



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:38 AM GMT

PDB ID : 2DSV  
Title : Interactions of protective signalling factor with chitin-like polysaccharide: Crystal structure of the complex between signalling protein from sheep (SPS-40) and a hexasaccharide at 2.5Å resolution  
Authors : Srivastava, D.B.; Ethayathulla, A.S.; Kumar, J.; Singh, N.; Sharma, S.; Bhushan, A.; Srinivasan, A.; Singh, T.P.  
Deposited on : 2006-07-07  
Resolution : 2.54 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) : 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

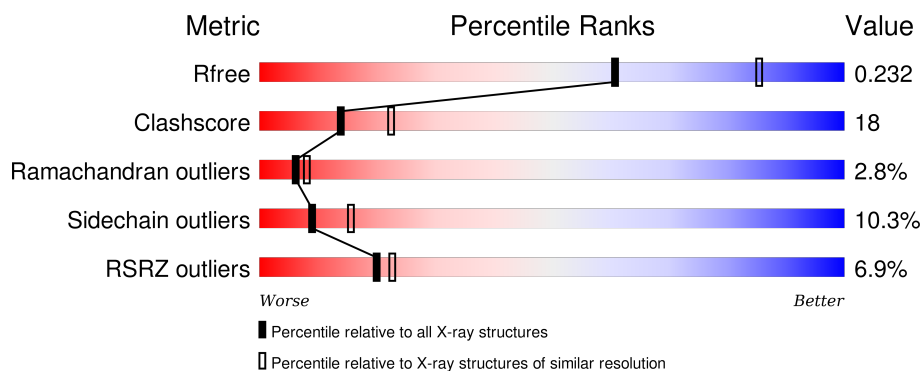
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.54 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	4549 (2.58-2.50)
Clashscore	102246	5292 (2.58-2.50)
Ramachandran outliers	100387	5194 (2.58-2.50)
Sidechain outliers	100360	5196 (2.58-2.50)
RSRZ outliers	91569	4561 (2.58-2.50)

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.

Mol	Chain	Length	Quality of chain
1	A	361	

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
2	NAG	A	363	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	NAG	A	364	X	-	-	-
2	MAN	A	365	X	-	-	-
2	MAN	A	366	X	-	-	-
3	NAG	A	367	X	-	X	X
3	NAG	A	369	X	-	-	-

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3143 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 Chitinase-3-like protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	361	Total	C	N	O	S	0	0	0
			2869	1832	499	529	9			

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

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	4	Total	C	N	O	0	0
			50	28	2	20		

- Molecule 3 is a polymer of unknown type called SUGAR (6-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	6	Total	C	N	O	0	0
			85	48	6	31		

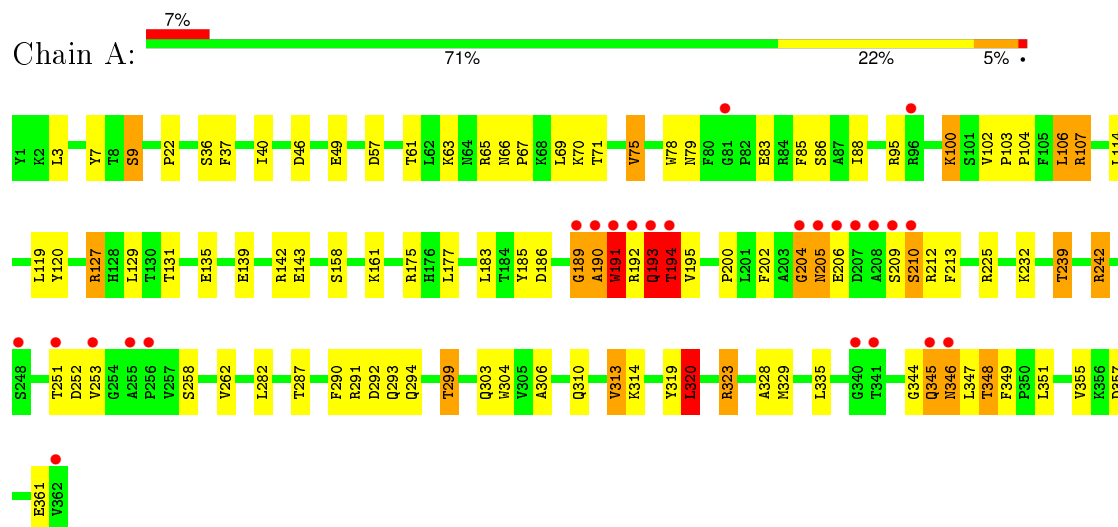
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	139	Total	O	0	0
			139	139		

### 3 Residue-property plots [i](#)

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: Chitinase-3-like protein 1



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	62.77Å 66.31Å 107.12Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	56.00 – 2.54 41.95 – 2.54	Depositor EDS
% Data completeness (in resolution range)	94.8 (56.00-2.54) 94.5 (41.95-2.54)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.12	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.40 (at 2.54Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.187 , 0.227 0.198 , 0.232	Depositor DCC
$R_{free}$ test set	314 reflections (2.17%)	DCC
Wilson B-factor (Å <sup>2</sup> )	29.7	Xtriage
Anisotropy	0.100	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 70.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 14469 reflections	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	3143	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.67% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NAG, MAN

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	0/2945	0.75	2/3995 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	A	3	0
3	A	2	0
All	All	5	0

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	A	320	LEU	CA-CB-CG	7.03	131.48	115.30
1	A	191	TRP	CA-CB-CG	6.73	126.49	113.70

All (5) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	364	NAG	C1
2	A	365	MAN	C1
2	A	366	MAN	C1
3	A	367	NAG	C1
3	A	369	NAG	C1

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	2869	0	2794	96	0
2	A	50	0	43	4	0
3	A	85	0	75	19	0
4	A	139	0	0	18	0
All	All	3143	0	2912	109	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 18.

All (109) 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:323:ARG:HG2	1:A:323:ARG:HH11	1.40	0.84
2:A:363:NAG:H83	2:A:363:NAG:H3	1.64	0.79
1:A:292:ASP:HB2	4:A:505:HOH:O	1.84	0.77
1:A:103:PRO:HB2	1:A:104:PRO:HD3	1.71	0.73
1:A:120:TYR:OH	3:A:367:NAG:H61	1.88	0.73
1:A:204:GLY:CA	1:A:292:ASP:HB3	2.21	0.71
1:A:323:ARG:NH1	1:A:323:ARG:HG2	2.04	0.71
1:A:78:TRP:CE2	3:A:367:NAG:O1	2.45	0.70
1:A:319:TYR:CZ	1:A:323:ARG:HD2	2.27	0.68
1:A:225:ARG:HG2	4:A:511:HOH:O	1.92	0.68
1:A:262:VAL:H	1:A:303:GLN:HE22	1.40	0.68
3:A:367:NAG:O4	3:A:368:NAG:H4	1.94	0.67
1:A:78:TRP:HZ3	1:A:120:TYR:CE1	2.12	0.67
1:A:262:VAL:H	1:A:303:GLN:NE2	1.93	0.67
1:A:183:LEU:HD21	3:A:368:NAG:H82	1.76	0.66
1:A:190:ALA:O	1:A:191:TRP:C	2.33	0.66
3:A:367:NAG:O3	3:A:368:NAG:H4	1.96	0.65
1:A:293:GLN:HG3	4:A:505:HOH:O	1.95	0.65
1:A:57:ASP:O	1:A:61:THR:HG23	1.95	0.65
2:A:365:MAN:H5	2:A:366:MAN:O2	1.97	0.65
1:A:120:TYR:CE2	1:A:158:SER:HB2	2.32	0.64
1:A:313:VAL:HG13	1:A:355:VAL:CG2	2.28	0.64
1:A:204:GLY:HA2	1:A:292:ASP:HB3	1.79	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:253:VAL:HG13	4:A:492:HOH:O	1.99	0.62
3:A:369:NAG:H62	3:A:370:NAG:C7	2.30	0.61
1:A:323:ARG:CG	1:A:323:ARG:HH11	2.10	0.61
1:A:204:GLY:HA3	1:A:292:ASP:HB3	1.83	0.60
2:A:363:NAG:H3	2:A:363:NAG:C8	2.32	0.60
1:A:189:GLY:O	1:A:191:TRP:N	2.36	0.59
1:A:239:THR:CG2	1:A:239:THR:O	2.50	0.58
1:A:239:THR:HG22	1:A:335:LEU:HB2	1.85	0.58
1:A:131:THR:HG23	4:A:506:HOH:O	2.03	0.58
3:A:367:NAG:H4	3:A:368:NAG:O6	2.05	0.57
1:A:78:TRP:CG	3:A:367:NAG:H3	2.40	0.56
1:A:177:LEU:O	1:A:232:LYS:HE2	2.06	0.56
1:A:40:ILE:HG21	1:A:88:ILE:CD1	2.36	0.56
1:A:40:ILE:HG21	1:A:88:ILE:HD13	1.89	0.55
1:A:120:TYR:OH	3:A:367:NAG:C6	2.53	0.55
1:A:348:THR:HG23	1:A:349:PHE:HD2	1.71	0.54
1:A:323:ARG:NH1	4:A:459:HOH:O	2.40	0.54
1:A:78:TRP:CZ2	3:A:367:NAG:O1	2.60	0.54
1:A:79:ASN:HB2	4:A:495:HOH:O	2.07	0.54
1:A:183:LEU:CD2	3:A:368:NAG:H82	2.38	0.53
1:A:310:GLN:O	1:A:314:LYS:HG3	2.09	0.53
1:A:186:ASP:OD1	3:A:367:NAG:H4	2.09	0.53
1:A:313:VAL:HG13	1:A:355:VAL:HG23	1.89	0.53
1:A:239:THR:HG23	1:A:239:THR:O	2.09	0.53
3:A:367:NAG:C1	4:A:400:HOH:O	2.57	0.52
1:A:205:ASN:HD22	1:A:206:GLU:N	2.07	0.52
1:A:313:VAL:CG1	1:A:355:VAL:HG23	2.40	0.52
3:A:370:NAG:H62	3:A:371:NAG:O7	2.12	0.50
1:A:323:ARG:NH2	4:A:406:HOH:O	2.45	0.50
1:A:239:THR:HG22	1:A:335:LEU:CB	2.42	0.50
1:A:107:ARG:NH1	1:A:143:GLU:OE1	2.45	0.49
3:A:367:NAG:H1	4:A:400:HOH:O	2.12	0.49
1:A:131:THR:HG22	1:A:135:GLU:OE2	2.13	0.49
1:A:194:THR:HG23	1:A:195:VAL:N	2.29	0.47
1:A:344:GLY:O	1:A:345:GLN:C	2.51	0.47
1:A:345:GLN:HG2	1:A:345:GLN:O	2.14	0.47
1:A:291:ARG:O	1:A:294:GLN:HG2	2.15	0.47
2:A:365:MAN:H3	2:A:366:MAN:O5	2.15	0.47
1:A:313:VAL:HG13	1:A:355:VAL:HG22	1.95	0.47
1:A:320:LEU:HD23	1:A:320:LEU:C	2.36	0.46
1:A:75:VAL:HG22	1:A:114:LEU:HD11	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:239:THR:CG2	1:A:335:LEU:HB2	2.46	0.46
1:A:95:ARG:NH2	1:A:131:THR:HG21	2.30	0.46
1:A:100:LYS:HE2	1:A:139:GLU:OE2	2.16	0.46
1:A:78:TRP:CD2	3:A:367:NAG:O1	2.66	0.46
1:A:186:ASP:CG	1:A:242:ARG:HH12	2.20	0.46
1:A:102:VAL:HB	1:A:103:PRO:HD3	1.98	0.45
1:A:185:TYR:HD1	1:A:186:ASP:OD2	2.00	0.45
1:A:344:GLY:O	1:A:346:ASN:N	2.50	0.45
1:A:287:THR:HG23	4:A:509:HOH:O	2.16	0.45
1:A:251:THR:HG23	1:A:299:THR:HG21	1.98	0.45
1:A:78:TRP:CE3	3:A:367:NAG:H5	2.52	0.45
1:A:70:LYS:HE2	4:A:445:HOH:O	2.17	0.45
1:A:210:SER:C	1:A:213:PHE:N	2.70	0.45
1:A:290:PHE:HB3	4:A:505:HOH:O	2.16	0.44
3:A:367:NAG:O3	3:A:368:NAG:C4	2.65	0.44
1:A:66:ASN:HA	1:A:67:PRO:HD3	1.81	0.44
1:A:205:ASN:HD22	1:A:205:ASN:C	2.21	0.44
1:A:46:ASP:HB2	4:A:500:HOH:O	2.16	0.44
1:A:142:ARG:HD3	4:A:421:HOH:O	2.17	0.44
1:A:9:SER:HB2	1:A:36:SER:O	2.18	0.43
1:A:210:SER:C	1:A:213:PHE:H	2.20	0.43
1:A:200:PRO:CB	1:A:293:GLN:HB3	2.48	0.43
1:A:78:TRP:CZ3	1:A:120:TYR:CE1	3.01	0.43
1:A:36:SER:HA	1:A:37:PHE:HA	1.80	0.43
1:A:7:TYR:CE2	1:A:22:PRO:HG3	2.53	0.43
1:A:202:PHE:HE2	1:A:294:GLN:NE2	2.16	0.43
1:A:63:LYS:HD2	1:A:71:THR:HG23	2.01	0.43
1:A:200:PRO:HB2	1:A:293:GLN:HB3	2.00	0.43
1:A:239:THR:CG2	1:A:335:LEU:C	2.88	0.43
1:A:88:ILE:HG13	4:A:489:HOH:O	2.18	0.42
1:A:100:LYS:HE2	1:A:139:GLU:CD	2.39	0.42
1:A:65:ARG:NE	4:A:497:HOH:O	2.46	0.42
1:A:119:LEU:HA	1:A:120:TYR:HA	1.80	0.42
1:A:304:TRP:CZ3	1:A:306:ALA:HB2	2.55	0.42
1:A:186:ASP:CG	1:A:242:ARG:NH1	2.73	0.41
1:A:193:GLN:O	1:A:194:THR:HB	2.20	0.41
1:A:9:SER:OG	1:A:49:GLU:HG3	2.19	0.41
1:A:186:ASP:OD2	1:A:242:ARG:NH1	2.54	0.41
1:A:320:LEU:HD11	1:A:328:ALA:HB2	2.01	0.41
1:A:357:ASP:O	1:A:361:GLU:HG3	2.21	0.41
1:A:319:TYR:OH	1:A:323:ARG:HD2	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:83:GLU:O	1:A:86:SER:HB3	2.21	0.40
1:A:102:VAL:HG12	1:A:106:LEU:HD22	2.03	0.40
3:A:371:NAG:H62	3:A:372:NAG:C8	2.51	0.40
1:A:127:ARG:NH1	4:A:494:HOH:O	2.54	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	359/361 (99%)	336 (94%)	13 (4%)	10 (3%)	<b>6</b> <b>8</b>

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	193	GLN
1	A	190	ALA
1	A	191	TRP
1	A	345	GLN
1	A	346	ASN
1	A	194	THR
1	A	209	SER
1	A	204	GLY
1	A	212	ARG
1	A	189	GLY

### 5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	302/302 (100%)	271 (90%)	31 (10%)	9	15

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LEU
1	A	9	SER
1	A	69	LEU
1	A	75	VAL
1	A	85	PHE
1	A	100	LYS
1	A	106	LEU
1	A	107	ARG
1	A	127	ARG
1	A	129	LEU
1	A	161	LYS
1	A	175	ARG
1	A	191	TRP
1	A	192	ARG
1	A	193	GLN
1	A	194	THR
1	A	205	ASN
1	A	210	SER
1	A	239	THR
1	A	242	ARG
1	A	252	ASP
1	A	258	SER
1	A	282	LEU
1	A	299	THR
1	A	313	VAL
1	A	320	LEU
1	A	323	ARG
1	A	329	MET
1	A	347	LEU
1	A	348	THR
1	A	351	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	60	ASN
1	A	79	ASN
1	A	150	GLN
1	A	205	ASN
1	A	294	GLN
1	A	303	GLN
1	A	315	ASN

### 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 ⓘ

10 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	$\# Z  > 2$
2	NAG	A	363	1,2	14,14,15	0.58	0	15,19,21	1.94	5 (33%)
2	NAG	A	364	2	14,14,15	1.01	0	15,19,21	1.19	2 (13%)
2	MAN	A	365	2	11,11,12	1.26	0	14,15,17	1.39	3 (21%)
2	MAN	A	366	2	11,11,12	0.71	0	14,15,17	1.28	1 (7%)
3	NAG	A	367	3	15,15,15	0.69	0	17,21,21	0.87	1 (5%)
3	NAG	A	368	3	14,14,15	1.01	1 (7%)	15,19,21	1.33	2 (13%)
3	NAG	A	369	3	14,14,15	1.01	1 (7%)	15,19,21	2.07	3 (20%)
3	NAG	A	370	3	14,14,15	0.68	0	15,19,21	1.37	2 (13%)
3	NAG	A	371	3	14,14,15	0.75	0	15,19,21	1.24	2 (13%)
3	NAG	A	372	3	14,14,15	0.67	0	15,19,21	1.42	2 (13%)

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	363	1,2	-	0/6/23/26	0/1/1/1
2	NAG	A	364	2	1/1/5/7	0/6/23/26	0/1/1/1
2	MAN	A	365	2	1/1/4/5	0/2/19/22	0/1/1/1
2	MAN	A	366	2	1/1/4/5	0/2/19/22	0/1/1/1
3	NAG	A	367	3	1/1/6/7	0/6/26/26	0/1/1/1
3	NAG	A	368	3	-	0/6/23/26	0/1/1/1
3	NAG	A	369	3	1/1/5/7	0/6/23/26	0/1/1/1
3	NAG	A	370	3	-	0/6/23/26	0/1/1/1
3	NAG	A	371	3	-	1/6/23/26	0/1/1/1
3	NAG	A	372	3	-	0/6/23/26	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	369	NAG	C1-C2	2.30	1.55	1.52
3	A	368	NAG	C1-C2	2.47	1.55	1.52

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	372	NAG	C4-C3-C2	-3.92	105.14	111.23
3	A	370	NAG	C2-N2-C7	-3.84	118.11	123.04
3	A	369	NAG	C4-C3-C2	-3.43	105.90	111.23
3	A	369	NAG	C2-N2-C7	-2.91	119.30	123.04
3	A	368	NAG	C4-C3-C2	-2.52	107.31	111.23
2	A	365	MAN	O4-C4-C3	-2.49	104.73	110.34
2	A	364	NAG	C2-N2-C7	-2.47	119.87	123.04
3	A	370	NAG	O4-C4-C5	-2.11	103.65	109.24
2	A	364	NAG	C3-C4-C5	-2.06	106.61	110.20
2	A	363	NAG	O7-C7-C8	-2.01	118.37	122.06
2	A	363	NAG	C3-C2-N2	2.06	115.50	110.56
3	A	371	NAG	C3-C4-C5	2.13	113.91	110.20
3	A	372	NAG	C2-N2-C7	2.18	125.84	123.04
2	A	365	MAN	C1-O5-C5	2.26	115.11	112.25
3	A	367	NAG	C4-C3-C2	2.28	113.59	110.43
2	A	363	NAG	C2-N2-C7	2.29	125.98	123.04
3	A	371	NAG	C2-N2-C7	2.61	126.39	123.04
2	A	363	NAG	C8-C7-N2	2.62	121.12	116.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	365	MAN	C1-C2-C3	3.05	113.14	109.54
2	A	366	MAN	O5-C5-C6	3.32	114.55	107.35
3	A	368	NAG	O4-C4-C3	3.35	117.88	110.34
2	A	363	NAG	C3-C4-C5	5.24	119.33	110.20
3	A	369	NAG	C1-O5-C5	6.20	120.11	112.25

All (5) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	A	369	NAG	C1
2	A	366	MAN	C1
3	A	367	NAG	C1
2	A	365	MAN	C1
2	A	364	NAG	C1

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	371	NAG	O7-C7-N2-C2

There are no ring outliers.

9 monomers are involved in 23 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	363	NAG	2	0
2	A	365	MAN	2	0
2	A	366	MAN	2	0
3	A	367	NAG	14	0
3	A	368	NAG	6	0
3	A	369	NAG	1	0
3	A	370	NAG	2	0
3	A	371	NAG	2	0
3	A	372	NAG	1	0

## 5.6 Ligand geometry [i](#)

There are no ligands in this entry.

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	361/361 (100%)	0.34	25 (6%)	20 23	11, 26, 58, 100	2 (0%)

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	208	ALA	8.8
1	A	207	ASP	7.2
1	A	255	ALA	7.0
1	A	209	SER	5.9
1	A	194	THR	5.5
1	A	362	VAL	4.9
1	A	205	ASN	4.8
1	A	191	TRP	4.8
1	A	210	SER	4.7
1	A	192	ARG	3.7
1	A	189	GLY	3.6
1	A	204	GLY	3.5
1	A	193	GLN	3.4
1	A	345	GLN	3.4
1	A	96	ARG	3.1
1	A	341	THR	3.0
1	A	346	ASN	2.9
1	A	256	PRO	2.6
1	A	340	GLY	2.6
1	A	190	ALA	2.5
1	A	206	GLU	2.4
1	A	251	THR	2.3
1	A	248	SER	2.3
1	A	81	GLY	2.1
1	A	253	VAL	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
3	NAG	A	367	15/15	0.80	0.34	3.32	40,45,48,59	0
2	NAG	A	363	14/15	0.81	0.28	2.55	17,23,29,39	14
3	NAG	A	370	14/15	0.91	0.18	1.82	42,46,51,51	0
3	NAG	A	368	14/15	0.90	0.21	0.75	42,47,53,54	0
3	NAG	A	369	14/15	0.91	0.16	0.04	41,44,46,48	0
2	MAN	A	366	11/12	0.35	0.52	-	58,59,60,60	11
2	MAN	A	365	11/12	0.22	0.52	-	56,58,58,60	11
3	NAG	A	372	14/15	0.75	0.51	-	40,44,47,48	0
2	NAG	A	364	14/15	0.77	0.40	-	43,48,52,52	14
3	NAG	A	371	14/15	0.82	0.30	-	43,46,48,49	0

## 6.4 Ligands [i](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.