584:HIS:ND1	2.41	0.44
1:A:416:PRO:HD3	1:A:620:THR:CG2	2.48	0.44
1:A:92:VAL:O	1:A:93:SER:C	2.55	0.44
1:A:242:ILE:HG21	1:A:242:ILE:HD13	1.77	0.44
1:A:123:GLN:CB	4:A:734:HOH:O	2.66	0.44
1:A:119:ASN:O	1:A:124:HIS:HE1	2.01	0.44
1:A:285:ASP:OD1	1:A:286:CYS:N	2.51	0.44
1:A:328:ILE:O	1:A:475:ARG:HD2	2.18	0.44
1:A:47:MET:CE	1:A:74:ASP:HA	2.49	0.43
1:A:17:THR:HG21	1:A:29:GLU:OE2	2.13	0.43
1:A:36:ARG:NE	1:A:358:GLU:OE1	2.39	0.43
1:A:92:VAL:CG2	1:A:116:LYS:HA	2.47	0.43
1:A:250:ASN:ND2	1:A:251:ARG:H	2.09	0.43
1:A:488:THR:OG1	1:A:625:GLN:NE2	2.46	0.43
1:A:213:ASP:OD2	1:A:216:ASN:HB3	2.17	0.43
1:A:81:GLN:O	1:A:82:ILE:HD13	2.19	0.43
1:A:20:PHE:CZ	1:A:88:ARG:NH2	2.87	0.43
1:A:460:GLY:O	1:A:590:PRO:HA	2.18	0.43
1:A:623:LEU:HD23	1:A:623:LEU:HA	1.74	0.43
1:A:499:SER:O	1:A:518:SER:HB3	2.19	0.43
1:A:68:VAL:HG13	1:A:83:ALA:HB2	2.01	0.42
1:A:450:VAL:HG22	1:A:453:GLY:O	2.19	0.42
1:A:177:VAL:HG12	1:A:178:GLY:O	2.20	0.42
1:A:322:MET:HE3	1:A:379:MET:CB	2.46	0.42
1:A:329:PHE:CE2	1:A:331:VAL:HA	2.54	0.42
1:A:420:LEU:C	1:A:420:LEU:HD23	2.39	0.42
1:A:360:ASN:HD21	1:A:363:ASP:CA	2.33	0.41
1:A:90:SER:OG	1:A:92:VAL:HG22	2.19	0.41
1:A:334:ILE:HB	1:A:372:PHE:HE2	1.81	0.41
1:A:555:ARG:HH11	1:A:575:ARG:HA	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:502:LEU:HD12	1:A:586:TYR:O	2.20	0.41
1:A:66:THR:HB	1:A:84:ILE:HD12	2.02	0.41
1:A:311:GLY:O	1:A:314:MET:HG2	2.21	0.41
1:A:121:TRP:CZ2	1:A:196:GLN:HG3	2.55	0.41
1:A:396:ILE:CG2	1:A:627:MET:HE3	2.51	0.41
1:A:506:LEU:CD2	1:A:514:LEU:HB2	2.50	0.41
1:A:511:ASP:N	1:A:511:ASP:OD2	2.53	0.41
1:A:188:ASN:HD21	1:A:190:ASN:ND2	2.17	0.41
1:A:329:PHE:HE2	1:A:331:VAL:HA	1.86	0.41
1:A:145:ASN:N	1:A:145:ASN:ND2	2.67	0.41
1:A:101:ILE:HG12	1:A:183:ALA:HB3	2.03	0.41
1:A:398:THR:HB	1:A:399:PRO:CD	2.51	0.41
1:A:130:TRP:CD2	1:A:175:GLU:HB3	2.57	0.40
1:A:90:SER:OG	1:A:91:SER:N	2.52	0.40
1:A:17:THR:HG23	1:A:29:GLU:OE2	2.15	0.40
1:A:532:ALA:CB	4:A:708:HOH:O	2.65	0.40
1:A:386:TRP:CD1	1:A:437:TYR:HD2	2.40	0.40
1:A:71:SER:OG	1:A:72:VAL:N	2.54	0.40
1:A:384:ARG:O	1:A:388:GLU:HG3	2.21	0.40
1:A:37:ILE:N	1:A:38:PRO:CD	2.84	0.40
1:A:60:ASP:OD1	3:A:700:DAN:H5	2.20	0.40
1:A:118:ARG:HG2	1:A:118:ARG:HH11	1.85	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	614/638 (96%)	544 (89%)	59 (10%)	11 (2%)	11 37

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	145	ASN
1	A	180	VAL
1	A	479	ALA
1	A	542	LEU
1	A	554	ASP
1	A	147	LYS
1	A	144	ALA
1	A	267	ALA
1	A	337	GLY
1	A	238	GLU
1	A	467	VAL

### 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	510/530 (96%)	448 (88%)	62 (12%)	6 18

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	LEU
1	A	8	VAL
1	A	17	THR
1	A	21	GLU
1	A	22	GLU
1	A	26	THR
1	A	27	ILE
1	A	77	THR
1	A	81	GLN
1	A	85	LYS
1	A	86	ASN
1	A	116	LYS
1	A	118	ARG
1	A	126	ASP
1	A	129	ASP
1	A	142	SER

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Mol	Chain	Res	Type
1	A	145	ASN
1	A	147	LYS
1	A	153	SER
1	A	160	LEU
1	A	163	LEU
1	A	169	ASP
1	A	197	ILE
1	A	216	ASN
1	A	219	LYS
1	A	222	GLU
1	A	226	LYS
1	A	250	ASN
1	A	258	SER
1	A	260	MET
1	A	262	LYS
1	A	270	THR
1	A	272	SER
1	A	277	ASN
1	A	281	SER
1	A	296	GLU
1	A	299	ARG
1	A	325	ASN
1	A	327	ARG
1	A	334	ILE
1	A	336	ILE
1	A	346	VAL
1	A	356	LEU
1	A	360	ASN
1	A	366	SER
1	A	382	VAL
1	A	427	SER
1	A	443	ASN
1	A	448	GLU
1	A	454	LEU
1	A	455	LYS
1	A	465	TRP
1	A	482	ARG
1	A	494	LEU
1	A	499	SER
1	A	513	LYS
1	A	528	LEU
1	A	535	SER

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Mol	Chain	Res	Type
1	A	551	THR
1	A	572	THR
1	A	579	LEU
1	A	599	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	86	ASN
1	A	124	HIS
1	A	145	ASN
1	A	196	GLN
1	A	250	ASN
1	A	277	ASN
1	A	305	HIS
1	A	333	GLN
1	A	357	HIS
1	A	360	ASN
1	A	377	GLN
1	A	443	ASN
1	A	472	GLN
1	A	547	HIS
1	A	605	ASN
1	A	625	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry

6 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	NAG	A	651	1	14,14,15	0.89	1 (7%)	15,19,21	1.19	1 (6%)
2	NAG	A	652	1	14,14,15	0.64	0	15,19,21	1.48	3 (20%)
2	NAG	A	653	1	14,14,15	0.85	1 (7%)	15,19,21	2.72	5 (33%)
2	NAG	A	654	1	14,14,15	0.65	0	15,19,21	1.83	3 (20%)
2	NAG	A	655	1	14,14,15	0.89	0	15,19,21	2.83	7 (46%)
3	DAN	A	700	-	16,20,20	3.82	4 (25%)	19,28,28	3.07	7 (36%)

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	NAG	A	651	1	-	0/6/23/26	0/1/1/1
2	NAG	A	652	1	-	0/6/23/26	0/1/1/1
2	NAG	A	653	1	-	0/6/23/26	0/1/1/1
2	NAG	A	654	1	-	0/6/23/26	0/1/1/1
2	NAG	A	655	1	-	0/6/23/26	0/1/1/1
3	DAN	A	700	-	-	0/14/34/34	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	700	DAN	O6-C2	-5.33	1.27	1.37
2	A	651	NAG	O5-C1	-2.34	1.39	1.43
3	A	700	DAN	O6-C6	-2.13	1.42	1.46
2	A	653	NAG	C1-C2	2.63	1.56	1.52
3	A	700	DAN	O10-C10	6.69	1.38	1.23
3	A	700	DAN	C3-C2	12.18	1.49	1.32

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	700	DAN	C4-C3-C2	-7.04	109.69	121.60
2	A	655	NAG	C4-C3-C2	-6.50	101.13	111.23
2	A	653	NAG	C4-C3-C2	-6.24	101.54	111.23
2	A	653	NAG	C3-C4-C5	-5.98	99.77	110.20
2	A	654	NAG	C2-N2-C7	-5.11	116.47	123.04
2	A	655	NAG	C3-C4-C5	-4.63	102.12	110.20
3	A	700	DAN	O7-C7-C6	-4.48	99.07	109.43
3	A	700	DAN	C6-C5-C4	-4.13	103.56	111.09
2	A	655	NAG	C2-N2-C7	-2.40	119.96	123.04
2	A	652	NAG	C4-C3-C2	-2.16	107.86	111.23
2	A	653	NAG	O5-C5-C6	2.03	111.73	107.35
2	A	652	NAG	O3-C3-C2	2.15	113.37	109.11
2	A	655	NAG	O3-C3-C2	2.50	114.06	109.11
2	A	653	NAG	O3-C3-C2	2.55	114.17	109.11
2	A	651	NAG	C3-C4-C5	2.73	114.96	110.20
2	A	654	NAG	C4-C3-C2	2.91	115.75	111.23
2	A	655	NAG	C6-C5-C4	2.96	120.31	113.02
2	A	654	NAG	C3-C4-C5	3.08	115.56	110.20
3	A	700	DAN	O9-C9-C8	3.32	118.31	111.10
2	A	655	NAG	C1-O5-C5	3.49	116.68	112.25
2	A	653	NAG	O4-C4-C5	3.63	118.86	109.24
2	A	652	NAG	C1-O5-C5	3.82	117.10	112.25
2	A	655	NAG	O3-C3-C4	3.84	118.99	110.34
3	A	700	DAN	C6-C5-N5	3.87	117.82	111.07
3	A	700	DAN	C9-C8-C7	4.54	123.12	112.48
3	A	700	DAN	C6-O6-C2	5.66	123.56	114.79

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	652	NAG	1	0
2	A	653	NAG	2	0
2	A	655	NAG	4	0
3	A	700	DAN	2	0

## 5.7 Other polymers

There are no such residues in this entry.

## 5.8 Polymer linkage issues

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section 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.