



# Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 06:57 AM GMT

PDB ID : 3D12  
Title : Crystal Structures of Nipah Virus G Attachment Glycoprotein in Complex with its Receptor Ephrin-B3  
Authors : Xu, K.; Rajashankar, K.R.; Chan, Y.P.; Himanen, P.; Broder, C.C.; Nikolov, D.B.  
Deposited on : 2008-05-02  
Resolution : 3.00 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

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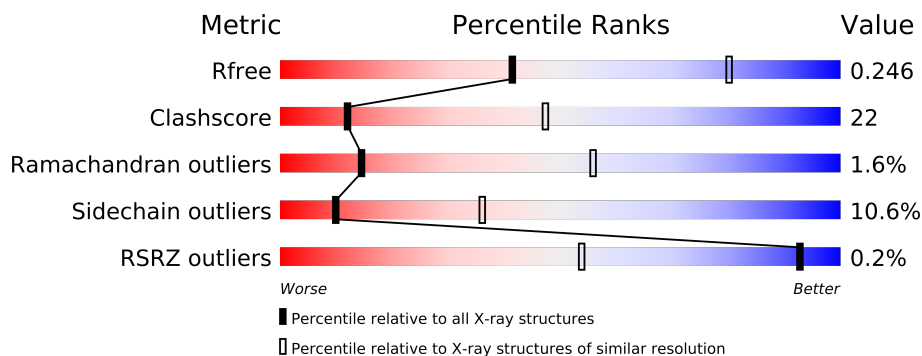
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 3.00 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))
$R_{free}$	66092	1216 (3.00-3.00)
Clashscore	79885	1594 (3.00-3.00)
Ramachandran outliers	78287	1537 (3.00-3.00)
Sidechain outliers	78261	1540 (3.00-3.00)
RSRZ outliers	66119	1217 (3.00-3.00)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	428	
1	D	428	
2	B	141	
2	E	141	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
5	NAG	A	1210	-	X
5	NAG	D	1310	-	X
7	SO4	A	1370	-	X
7	SO4	A	1373	-	X

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Mol	Type	Chain	Res	Geometry	Electron density
7	SO4	A	1374	-	X
7	SO4	A	1375	-	X
7	SO4	B	202	-	X
7	SO4	B	203	-	X
7	SO4	D	1471	-	X
7	SO4	D	1472	-	X
7	SO4	D	1473	-	X
7	SO4	D	1474	-	X
7	SO4	D	1477	-	X

## 2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 9598 atoms, of which 0 are hydrogen and 0 are deuterium.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Hemagglutinin-neuraminidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	428	Total	C	N	O	S	0	0	0
			3370	2146	568	635	21			
1	D	428	Total	C	N	O	S	0	0	0
			3370	2146	568	635	21			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	603	ALA	-	EXPRESSION TAG	UNP Q9IH62
D	603	ALA	-	EXPRESSION TAG	UNP Q9IH62

- Molecule 2 is a protein called Ephrin-B3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	141	Total	C	N	O	S	0	0	0
			1132	720	201	206	5			
2	E	141	Total	C	N	O	S	0	0	0
			1132	720	201	206	5			

- Molecule 3 is a polymer of unknown type called SUGAR (5-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	5	Total	C	N	O	0	0
			61	34	2	25		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	603	ALA	-	EXPRESSION TAG	UNP Q9IH62

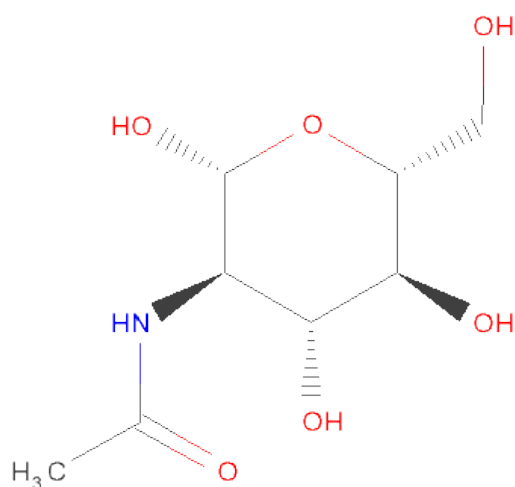
- Molecule 4 is a polymer of unknown type called SUGAR (7-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	7	Total	C	N	O	0	0
			83	46	2	35		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	603	ALA	-	EXPRESSION TAG	UNP Q9IH62

- Molecule 5 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula:  $C_8H_{15}NO_6$ ).



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

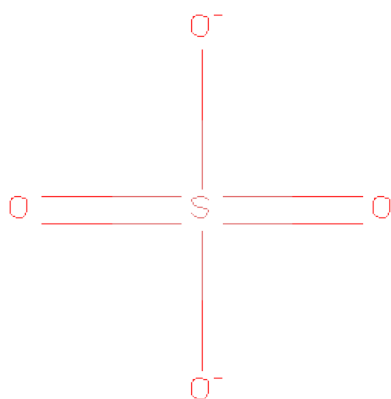
- Molecule 6 is a polymer of unknown type called SUGAR (2-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	2	Total	C	N	O	0	0
			28	16	2	10		
6	D	2	Total	C	N	O	0	0
			28	16	2	10		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	603	ALA	-	EXPRESSION TAG	UNP Q9IH62
D	603	ALA	-	EXPRESSION TAG	UNP Q9IH62

- Molecule 7 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	O	S	0	0
			5	4	1		
7	A	1	Total	O	S	0	0
			5	4	1		
7	A	1	Total	O	S	0	0
			5	4	1		
7	A	1	Total	O	S	0	0
			5	4	1		
7	A	1	Total	O	S	0	0
			5	4	1		
7	A	1	Total	O	S	0	0
			5	4	1		
7	A	1	Total	O	S	0	0
			5	4	1		
7	B	1	Total	O	S	0	0
			5	4	1		
7	B	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	D	1	Total	O	S	0	0
			5	4	1		
7	D	1	Total	O	S	0	0
			5	4	1		
7	D	1	Total	O	S	0	0
			5	4	1		
7	D	1	Total	O	S	0	0
			5	4	1		
7	D	1	Total	O	S	0	0
			5	4	1		
7	D	1	Total	O	S	0	0
			5	4	1		
7	D	1	Total	O	S	0	0
			5	4	1		
7	E	1	Total	O	S	0	0
			5	4	1		

- Molecule 8 is a polymer of unknown type called SUGAR (5-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	D	5	Total	C	N	O	0	0
			61	34	2	25		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	603	ALA	-	EXPRESSION TAG	UNP Q9IH62

- Molecule 9 is a polymer of unknown type called SUGAR (7-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
9	D	7	Total	C	N	O	0	0
			83	46	2	35		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	603	ALA	-	EXPRESSION TAG	UNP Q9IH62

- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	46	Total	O	0	0
			46	46		
10	B	9	Total	O	0	0
			9	9		
10	D	44	Total	O	0	0
			44	44		
10	E	13	Total	O	0	0
			13	13		

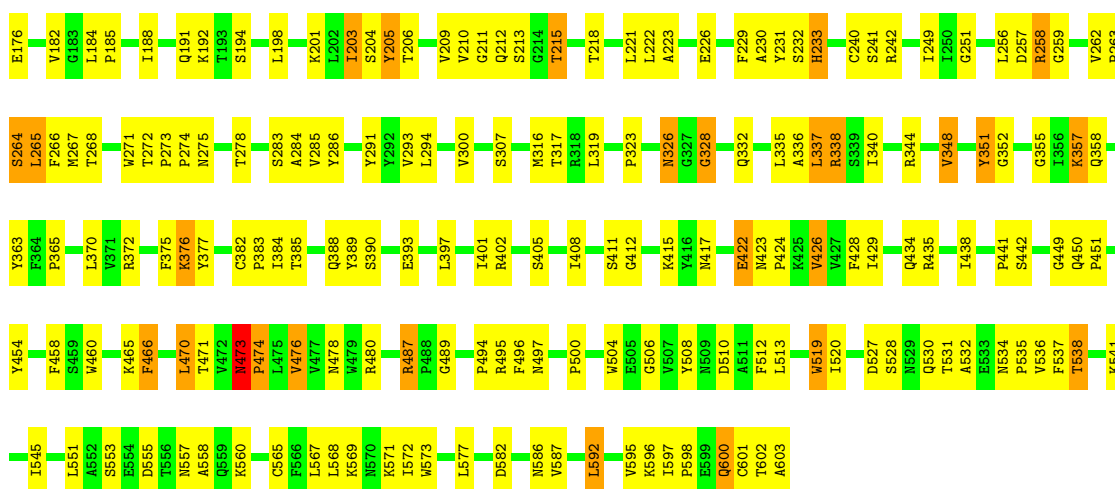


### 3 Residue-property plots

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.

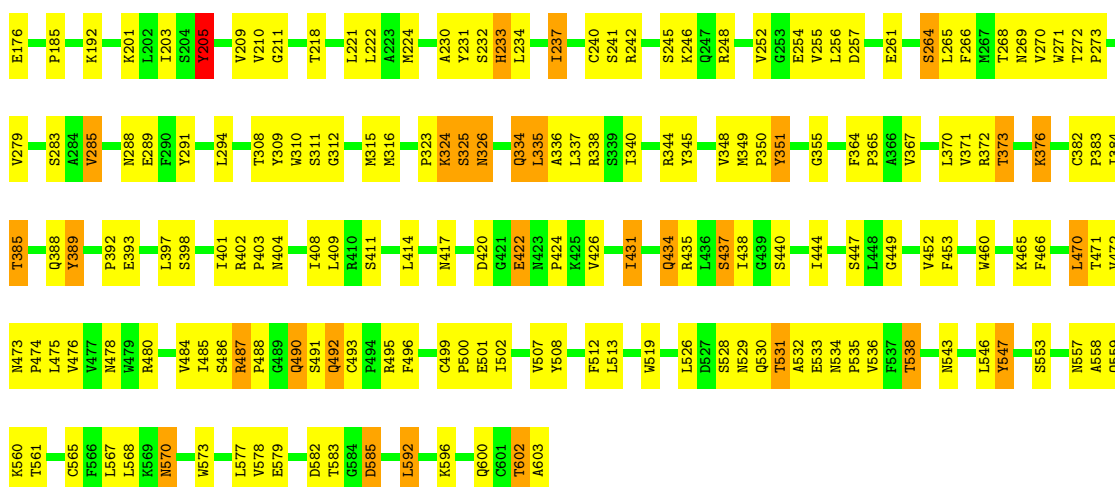
#### • Molecule 1: Hemagglutinin-neuraminidase

Chain A:



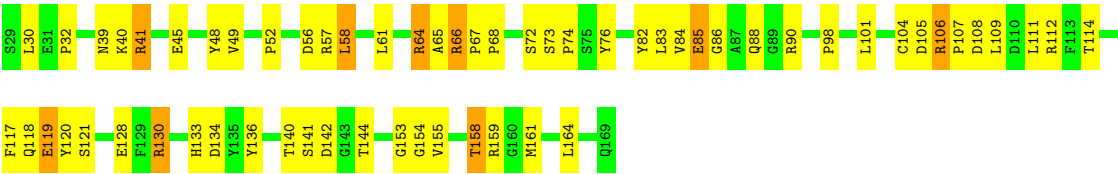
#### • Molecule 1: Hemagglutinin-neuraminidase

Chain D:



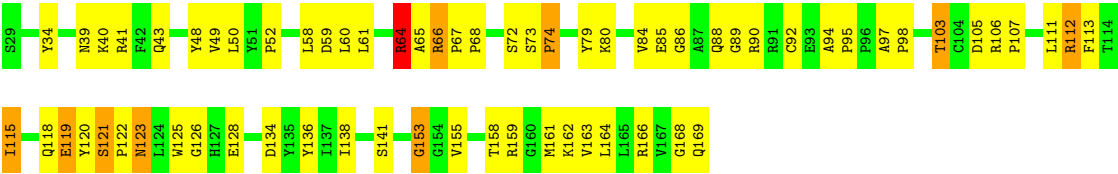
#### • Molecule 2: Ephrin-B3

Chain B:



• Molecule 2: Ephrin-B3

Chain E:



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	189.49Å 189.49Å 277.05Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	39.34 – 3.00 39.34 – 3.00	Depositor EDS
% Data completeness (in resolution range)	91.3 (39.34-3.00) 91.4 (39.34-3.00)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.57 (at 3.01Å)	Xtriage
Refinement program	PHENIX	Depositor
R, $R_{free}$	0.221 , 0.265 0.207 , 0.246	Depositor DCC
$R_{free}$ test set	2673 reflections (4.97%)	DCC
Wilson B-factor (Å <sup>2</sup> )	62.2	Xtriage
Anisotropy	0.466	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 37.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 58204 reflections (0.002%)	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	9598	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	66.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.63% of the height of the origin peak. No significant pseudotranslation is detected.*

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

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: BGC, NAG, GXL, GLC, LXZ, NGA, BMA, SO4, GL0, MAN, NGZ, LXB

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	0.46	1/3451 (0.0%)	0.54	0/4694
1	D	0.45	1/3451 (0.0%)	0.54	0/4694
2	B	0.37	0/1165	0.54	0/1582
2	E	0.39	0/1165	0.54	0/1582
All	All	0.44	2/9232 (0.0%)	0.54	0/12552

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
1	A	0	1
1	D	0	1
4	A	17	0
6	A	2	0
6	D	3	0
9	D	9	0
All	All	31	2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	176	GLU	CD-OE2	7.35	1.33	1.25
1	D	176	GLU	CD-OE2	7.19	1.33	1.25

There are no bond angle outliers.

All (31) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
4	A	1366	BGC	C5
4	A	1362	NGA	C2,C5,C3,C1,C4
4	A	1363	GL0	C1
4	A	1365	GLC	C2,C5,C3,C1,C4
4	A	1367	GLC	C2,C5,C3,C1,C4
6	A	1313	NAG	C1
6	A	1314	NAG	C1
9	D	1466	BGC	C5
9	D	1462	NGZ	C1
9	D	1463	GL0	C1
9	D	1465	GLC	C2,C5,C3,C1,C4
9	D	1467	GXL	C4
6	D	1413	NAG	C2,C1
6	D	1414	NAG	C1

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	473	ASN	Peptide
1	D	473	ASN	Peptide

## 5.2 Close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3370	0	3313	145	0
1	D	3370	0	3313	143	0
2	B	1132	0	1094	53	0
2	E	1132	0	1094	66	0
3	A	61	0	52	1	0
4	A	83	0	70	1	0
5	A	14	0	13	0	0
5	D	14	0	13	0	0
6	A	28	0	25	2	0
6	D	28	0	25	1	0
7	A	45	0	0	4	0
7	B	10	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	D	50	0	0	0	0
7	E	5	0	0	1	0
8	D	61	0	52	3	0
9	D	83	0	70	5	0
10	A	46	0	0	2	0
10	B	9	0	0	0	0
10	D	44	0	0	2	0
10	E	13	0	0	1	0
All	All	9598	0	9134	401	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 22.

All (401) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:376:LYS:H	1:A:376:LYS:HD3	1.26	0.97
1:D:403:PRO:HG3	1:D:502:ILE:HD11	1.44	0.97
1:A:376:LYS:N	1:A:376:LYS:HD3	1.81	0.94
1:D:531:THR:HG21	1:D:533:GLU:OE2	1.71	0.90
2:B:86:GLY:HA2	2:B:136:TYR:HD1	1.37	0.90
2:B:66:ARG:HB3	2:B:67:PRO:HD3	1.54	0.89
1:A:602:THR:HG23	1:A:603:ALA:H	1.40	0.87
2:E:118:GLN:HE22	2:E:120:TYR:HB2	1.40	0.86
2:B:118:GLN:NE2	2:B:120:TYR:HB2	1.89	0.86
1:D:376:LYS:HD3	1:D:376:LYS:H	1.42	0.84
2:E:66:ARG:HB3	2:E:67:PRO:HD3	1.57	0.84
2:B:66:ARG:HB3	2:B:67:PRO:CD	2.07	0.83
1:D:376:LYS:N	1:D:376:LYS:HD3	1.93	0.83
2:B:86:GLY:HA2	2:B:136:TYR:CD1	2.14	0.82
2:E:113:PHE:CE2	2:E:115:ILE:HG12	2.14	0.82
1:A:210:VAL:HG12	1:A:211:GLY:H	1.43	0.82
2:B:118:GLN:HG2	2:B:119:GLU:N	1.96	0.81
1:D:471:THR:HG23	1:D:474:PRO:O	1.81	0.81
1:D:376:LYS:H	1:D:376:LYS:CD	1.95	0.79
2:E:113:PHE:HE2	2:E:115:ILE:HG12	1.45	0.78
1:A:363:TYR:CE2	1:A:415:LYS:HG3	2.18	0.77
2:E:141:SER:O	2:E:155:VAL:HG23	1.85	0.77
1:A:557:ASN:HB3	2:B:118:GLN:NE2	1.99	0.76
1:D:422:GLU:CD	1:D:422:GLU:H	1.88	0.76
2:E:118:GLN:NE2	2:E:120:TYR:HB2	1.99	0.76
1:D:402:ARG:HH12	1:D:404:ASN:HB2	1.50	0.76
2:B:130:ARG:HG2	2:B:130:ARG:HH11	1.51	0.75

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
8:D:1240:BMA:H2	8:D:1241:BGC:H2	1.69	0.74
2:B:39:ASN:OD1	2:B:41:ARG:HB2	1.86	0.74
2:E:118:GLN:HG2	2:E:119:GLU:H	1.50	0.74
1:A:535:PRO:HG3	1:A:558:ALA:HB2	1.68	0.74
1:D:585:ASP:HB3	10:D:1515:HOH:O	1.88	0.73
1:D:557:ASN:HB3	2:E:118:GLN:NE2	2.03	0.72
1:D:531:THR:CG2	1:D:533:GLU:HG2	2.20	0.72
2:E:66:ARG:O	2:E:68:PRO:HD3	1.89	0.72
1:D:384:ILE:HD11	1:D:392:PRO:HA	1.72	0.71
1:A:535:PRO:HG3	1:A:558:ALA:CB	2.21	0.71
2:B:66:ARG:O	2:B:68:PRO:HD3	1.91	0.70
2:B:64:ARG:HD2	2:B:67:PRO:HD2	1.73	0.70
2:E:118:GLN:HG2	2:E:119:GLU:N	2.03	0.70
2:E:39:ASN:O	2:E:40:LYS:HB3	1.91	0.70
1:D:403:PRO:CG	1:D:502:ILE:HD11	2.21	0.70
2:E:79:TYR:HE2	2:E:155:VAL:HG21	1.56	0.70
2:E:66:ARG:HB3	2:E:67:PRO:CD	2.22	0.69
2:E:64:ARG:CD	2:E:67:PRO:HD2	2.23	0.69
2:B:73:SER:HB2	2:B:74:PRO:HD2	1.73	0.69
1:D:242:ARG:HD3	2:E:126:GLY:O	1.92	0.68
8:D:1240:BMA:C2	8:D:1241:BGC:H2	2.20	0.68
1:D:471:THR:HG22	1:D:476:VAL:CG1	2.23	0.68
1:A:223:ALA:HB3	1:A:230:ALA:HB3	1.75	0.68
1:A:471:THR:HG22	1:A:476:VAL:HG13	1.77	0.67
2:E:88:GLN:NE2	2:E:95:PRO:HG3	2.10	0.67
1:D:531:THR:HG22	1:D:533:GLU:HG2	1.77	0.67
1:D:370:LEU:HB3	1:D:408:ILE:HD11	1.76	0.67
1:A:256:LEU:HD12	1:A:266:PHE:CD2	2.30	0.67
1:D:536:VAL:HG12	1:D:538:THR:HG22	1.77	0.66
1:D:240:CYS:SG	2:E:119:GLU:HG2	2.36	0.66
2:E:155:VAL:HG12	2:E:159:ARG:HB2	1.77	0.66
2:E:49:VAL:HG13	2:E:164:LEU:HD23	1.78	0.66
2:E:89:GLY:HA3	2:E:136:TYR:CD1	2.31	0.66
1:A:471:THR:HG23	1:A:474:PRO:O	1.96	0.66
2:E:85:GLU:HB2	2:E:88:GLN:HG3	1.78	0.66
1:A:557:ASN:HB3	2:B:118:GLN:HE21	1.61	0.65
1:A:215:THR:HG22	1:A:587:VAL:HG23	1.78	0.65
1:D:558:ALA:HA	1:D:579:GLU:O	1.96	0.65
1:A:206:THR:HG21	1:A:264:SER:OG	1.96	0.65
2:E:84:VAL:HG13	2:E:88:GLN:HB2	1.78	0.65
1:A:226:GLU:OE1	1:A:286:TYR:HE2	1.80	0.65
1:D:471:THR:HG22	1:D:476:VAL:HG12	1.78	0.65

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:437:SER:HB2	1:D:465:LYS:HZ2	1.62	0.65
1:A:508:TYR:HB3	1:A:560:LYS:HD3	1.77	0.64
1:D:372:ARG:NH2	1:D:397:LEU:HA	2.12	0.64
2:B:66:ARG:HB2	2:B:107:PRO:O	1.98	0.64
2:B:105:ASP:HB2	2:B:106:ARG:HE	1.63	0.64
1:D:508:TYR:HB3	1:D:560:LYS:HD3	1.80	0.64
1:D:557:ASN:HB3	2:E:118:GLN:HE21	1.63	0.64
1:A:388:GLN:O	1:A:389:TYR:HB2	1.99	0.63
2:E:88:GLN:HE22	2:E:95:PRO:HG3	1.64	0.63
2:E:105:ASP:HB2	2:E:106:ARG:HE	1.64	0.63
1:D:398:SER:OG	1:D:502:ILE:HA	1.99	0.63
1:D:535:PRO:HG3	1:D:558:ALA:HB2	1.81	0.62
1:D:437:SER:HB2	1:D:465:LYS:NZ	2.14	0.62
2:B:64:ARG:HD3	2:B:108:ASP:O	2.00	0.62
2:E:65:ALA:HB2	2:E:72:SER:HB2	1.79	0.62
2:B:118:GLN:HE22	2:B:120:TYR:HB2	1.64	0.62
2:E:80:LYS:HE2	2:E:103:THR:HG23	1.81	0.62
1:A:344:ARG:NH2	1:A:376:LYS:HD2	2.15	0.61
2:E:79:TYR:CE2	2:E:155:VAL:HG21	2.35	0.61
2:B:104:CYS:HA	2:B:111:LEU:HD12	1.81	0.61
1:D:388:GLN:O	1:D:389:TYR:HB2	1.98	0.61
1:D:512:PHE:CD2	1:D:565:CYS:HB2	2.36	0.61
1:D:210:VAL:HG12	1:D:211:GLY:H	1.65	0.61
1:D:310:TRP:CZ3	1:D:312:GLY:HA2	2.36	0.61
1:A:240:CYS:SG	2:B:119:GLU:HG2	2.41	0.60
9:D:1461:LXB:O6	9:D:1461:LXB:O4	2.19	0.59
1:D:471:THR:CG2	1:D:476:VAL:HG12	2.32	0.59
1:A:402:ARG:HD2	7:A:1368:SO4:O2	2.03	0.59
1:A:351:TYR:CE2	1:A:441:PRO:HD3	2.37	0.59
1:A:405:SER:HA	7:A:1368:SO4:O2	2.02	0.59
2:E:86:GLY:HA2	2:E:136:TYR:HD1	1.67	0.59
1:A:372:ARG:HH21	1:A:397:LEU:HA	1.67	0.59
1:A:495:ARG:O	1:A:496:PHE:HB2	2.03	0.59
1:A:489:GLY:O	1:A:530:GLN:HA	2.02	0.58
2:B:39:ASN:O	2:B:40:LYS:HB3	2.03	0.58
1:A:212:GLN:H	1:A:212:GLN:CD	2.07	0.58
1:A:242:ARG:HE	2:B:128:GLU:HG2	1.68	0.58
1:A:537:PHE:HB2	1:A:551:LEU:HD21	1.86	0.58
1:A:210:VAL:HG12	1:A:211:GLY:N	2.17	0.58
1:A:231:TYR:OH	1:A:233:HIS:HD2	1.87	0.58
1:D:492:GLN:CA	1:D:492:GLN:HE21	2.17	0.58
1:A:205:TYR:CE1	1:A:206:THR:HG23	2.39	0.57

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:490:GLN:OE1	2:E:123:ASN:HB2	2.04	0.57
1:D:577:LEU:HD12	1:D:592:LEU:HD11	1.86	0.57
1:A:352:GLY:HA3	1:A:442:SER:O	2.03	0.57
1:A:348:VAL:HG21	1:A:428:PHE:HB2	1.86	0.56
1:A:184:LEU:HD12	1:D:472:VAL:HG23	1.86	0.56
1:D:484:VAL:HG23	1:D:485:ILE:HG12	1.87	0.56
1:A:528:SER:OG	1:A:531:THR:HG22	2.05	0.56
1:D:348:VAL:HG11	1:D:414:LEU:HD21	1.87	0.56
9:D:1463:GL0:H4	9:D:1464:BGC:O5	2.05	0.56
1:A:372:ARG:NH2	1:A:397:LEU:HA	2.20	0.56
1:A:553:SER:OG	10:A:48:HOH:O	2.18	0.56
1:A:232:SER:HA	1:A:249:ILE:O	2.07	0.55
9:D:1461:LXB:H8	10:D:1511:HOH:O	2.07	0.55
2:B:30:LEU:HD11	2:B:52:PRO:HB3	1.87	0.55
1:A:600:GLN:HB3	1:A:602:THR:HG22	1.89	0.55
6:D:1414:NAG:H82	6:D:1414:NAG:O3	2.07	0.54
2:E:34:TYR:CE2	2:E:61:LEU:HD12	2.42	0.54
1:D:271:TRP:CZ3	1:D:273:PRO:HG3	2.42	0.54
1:A:251:GLY:HA3	1:A:267:MET:HE1	1.90	0.54
1:D:370:LEU:HB3	1:D:408:ILE:CD1	2.37	0.54
1:A:417:ASN:O	1:A:424:PRO:HB3	2.08	0.54
1:D:417:ASN:HB3	1:D:420:ASP:HB2	1.90	0.54
1:A:536:VAL:O	1:A:538:THR:HG22	2.08	0.54
2:E:134:ASP:HB3	2:E:164:LEU:HD11	1.90	0.54
2:E:80:LYS:HG2	2:E:103:THR:HG23	1.90	0.54
1:A:338:ARG:HH11	1:A:423:ASN:HB3	1.72	0.53
1:A:577:LEU:HD23	1:A:577:LEU:C	2.29	0.53
1:D:344:ARG:NH2	1:D:376:LYS:HD2	2.23	0.53
1:A:283:SER:OG	1:A:365:PRO:HD2	2.08	0.53
1:D:231:TYR:OH	1:D:233:HIS:HD2	1.92	0.53
1:A:372:ARG:NH1	1:A:405:SER:OG	2.42	0.53
1:D:397:LEU:HB3	1:D:502:ILE:HD12	1.91	0.53
1:D:269:ASN:HD22	1:D:323:PRO:HG3	1.73	0.53
1:D:470:LEU:CD1	1:D:478:ASN:HB2	2.39	0.53
1:A:294:LEU:HD13	1:A:316:MET:HE3	1.91	0.52
2:E:50:LEU:C	2:E:52:PRO:HD3	2.30	0.52
2:E:64:ARG:HD3	2:E:67:PRO:HD2	1.89	0.52
1:D:309:TYR:CD2	8:D:1238:NAG:H5	2.44	0.52
1:D:264:SER:HB3	1:D:266:PHE:CE2	2.44	0.52
2:B:141:SER:O	2:B:155:VAL:HG23	2.08	0.52
1:A:450:GLN:HG3	1:A:451:PRO:HD2	1.90	0.52
1:D:230:ALA:HA	1:D:252:VAL:HG12	1.92	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:340:ILE:HA	1:A:426:VAL:HG23	1.90	0.52
2:B:130:ARG:CG	2:B:130:ARG:HH11	2.22	0.52
1:D:408:ILE:O	1:D:408:ILE:HD12	2.10	0.52
2:B:84:VAL:HG13	2:B:88:GLN:HB2	1.90	0.52
1:A:258:ARG:HG2	1:A:262:VAL:O	2.10	0.52
1:A:401:ILE:HG22	1:A:504:TRP:CE2	2.45	0.52
2:E:50:LEU:O	2:E:52:PRO:HD3	2.09	0.52
1:A:408:ILE:HD12	1:A:408:ILE:C	2.30	0.52
1:D:491:SER:HB3	2:E:113:PHE:HB2	1.91	0.52
1:D:561:THR:HG22	1:D:578:VAL:HG13	1.92	0.52
1:D:351:TYR:HB2	1:D:367:VAL:HG12	1.92	0.51
1:A:449:GLY:HA3	1:D:449:GLY:HA3	1.92	0.51
1:A:328:GLY:HA2	10:A:15:HOH:O	2.11	0.51
2:E:106:ARG:HG2	2:E:111:LEU:HD11	1.92	0.51
1:D:480:ARG:HD3	1:D:519:TRP:CH2	2.46	0.51
1:D:237:ILE:HD11	1:D:246:LYS:HB2	1.91	0.51
1:D:397:LEU:CB	1:D:502:ILE:HD12	2.40	0.51
2:E:123:ASN:ND2	2:E:125:TRP:H	2.08	0.51
1:D:348:VAL:CG1	1:D:414:LEU:HD21	2.41	0.51
1:A:188:ILE:HG21	1:A:572:ILE:HD12	1.93	0.51
1:D:531:THR:HG21	1:D:533:GLU:HG2	1.92	0.51
2:E:66:ARG:HB2	2:E:107:PRO:O	2.11	0.51
2:B:118:GLN:HG2	2:B:120:TYR:H	1.75	0.51
1:D:335:LEU:N	1:D:335:LEU:HD12	2.23	0.51
2:B:118:GLN:HG2	2:B:119:GLU:H	1.72	0.51
2:B:155:VAL:HA	2:B:158:THR:HB	1.93	0.51
1:D:294:LEU:HD13	1:D:316:MET:HE2	1.92	0.50
1:D:422:GLU:O	1:D:424:PRO:HD3	2.10	0.50
1:A:233:HIS:C	1:A:233:HIS:ND1	2.65	0.50
1:D:372:ARG:HH21	1:D:397:LEU:HA	1.76	0.50
1:D:431:ILE:HA	1:D:475:LEU:HD23	1.94	0.50
1:D:401:ILE:HG13	1:D:402:ARG:HG3	1.91	0.50
6:A:1314:NAG:H5	7:A:1371:SO4:O2	2.11	0.50
1:A:264:SER:HB3	1:A:266:PHE:CE2	2.46	0.50
2:E:80:LYS:HE2	2:E:103:THR:CG2	2.41	0.50
4:A:1362:NGA:H62	4:A:1367:GLC:H2	1.94	0.50
1:D:224:MET:HE3	1:D:255:VAL:HG21	1.93	0.50
1:A:388:GLN:O	1:A:389:TYR:CB	2.59	0.50
1:A:429:ILE:HD13	1:A:473:ASN:HA	1.93	0.50
1:D:496:PHE:HD2	9:D:1461:LXB:H5	1.76	0.49
1:D:559:GLN:HB3	1:D:579:GLU:OE1	2.12	0.49
1:A:185:PRO:HB3	1:A:567:LEU:HD21	1.93	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:E:138:ILE:HG22	2:E:162:LYS:HB3	1.93	0.49
1:D:460:TRP:CB	1:D:501:GLU:H	2.26	0.49
1:A:257:ASP:C	1:A:259:GLY:H	2.16	0.49
2:B:65:ALA:HB2	2:B:72:SER:HB2	1.93	0.49
1:A:221:LEU:C	1:A:221:LEU:HD23	2.33	0.49
1:A:389:TYR:CG	2:B:106:ARG:HD2	2.48	0.49
1:A:423:ASN:N	1:A:424:PRO:HD3	2.28	0.49
2:B:155:VAL:HG11	2:B:161:MET:SD	2.52	0.49
2:E:138:ILE:HG22	2:E:162:LYS:CB	2.43	0.49
1:D:340:ILE:HG12	1:D:426:VAL:CG2	2.43	0.49
1:D:447:SER:HB3	1:D:513:LEU:HD23	1.95	0.49
1:A:376:LYS:H	1:A:376:LYS:CD	2.13	0.49
1:D:486:SER:HB3	1:D:495:ARG:HG3	1.94	0.48
1:A:192:LYS:HD2	1:A:545:ILE:O	2.13	0.48
1:D:324:LYS:O	1:D:326:ASN:N	2.43	0.48
1:A:602:THR:HG23	1:A:603:ALA:N	2.19	0.48
1:D:349:MET:HB2	1:D:350:PRO:HD2	1.95	0.48
2:B:32:PRO:HB3	2:B:61:LEU:HD11	1.95	0.48
1:D:285:VAL:HG13	1:D:364:PHE:CZ	2.48	0.48
3:A:1140:BMA:H2	3:A:1141:BMA:O2	2.13	0.48
2:B:76:TYR:CD1	2:B:107:PRO:HA	2.49	0.48
1:A:218:THR:HG21	1:A:240:CYS:HB3	1.95	0.48
2:E:86:GLY:HA2	2:E:136:TYR:CD1	2.47	0.48
2:B:104:CYS:HA	2:B:111:LEU:CD1	2.43	0.48
1:D:257:ASP:OD2	1:D:568:LEU:HD21	2.14	0.48
1:D:257:ASP:OD2	1:D:261:GLU:HA	2.13	0.48
1:D:417:ASN:O	1:D:424:PRO:HB3	2.14	0.48
2:E:85:GLU:OE1	2:E:85:GLU:HA	2.14	0.48
1:A:348:VAL:HG21	1:A:428:PHE:CB	2.43	0.48
1:A:600:GLN:NE2	1:A:600:GLN:HA	2.29	0.48
1:A:212:GLN:O	1:A:215:THR:HB	2.13	0.48
1:D:486:SER:HB3	1:D:495:ARG:CG	2.44	0.48
1:D:460:TRP:HB3	1:D:501:GLU:H	1.78	0.48
1:D:294:LEU:HB2	1:D:316:MET:HE2	1.95	0.48
1:A:470:LEU:HD13	1:A:478:ASN:HB2	1.94	0.48
1:D:234:LEU:HD11	1:D:245:SER:HB3	1.96	0.47
1:D:495:ARG:O	1:D:496:PHE:HB2	2.14	0.47
1:A:340:ILE:HG12	1:A:426:VAL:HG21	1.96	0.47
1:D:254:GLU:HG3	1:D:256:LEU:HD11	1.96	0.47
2:E:84:VAL:HG13	2:E:85:GLU:N	2.30	0.47
1:D:492:GLN:NE2	1:D:492:GLN:HA	2.28	0.47
1:D:535:PRO:HG3	1:D:558:ALA:CB	2.44	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:345:TYR:CD1	1:D:370:LEU:HB2	2.50	0.47
1:A:597:ILE:HA	1:A:598:PRO:HD3	1.70	0.47
1:A:573:TRP:CZ2	1:A:596:LYS:HB2	2.49	0.47
1:A:460:TRP:CD2	1:A:500:PRO:HA	2.49	0.47
2:E:158:THR:HG22	2:E:159:ARG:HG3	1.97	0.47
1:A:376:LYS:N	1:A:376:LYS:CD	2.63	0.47
1:D:283:SER:OG	1:D:365:PRO:HD2	2.15	0.47
2:E:159:ARG:HB2	2:E:161:MET:HG3	1.96	0.47
1:D:491:SER:HA	1:D:530:GLN:NE2	2.30	0.47
1:D:547:TYR:N	1:D:547:TYR:CD2	2.83	0.47
2:E:85:GLU:HB2	2:E:88:GLN:CG	2.45	0.46
1:D:480:ARG:HD3	1:D:519:TRP:CZ3	2.49	0.46
1:A:471:THR:CG2	1:A:476:VAL:HG13	2.45	0.46
2:E:92:CYS:O	2:E:153:GLY:N	2.47	0.46
1:D:488:PRO:HA	1:D:529:ASN:O	2.15	0.46
1:A:348:VAL:HG21	1:A:428:PHE:CG	2.51	0.46
1:A:203:ILE:HG12	1:A:592:LEU:HB3	1.97	0.46
2:B:64:ARG:CD	2:B:67:PRO:HD2	2.43	0.46
2:B:85:GLU:HB2	2:B:88:GLN:CD	2.36	0.46
1:D:532:ALA:HB1	1:D:558:ALA:O	2.15	0.46
1:D:334:GLN:HE21	1:D:334:GLN:HB3	1.52	0.46
1:A:422:GLU:CD	1:A:422:GLU:H	2.20	0.46
1:D:585:ASP:N	1:D:585:ASP:OD1	2.45	0.46
1:A:229:PHE:CD2	1:A:265:LEU:HD13	2.51	0.46
1:D:388:GLN:O	1:D:389:TYR:CB	2.64	0.46
1:D:408:ILE:C	1:D:408:ILE:HD12	2.37	0.45
1:A:268:THR:O	1:A:323:PRO:HG2	2.17	0.45
1:A:274:PRO:HG2	1:A:275:ASN:H	1.81	0.45
1:A:454:TYR:CE1	1:A:510:ASP:HA	2.51	0.45
1:D:492:GLN:CA	1:D:492:GLN:NE2	2.79	0.45
1:D:248:ARG:O	1:D:272:THR:HG22	2.16	0.45
1:D:470:LEU:HD11	1:D:478:ASN:HB2	1.99	0.45
1:D:221:LEU:HD23	1:D:221:LEU:C	2.37	0.45
1:A:532:ALA:HB1	1:A:558:ALA:O	2.16	0.45
1:A:275:ASN:ND2	1:A:278:THR:HG23	2.31	0.45
2:B:56:ASP:O	2:B:117:PHE:HD1	2.00	0.45
1:A:294:LEU:HB2	1:A:316:MET:CE	2.47	0.45
1:D:460:TRP:CG	1:D:500:PRO:HA	2.52	0.45
1:D:434:GLN:O	1:D:435:ARG:C	2.55	0.45
1:D:192:LYS:HA	1:D:546:LEU:HA	1.98	0.45
1:A:351:TYR:CZ	1:A:441:PRO:HD3	2.52	0.45
1:D:222:LEU:HD12	1:D:230:ALA:O	2.17	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:494:PRO:HD2	1:A:497:ASN:HB2	1.98	0.45
2:B:140:THR:HA	2:B:154:GLY:H	1.82	0.45
1:A:513:LEU:HD12	1:A:520:ILE:O	2.17	0.45
2:B:130:ARG:CG	2:B:130:ARG:NH1	2.80	0.44
1:D:496:PHE:CD2	9:D:1461:LXB:H5	2.50	0.44
1:A:213:SER:OG	1:A:586:ASN:HB3	2.17	0.44
1:D:355:GLY:HA2	1:D:444:ILE:HG23	1.99	0.44
1:A:527:ASP:HB3	1:A:534:ASN:ND2	2.32	0.44
1:D:370:LEU:HD23	1:D:408:ILE:HD11	1.99	0.44
1:A:231:TYR:OH	1:A:233:HIS:CD2	2.68	0.44
1:D:460:TRP:CD2	1:D:500:PRO:HA	2.53	0.44
2:B:82:TYR:CD2	2:B:98:PRO:HB2	2.52	0.44
1:D:315:MET:HG2	1:D:316:MET:N	2.31	0.44
2:B:90:ARG:HD2	7:B:203:SO4:O4	2.18	0.44
1:D:487:ARG:NH2	1:D:493:CYS:O	2.44	0.44
1:A:592:LEU:HA	1:A:592:LEU:HD12	1.89	0.44
1:D:600:GLN:HA	1:D:600:GLN:NE2	2.32	0.44
1:A:422:GLU:C	1:A:424:PRO:HD3	2.38	0.44
1:A:412:GLY:HA3	1:A:429:ILE:O	2.17	0.44
1:A:569:LYS:O	1:A:571:LYS:HG3	2.18	0.44
1:A:294:LEU:HB2	1:A:316:MET:HE1	2.00	0.44
1:D:499:CYS:HA	1:D:500:PRO:HD3	1.91	0.44
1:A:271:TRP:CZ3	1:A:273:PRO:HG3	2.53	0.44
1:D:294:LEU:HD13	1:D:316:MET:CE	2.48	0.44
2:E:66:ARG:CB	2:E:67:PRO:HD3	2.40	0.44
2:B:83:LEU:HB2	2:B:101:LEU:HD11	2.00	0.44
1:A:335:LEU:HG	1:A:337:LEU:H	1.82	0.44
1:D:531:THR:HG21	1:D:533:GLU:CD	2.38	0.44
2:E:158:THR:HG21	7:E:190:SO4:O3	2.18	0.44
1:A:458:PHE:O	1:A:504:TRP:HA	2.18	0.44
1:D:528:SER:OG	1:D:531:THR:HB	2.17	0.43
1:D:409:LEU:HD12	1:D:438:ILE:HD11	1.99	0.43
1:A:401:ILE:HG13	1:A:402:ARG:HG3	2.00	0.43
1:A:375:PHE:CE2	1:A:377:TYR:HB2	2.53	0.43
2:E:94:ALA:HA	2:E:95:PRO:HD3	1.85	0.43
1:A:191:GLN:HB2	1:A:601:CYS:SG	2.57	0.43
1:A:229:PHE:CG	1:A:265:LEU:HD13	2.53	0.43
1:D:310:TRP:CE3	1:D:312:GLY:HA2	2.53	0.43
2:B:49:VAL:HG22	2:B:164:LEU:HB3	2.00	0.43
2:E:169:GLN:HG2	10:E:200:HOH:O	2.18	0.43
1:A:510:ASP:C	1:A:510:ASP:OD1	2.56	0.43
1:D:376:LYS:CE	1:D:376:LYS:H	2.31	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:E:105:ASP:HB2	2:E:106:ARG:NE	2.32	0.43
1:A:338:ARG:NH1	1:A:423:ASN:HB3	2.33	0.43
1:A:336:ALA:O	1:A:337:LEU:HB2	2.18	0.43
1:D:573:TRP:CZ2	1:D:596:LYS:HB2	2.54	0.43
1:A:205:TYR:N	1:A:205:TYR:CD1	2.86	0.43
1:D:470:LEU:HD13	1:D:478:ASN:HB2	2.01	0.43
1:A:319:LEU:HD22	1:A:332:GLN:HG2	2.01	0.43
1:D:255:VAL:HG22	1:D:265:LEU:HD22	2.01	0.43
1:D:271:TRP:CZ2	1:D:334:GLN:OE1	2.72	0.42
1:D:602:THR:OG1	1:D:603:ALA:N	2.51	0.42
2:B:66:ARG:CB	2:B:67:PRO:CD	2.91	0.42
1:D:420:ASP:HB3	1:D:424:PRO:HB3	2.01	0.42
1:A:358:GLN:HG3	1:D:570:ASN:ND2	2.34	0.42
2:B:58:LEU:O	2:B:114:THR:HA	2.19	0.42
1:A:326:ASN:HD22	1:A:326:ASN:HA	1.65	0.42
1:A:294:LEU:HD13	1:A:316:MET:CE	2.48	0.42
2:B:159:ARG:HD2	2:B:161:MET:HE3	2.02	0.42
1:A:382:CYS:HA	1:A:383:PRO:HD3	1.78	0.42
1:D:336:ALA:HB3	1:D:338:ARG:CZ	2.49	0.42
1:A:466:PHE:CZ	1:A:519:TRP:HZ3	2.38	0.42
1:A:471:THR:HG22	1:A:476:VAL:CG1	2.47	0.42
2:B:76:TYR:CE1	2:B:107:PRO:HA	2.54	0.42
1:D:384:ILE:O	1:D:384:ILE:HG22	2.19	0.42
1:A:370:LEU:HB3	1:A:408:ILE:HD11	2.01	0.42
1:A:434:GLN:HE21	1:A:434:GLN:HB3	1.65	0.42
1:D:185:PRO:HB3	1:D:567:LEU:HD21	2.01	0.42
1:A:223:ALA:O	1:A:229:PHE:HA	2.20	0.42
2:E:60:LEU:O	2:E:112:ARG:HA	2.19	0.42
1:A:402:ARG:NH1	7:A:1368:SO4:O1	2.53	0.42
1:A:512:PHE:CD2	1:A:565:CYS:HB2	2.54	0.42
2:B:142:ASP:OD1	2:B:144:THR:HG23	2.20	0.42
1:A:555:ASP:HB3	2:B:57:ARG:HH22	1.85	0.42
2:E:64:ARG:HD2	2:E:67:PRO:HD2	1.99	0.41
1:A:508:TYR:CZ	1:A:510:ASP:HB3	2.55	0.41
1:D:452:VAL:HG12	1:D:453:PHE:N	2.35	0.41
1:A:435:ARG:HG2	1:A:465:LYS:HE3	2.02	0.41
1:D:268:THR:O	1:D:323:PRO:HG2	2.20	0.41
1:A:600:GLN:HA	1:A:600:GLN:HE21	1.85	0.41
1:D:248:ARG:HB3	1:D:273:PRO:HD2	2.01	0.41
1:A:384:ILE:HG13	1:A:390:SER:HB3	2.02	0.41
1:A:470:LEU:CD1	1:A:478:ASN:HB2	2.51	0.41
1:A:465:LYS:HE2	1:A:465:LYS:HB3	1.85	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:198:LEU:HD22	1:A:595:VAL:HG11	2.01	0.41
1:A:335:LEU:HD23	1:A:337:LEU:HD22	2.03	0.41
2:E:97:ALA:N	2:E:98:PRO:HD3	2.36	0.41
1:A:265:LEU:HD23	1:A:265:LEU:N	2.34	0.41
1:A:284:ALA:HB2	1:A:293:VAL:HG22	2.01	0.41
1:D:288:ASN:O	1:D:289:GLU:HB2	2.20	0.41
1:A:355:GLY:HA3	1:A:363:TYR:O	2.21	0.41
1:A:286:TYR:O	1:A:357:LYS:HG2	2.21	0.41
2:E:50:LEU:HD12	2:E:163:VAL:HG11	2.03	0.41
2:E:90:ARG:C	2:E:92:CYS:H	2.23	0.41
1:A:300:VAL:O	1:A:300:VAL:HG23	2.21	0.41
1:D:218:THR:HG21	1:D:240:CYS:HB3	2.03	0.41
2:E:155:VAL:HG11	2:E:161:MET:SD	2.60	0.41
2:E:43:GLN:OE1	2:E:159:ARG:HA	2.20	0.41
1:D:512:PHE:CE2	1:D:565:CYS:HB2	2.55	0.41
1:A:221:LEU:HD23	1:A:222:LEU:N	2.36	0.41
1:A:316:MET:HG3	1:A:317:THR:N	2.35	0.41
2:E:73:SER:HB2	2:E:74:PRO:HD2	2.02	0.41
1:D:526:LEU:HA	1:D:534:ASN:O	2.21	0.41
1:A:487:ARG:HH11	1:A:506:GLY:HA3	1.85	0.41
1:A:263:PRO:HD3	1:A:568:LEU:HD23	2.03	0.41
2:E:59:ASP:HA	2:E:113:PHE:O	2.21	0.41
1:D:221:LEU:HB3	1:D:232:SER:HB3	2.03	0.41
2:E:121:SER:HA	2:E:122:PRO:HD3	1.94	0.41
1:A:438:ILE:HD13	1:A:504:TRP:CD1	2.56	0.40
1:D:371:VAL:HG12	1:D:373:THR:HG23	2.02	0.40
1:A:454:TYR:CZ	1:A:510:ASP:HA	2.56	0.40
2:B:130:ARG:HG2	2:B:130:ARG:NH1	2.26	0.40
1:D:508:TYR:O	1:D:508:TYR:HD2	2.04	0.40
1:D:382:CYS:HA	1:D:383:PRO:HD3	1.70	0.40
1:D:203:ILE:HA	1:D:205:TYR:CE1	2.56	0.40
2:B:84:VAL:CG1	2:B:88:GLN:HB2	2.50	0.40
2:E:134:ASP:OD1	2:E:166:ARG:HG2	2.22	0.40
1:A:221:LEU:HB3	1:A:232:SER:HB3	2.03	0.40
1:D:269:ASN:ND2	1:D:323:PRO:HG3	2.36	0.40
2:B:155:VAL:HG12	2:B:161:MET:HB2	2.02	0.40
2:B:134:ASP:HB3	2:B:164:LEU:HD11	2.03	0.40
1:A:480:ARG:HA	6:A:1313:NAG:H82	2.04	0.40

There are no symmetry-related clashes.



## 5.3 Torsion angles

### 5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	426/428 (100%)	382 (90%)	40 (9%)	4 (1%)	25	73
1	D	426/428 (100%)	372 (87%)	50 (12%)	4 (1%)	25	73
2	B	139/141 (99%)	120 (86%)	13 (9%)	6 (4%)	4	23
2	E	139/141 (99%)	118 (85%)	17 (12%)	4 (3%)	7	35
All	All	1130/1138 (99%)	992 (88%)	120 (11%)	18 (2%)	14	56

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	64	ARG
2	E	64	ARG
1	A	338	ARG
2	B	45	GLU
2	B	153	GLY
2	B	158	THR
1	D	205	TYR
1	D	325	SER
2	E	153	GLY
1	A	258	ARG
2	B	66	ARG
1	A	474	PRO
1	D	385	THR
1	D	389	TYR
2	B	85	GLU
2	E	168	GLY
1	A	328	GLY
2	E	74	PRO

### 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	382/382 (100%)	344 (90%)	38 (10%)	11	40
1	D	382/382 (100%)	335 (88%)	47 (12%)	7	28
2	B	121/121 (100%)	111 (92%)	10 (8%)	16	52
2	E	121/121 (100%)	109 (90%)	12 (10%)	11	40
All	All	1006/1006 (100%)	899 (89%)	107 (11%)	10	36

All (107) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	182	VAL
1	A	194	SER
1	A	201	LYS
1	A	203	ILE
1	A	204	SER
1	A	205	TYR
1	A	209	VAL
1	A	215	THR
1	A	233	HIS
1	A	241	SER
1	A	264	SER
1	A	265	LEU
1	A	272	THR
1	A	285	VAL
1	A	291	TYR
1	A	307	SER
1	A	326	ASN
1	A	337	LEU
1	A	348	VAL
1	A	351	TYR
1	A	357	LYS
1	A	376	LYS
1	A	385	THR
1	A	393	GLU
1	A	411	SER
1	A	422	GLU
1	A	426	VAL
1	A	466	PHE
1	A	470	LEU
1	A	473	ASN
1	A	476	VAL

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Mol	Chain	Res	Type
1	A	487	ARG
1	A	519	TRP
1	A	538	THR
1	A	541	LYS
1	A	582	ASP
1	A	592	LEU
1	A	600	GLN
2	B	41	ARG
2	B	48	TYR
2	B	58	LEU
2	B	106	ARG
2	B	109	LEU
2	B	112	ARG
2	B	119	GLU
2	B	121	SER
2	B	130	ARG
2	B	133	HIS
1	D	201	LYS
1	D	205	TYR
1	D	209	VAL
1	D	233	HIS
1	D	237	ILE
1	D	241	SER
1	D	264	SER
1	D	270	VAL
1	D	279	VAL
1	D	285	VAL
1	D	291	TYR
1	D	308	THR
1	D	311	SER
1	D	324	LYS
1	D	325	SER
1	D	326	ASN
1	D	334	GLN
1	D	335	LEU
1	D	337	LEU
1	D	351	TYR
1	D	373	THR
1	D	376	LYS
1	D	385	THR
1	D	393	GLU
1	D	411	SER

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Mol	Chain	Res	Type
1	D	422	GLU
1	D	431	ILE
1	D	434	GLN
1	D	437	SER
1	D	440	SER
1	D	466	PHE
1	D	470	LEU
1	D	487	ARG
1	D	490	GLN
1	D	492	GLN
1	D	507	VAL
1	D	531	THR
1	D	538	THR
1	D	543	ASN
1	D	547	TYR
1	D	553	SER
1	D	570	ASN
1	D	582	ASP
1	D	583	THR
1	D	585	ASP
1	D	592	LEU
1	D	602	THR
2	E	41	ARG
2	E	48	TYR
2	E	58	LEU
2	E	64	ARG
2	E	66	ARG
2	E	103	THR
2	E	112	ARG
2	E	115	ILE
2	E	119	GLU
2	E	121	SER
2	E	123	ASN
2	E	128	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (25) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	233	HIS
1	A	326	ASN
1	A	334	GLN
1	A	434	GLN

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Mol	Chain	Res	Type
1	A	473	ASN
1	A	492	GLN
1	A	497	ASN
2	B	118	GLN
2	B	127	HIS
1	D	233	HIS
1	D	269	ASN
1	D	326	ASN
1	D	332	GLN
1	D	334	GLN
1	D	423	ASN
1	D	434	GLN
1	D	450	GLN
1	D	473	ASN
1	D	492	GLN
1	D	497	ASN
1	D	586	ASN
1	D	600	GLN
2	E	118	GLN
2	E	123	ASN
2	E	132	HIS

### 5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

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

### 5.5 Carbohydrates ⓘ

28 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
3	NAG	A	1138	1,3	12,14,15	1.90	4 (33%)	15,19,21	1.78	4 (26%)
3	NAG	A	1139	3	12,14,15	2.10	4 (33%)	15,19,21	1.50	3 (20%)
3	BMA	A	1140	3	10,11,12	1.64	3 (30%)	11,15,17	1.49	2 (18%)
3	BMA	A	1141	3	10,11,12	1.02	0	11,15,17	2.45	4 (36%)
3	MAN	A	1142	3	10,11,12	0.96	0	11,15,17	1.09	1 (9%)
6	NAG	A	1313	1,6	12,14,15	1.84	5 (41%)	15,19,21	1.77	4 (26%)
6	NAG	A	1314	6	12,14,15	1.54	1 (8%)	15,19,21	1.73	4 (26%)
4	LXZ	A	1361	1,4	12,14,15	2.43	5 (41%)	15,19,21	1.90	3 (20%)
4	NGA	A	1362	4	12,14,15	1.77	4 (33%)	15,19,21	1.88	4 (26%)
4	GL0	A	1363	4	10,11,12	1.40	2 (20%)	11,15,17	2.81	4 (36%)
4	BGC	A	1364	4	10,11,12	0.94	0	11,15,17	1.54	2 (18%)
4	GLC	A	1365	4	10,11,12	1.13	0	11,15,17	1.29	2 (18%)
4	BGC	A	1366	4	10,11,12	0.93	0	11,15,17	2.55	2 (18%)
4	GLC	A	1367	4	10,11,12	0.85	0	11,15,17	1.86	2 (18%)
8	NAG	D	1238	1,8	12,14,15	2.00	5 (41%)	15,19,21	1.26	2 (13%)
8	NAG	D	1239	8	12,14,15	2.03	5 (41%)	15,19,21	1.79	4 (26%)
8	BMA	D	1240	8	10,11,12	1.76	3 (30%)	11,15,17	1.54	4 (36%)
8	BGC	D	1241	8	10,11,12	1.14	1 (10%)	11,15,17	2.67	4 (36%)
8	MAN	D	1242	8	10,11,12	1.02	0	11,15,17	1.15	1 (9%)
6	NAG	D	1413	1,6	12,14,15	1.79	4 (33%)	15,19,21	1.78	3 (20%)
6	NAG	D	1414	6	12,14,15	1.61	2 (16%)	15,19,21	1.38	3 (20%)
9	LXB	D	1461	9,1	12,14,15	2.41	4 (33%)	15,19,21	2.03	4 (26%)
9	NGZ	D	1462	9	12,14,15	1.78	4 (33%)	15,19,21	1.62	2 (13%)
9	GL0	D	1463	9	10,11,12	1.34	2 (20%)	11,15,17	2.84	4 (36%)
9	BGC	D	1464	9	10,11,12	0.92	0	11,15,17	1.22	1 (9%)
9	GLC	D	1465	9	10,11,12	1.09	0	11,15,17	1.35	2 (18%)
9	BGC	D	1466	9	10,11,12	0.92	0	11,15,17	1.65	1 (9%)
9	GXL	D	1467	9	10,11,12	0.81	0	11,15,17	1.48	1 (9%)

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
3	NAG	A	1138	1,3	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	A	1139	3	-	0/6/23/26	0/1/1/1
3	BMA	A	1140	3	-	0/2/19/22	0/1/1/1
3	BMA	A	1141	3	-	0/2/19/22	0/1/1/1
3	MAN	A	1142	3	-	0/2/19/22	0/1/1/1
6	NAG	A	1313	1,6	1/1/5/7	0/6/23/26	0/1/1/1
6	NAG	A	1314	6	1/1/5/7	0/6/23/26	0/1/1/1
4	LXZ	A	1361	1,4	-	0/6/23/26	0/1/1/1
4	NGA	A	1362	4	5/5/5/7	0/6/23/26	0/1/1/1
4	GL0	A	1363	4	1/1/4/5	0/2/19/22	0/1/1/1
4	BGC	A	1364	4	-	0/2/19/22	0/1/1/1
4	GLC	A	1365	4	5/5/4/5	0/2/19/22	0/1/1/1
4	BGC	A	1366	4	1/1/4/5	0/2/19/22	0/1/1/1
4	GLC	A	1367	4	5/5/4/5	0/2/19/22	0/1/1/1
8	NAG	D	1238	1,8	-	0/6/23/26	0/1/1/1
8	NAG	D	1239	8	-	0/6/23/26	0/1/1/1
8	BMA	D	1240	8	-	0/2/19/22	0/1/1/1
8	BGC	D	1241	8	-	0/2/19/22	0/1/1/1
8	MAN	D	1242	8	-	0/2/19/22	0/1/1/1
6	NAG	D	1413	1,6	2/2/5/7	0/6/23/26	0/1/1/1
6	NAG	D	1414	6	1/1/5/7	0/6/23/26	0/1/1/1
9	LXB	D	1461	9,1	-	0/6/23/26	0/1/1/1
9	NGZ	D	1462	9	1/1/5/7	0/6/23/26	0/1/1/1
9	GL0	D	1463	9	1/1/4/5	0/2/19/22	0/1/1/1
9	BGC	D	1464	9	-	0/2/19/22	0/1/1/1
9	GLC	D	1465	9	5/5/4/5	0/2/19/22	0/1/1/1
9	BGC	D	1466	9	1/1/4/5	0/2/19/22	0/1/1/1
9	GXL	D	1467	9	1/1/4/5	0/2/19/22	0/1/1/1

All (58) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	D	1461	LXB	C4-C5	4.55	1.63	1.53
4	A	1361	LXZ	C4-C5	4.48	1.62	1.53
9	D	1461	LXB	C7-N2	3.84	1.50	1.34
3	A	1139	NAG	C7-N2	3.82	1.49	1.34
4	A	1361	LXZ	C7-N2	3.82	1.49	1.34
4	A	1361	LXZ	C4-C3	3.77	1.62	1.52
3	A	1139	NAG	C2-N2	3.71	1.50	1.46
8	D	1239	NAG	C7-N2	3.69	1.49	1.34
6	A	1313	NAG	C7-N2	3.67	1.49	1.34
9	D	1461	LXB	C3-C4	3.64	1.62	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	D	1413	NAG	C7-N2	3.56	1.48	1.34
8	D	1238	NAG	C3-C2	3.49	1.59	1.52
6	D	1414	NAG	C7-N2	3.47	1.48	1.34
4	A	1362	NGA	C7-N2	3.38	1.48	1.34
6	A	1314	NAG	C7-N2	3.27	1.47	1.34
8	D	1239	NAG	C2-N2	3.26	1.50	1.46
8	D	1238	NAG	C7-N2	3.23	1.47	1.34
3	A	1138	NAG	C3-C2	3.23	1.59	1.52
9	D	1462	NGZ	C7-N2	3.22	1.47	1.34
3	A	1138	NAG	C7-N2	3.11	1.47	1.34
8	D	1240	BMA	O5-C5	2.87	1.50	1.45
8	D	1240	BMA	C4-C5	2.79	1.59	1.53
8	D	1238	NAG	C4-C3	2.76	1.59	1.52
3	A	1139	NAG	C3-C2	2.74	1.58	1.52
3	A	1138	NAG	C4-C3	2.66	1.59	1.52
9	D	1462	NGZ	C3-C2	2.64	1.58	1.52
3	A	1139	NAG	C8-C7	2.63	1.56	1.50
3	A	1140	BMA	C4-C5	2.61	1.58	1.53
4	A	1361	LXZ	C2-N2	2.60	1.49	1.46
8	D	1239	NAG	C8-C7	2.59	1.56	1.50
6	D	1413	NAG	C2-N2	2.56	1.49	1.46
3	A	1140	BMA	O5-C5	2.50	1.50	1.45
8	D	1238	NAG	C4-C5	2.48	1.58	1.53
4	A	1363	GL0	O5-C5	2.41	1.49	1.45
4	A	1362	NGA	C3-C2	2.37	1.57	1.52
6	D	1413	NAG	C3-C2	2.36	1.57	1.52
9	D	1461	LXB	C2-N2	2.29	1.49	1.46
6	A	1313	NAG	C2-N2	2.29	1.49	1.46
6	A	1313	NAG	C8-C7	2.27	1.55	1.50
3	A	1138	NAG	C4-C5	2.21	1.57	1.53
8	D	1239	NAG	C3-C2	2.21	1.57	1.52
4	A	1361	LXZ	C8-C7	2.20	1.55	1.50
9	D	1463	GL0	O5-C5	2.18	1.49	1.45
4	A	1363	GL0	C3-C2	2.16	1.57	1.52
8	D	1240	BMA	C4-C3	2.14	1.58	1.52
9	D	1462	NGZ	O4-C4	2.12	1.48	1.43
9	D	1462	NGZ	C8-C7	2.11	1.55	1.50
6	D	1414	NAG	C2-N2	2.09	1.48	1.46
8	D	1241	BGC	C3-C2	2.09	1.57	1.52
9	D	1463	GL0	C3-C2	2.09	1.57	1.52
8	D	1238	NAG	O4-C4	2.09	1.48	1.43
3	A	1140	BMA	C4-C3	2.06	1.57	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	1313	NAG	C4-C3	2.05	1.57	1.52
4	A	1362	NGA	C8-C7	2.04	1.54	1.50
6	D	1413	NAG	C8-C7	2.04	1.54	1.50
4	A	1362	NGA	C2-N2	2.03	1.48	1.46
8	D	1239	NAG	C4-C5	2.02	1.57	1.53
6	A	1313	NAG	C3-C2	2.01	1.56	1.52

All (77) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1366	BGC	O5-C5-C6	7.48	114.83	106.98
4	A	1363	GL0	O5-C5-C6	6.48	113.78	106.98
9	D	1463	GL0	O5-C5-C6	6.19	113.48	106.98
4	A	1361	LXZ	O5-C5-C4	5.70	117.88	110.65
9	D	1463	GL0	C4-C3-C2	5.40	117.75	110.50
8	D	1241	BGC	O5-C5-C4	5.02	117.03	110.65
3	A	1141	BMA	O5-C5-C4	4.96	116.95	110.65
9	D	1466	BGC	O5-C5-C6	4.75	111.96	106.98
8	D	1241	BGC	O5-C5-C6	4.60	111.80	106.98
8	D	1239	NAG	O5-C5-C4	4.56	116.44	110.65
9	D	1461	LXB	O5-C5-C6	4.52	111.72	106.98
3	A	1141	BMA	C3-C4-C5	4.42	118.09	110.20
4	A	1367	GLC	O5-C5-C4	4.34	116.17	110.65
9	D	1462	NGZ	O5-C5-C4	4.34	116.16	110.65
6	D	1413	NAG	C3-C2-N2	-4.19	105.37	111.76
4	A	1363	GL0	C4-C3-C2	4.14	116.06	110.50
9	D	1461	LXB	O5-C5-C4	3.96	115.68	110.65
4	A	1362	NGA	O5-C5-C4	3.88	115.58	110.65
4	A	1362	NGA	C3-C4-C5	3.79	116.97	110.20
8	D	1241	BGC	C3-C4-C5	3.75	116.90	110.20
8	D	1241	BGC	C4-C3-C2	3.72	115.50	110.50
4	A	1363	GL0	O5-C5-C4	3.60	115.22	110.65
4	A	1363	GL0	C3-C4-C5	3.59	116.61	110.20
3	A	1138	NAG	O5-C5-C4	3.46	115.05	110.65
6	D	1413	NAG	O5-C5-C6	3.46	110.61	106.98
9	D	1465	GLC	O5-C5-C6	3.43	110.58	106.98
3	A	1138	NAG	C6-C5-C4	-3.43	104.72	113.00
6	A	1313	NAG	O5-C5-C6	3.43	110.58	106.98
3	A	1138	NAG	C2-N2-C7	-3.42	117.35	123.09
6	A	1314	NAG	O5-C5-C4	3.35	114.90	110.65
3	A	1140	BMA	O5-C5-C4	3.34	114.89	110.65
3	A	1141	BMA	C4-C3-C2	3.29	114.93	110.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	D	1467	GXL	O5-C5-C4	3.21	114.73	110.65
4	A	1364	BGC	O5-C5-C6	3.14	110.28	106.98
9	D	1463	GL0	C3-C4-C5	3.07	115.69	110.20
6	A	1313	NAG	O5-C5-C4	3.06	114.53	110.65
4	A	1361	LXZ	C3-C4-C5	3.04	115.63	110.20
6	D	1414	NAG	C8-C7-N2	2.99	121.95	116.11
9	D	1461	LXB	C3-C4-C5	2.95	115.47	110.20
3	A	1139	NAG	O5-C5-C6	2.92	110.04	106.98
8	D	1242	MAN	O5-C5-C6	2.90	110.02	106.98
4	A	1362	NGA	C3-C2-N2	-2.82	107.47	111.76
6	D	1413	NAG	C2-N2-C7	2.74	127.68	123.09
3	A	1139	NAG	C8-C7-N2	2.73	121.44	116.11
9	D	1463	GL0	O5-C5-C4	2.72	114.10	110.65
8	D	1238	NAG	O5-C5-C4	2.71	114.10	110.65
4	A	1364	BGC	O5-C5-C4	2.70	114.08	110.65
6	A	1314	NAG	O5-C5-C6	2.67	109.78	106.98
8	D	1238	NAG	C2-N2-C7	-2.62	118.69	123.09
3	A	1141	BMA	O5-C5-C6	2.61	109.72	106.98
6	A	1313	NAG	C2-N2-C7	-2.59	118.73	123.09
4	A	1365	GLC	O5-C5-C6	2.57	109.67	106.98
8	D	1240	BMA	O3-C3-C2	2.56	114.62	109.94
4	A	1365	GLC	O5-C5-C4	2.49	113.81	110.65
8	D	1240	BMA	O3-C3-C4	2.47	115.89	110.35
6	A	1313	NAG	C8-C7-N2	2.46	120.92	116.11
6	D	1414	NAG	O5-C5-C4	2.46	113.77	110.65
4	A	1367	GLC	C6-C5-C4	-2.45	107.09	113.00
8	D	1240	BMA	O5-C5-C4	2.44	113.75	110.65
3	A	1142	MAN	O5-C5-C4	2.40	113.70	110.65
3	A	1140	BMA	C3-C4-C5	2.36	114.42	110.20
8	D	1239	NAG	C2-N2-C7	-2.35	119.14	123.09
9	D	1465	GLC	C4-C3-C2	2.28	113.57	110.50
4	A	1362	NGA	C4-C3-C2	2.28	116.89	111.32
8	D	1240	BMA	O5-C5-C6	2.25	109.34	106.98
8	D	1239	NAG	O5-C5-C6	2.25	109.34	106.98
4	A	1366	BGC	O2-C2-C3	-2.24	105.33	110.18
4	A	1361	LXZ	C8-C7-N2	2.21	120.44	116.11
3	A	1139	NAG	C2-N2-C7	-2.20	119.40	123.09
6	A	1314	NAG	C3-C2-N2	-2.19	108.42	111.76
6	D	1414	NAG	O5-C5-C6	2.16	109.25	106.98
9	D	1461	LXB	C4-C3-C2	2.11	116.47	111.32
9	D	1464	BGC	O5-C5-C6	2.08	109.17	106.98
3	A	1138	NAG	C3-C4-C5	2.08	113.92	110.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	1314	NAG	C8-C7-N2	2.08	120.17	116.11
8	D	1239	NAG	C8-C7-N2	2.05	120.11	116.11
9	D	1462	NGZ	O4-C4-C5	-2.05	103.88	109.28

All (31) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
9	D	1465	GLC	C2
9	D	1465	GLC	C5
9	D	1465	GLC	C3
9	D	1465	GLC	C1
9	D	1465	GLC	C4
9	D	1466	BGC	C5
9	D	1467	GXL	C4
4	A	1362	NGA	C2
4	A	1362	NGA	C5
4	A	1362	NGA	C3
4	A	1362	NGA	C1
4	A	1362	NGA	C4
6	D	1414	NAG	C1
4	A	1363	GL0	C1
9	D	1462	NGZ	C1
4	A	1365	GLC	C2
4	A	1365	GLC	C5
4	A	1365	GLC	C3
4	A	1365	GLC	C1
4	A	1365	GLC	C4
6	A	1313	NAG	C1
4	A	1366	BGC	C5
6	A	1314	NAG	C1
4	A	1367	GLC	C2
4	A	1367	GLC	C5
4	A	1367	GLC	C3
4	A	1367	GLC	C1
4	A	1367	GLC	C4
9	D	1463	GL0	C1
6	D	1413	NAG	C2
6	D	1413	NAG	C1

There are no torsion outliers.

There are no ring outliers.

## 5.6 Ligand geometry ⓘ

24 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	$\# Z  > 2$
5	NAG	A	1210	1	12,14,15	1.46	2 (16%)	15,19,21	1.41	3 (20%)
7	SO4	A	1368	-	4,4,4	0.30	0	6,6,6	0.30	0
7	SO4	A	1369	-	4,4,4	0.29	0	6,6,6	0.26	0
7	SO4	A	1370	-	4,4,4	0.10	0	6,6,6	0.17	0
7	SO4	A	1371	-	4,4,4	0.04	0	6,6,6	0.10	0
7	SO4	A	1372	-	4,4,4	0.21	0	6,6,6	0.09	0
7	SO4	A	1373	-	4,4,4	0.15	0	6,6,6	0.18	0
7	SO4	A	1374	-	4,4,4	0.12	0	6,6,6	0.19	0
7	SO4	A	1375	-	4,4,4	0.28	0	6,6,6	0.38	0
7	SO4	A	1376	-	4,4,4	0.11	0	6,6,6	0.19	0
7	SO4	B	202	-	4,4,4	0.09	0	6,6,6	0.11	0
7	SO4	B	203	-	4,4,4	0.06	0	6,6,6	0.13	0
5	NAG	D	1310	1	12,14,15	1.47	2 (16%)	15,19,21	1.26	2 (13%)
7	SO4	D	1468	-	4,4,4	0.15	0	6,6,6	0.18	0
7	SO4	D	1469	-	4,4,4	0.10	0	6,6,6	0.07	0
7	SO4	D	1470	-	4,4,4	0.08	0	6,6,6	0.09	0
7	SO4	D	1471	-	4,4,4	0.15	0	6,6,6	0.11	0
7	SO4	D	1472	-	4,4,4	0.09	0	6,6,6	0.13	0
7	SO4	D	1473	-	4,4,4	0.07	0	6,6,6	0.11	0
7	SO4	D	1474	-	4,4,4	0.24	0	6,6,6	0.38	0
7	SO4	D	1475	-	4,4,4	0.13	0	6,6,6	0.09	0
7	SO4	D	1476	-	4,4,4	0.09	0	6,6,6	0.20	0
7	SO4	D	1477	-	4,4,4	0.03	0	6,6,6	0.16	0
7	SO4	E	190	-	4,4,4	0.08	0	6,6,6	0.22	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	A	1210	1	-	0/6/23/26	0/1/1/1
7	SO4	A	1368	-	-	0/0/0/0	0/0/0/0
7	SO4	A	1369	-	-	0/0/0/0	0/0/0/0
7	SO4	A	1370	-	-	0/0/0/0	0/0/0/0
7	SO4	A	1371	-	-	0/0/0/0	0/0/0/0
7	SO4	A	1372	-	-	0/0/0/0	0/0/0/0
7	SO4	A	1373	-	-	0/0/0/0	0/0/0/0
7	SO4	A	1374	-	-	0/0/0/0	0/0/0/0
7	SO4	A	1375	-	-	0/0/0/0	0/0/0/0
7	SO4	A	1376	-	-	0/0/0/0	0/0/0/0
7	SO4	B	202	-	-	0/0/0/0	0/0/0/0
7	SO4	B	203	-	-	0/0/0/0	0/0/0/0
5	NAG	D	1310	1	1/1/5/7	0/6/23/26	0/1/1/1
7	SO4	D	1468	-	-	0/0/0/0	0/0/0/0
7	SO4	D	1469	-	-	0/0/0/0	0/0/0/0
7	SO4	D	1470	-	-	0/0/0/0	0/0/0/0
7	SO4	D	1471	-	-	0/0/0/0	0/0/0/0
7	SO4	D	1472	-	-	0/0/0/0	0/0/0/0
7	SO4	D	1473	-	-	0/0/0/0	0/0/0/0
7	SO4	D	1474	-	-	0/0/0/0	0/0/0/0
7	SO4	D	1475	-	-	0/0/0/0	0/0/0/0
7	SO4	D	1476	-	-	0/0/0/0	0/0/0/0
7	SO4	D	1477	-	-	0/0/0/0	0/0/0/0
7	SO4	E	190	-	-	0/0/0/0	0/0/0/0

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	1210	NAG	C7-N2	3.27	1.47	1.34
5	D	1310	NAG	C7-N2	3.14	1.47	1.34
5	D	1310	NAG	C3-C2	2.07	1.56	1.52
5	A	1210	NAG	C8-C7	2.03	1.54	1.50

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1210	NAG	C8-C7-N2	2.95	121.87	116.11
5	D	1310	NAG	O5-C5-C6	2.48	109.58	106.98
5	A	1210	NAG	O5-C5-C4	2.36	113.65	110.65
5	A	1210	NAG	C6-C5-C4	-2.17	107.76	113.00
5	D	1310	NAG	O5-C5-C4	2.09	113.30	110.65

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
5	D	1310	NAG	C1

There are no torsion outliers.

There are no ring outliers.

## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

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	428/428 (100%)	-0.21	0 100 100	43, 59, 85, 110	0
1	D	428/428 (100%)	-0.23	0 100 100	45, 61, 86, 104	0
2	B	141/141 (100%)	-0.18	0 100 100	45, 72, 94, 106	0
2	E	141/141 (100%)	-0.23	0 100 100	48, 70, 97, 99	0
All	All	1138/1138 (100%)	-0.21	0 93 100	43, 63, 90, 110	0

There are no RSRZ outliers to report.

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

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

### 6.3 Carbohydrates ⓘ

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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
8	BGC	D	1241	11/12	0.55	7.94	75,98,105,106	0
6	NAG	A	1313	14/15	0.23	7.07	62,79,89,97	0
6	NAG	D	1414	14/15	0.33	3.61	89,102,109,110	0
6	NAG	A	1314	14/15	0.35	2.95	85,95,100,100	0
4	GLC	A	1367	11/12	0.20	2.37	69,75,82,90	0
9	GXL	D	1467	11/12	0.19	0.80	69,79,90,91	0
8	NAG	D	1238	14/15	0.27	0.79	85,90,100,104	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
6	NAG	D	1413	14/15	0.17	0.16	72,83,92,104	0
4	LXZ	A	1361	14/15	0.15	-1.12	65,70,80,81	0
3	NAG	A	1138	14/15	0.14	-1.44	55,59,65,67	0
9	LXB	D	1461	14/15	0.13	-1.55	73,76,80,85	0
3	NAG	A	1139	14/15	0.13	-1.80	60,67,72,73	0
3	BMA	A	1140	11/12	0.15	-2.63	75,81,86,89	0
4	GLC	A	1365	11/12	0.38	-	87,97,101,101	0
9	GLC	D	1465	11/12	0.36	-	98,107,112,113	0
8	MAN	D	1242	11/12	0.71	-	108,118,121,122	0
3	MAN	A	1142	11/12	0.33	-	87,91,96,96	0
8	NAG	D	1239	14/15	0.30	-	90,103,113,116	0
4	BGC	A	1366	11/12	0.20	-	75,81,87,90	0
4	NGA	A	1362	14/15	0.17	-	71,80,84,91	0
9	BGC	D	1466	11/12	0.18	-	79,82,87,89	0
3	BMA	A	1141	11/12	0.25	-	86,90,93,93	0
4	GL0	A	1363	11/12	0.25	-	93,97,99,101	0
4	BGC	A	1364	11/12	0.38	-	93,103,106,106	0
9	GL0	D	1463	11/12	0.33	-	110,116,119,121	0
8	BMA	D	1240	11/12	0.45	-	101,110,116,117	0
9	BGC	D	1464	11/12	0.38	-	113,120,124,125	0
9	NGZ	D	1462	14/15	0.21	-	70,92,99,106	0

## 6.4 Ligands ⓘ

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	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
5	NAG	D	1310	14/15	0.34	48.33	95,105,108,108	0
7	SO4	A	1373	5/5	0.28	14.38	93,99,118,124	0
7	SO4	D	1472	5/5	0.36	6.11	98,103,119,121	0
7	SO4	D	1473	5/5	0.44	5.65	105,111,123,131	0
7	SO4	D	1474	5/5	0.39	5.28	90,95,109,114	0
7	SO4	A	1370	5/5	0.26	4.20	116,121,126,139	0
5	NAG	A	1210	14/15	0.30	3.31	88,92,97,97	0
7	SO4	B	203	5/5	0.34	2.92	108,108,125,130	0
7	SO4	A	1375	5/5	0.38	2.90	82,83,103,104	0
7	SO4	D	1471	5/5	0.55	2.73	84,97,109,120	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
7	SO4	B	202	5/5	0.34	2.46	102,106,112,126	0
7	SO4	A	1374	5/5	0.34	2.44	92,99,110,121	0
7	SO4	D	1477	5/5	0.33	2.06	105,107,117,130	0
7	SO4	A	1372	5/5	0.33	1.82	81,92,102,116	0
7	SO4	D	1469	5/5	0.30	1.73	124,128,134,144	0
7	SO4	E	190	5/5	0.36	1.32	105,109,123,130	0
7	SO4	A	1368	5/5	0.20	0.81	58,60,64,67	0
7	SO4	A	1376	5/5	0.21	0.11	86,88,100,108	0
7	SO4	A	1369	5/5	0.15	0.09	57,66,73,79	0
7	SO4	D	1475	5/5	0.22	-0.14	88,88,100,108	0
7	SO4	D	1476	5/5	0.12	-1.70	86,89,99,105	0
7	SO4	D	1468	5/5	0.12	-1.76	81,86,90,97	0
7	SO4	D	1470	5/5	0.30	-	109,117,128,134	0
7	SO4	A	1371	5/5	0.31	-	108,110,117,129	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.