Full wwPDB X-ray Structure Validation Report (i)

Sep 12, 2023 – 02:19 PM EDT

PDB ID	:	4NZQ
Title	:	Crystal structure of Ca2+-free prothrombin deletion mutant residues 146-167
Authors	:	Pozzi, N.; Chen, Z.; Shropshire, D.B.; Pelc, L.A.; Di Cera, E.
Deposited on	:	2013-12-12
Resolution	:	2.81 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*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

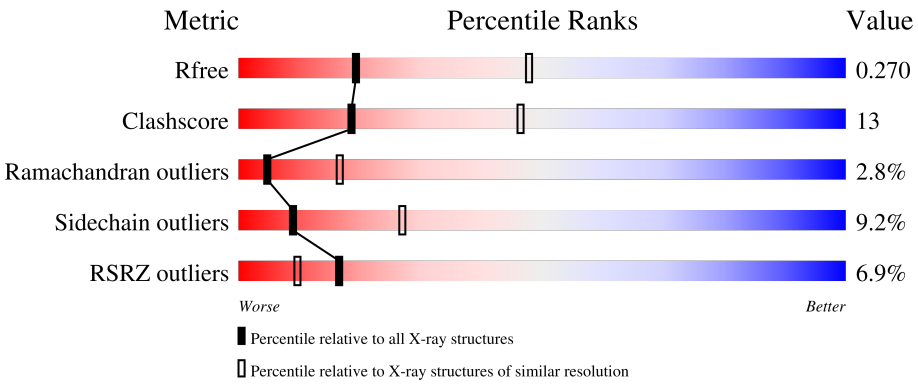
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.81 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}	130704	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. 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 <=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	557	6% 61% 26% • 9%

2 Entry composition (i)

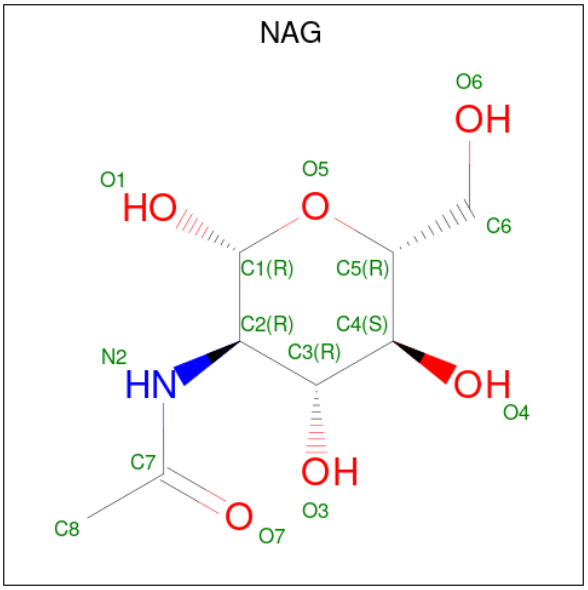
There are 2 unique types of molecules in this entry. The entry contains 4084 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Prothrombin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	508	4042	2533	714	766	29	0	0	0

- Molecule 2 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



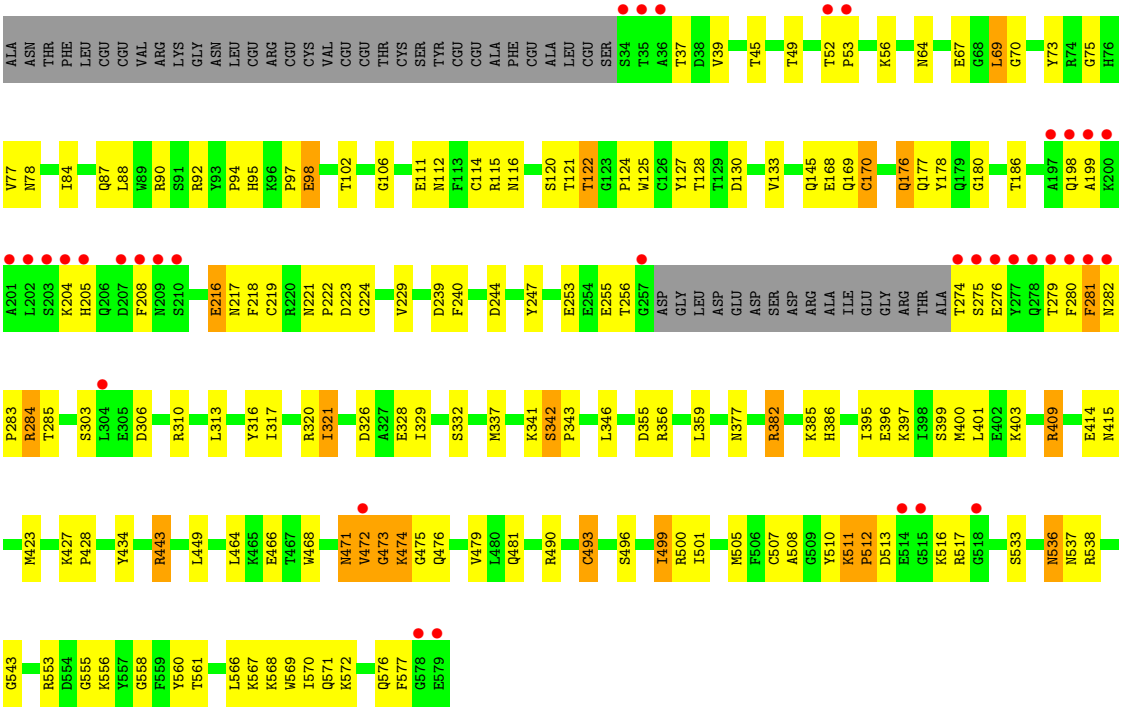
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	A	1	14	8	1	5	0	0
2	A	1	14	8	1	5	0	0
2	A	1	14	8	1	5	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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: Prothrombin

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α, β, γ	88.75Å 171.74Å 141.25Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.23 – 2.81 30.23 – 2.81	Depositor EDS
% Data completeness (in resolution range)	98.0 (30.23-2.81) 98.2 (30.23-2.81)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.56 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.226 , 0.279 0.223 , 0.270	Depositor DCC
R_{free} test set	1330 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å^2)	53.0	Xtriage
Anisotropy	0.076	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 35.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.44$, $< L^2 > = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.91	EDS
Total number of atoms	4084	wwPDB-VP
Average B, all atoms (Å^2)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.64% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.78	1/4144 (0.0%)	0.82	2/5614 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	170	CYS	CB-SG	-5.83	1.72	1.81

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	382	ARG	NE-CZ-NH2	7.54	124.07	120.30
1	A	382	ARG	NE-CZ-NH1	-6.17	117.21	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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	4042	0	3855	105	0
2	A	42	0	39	3	0
All	All	4084	0	3894	105	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 13.

All (105) 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:474:LYS:NZ	1:A:475:GLY:O	1.88	1.05
1:A:84:ILE:HG21	1:A:128:THR:CG2	1.93	0.99
1:A:52:THR:HG23	1:A:53:PRO:HD3	1.45	0.96
1:A:176:GLN:HE22	1:A:224:GLY:HA2	1.32	0.93
1:A:280:PHE:CE1	1:A:567:LYS:HB3	2.17	0.80
1:A:52:THR:HG23	1:A:53:PRO:CD	2.15	0.76
1:A:53:PRO:HG2	1:A:169:GLN:CD	2.07	0.74
1:A:490:ARG:HH12	1:A:501:ILE:HG22	1.52	0.73
1:A:355:ASP:OD1	1:A:356:ARG:HB2	1.90	0.71
1:A:281:PHE:CZ	1:A:567:LYS:HD2	2.28	0.69
1:A:84:ILE:HG21	1:A:128:THR:HG22	1.74	0.68
1:A:415:ASN:HD22	1:A:499:ILE:HB	1.58	0.68
1:A:69:LEU:HD22	1:A:121:THR:HA	1.77	0.67
1:A:87:GLN:HE22	1:A:95:HIS:H	1.44	0.65
1:A:53:PRO:HG2	1:A:169:GLN:OE1	1.96	0.65
1:A:386:HIS:NE2	1:A:479:VAL:HG13	2.13	0.64
1:A:128:THR:HB	1:A:133:VAL:O	1.96	0.64
1:A:382:ARG:HH21	1:A:385:LYS:HD2	1.62	0.63
1:A:490:ARG:NH1	1:A:501:ILE:HG22	2.14	0.62
1:A:471:ASN:O	1:A:473:GLY:N	2.33	0.62
1:A:382:ARG:NH2	1:A:385:LYS:HD2	2.14	0.61
1:A:84:ILE:HG21	1:A:128:THR:HG23	1.79	0.61
1:A:145:GLN:NE2	1:A:145:GLN:HA	2.14	0.61
1:A:49:THR:OG1	1:A:67:GLU:OE1	2.20	0.59
1:A:536:ASN:C	1:A:536:ASN:HD22	2.05	0.59
1:A:566:LEU:O	1:A:570:ILE:HG13	2.02	0.59
1:A:510:TYR:O	1:A:512:PRO:HD3	2.04	0.58
1:A:114:CYS:O	1:A:115:ARG:HD3	2.04	0.57
1:A:88:LEU:HD23	1:A:112:ASN:HB3	1.85	0.57
1:A:533:SER:HB3	1:A:536:ASN:ND2	2.18	0.57
1:A:102:THR:HG21	2:A:703:NAG:H82	1.87	0.56
1:A:320:ARG:O	1:A:321:ILE:HB	2.04	0.56
1:A:84:ILE:CG2	1:A:128:THR:CG2	2.79	0.56
1:A:508:ALA:HB3	1:A:560:TYR:CE2	2.41	0.55
1:A:427:LYS:HD3	1:A:428:PRO:O	2.08	0.54
1:A:77:VAL:HG23	1:A:114:CYS:SG	2.48	0.53
1:A:576:GLN:O	1:A:577:PHE:CD1	2.60	0.53
1:A:341:LYS:NZ	1:A:377:ASN:O	2.27	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:178:TYR:CZ	1:A:180:GLY:HA3	2.44	0.52
1:A:513:ASP:N	1:A:516:LYS:O	2.37	0.52
1:A:145:GLN:HA	1:A:145:GLN:HE21	1.75	0.52
1:A:401:LEU:HD22	1:A:423:MET:HB3	1.92	0.52
1:A:95:HIS:O	1:A:97:PRO:HD3	2.10	0.52
1:A:337:MET:HG2	1:A:346:LEU:HD13	1.91	0.52
1:A:415:ASN:ND2	1:A:499:ILE:HB	2.23	0.51
1:A:501:ILE:HG23	1:A:505:MET:HB2	1.93	0.51
1:A:464:LEU:HD12	1:A:468:TRP:CD1	2.46	0.50
1:A:511:LYS:HE2	1:A:512:PRO:HD2	1.93	0.50
1:A:98:GLU:OE1	1:A:127:TYR:OH	2.25	0.50
1:A:87:GLN:HE21	1:A:94:PRO:HD2	1.76	0.50
1:A:536:ASN:C	1:A:536:ASN:ND2	2.65	0.49
1:A:255:GLU:OE2	1:A:443:ARG:NH1	2.46	0.49
1:A:326:ASP:HA	1:A:481:GLN:NE2	2.27	0.49
1:A:78:ASN:ND2	2:A:702:NAG:C7	2.75	0.49
1:A:84:ILE:CG2	1:A:128:THR:HG23	2.41	0.49
1:A:513:ASP:HB2	1:A:516:LYS:HB3	1.94	0.49
1:A:186:THR:HG22	1:A:244:ASP:HB2	1.96	0.48
1:A:386:HIS:NE2	1:A:479:VAL:CG1	2.75	0.48
1:A:73:TYR:CZ	1:A:75:GLY:HA3	2.49	0.47
1:A:553:ARG:HB2	1:A:556:LYS:HG3	1.94	0.47
1:A:53:PRO:O	1:A:56:LYS:HB3	2.14	0.47
1:A:472:VAL:HG12	1:A:472:VAL:O	2.15	0.47
1:A:247:TYR:HE2	1:A:409:ARG:HH22	1.62	0.47
1:A:219:CYS:HB3	1:A:229:VAL:HG23	1.97	0.47
1:A:414:GLU:OE2	1:A:500:ARG:NH2	2.48	0.46
1:A:84:ILE:HD13	1:A:128:THR:HG21	1.97	0.46
1:A:281:PHE:HZ	1:A:567:LYS:HD2	1.78	0.46
1:A:128:THR:HG22	1:A:130:ASP:H	1.80	0.46
1:A:536:ASN:ND2	1:A:538:ARG:H	2.12	0.46
1:A:87:GLN:NE2	1:A:94:PRO:HD2	2.30	0.45
1:A:326:ASP:HA	1:A:481:GLN:HE22	1.81	0.45
1:A:508:ALA:HB3	1:A:560:TYR:HE2	1.81	0.45
1:A:329:ILE:HD13	1:A:386:HIS:CE1	2.52	0.45
1:A:130:ASP:OD1	1:A:130:ASP:C	2.55	0.45
1:A:223:ASP:OD2	1:A:572:LYS:NZ	2.48	0.45
1:A:466:GLU:O	1:A:466:GLU:HG3	2.17	0.44
1:A:316:TYR:CD1	1:A:317:ILE:HG13	2.52	0.44
1:A:423:MET:HB2	1:A:423:MET:HE2	1.93	0.44
1:A:116:ASN:HB2	1:A:124:PRO:HA	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:114:CYS:HA	1:A:125:TRP:O	2.18	0.44
1:A:342:SER:HA	1:A:343:PRO:C	2.39	0.44
1:A:464:LEU:HD12	1:A:468:TRP:HD1	1.83	0.44
1:A:320:ARG:O	1:A:321:ILE:CB	2.66	0.43
1:A:449:LEU:HD23	1:A:449:LEU:HA	1.82	0.43
1:A:198:GLN:HE21	1:A:199:ALA:HB2	1.83	0.43
1:A:306:ASP:HB2	1:A:328:GLU:OE2	2.18	0.43
1:A:543:GLY:HA2	1:A:561:THR:O	2.19	0.43
1:A:568:LYS:HA	1:A:571:GLN:HB2	2.01	0.43
1:A:216:GLU:HB3	1:A:217:ASN:H	1.59	0.43
1:A:221:ASN:HB2	1:A:229:VAL:HA	2.01	0.42
1:A:208:PHE:HB3	1:A:222:PRO:HB2	2.01	0.42
1:A:471:ASN:C	1:A:473:GLY:N	2.73	0.42
1:A:512:PRO:HG2	1:A:555:GLY:HA2	2.01	0.42
1:A:120:SER:C	1:A:122:THR:H	2.23	0.42
1:A:382:ARG:NH2	1:A:396:GLU:OE2	2.53	0.42
1:A:395:ILE:HG23	1:A:434:TYR:CD2	2.55	0.41
1:A:507:CYS:HA	1:A:558:GLY:O	2.20	0.41
1:A:240:PHE:HZ	1:A:569:TRP:HD1	1.68	0.41
1:A:536:ASN:HD22	1:A:537:ASN:N	2.19	0.41
1:A:282:ASN:C	1:A:284:ARG:H	2.24	0.41
1:A:409:ARG:HG3	1:A:409:ARG:HH11	1.85	0.41
1:A:409:ARG:HG3	1:A:409:ARG:NH1	2.36	0.40
1:A:102:THR:CB	2:A:703:NAG:H82	2.51	0.40
1:A:218:PHE:N	1:A:218:PHE:CD1	2.87	0.40
1:A:493:CYS:O	1:A:496:SER:OG	2.29	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	504/557 (90%)	449 (89%)	41 (8%)	14 (3%)	5 17

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	321	ILE
1	A	472	VAL
1	A	70	GLY
1	A	275	SER
1	A	473	GLY
1	A	106	GLY
1	A	256	THR
1	A	69	LEU
1	A	170	CYS
1	A	512	PRO
1	A	111	GLU
1	A	303	SER
1	A	471	ASN
1	A	283	PRO

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	435/466 (93%)	395 (91%)	40 (9%)	9 27

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	THR
1	A	39	VAL
1	A	45	THR
1	A	64	ASN
1	A	90	ARG
1	A	92	ARG
1	A	98	GLU

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Mol	Chain	Res	Type
1	A	122	THR
1	A	168	GLU
1	A	176	GLN
1	A	177	GLN
1	A	204	LYS
1	A	205	HIS
1	A	216	GLU
1	A	239	ASP
1	A	253	GLU
1	A	274	THR
1	A	276	GLU
1	A	279	THR
1	A	281	PHE
1	A	284	ARG
1	A	285	THR
1	A	310	ARG
1	A	313	LEU
1	A	332	SER
1	A	342	SER
1	A	359	LEU
1	A	397	LYS
1	A	399	SER
1	A	400	MET
1	A	403	LYS
1	A	409	ARG
1	A	443	ARG
1	A	474	LYS
1	A	476	GLN
1	A	493	CYS
1	A	499	ILE
1	A	511	LYS
1	A	517	ARG
1	A	536	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	87	GLN
1	A	104	HIS
1	A	145	GLN
1	A	176	GLN
1	A	179	GLN

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Mol	Chain	Res	Type
1	A	198	GLN
1	A	415	ASN
1	A	476	GLN
1	A	481	GLN
1	A	536	ASN
1	A	541	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

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	NAG	A	701	1	14,14,15	0.63	0	17,19,21	1.92	6 (35%)
2	NAG	A	702	1	14,14,15	0.96	1 (7%)	17,19,21	1.52	3 (17%)
2	NAG	A	703	1	14,14,15	1.18	1 (7%)	17,19,21	1.98	5 (29%)

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	701	1	-	0/6/23/26	0/1/1/1
2	NAG	A	702	1	-	3/6/23/26	0/1/1/1
2	NAG	A	703	1	-	2/6/23/26	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	703	NAG	C1-C2	3.45	1.57	1.52
2	A	702	NAG	C4-C5	2.44	1.58	1.53

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	703	NAG	C2-N2-C7	4.84	129.79	122.90
2	A	701	NAG	C1-O5-C5	3.77	117.31	112.19
2	A	702	NAG	C4-C3-C2	-3.17	106.36	111.02
2	A	703	NAG	O5-C5-C4	-3.06	103.39	110.83
2	A	703	NAG	C4-C3-C2	3.04	115.48	111.02
2	A	701	NAG	C4-C3-C2	2.97	115.37	111.02
2	A	701	NAG	O7-C7-C8	-2.68	117.08	122.06
2	A	701	NAG	O5-C5-C6	2.58	111.24	107.20
2	A	703	NAG	O7-C7-C8	-2.51	117.40	122.06
2	A	703	NAG	O7-C7-N2	2.44	126.43	121.95
2	A	701	NAG	C2-N2-C7	2.35	126.25	122.90
2	A	701	NAG	O3-C3-C4	-2.31	105.02	110.35
2	A	702	NAG	C1-O5-C5	-2.27	109.11	112.19
2	A	702	NAG	O4-C4-C5	2.24	114.85	109.30

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	703	NAG	C4-C5-C6-O6
2	A	703	NAG	O5-C5-C6-O6
2	A	702	NAG	C4-C5-C6-O6
2	A	702	NAG	O5-C5-C6-O6
2	A	702	NAG	C3-C2-N2-C7

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	702	NAG	1	0
2	A	703	NAG	2	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 (i)

6.1 Protein, DNA and RNA chains (i)

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å^2)	Q<0.9
1	A	508/557 (91%)	-0.00	35 (6%) 16 10	25, 44, 88, 114	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	279	THR	9.5
1	A	275	SER	6.9
1	A	282	ASN	6.4
1	A	52	THR	6.3
1	A	514	GLU	6.2
1	A	278	GLN	6.2
1	A	276	GLU	6.1
1	A	203	SER	6.0
1	A	205	HIS	6.0
1	A	197	ALA	5.9
1	A	202	LEU	5.6
1	A	518	GLY	5.2
1	A	515	GLY	5.2
1	A	274	THR	5.0
1	A	579	GLU	4.9
1	A	472	VAL	4.3
1	A	277	TYR	3.9
1	A	210	SER	3.9
1	A	199	ALA	3.8
1	A	207	ASP	3.7
1	A	208	PHE	3.6
1	A	198	GLN	3.5
1	A	280	PHE	3.5
1	A	53	PRO	3.5
1	A	201	ALA	3.4
1	A	281	PHE	3.3
1	A	578	GLY	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	204	LYS	2.9
1	A	200	LYS	2.8
1	A	257	GLY	2.7
1	A	209	ASN	2.6
1	A	34	SER	2.4
1	A	35	THR	2.3
1	A	304	LEU	2.2
1	A	36	ALA	2.1

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 monosaccharides 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. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	NAG	A	703	14/15	0.68	0.32	88,90,92,92	0
2	NAG	A	702	14/15	0.83	0.24	66,72,75,75	0
2	NAG	A	701	14/15	0.90	0.25	62,67,71,72	0

6.5 Other polymers (i)

There are no such residues in this entry.