



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 06:29 PM GMT

PDB ID : 1AZX  
Title : ANTITHROMBIN/PENTASACCHARIDE COMPLEX  
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Deposited on : 1997-11-23  
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) : 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

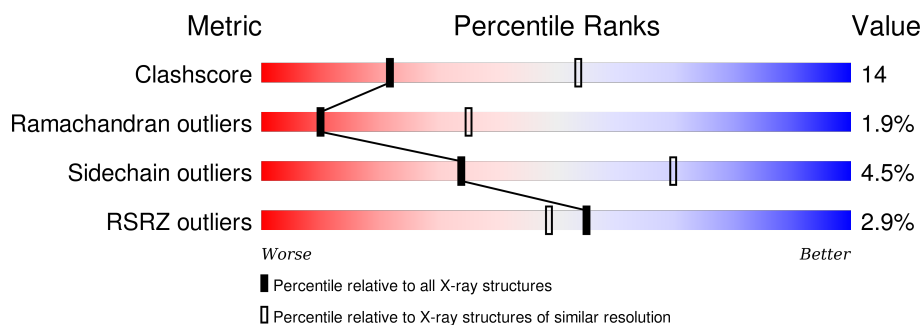
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.90 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)
RSRZ outliers	91569	1456 (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.

Mol	Chain	Length	Quality of chain
1	I	432	<div> <div>4%</div> <div>65%</div> <div>29%</div> <div>• •</div> </div>
1	L	432	<div> <div>2%</div> <div>64%</div> <div>31%</div> <div>• •</div> </div>

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 6881 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 ANTITHROMBIN.

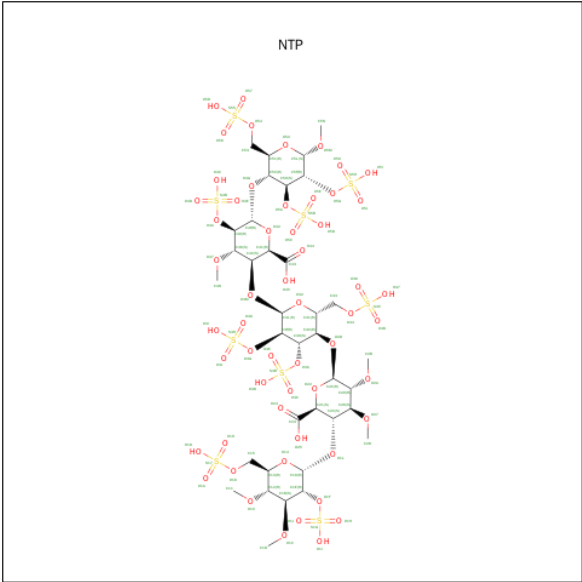
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	I	417	Total	C	N	O	S	0	0	0
			3328	2121	562	627	18			
1	L	414	Total	C	N	O	S	0	0	0
			3311	2111	559	623	18			

- Molecule 2 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula:  $C_8H_{15}NO_6$ ).



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

- Molecule 3 is HEPARIN PENTASACCHARIDE (three-letter code: NTP) (formula:  $C_{36}H_{60}O_{55}S_9$ ).

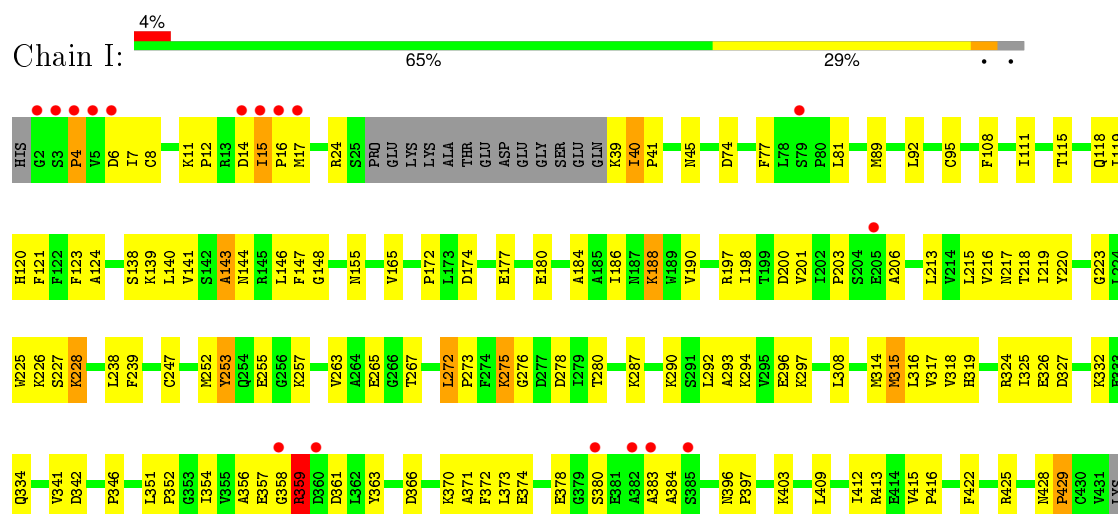


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	I	1	Total	C	O	S	0	0
			100	36	55	9		
3	L	1	Total	C	O	S	0	0
			100	36	55	9		

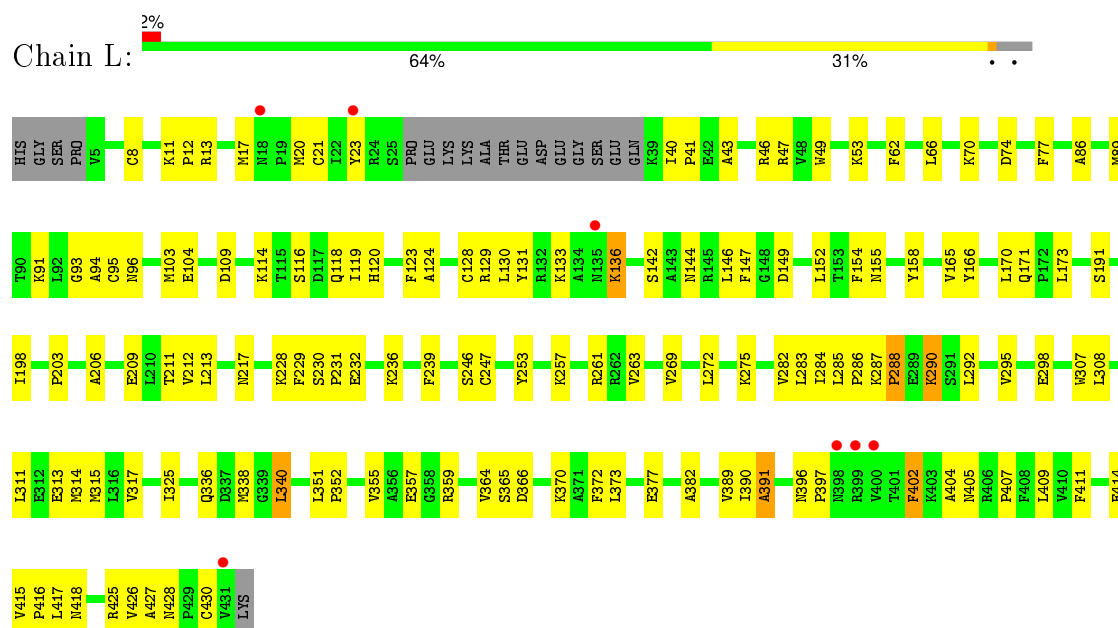
### 3 Residue-property plots

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: ANTITHROMBIN



#### • Molecule 1: ANTITHROMBIN



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	70.44Å 86.97Å 97.22Å 90.00° 108.88° 90.00°	Depositor
Resolution (Å)	30.00 – 2.90 27.65 – 2.90	Depositor EDS
% Data completeness (in resolution range)	94.2 (30.00-2.90) 94.5 (27.65-2.90)	Depositor EDS
$R_{merge}$	0.12	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.01 (at 2.90Å)	Xtriage
Refinement program	TNT V. 5-D	Depositor
R, $R_{free}$	0.195 , 0.280 0.208 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	45.4	Xtriage
Anisotropy	0.356	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 94.5	EDS
Estimated twinning fraction	0.023 for h,-k,-h-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 23424 reflections	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	6881	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.65% 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: NTP, 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	I	0.34	0/3394	0.59	0/4582
1	L	0.35	0/3376	0.58	0/4557
All	All	0.35	0/6770	0.59	0/9139

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts

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	I	3328	0	3333	90	0
1	L	3311	0	3323	91	0
2	I	14	0	13	5	0
2	L	28	0	26	3	0
3	I	100	0	49	6	0
3	L	100	0	49	9	0
All	All	6881	0	6793	189	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 14.

All (189) 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:I:356:ALA:O	2:I:501:NAG:H81	1.60	1.01
1:L:129:ARG:HD3	1:L:417:LEU:HD21	1.56	0.87
1:L:20:MET:HA	2:L:501:NAG:H61	1.55	0.87
1:I:17:MET:HB2	1:I:120:HIS:HE1	1.39	0.86
1:I:144:ASN:HB3	1:I:217:ASN:HA	1.67	0.77
1:I:324:ARG:HG2	1:I:374:GLU:HG3	1.67	0.75
3:L:503:NTP:O44	3:L:503:NTP:H31	1.91	0.71
1:I:319:HIS:HB2	1:I:403:LYS:HA	1.73	0.69
1:I:155:ASN:HD22	2:I:501:NAG:H83	1.55	0.69
3:L:503:NTP:H283	3:L:503:NTP:H2B2	1.75	0.69
1:I:356:ALA:O	2:I:501:NAG:C8	2.41	0.68
1:I:108:PHE:O	1:I:111:ILE:HG12	1.94	0.67
1:I:358:GLY:O	1:I:359:ARG:HD3	1.94	0.66
3:L:503:NTP:H151	3:L:503:NTP:H112	1.76	0.66
1:L:20:MET:CA	2:L:501:NAG:H61	2.26	0.66
1:L:355:VAL:HG21	1:L:359:ARG:HB2	1.78	0.65
1:I:17:MET:HB2	1:I:120:HIS:CE1	2.28	0.65
1:I:257:LYS:HA	1:I:314:MET:O	1.96	0.65
1:I:77:PHE:CZ	1:I:373:LEU:HB2	2.33	0.64
1:L:47:ARG:NH1	1:L:114:LYS:HB3	2.11	0.64
1:I:308:LEU:HD13	1:I:413:ARG:NH1	2.14	0.63
1:I:255:GLU:HG2	1:I:317:VAL:HG22	1.80	0.62
1:I:415:VAL:HB	1:I:416:PRO:HD3	1.81	0.62
1:L:415:VAL:HB	1:L:416:PRO:HD3	1.80	0.62
1:L:229:PHE:O	1:L:377:GLU:HB3	2.00	0.62
1:I:17:MET:SD	1:I:121:PHE:HB2	2.40	0.61
3:L:503:NTP:H3L	3:L:503:NTP:O3K	2.00	0.61
1:L:286:PRO:HD3	1:L:292:LEU:HD13	1.81	0.61
1:L:257:LYS:HG2	1:L:315:MET:HG2	1.82	0.61
1:L:247:CYS:HG	1:L:430:CYS:CB	2.13	0.60
1:I:239:PHE:HB3	1:I:247:CYS:HB3	1.83	0.60
1:I:412:ILE:HB	1:I:422:PHE:HB2	1.84	0.59
1:L:152:LEU:HD13	1:L:212:VAL:HG13	1.83	0.59
1:I:139:LYS:HZ3	1:I:197:ARG:HH22	1.48	0.59
1:L:211:THR:HA	1:L:391:ALA:O	2.01	0.59
1:L:407:PRO:HB3	1:L:427:ALA:HA	1.85	0.59
1:I:265:GLU:O	1:I:287:LYS:HD2	2.03	0.58
1:I:140:LEU:HD11	1:I:219:ILE:HD11	1.85	0.58
1:I:228:LYS:HB3	1:I:228:LYS:HZ2	1.69	0.58
1:I:198:ILE:HG23	1:I:370:LYS:HD3	1.87	0.57
1:I:228:LYS:HZ2	1:I:378:GLU:HB2	1.69	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:272:LEU:HD11	1:L:283:LEU:HD11	1.87	0.57
1:L:313:GLU:CD	1:L:313:GLU:H	2.08	0.57
1:L:364:VAL:HG22	1:L:390:ILE:HD13	1.88	0.56
1:L:21:CYS:SG	1:L:95:CYS:CB	2.94	0.56
1:I:4:PRO:HB2	1:I:8:CYS:SG	2.46	0.55
1:L:308:LEU:HA	1:L:311:LEU:HG	1.88	0.55
1:I:138:SER:HB3	1:I:223:GLY:HA2	1.88	0.55
1:L:355:VAL:CG2	1:L:359:ARG:HB2	2.36	0.55
1:L:116:SER:O	1:L:119:ILE:HG12	2.07	0.55
1:L:95:CYS:SG	1:L:96:ASN:ND2	2.80	0.55
1:I:143:ALA:HB3	1:I:218:THR:OG1	2.07	0.55
1:L:21:CYS:SG	1:L:95:CYS:HB2	2.47	0.54
1:L:236:LYS:NZ	1:L:236:LYS:HB2	2.22	0.54
1:I:356:ALA:O	2:I:501:NAG:O7	2.26	0.53
1:L:12:PRO:HB3	1:L:118:GLN:NE2	2.24	0.53
1:I:238:LEU:HA	1:I:247:CYS:O	2.10	0.52
1:L:253:TYR:CE1	1:L:317:VAL:HG13	2.45	0.52
1:L:40:ILE:HB	1:L:41:PRO:HD2	1.92	0.52
1:I:146:LEU:HG	1:I:215:LEU:HG	1.91	0.52
1:L:130:LEU:HD23	1:L:414:GLU:OE1	2.11	0.51
1:L:365:SER:HB3	1:L:389:VAL:HG23	1.91	0.51
3:I:502:NTP:H283	3:I:502:NTP:H2B2	1.92	0.51
1:L:20:MET:HA	2:L:501:NAG:C6	2.36	0.51
1:L:257:LYS:HA	1:L:314:MET:O	2.10	0.51
1:I:147:PHE:CE2	1:I:186:ILE:HG23	2.45	0.51
1:L:370:LYS:HB2	1:L:370:LYS:NZ	2.25	0.51
1:L:287:LYS:HB3	1:L:290:LYS:HG3	1.91	0.51
1:I:15:ILE:HD13	1:I:15:ILE:N	2.26	0.51
1:I:174:ASP:OD2	1:I:177:GLU:HB2	2.11	0.51
1:I:40:ILE:HG13	1:I:41:PRO:HD2	1.94	0.50
1:L:23:TYR:CD1	1:L:103:MET:HG3	2.47	0.50
1:I:148:GLY:O	1:I:172:PRO:HA	2.12	0.50
1:I:184:ALA:O	1:I:188:LYS:HD2	2.12	0.50
1:L:402:PHE:HD1	1:L:402:PHE:O	1.95	0.50
1:I:253:TYR:HA	1:I:318:VAL:O	2.12	0.50
1:L:77:PHE:CE2	1:L:373:LEU:HB2	2.46	0.50
1:I:220:TYR:HD1	1:I:372:PHE:HB3	1.77	0.49
1:L:86:ALA:O	1:L:89:MET:HB2	2.12	0.49
1:I:380:SER:HB2	1:I:384:ALA:HB2	1.94	0.49
1:I:314:MET:HG2	1:I:315:MET:N	2.26	0.49
1:I:290:LYS:HD2	1:I:294:LYS:HD3	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:120:HIS:HD2	1:L:166:TYR:OH	1.96	0.49
1:I:17:MET:HG3	1:I:118:GLN:HE22	1.77	0.49
1:L:93:GLY:O	1:L:351:LEU:HA	2.13	0.49
1:L:91:LYS:NZ	1:L:120:HIS:NE2	2.61	0.49
1:I:326:GLU:HA	1:I:371:ALA:O	2.13	0.49
1:I:396:ASN:OD1	1:I:397:PRO:HD2	2.13	0.48
1:L:149:ASP:HA	1:L:173:LEU:O	2.13	0.48
1:L:21:CYS:HG	1:L:95:CYS:HG	0.48	0.48
1:L:203:PRO:HG2	1:L:206:ALA:HB2	1.94	0.48
1:I:12:PRO:HD2	3:I:502:NTP:O3J	2.14	0.48
1:L:155:ASN:HB3	1:L:158:TYR:HB3	1.96	0.48
1:L:290:LYS:NZ	1:L:290:LYS:HB3	2.28	0.48
1:L:77:PHE:CE1	1:L:373:LEU:HD22	2.48	0.48
1:L:370:LYS:HZ2	1:L:370:LYS:HB2	1.79	0.48
3:I:502:NTP:C28	3:I:502:NTP:H1K	2.44	0.47
1:I:272:LEU:N	1:I:272:LEU:HD12	2.29	0.47
1:I:141:VAL:HG23	1:I:220:TYR:HB3	1.94	0.47
1:L:131:TYR:CE1	1:L:142:SER:HB2	2.49	0.47
1:L:144:ASN:OD1	1:L:217:ASN:ND2	2.47	0.47
1:L:239:PHE:O	1:L:246:SER:HA	2.13	0.47
1:L:62:PHE:HA	1:L:338:MET:HE1	1.96	0.47
3:I:502:NTP:H282	3:I:502:NTP:H1K	1.97	0.47
1:L:407:PRO:HB2	1:L:425:ARG:HG2	1.97	0.47
1:L:247:CYS:SG	1:L:430:CYS:CB	3.03	0.47
1:L:154:PHE:HA	1:L:355:VAL:HA	1.96	0.47
1:L:47:ARG:HH12	1:L:114:LYS:HB3	1.78	0.47
1:L:43:ALA:HB1	3:L:503:NTP:H113	1.97	0.46
1:I:188:LYS:N	1:I:188:LYS:HE3	2.30	0.46
1:L:8:CYS:HG	1:L:128:CYS:CB	2.26	0.46
1:I:119:ILE:O	1:I:123:PHE:HD2	1.98	0.46
1:L:13:ARG:NH1	3:L:503:NTP:O3I	2.48	0.46
1:I:293:ALA:O	1:I:297:LYS:HG3	2.16	0.46
1:L:229:PHE:HB2	1:L:377:GLU:HA	1.96	0.46
1:I:273:PRO:HA	1:I:280:THR:HG22	1.97	0.46
1:L:47:ARG:CZ	1:L:114:LYS:HD3	2.46	0.46
1:L:282:VAL:HB	1:L:411:PHE:HB2	1.97	0.46
1:I:292:LEU:O	1:I:296:GLU:HG3	2.16	0.45
1:L:191:SER:HA	1:L:198:ILE:O	2.16	0.45
1:I:292:LEU:HD11	1:I:409:LEU:HG	1.97	0.45
1:I:141:VAL:CG2	1:I:220:TYR:HB3	2.47	0.45
1:I:147:PHE:O	1:I:213:LEU:HD12	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:324:ARG:HA	1:I:373:LEU:O	2.17	0.44
1:L:275:LYS:HD2	1:L:275:LYS:HA	1.79	0.44
1:L:133:LYS:O	1:L:136:LYS:HE2	2.17	0.44
1:L:261:ARG:HB2	1:L:311:LEU:HA	1.99	0.44
1:I:95:CYS:SG	1:I:352:PRO:CG	3.05	0.44
1:L:124:ALA:HB2	1:L:165:VAL:HG13	2.00	0.44
1:L:295:VAL:HG13	1:L:307:TRP:HH2	1.83	0.44
1:I:358:GLY:C	1:I:359:ARG:HD3	2.37	0.44
1:I:124:ALA:HB2	1:I:165:VAL:HG13	2.00	0.44
1:L:66:LEU:O	1:L:70:LYS:HG3	2.17	0.44
1:I:325:ILE:O	1:I:372:PHE:HA	2.17	0.44
1:I:201:VAL:HG11	1:I:216:VAL:HG13	1.99	0.44
1:I:252:MET:O	1:I:319:HIS:HA	2.18	0.43
1:I:357:GLU:HB2	1:I:359:ARG:NH1	2.34	0.43
1:I:74:ASP:O	1:I:425:ARG:NH1	2.50	0.43
1:L:239:PHE:HB3	1:L:430:CYS:SG	2.58	0.43
1:L:287:LYS:HB3	1:L:290:LYS:HB2	2.00	0.43
1:L:336:GLN:HA	1:L:340:LEU:O	2.18	0.43
1:I:341:VAL:HG23	1:I:342:ASP:H	1.83	0.43
1:I:155:ASN:ND2	2:I:501:NAG:H83	2.28	0.43
1:L:407:PRO:HB3	1:L:427:ALA:CA	2.48	0.43
1:L:325:ILE:HD13	1:L:426:VAL:HG22	2.01	0.43
1:I:198:ILE:HG21	1:I:218:THR:HB	2.01	0.42
1:L:13:ARG:NH1	3:L:503:NTP:H482	2.34	0.42
1:I:319:HIS:CB	1:I:403:LYS:HG3	2.49	0.42
1:I:11:LYS:HB3	3:I:502:NTP:O3J	2.19	0.42
1:L:147:PHE:HA	1:L:171:GLN:O	2.20	0.42
1:I:17:MET:HG3	1:I:118:GLN:OE1	2.20	0.42
1:I:203:PRO:HG2	1:I:206:ALA:HB2	2.00	0.42
1:L:124:ALA:O	1:L:128:CYS:SG	2.78	0.42
1:I:139:LYS:HA	1:I:139:LYS:HD2	1.92	0.42
1:I:294:LYS:HZ2	1:I:294:LYS:HB3	1.84	0.42
1:I:14:ASP:HB2	1:I:15:ILE:HD13	2.02	0.42
1:L:285:LEU:HD12	1:L:285:LEU:N	2.35	0.42
1:I:190:VAL:HG21	1:I:201:VAL:HG21	2.02	0.42
1:L:230:SER:HA	1:L:231:PRO:HD3	1.90	0.42
1:L:292:LEU:HD21	1:L:409:LEU:HG	2.01	0.41
1:L:239:PHE:HZ	1:L:405:ASN:O	2.02	0.41
1:I:39:LYS:O	1:I:40:ILE:HB	2.20	0.41
1:L:20:MET:HG2	1:L:352:PRO:HG2	2.01	0.41
1:I:308:LEU:HD13	1:I:413:ARG:HH11	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:327:ASP:O	1:I:370:LYS:HA	2.20	0.41
1:L:366:ASP:O	1:L:389:VAL:HG22	2.21	0.41
1:I:45:ASN:HA	3:I:502:NTP:H282	2.02	0.41
1:I:351:LEU:HD13	1:I:354:ILE:HD12	2.01	0.41
3:L:503:NTP:H282	3:L:503:NTP:H1K	2.01	0.41
1:I:228:LYS:NZ	1:I:378:GLU:HB2	2.34	0.41
1:L:372:PHE:O	1:L:382:ALA:HA	2.20	0.41
1:L:11:LYS:HD3	3:L:503:NTP:H3F	2.02	0.41
1:I:276:GLY:C	1:I:278:ASP:H	2.24	0.41
1:L:74:ASP:O	1:L:425:ARG:HD2	2.21	0.41
1:I:314:MET:HG2	1:I:315:MET:H	1.85	0.41
1:L:147:PHE:O	1:L:213:LEU:HD12	2.21	0.41
1:I:263:VAL:CG2	1:I:267:THR:HB	2.51	0.41
1:I:15:ILE:HA	1:I:16:PRO:HD2	1.87	0.40
1:L:94:ALA:HA	1:L:351:LEU:HD23	2.03	0.40
1:L:230:SER:HB3	1:L:232:GLU:CD	2.41	0.40
1:L:146:LEU:HB2	1:L:170:LEU:HA	2.04	0.40
1:I:346:PRO:HG3	1:I:363:TYR:CE2	2.56	0.40
1:I:17:MET:HG3	1:I:118:GLN:NE2	2.35	0.40
1:L:269:VAL:HG13	1:L:284:ILE:HD13	2.04	0.40
1:I:89:MET:O	1:I:92:LEU:HB3	2.21	0.40
1:L:49:TRP:CZ2	1:L:53:LYS:HD3	2.56	0.40
1:I:226:LYS:HA	1:I:275:LYS:HD2	2.02	0.40
1:I:115:THR:O	1:I:119:ILE:N	2.54	0.40
1:L:104:GLU:HA	1:L:109:ASP:OD1	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	I	413/432 (96%)	367 (89%)	36 (9%)	10 (2%)	7 29

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	L	410/432 (95%)	377 (92%)	27 (7%)	6 (2%)	13	42
All	All	823/864 (95%)	744 (90%)	63 (8%)	16 (2%)	10	35

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	I	4	PRO
1	I	200	ASP
1	I	359	ARG
1	L	404	ALA
1	L	288	PRO
1	I	6	ASP
1	I	383	ALA
1	I	429	PRO
1	L	357	GLU
1	I	143	ALA
1	I	361	ASP
1	L	391	ALA
1	I	332	LYS
1	I	40	ILE
1	L	397	PRO
1	L	263	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	I	370/383 (97%)	351 (95%)	19 (5%)	29	65
1	L	368/383 (96%)	354 (96%)	14 (4%)	40	76
All	All	738/766 (96%)	705 (96%)	33 (4%)	34	70

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	I	7	ILE
1	I	15	ILE
1	I	24	ARG
1	I	81	LEU
1	I	180	GLU
1	I	188	LYS
1	I	225	TRP
1	I	227	SER
1	I	228	LYS
1	I	253	TYR
1	I	272	LEU
1	I	275	LYS
1	I	315	MET
1	I	316	LEU
1	I	334	GLN
1	I	359	ARG
1	I	366	ASP
1	I	428	ASN
1	I	429	PRO
1	L	17	MET
1	L	46	ARG
1	L	123	PHE
1	L	136	LYS
1	L	209	GLU
1	L	228	LYS
1	L	288	PRO
1	L	290	LYS
1	L	298	GLU
1	L	340	LEU
1	L	396	ASN
1	L	402	PHE
1	L	418	ASN
1	L	428	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	I	120	HIS
1	I	127	ASN
1	I	233	ASN
1	I	428	ASN
1	L	96	ASN
1	L	118	GLN

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Mol	Chain	Res	Type
1	L	120	HIS
1	L	396	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 ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

5 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	I	501	1	14,14,15	0.84	0	15,19,21	1.54	3 (20%)
3	NTP	I	502	-	98,104,104	1.00	1 (1%)	140,162,162	0.84	3 (2%)
2	NAG	L	501	1	14,14,15	0.84	0	15,19,21	1.54	3 (20%)
2	NAG	L	502	1	14,14,15	0.84	0	15,19,21	1.54	3 (20%)
3	NTP	L	503	-	98,104,104	0.95	2 (2%)	140,162,162	0.90	6 (4%)

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	I	501	1	-	0/6/23/26	0/1/1/1
3	NTP	I	502	-	-	0/76/184/184	0/5/5/5
2	NAG	L	501	1	-	0/6/23/26	0/1/1/1
2	NAG	L	502	1	-	0/6/23/26	0/1/1/1
3	NTP	L	503	-	-	0/76/184/184	0/5/5/5

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	I	502	NTP	O3A-C39	-2.33	1.41	1.46
3	L	503	NTP	O3A-C39	-2.09	1.41	1.46
3	L	503	NTP	C40-C41	2.64	1.57	1.52

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	L	501	NAG	O4-C4-C3	-3.10	103.37	110.34
2	L	502	NAG	O4-C4-C3	-3.09	103.37	110.34
2	I	501	NAG	O4-C4-C3	-3.09	103.38	110.34
2	L	501	NAG	O7-C7-C8	-2.24	117.95	122.06
2	L	502	NAG	O7-C7-C8	-2.24	117.95	122.06
2	I	501	NAG	O7-C7-C8	-2.21	118.01	122.06
3	L	503	NTP	O34-C33-C31	-2.20	103.67	107.90
3	L	503	NTP	O5G-C5F-C5L	2.03	110.41	107.65
3	L	503	NTP	O3G-C3F-C3L	2.03	110.42	107.65
3	L	503	NTP	O42-C4F-C49	2.07	113.65	109.47
3	L	503	NTP	C4F-O42-C41	2.09	115.42	112.17
3	I	502	NTP	O3G-C3F-C3L	2.10	110.50	107.65
3	I	502	NTP	O5G-C5F-C5L	2.13	110.54	107.65
3	L	503	NTP	O5M-C5L-C5F	2.13	111.45	107.82
3	I	502	NTP	O4A-C49-C4F	2.29	110.77	107.65
2	L	501	NAG	C1-O5-C5	3.17	116.27	112.25
2	I	501	NAG	C1-O5-C5	3.18	116.28	112.25
2	L	502	NAG	C1-O5-C5	3.18	116.28	112.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 23 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	I	501	NAG	5	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	I	502	NTP	6	0
2	L	501	NAG	3	0
3	L	503	NTP	9	0

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	I	417/432 (96%)	-0.22	17 (4%) 41 34	2, 35, 78, 114	0
1	L	414/432 (95%)	-0.32	7 (1%) 73 70	2, 27, 70, 97	0
All	All	831/864 (96%)	-0.27	24 (2%) 55 49	2, 32, 74, 114	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	I	3	SER	7.4
1	I	2	GLY	5.9
1	I	5	VAL	4.4
1	I	4	PRO	4.3
1	L	18	ASN	3.9
1	I	385	SER	3.8
1	I	382	ALA	3.6
1	I	15	ILE	3.5
1	L	399	ARG	3.3
1	I	6	ASP	3.0
1	I	383	ALA	2.9
1	L	135	ASN	2.9
1	I	205	GLU	2.7
1	I	14	ASP	2.7
1	L	23	TYR	2.7
1	L	398	ASN	2.4
1	L	400	VAL	2.4
1	I	79	SER	2.3
1	I	360	ASP	2.3
1	L	431	VAL	2.3
1	I	16	PRO	2.2
1	I	358	GLY	2.1
1	I	17	MET	2.1
1	I	380	SER	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](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 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( $\text{\AA}^2$ )	Q<0.9
2	NAG	L	501	14/15	0.55	0.38	1.83	30,60,78,82	0
2	NAG	I	501	14/15	0.86	0.30	1.68	35,55,55,65	0
3	NTP	L	503	100/100	0.95	0.13	-0.69	21,31,56,74	0
3	NTP	I	502	100/100	0.94	0.12	-1.07	27,50,65,71	0
2	NAG	L	502	14/15	0.65	0.43	-	30,60,78,82	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.