



Full wwPDB X-ray Structure Validation Report

Feb 26, 2014 – 08:24 PM GMT

PDB ID : 3PPS
Title : Crystal structure of an ascomycete fungal laccase from Thielavia arenaria
Authors : Kallio, J.P.; Rouvinen, J.; Hakulinen, N.
Deposited on : 2010-11-25
Resolution : 2.50 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

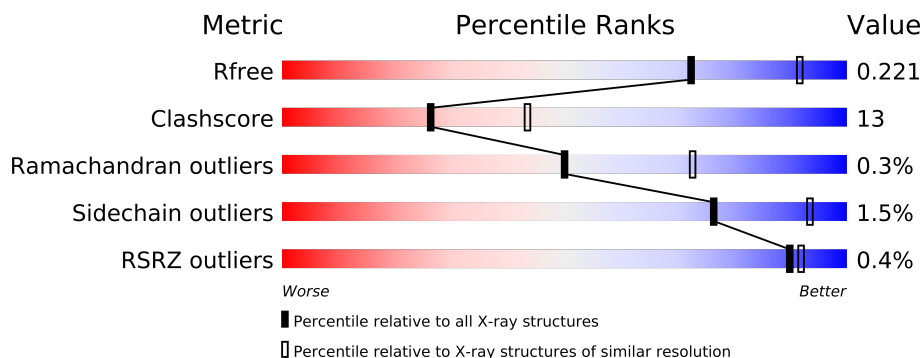
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.50 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	2784 (2.50-2.50)
Clashscore	79885	3562 (2.50-2.50)
Ramachandran outliers	78287	3480 (2.50-2.50)
Sidechain outliers	78261	3482 (2.50-2.50)
RSRZ outliers	66119	2785 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 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	604	
1	B	604	
1	C	604	
1	D	604	

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

Mol	Type	Chain	Res	Geometry	Electron density
6	NAG	A	730	-	X
6	NAG	C	730	-	X

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 18812 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 Laccase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	564	Total	C	N	O	S	0	0	0
			4406	2793	776	821	16			
1	B	564	Total	C	N	O	S	0	0	0
			4406	2793	776	821	16			
1	C	564	Total	C	N	O	S	0	0	0
			4406	2793	776	821	16			
1	D	564	Total	C	N	O	S	0	0	0
			4406	2793	776	821	16			

- Molecule 2 is COPPER (II) ION (three-letter code: CU) (formula: Cu).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	4	Total	Cu	0	0
			4	4		
2	A	4	Total	Cu	0	0
			4	4		
2	D	4	Total	Cu	0	0
			4	4		
2	C	4	Total	Cu	0	0
			4	4		

- Molecule 3 is OXYGEN MOLECULE (three-letter code: OXY) (formula: O₂).

○ ○

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O 2 2	0	0

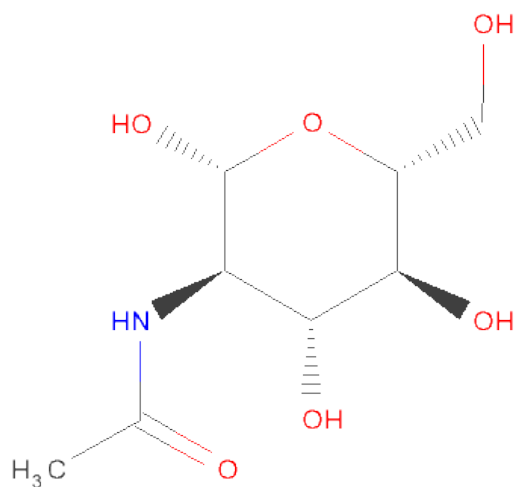
- Molecule 4 is a polymer of unknown type called SUGAR (2-MER).

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

- Molecule 5 is a polymer of unknown type called SUGAR (6-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	6	Total	C	N	O	0	0
			72	40	2	30		

- Molecule 6 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



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

- Molecule 7 is a polymer of unknown type called SUGAR (5-MER).

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

- Molecule 8 is a polymer of unknown type called SUGAR (3-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	C	3	Total	C	N	O	0	0
			39	22	2	15		
8	D	3	Total	C	N	O	0	0
			39	22	2	15		

- Molecule 9 is water.

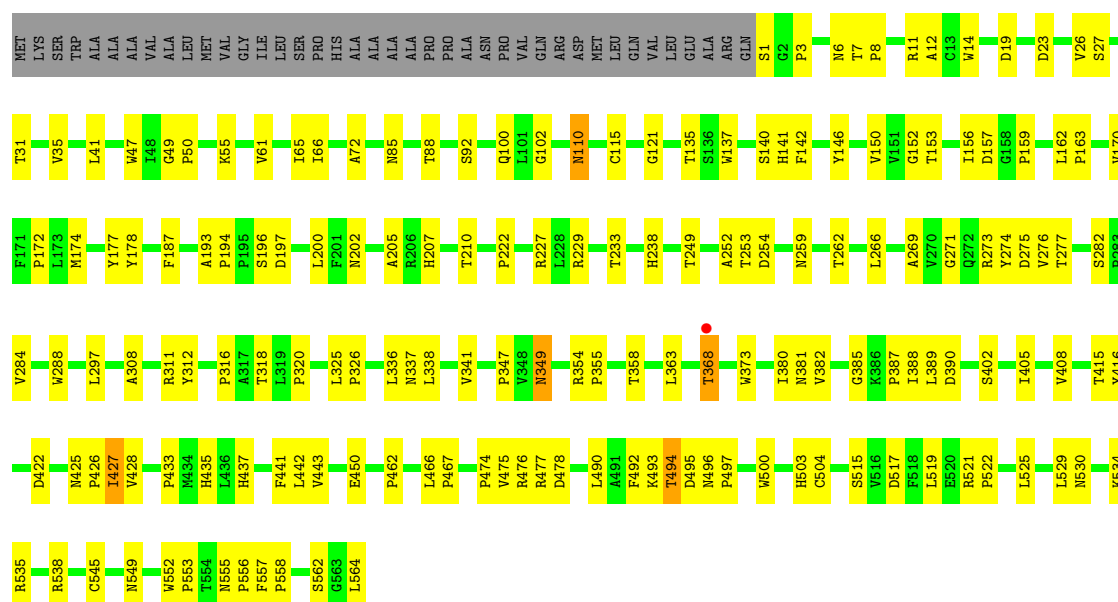
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	159	Total	O	0	0
			159	159		
9	B	138	Total	O	0	0
			138	138		
9	C	124	Total	O	0	0
			124	124		
9	D	141	Total	O	0	0
			141	141		

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.

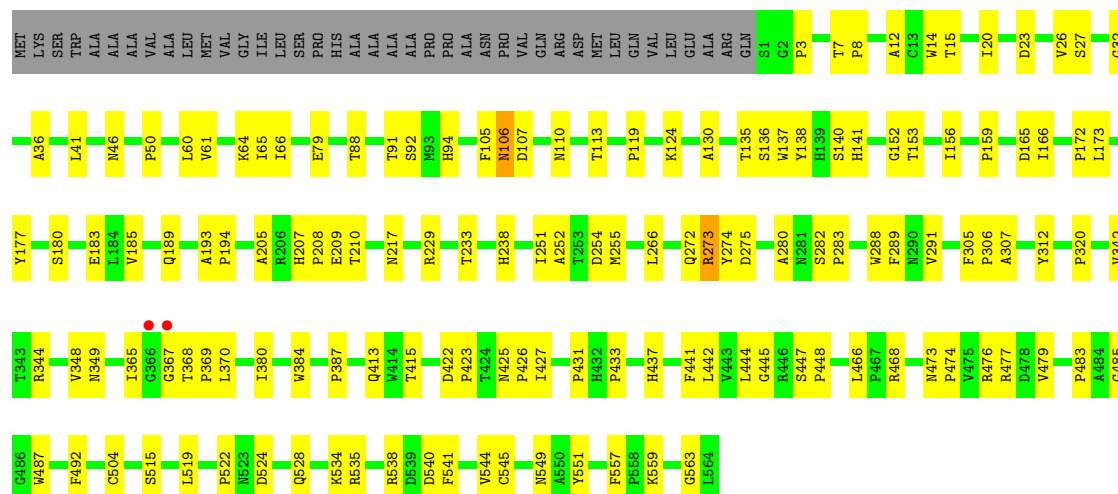
• Molecule 1: Laccase

Chain A:



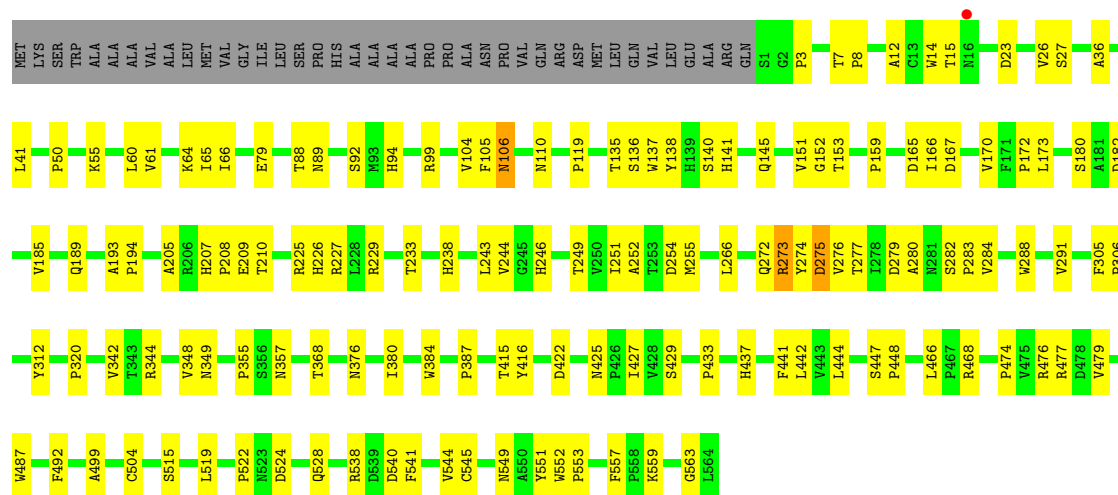
• Molecule 1: Laccase

Chain B:



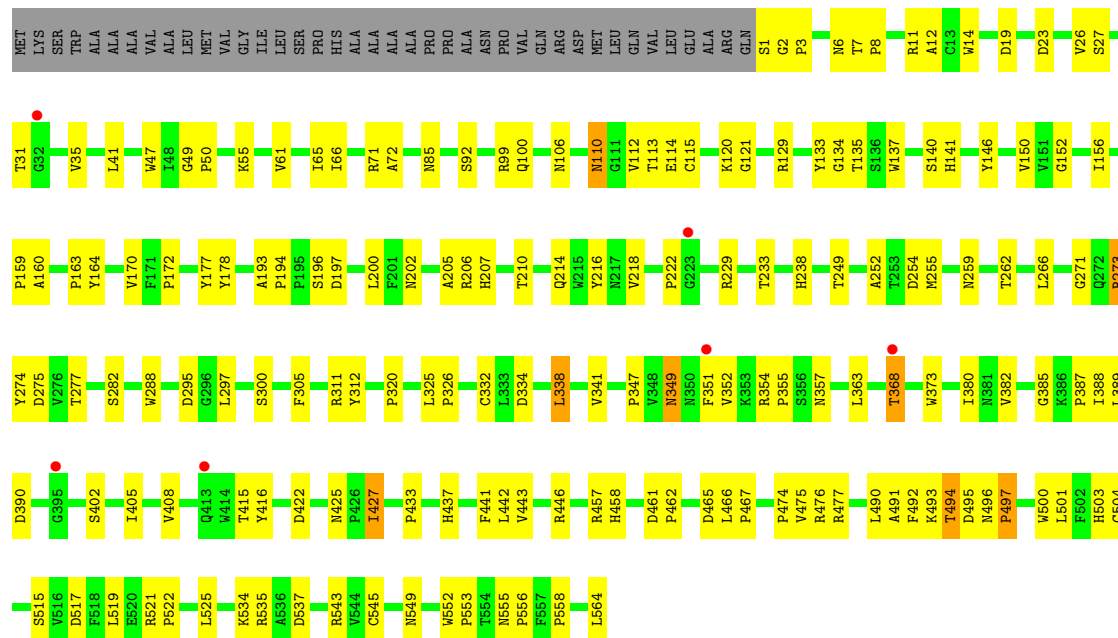
- Molecule 1: Laccase

Chain C:



- Molecule 1: Laccase

Chain D:



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	61.36Å 178.95Å 118.13Å 90.00° 90.26° 90.00°	Depositor
Resolution (Å)	42.65 – 2.50 42.65 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.9 (42.65-2.50) 98.6 (42.65-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.63 (at 2.51Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.5_2)	Depositor
R, R_{free}	0.182 , 0.222 0.179 , 0.221	Depositor DCC
R_{free} test set	4405 reflections (5.02%)	DCC
Wilson B-factor (Å ²)	12.6	Xtriage
Anisotropy	1.837	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 8.0	EDS
Estimated twinning fraction	0.319 for h,-k,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.39$, $\langle L^2 \rangle = 0.21$	Xtriage
Outliers	0 of 87766 reflections	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	18812	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 37.22 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 4.4366e-04.*

¹Intensities 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: MAN, BMA, NAG, CU, OXY

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	0.36	0/4539	0.53	0/6219
1	B	0.36	0/4539	0.53	0/6219
1	C	0.35	0/4539	0.53	0/6219
1	D	0.35	0/4539	0.53	0/6219
All	All	0.35	0/18156	0.53	0/24876

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	4406	0	4223	129	0
1	B	4406	0	4223	105	0
1	C	4406	0	4224	103	0
1	D	4406	0	4225	136	0
2	A	4	0	0	0	0
2	B	4	0	0	0	0
2	C	4	0	0	0	0
2	D	4	0	0	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	2	0	0	0	0
4	A	56	0	50	4	0
4	B	56	0	50	2	0
4	C	84	0	75	9	0
4	D	28	0	25	0	0
5	A	72	0	61	1	0
6	A	42	0	39	1	0
6	B	42	0	39	0	0
6	C	14	0	13	0	0
6	D	14	0	13	1	0
7	B	61	0	52	2	0
7	D	61	0	52	4	0
8	C	39	0	34	3	0
8	D	39	0	34	1	0
9	A	159	0	0	12	0
9	B	138	0	0	15	0
9	C	124	0	0	14	0
9	D	141	0	0	28	0
All	All	18812	0	17432	473	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 13.

All (473) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:376:ASN:HD22	4:C:760:NAG:H83	1.35	0.92
1:A:253:THR:HG21	1:A:338:LEU:HD12	1.59	0.83
1:D:207:HIS:HD2	1:D:210:THR:H	1.27	0.82
1:A:207:HIS:HD2	1:A:210:THR:H	1.28	0.81
1:C:104:VAL:HG11	9:C:694:HOH:O	1.79	0.81
1:D:442:LEU:HB2	9:D:663:HOH:O	1.80	0.79
1:D:552:TRP:HB3	1:D:553:PRO:HD3	1.65	0.79
1:A:14:TRP:HB2	1:A:159:PRO:HG3	1.65	0.78
1:D:252:ALA:HB3	1:D:275:ASP:HB2	1.66	0.77
1:D:14:TRP:HB2	1:D:159:PRO:HG3	1.66	0.77
1:B:422:ASP:OD1	1:B:425:ASN:HB2	1.86	0.76
1:C:422:ASP:OD1	1:C:425:ASN:HB2	1.86	0.75
1:A:552:TRP:HB3	1:A:553:PRO:HD3	1.67	0.75
1:D:114:GLU:HB3	9:D:582:HOH:O	1.86	0.74
1:C:167:ASP:OD1	9:C:654:HOH:O	2.05	0.73
1:C:429:SER:HB2	9:C:663:HOH:O	1.89	0.73
1:D:352:VAL:HB	9:D:665:HOH:O	1.87	0.72

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:50:PRO:HB2	1:A:205:ALA:HB3	1.72	0.71
1:A:349:ASN:ND2	9:A:635:HOH:O	2.22	0.71
1:C:14:TRP:HB2	1:C:159:PRO:HG3	1.72	0.70
1:D:238:HIS:NE2	1:D:433:PRO:HG3	2.07	0.70
1:D:50:PRO:HB2	1:D:205:ALA:HB3	1.73	0.70
1:C:244:VAL:O	9:C:651:HOH:O	2.08	0.70
1:A:503:HIS:HB3	1:A:515:SER:OG	1.91	0.69
1:B:14:TRP:HB2	1:B:159:PRO:HG3	1.72	0.69
1:A:102:GLY:HA2	9:A:705:HOH:O	1.91	0.69
1:B:487:TRP:O	9:B:631:HOH:O	2.09	0.69
1:A:238:HIS:NE2	1:A:433:PRO:HG3	2.08	0.69
1:C:540:ASP:O	1:C:544:VAL:HG23	1.93	0.68
1:D:462:PRO:HA	1:D:466:LEU:HD13	1.74	0.68
1:B:540:ASP:O	1:B:544:VAL:HG23	1.94	0.67
1:A:426:PRO:HG2	1:B:370:LEU:HD21	1.77	0.67
1:B:14:TRP:CD1	1:B:15:THR:HG23	2.30	0.67
1:A:252:ALA:HB3	1:A:275:ASP:HB2	1.75	0.67
1:D:503:HIS:HB3	1:D:515:SER:OG	1.95	0.66
1:C:433:PRO:HD2	1:C:504:CYS:SG	2.35	0.66
1:D:338:LEU:HB2	1:D:475:VAL:HG22	1.78	0.66
1:A:462:PRO:HA	1:A:466:LEU:HD13	1.76	0.66
1:C:376:ASN:ND2	4:C:760:NAG:H83	2.09	0.66
1:B:3:PRO:HG3	1:B:14:TRP:CH2	2.32	0.65
1:B:92:SER:O	1:B:140:SER:HA	1.95	0.65
1:D:85:ASN:OD1	1:D:121:GLY:HA2	1.96	0.65
1:C:448:PRO:HD3	1:C:468:ARG:NH2	2.12	0.65
1:A:254:ASP:OD2	1:A:477:ARG:HB2	1.97	0.65
1:B:32:GLY:N	9:B:642:HOH:O	2.27	0.65
1:D:254:ASP:OD2	1:D:477:ARG:HB2	1.96	0.65
1:C:3:PRO:HG3	1:C:14:TRP:CH2	2.33	0.64
1:A:557:PHE:HB2	4:A:700:NAG:O7	1.96	0.64
1:B:106:ASN:HD22	1:B:106:ASN:N	1.95	0.64
1:C:106:ASN:N	1:C:106:ASN:HD22	1.95	0.64
1:B:94:HIS:CE1	1:B:437:HIS:CE1	2.86	0.64
1:C:14:TRP:CD1	1:C:15:THR:HG23	2.33	0.64
1:C:182:ASP:HB2	4:C:701:NAG:H82	1.78	0.64
1:D:446:ARG:NH2	9:D:656:HOH:O	2.31	0.63
1:B:433:PRO:HD2	1:B:504:CYS:SG	2.38	0.63
1:C:442:LEU:HD23	1:C:476:ARG:HB3	1.81	0.63
1:C:477:ARG:NH1	9:C:579:HOH:O	2.32	0.62
1:D:135:THR:HG21	1:D:229:ARG:HB3	1.80	0.62
1:A:100:GLN:OE1	1:A:100:GLN:HA	1.98	0.62

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:408:VAL:HG11	1:A:494:THR:HG21	1.81	0.62
1:A:380:ILE:HG21	1:A:515:SER:HB3	1.81	0.62
1:C:92:SER:O	1:C:140:SER:HA	1.99	0.62
1:D:100:GLN:HA	1:D:100:GLN:OE1	2.00	0.61
1:C:244:VAL:C	9:C:651:HOH:O	2.37	0.61
1:B:138:TYR:CZ	1:B:152:GLY:HA3	2.34	0.61
1:D:519:LEU:HG	1:D:522:PRO:HG3	1.83	0.61
1:D:368:THR:O	1:D:368:THR:HG22	2.01	0.61
1:D:380:ILE:HG21	1:D:515:SER:HB3	1.83	0.60
1:A:92:SER:OG	1:A:110:ASN:HB3	2.01	0.60
1:C:138:TYR:CZ	1:C:152:GLY:HA3	2.36	0.60
1:A:85:ASN:OD1	1:A:121:GLY:HA2	2.02	0.60
1:D:408:VAL:HG11	1:D:494:THR:HG21	1.82	0.60
1:A:519:LEU:HG	1:A:522:PRO:HG3	1.82	0.60
1:B:185:VAL:O	1:B:189:GLN:HG3	2.02	0.60
1:D:92:SER:OG	1:D:110:ASN:HB3	2.02	0.59
1:C:185:VAL:O	1:C:189:GLN:HG3	2.02	0.59
1:A:135:THR:HG21	1:A:229:ARG:HB3	1.84	0.59
1:A:12:ALA:O	1:A:19:ASP:HB2	2.03	0.59
1:D:491:ALA:HB3	9:D:663:HOH:O	2.02	0.59
1:D:249:THR:HB	1:D:277:THR:OG1	2.03	0.59
1:B:425:ASN:OD1	1:B:427:ILE:HG22	2.03	0.59
1:C:94:HIS:CE1	1:C:437:HIS:CE1	2.90	0.59
1:D:2:GLY:N	9:D:615:HOH:O	2.35	0.59
1:D:214:GLN:CD	6:D:720:NAG:H83	2.23	0.58
1:C:254:ASP:OD2	1:C:477:ARG:HB2	2.02	0.58
1:C:238:HIS:NE2	1:C:433:PRO:HG3	2.18	0.58
1:B:442:LEU:HD23	1:B:476:ARG:HB3	1.86	0.58
1:C:425:ASN:OD1	1:C:427:ILE:HG22	2.04	0.58
1:A:88:THR:HG22	9:A:573:HOH:O	2.04	0.58
1:A:92:SER:O	1:A:140:SER:HA	2.04	0.57
1:B:448:PRO:HD3	1:B:468:ARG:NH2	2.19	0.57
1:B:282:SER:HB3	1:B:283:PRO:HD2	1.86	0.57
1:D:357:ASN:ND2	9:D:699:HOH:O	2.32	0.57
1:B:266:LEU:HD21	1:B:274:TYR:CG	2.39	0.57
1:D:415:THR:HG22	1:D:492:PHE:O	2.05	0.57
1:C:266:LEU:HD21	1:C:274:TYR:CG	2.40	0.57
1:C:415:THR:HG22	1:C:492:PHE:O	2.04	0.57
1:C:448:PRO:HD3	1:C:468:ARG:HH21	1.70	0.57
1:A:368:THR:CG2	1:A:368:THR:O	2.52	0.57
1:A:249:THR:HB	1:A:277:THR:OG1	2.04	0.57
1:A:253:THR:CG2	1:A:338:LEU:HD12	2.32	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:238:HIS:NE2	1:B:433:PRO:HG3	2.19	0.56
1:D:457:ARG:C	9:D:671:HOH:O	2.43	0.56
1:D:352:VAL:N	9:D:670:HOH:O	2.24	0.56
1:D:92:SER:O	1:D:140:SER:HA	2.06	0.56
1:D:368:THR:CG2	1:D:368:THR:O	2.53	0.56
1:A:170:VAL:HA	1:A:229:ARG:HB2	1.88	0.56
1:C:170:VAL:O	8:C:710:NAG:N2	2.37	0.56
1:D:170:VAL:HA	1:D:229:ARG:HB2	1.88	0.55
1:A:65:ILE:O	1:A:66:ILE:HD13	2.07	0.55
1:A:415:THR:HG22	1:A:492:PHE:O	2.05	0.55
1:B:254:ASP:OD2	1:B:477:ARG:HB2	2.06	0.55
1:A:405:ILE:HD13	1:A:517:ASP:HB3	1.89	0.55
1:B:483:PRO:HG2	9:B:631:HOH:O	2.07	0.55
1:A:308:ALA:HA	9:A:617:HOH:O	2.05	0.55
1:B:415:THR:HG22	1:B:492:PHE:O	2.06	0.55
1:C:7:THR:HB	1:C:8:PRO:HD2	1.88	0.55
1:C:243:LEU:HG	9:C:651:HOH:O	2.07	0.55
1:C:180:SER:OG	4:C:701:NAG:H81	2.07	0.55
1:B:368:THR:HG21	9:B:622:HOH:O	2.06	0.55
1:D:65:ILE:O	1:D:66:ILE:HD13	2.06	0.55
1:C:282:SER:HB3	1:C:283:PRO:HD2	1.87	0.54
1:A:23:ASP:HB3	1:A:26:VAL:HG22	1.89	0.54
1:A:338:LEU:HB2	1:A:475:VAL:HG22	1.87	0.54
1:D:71:ARG:HH12	7:D:711:NAG:H4	1.71	0.54
1:C:65:ILE:O	1:C:66:ILE:HD13	2.08	0.54
1:A:337:ASN:O	1:A:338:LEU:HD23	2.08	0.54
1:D:23:ASP:HB3	1:D:26:VAL:HG22	1.90	0.54
1:D:363:LEU:HB2	1:D:373:TRP:CZ3	2.43	0.54
1:D:200:LEU:HD23	1:D:205:ALA:HB2	1.90	0.54
1:B:23:ASP:HB3	1:B:26:VAL:HG22	1.90	0.54
1:D:491:ALA:N	9:D:663:HOH:O	2.40	0.54
1:C:524:ASP:O	1:C:528:GLN:HG3	2.08	0.54
1:D:405:ILE:HD13	1:D:517:ASP:HB3	1.90	0.54
1:A:1:SER:HB2	1:A:31:THR:HG22	1.90	0.53
1:A:282:SER:HB2	1:A:312:TYR:OH	2.08	0.53
1:A:368:THR:HG23	1:A:368:THR:O	2.09	0.53
1:B:207:HIS:CD2	1:B:209:GLU:H	2.25	0.53
1:C:291:VAL:HB	1:C:306:PRO:HB2	1.91	0.53
1:A:450:GLU:HB2	9:A:670:HOH:O	2.06	0.53
1:A:200:LEU:HD23	1:A:205:ALA:HB2	1.91	0.53
1:C:246:HIS:N	9:C:651:HOH:O	2.41	0.53
1:D:288:TRP:CE2	1:D:320:PRO:HB2	2.44	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:23:ASP:HB3	1:C:26:VAL:CG2	2.39	0.53
1:C:23:ASP:HB3	1:C:26:VAL:HG22	1.91	0.53
1:D:282:SER:HB2	1:D:312:TYR:OH	2.08	0.53
1:D:12:ALA:O	1:D:19:ASP:HB2	2.08	0.53
1:A:427:ILE:HD13	1:B:365:ILE:HD11	1.91	0.53
1:B:7:THR:HB	1:B:8:PRO:HD2	1.90	0.53
1:B:485:GLY:O	9:B:661:HOH:O	2.19	0.52
1:D:461:ASP:N	9:D:650:HOH:O	2.42	0.52
1:C:135:THR:HG21	1:C:229:ARG:HB3	1.92	0.52
1:D:129:ARG:HD3	9:D:661:HOH:O	2.07	0.52
1:B:535:ARG:NH2	9:B:569:HOH:O	2.41	0.52
1:A:194:PRO:HD2	9:A:708:HOH:O	2.09	0.52
1:B:291:VAL:HB	1:B:306:PRO:HB2	1.92	0.52
1:B:64:LYS:HB2	9:B:688:HOH:O	2.09	0.52
1:C:207:HIS:CD2	1:C:209:GLU:H	2.27	0.52
1:A:363:LEU:HB2	1:A:373:TRP:CZ3	2.44	0.52
8:C:711:NAG:O3	8:C:711:NAG:H83	2.09	0.52
1:D:216:TYR:HE2	7:D:710:NAG:H82	1.74	0.52
1:B:448:PRO:HD3	1:B:468:ARG:HH21	1.75	0.52
1:D:305:PHE:HA	9:D:703:HOH:O	2.10	0.52
1:D:146:TYR:HB2	1:D:150:VAL:O	2.09	0.51
1:A:385:GLY:O	1:A:387:PRO:HD3	2.11	0.51
1:D:120:LYS:HA	9:D:595:HOH:O	2.09	0.51
1:B:519:LEU:HG	1:B:522:PRO:HG3	1.92	0.51
1:A:557:PHE:CD2	4:A:701:NAG:H62	2.45	0.51
1:A:61:VAL:N	9:A:577:HOH:O	2.19	0.51
1:D:385:GLY:O	1:D:387:PRO:HD3	2.11	0.51
1:D:1:SER:HB2	1:D:31:THR:HG22	1.92	0.51
1:A:3:PRO:HG3	1:A:14:TRP:CH2	2.46	0.51
1:B:23:ASP:HB3	1:B:26:VAL:CG2	2.41	0.51
1:B:524:ASP:O	1:B:528:GLN:HG3	2.10	0.51
1:D:47:TRP:CZ3	1:D:49:GLY:HA2	2.46	0.51
1:C:119:PRO:HG2	1:C:551:TYR:CZ	2.46	0.51
1:D:6:ASN:HD22	1:D:14:TRP:HE3	1.59	0.51
1:D:26:VAL:HG23	1:D:27:SER:N	2.26	0.51
1:D:325:LEU:HB3	1:D:326:PRO:CD	2.41	0.51
1:C:384:TRP:CZ3	1:C:559:LYS:HD2	2.45	0.51
1:C:41:LEU:HA	1:C:60:LEU:O	2.11	0.51
1:D:354:ARG:NH1	9:D:665:HOH:O	2.43	0.50
1:D:433:PRO:HD2	1:D:504:CYS:SG	2.52	0.50
1:D:218:VAL:HG22	7:D:710:NAG:H81	1.93	0.50
1:B:153:THR:HG21	1:B:172:PRO:HG3	1.92	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:47:TRP:CZ3	1:A:49:GLY:HA2	2.46	0.50
1:C:380:ILE:HG21	1:C:515:SER:HB3	1.94	0.50
1:D:3:PRO:HG3	1:D:14:TRP:CH2	2.47	0.50
1:C:172:PRO:O	1:C:173:LEU:HD23	2.11	0.50
1:B:65:ILE:O	1:B:66:ILE:HD13	2.12	0.50
1:A:41:LEU:CD2	1:A:61:VAL:HG22	2.42	0.50
1:D:437:HIS:O	1:D:500:TRP:HB3	2.11	0.50
1:B:384:TRP:CZ2	1:B:563:GLY:HA2	2.46	0.50
1:A:207:HIS:CD2	1:A:210:THR:H	2.19	0.50
1:C:104:VAL:CG1	9:C:694:HOH:O	2.50	0.50
1:B:384:TRP:CZ3	1:B:559:LYS:HD2	2.46	0.50
1:C:519:LEU:HG	1:C:522:PRO:HG3	1.93	0.50
1:B:305:PHE:CZ	4:B:741:NAG:H61	2.47	0.50
1:A:288:TRP:CE2	1:A:320:PRO:HB2	2.47	0.49
1:C:384:TRP:CE3	1:C:559:LYS:HD2	2.47	0.49
1:B:272:GLN:HG3	1:B:479:VAL:CG1	2.43	0.49
1:B:41:LEU:HD23	1:B:61:VAL:HG22	1.94	0.49
1:C:41:LEU:HD23	1:C:61:VAL:HG22	1.94	0.49
1:A:422:ASP:OD1	1:A:425:ASN:HB2	2.13	0.49
1:B:172:PRO:O	1:B:173:LEU:HD23	2.12	0.49
1:B:384:TRP:CE3	1:B:559:LYS:HD2	2.47	0.49
1:D:177:TYR:CE2	1:D:196:SER:HA	2.48	0.49
1:B:413:GLN:HB2	9:B:580:HOH:O	2.12	0.49
1:C:384:TRP:CZ2	1:C:563:GLY:HA2	2.48	0.49
1:A:354:ARG:HB2	1:A:355:PRO:HD2	1.95	0.49
1:D:347:PRO:HD2	1:D:416:TYR:OH	2.13	0.49
1:B:280:ALA:HB1	1:B:312:TYR:CE2	2.48	0.49
1:C:105:PHE:C	1:C:106:ASN:HD22	2.15	0.48
1:C:50:PRO:HB2	1:C:205:ALA:HB3	1.94	0.48
1:B:105:PHE:C	1:B:106:ASN:HD22	2.14	0.48
1:A:26:VAL:HG23	1:A:27:SER:N	2.27	0.48
1:D:216:TYR:CE2	7:D:710:NAG:H82	2.48	0.48
1:B:380:ILE:HG21	1:B:515:SER:HB3	1.96	0.48
1:C:12:ALA:O	1:C:159:PRO:HB3	2.12	0.48
1:D:338:LEU:HB2	1:D:475:VAL:CG2	2.43	0.48
1:B:46:ASN:ND2	9:B:685:HOH:O	2.36	0.48
1:D:493:LYS:O	1:D:495:ASP:N	2.44	0.48
1:D:354:ARG:HB2	1:D:355:PRO:HD2	1.96	0.48
1:D:461:ASP:C	9:D:650:HOH:O	2.51	0.48
1:C:280:ALA:HB1	1:C:312:TYR:CE2	2.49	0.48
1:B:119:PRO:HG2	1:B:551:TYR:CZ	2.49	0.48
1:B:135:THR:HG21	1:B:229:ARG:HB3	1.96	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:382:VAL:CG1	1:D:564:LEU:HD11	2.43	0.48
1:A:6:ASN:HD22	1:A:14:TRP:HE3	1.62	0.48
1:C:557:PHE:CD2	4:C:701:NAG:H62	2.49	0.48
1:B:124:LYS:NZ	9:B:626:HOH:O	2.46	0.48
1:A:146:TYR:HB2	1:A:150:VAL:O	2.14	0.48
1:D:172:PRO:HD2	1:D:202:ASN:HA	1.95	0.48
1:C:26:VAL:HG23	1:C:27:SER:N	2.30	0.47
1:D:255:MET:HB2	9:D:566:HOH:O	2.13	0.47
1:D:442:LEU:HD23	1:D:476:ARG:HB3	1.95	0.47
1:A:433:PRO:HD2	1:A:504:CYS:SG	2.54	0.47
1:D:382:VAL:HG13	1:D:564:LEU:HD11	1.96	0.47
1:C:368:THR:HG23	1:C:368:THR:O	2.14	0.47
1:A:466:LEU:N	1:A:467:PRO:CD	2.77	0.47
1:D:266:LEU:HD21	1:D:274:TYR:CG	2.50	0.47
1:A:530:ASN:HA	9:A:653:HOH:O	2.13	0.47
1:D:334:ASP:HB2	9:D:664:HOH:O	2.15	0.47
1:B:538:ARG:O	1:B:541:PHE:HB3	2.15	0.47
1:A:55:LYS:HB3	1:A:55:LYS:HE3	1.60	0.47
1:A:325:LEU:HB3	1:A:326:PRO:CD	2.44	0.47
1:A:408:VAL:CG1	1:A:494:THR:HG21	2.44	0.47
1:D:249:THR:HA	1:D:262:THR:HA	1.97	0.47
1:C:251:ILE:O	1:C:252:ALA:HB2	2.15	0.47
1:B:217:ASN:HB2	9:B:612:HOH:O	2.14	0.47
1:A:227:ARG:NH2	9:A:678:HOH:O	2.46	0.47
1:D:252:ALA:HB1	9:D:596:HOH:O	2.15	0.47
1:D:55:LYS:HB3	1:D:55:LYS:HE3	1.66	0.47
1:C:441:PHE:O	1:C:476:ARG:HB2	2.15	0.47
1:C:266:LEU:HD21	1:C:274:TYR:CD1	2.50	0.47
1:D:26:VAL:HA	9:D:625:HOH:O	2.15	0.47
1:D:26:VAL:HG23	1:D:27:SER:H	1.79	0.47
1:D:425:ASN:HA	9:D:589:HOH:O	2.13	0.47
1:C:88:THR:HG21	1:C:557:PHE:HE1	1.80	0.47
1:C:153:THR:HG21	1:C:172:PRO:HG3	1.96	0.46
1:C:99:ARG:HG3	9:C:676:HOH:O	2.14	0.46
1:A:177:TYR:CE2	1:A:196:SER:HA	2.50	0.46
1:C:545:CYS:O	1:C:549:ASN:ND2	2.44	0.46
1:A:557:PHE:HA	1:A:558:PRO:HD3	1.80	0.46
1:C:165:ASP:HB2	1:C:166:ILE:HD12	1.97	0.46
1:C:474:PRO:HA	9:C:598:HOH:O	2.15	0.46
1:B:445:GLY:HA2	9:B:695:HOH:O	2.14	0.46
1:A:249:THR:HA	1:A:262:THR:HA	1.96	0.46
1:D:458:HIS:N	9:D:671:HOH:O	2.48	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:88:THR:HG21	1:B:557:PHE:HE1	1.80	0.46
1:D:207:HIS:CD2	1:D:210:THR:H	2.18	0.46
1:D:388:ILE:HD12	1:D:405:ILE:HD11	1.98	0.46
1:A:380:ILE:O	1:A:562:SER:HB2	2.16	0.46
1:B:431:PRO:HB3	9:B:629:HOH:O	2.16	0.46
4:C:760:NAG:H2	4:C:760:NAG:H82	1.74	0.46
1:A:26:VAL:HG23	1:A:27:SER:H	1.80	0.46
1:B:41:LEU:HA	1:B:60:LEU:O	2.15	0.46
1:A:442:LEU:HD23	1:A:476:ARG:HB3	1.97	0.46
1:C:272:GLN:HG3	1:C:479:VAL:CG1	2.46	0.46
1:B:130:ALA:HB1	1:B:156:ILE:HD13	1.98	0.46
1:D:193:ALA:HA	1:D:194:PRO:HD3	1.80	0.46
1:A:428:VAL:CG1	1:B:427:ILE:HD11	2.46	0.46
1:A:142:PHE:HZ	4:A:700:NAG:H83	1.81	0.46
1:D:295:ASP:HB3	1:D:332:CYS:SG	2.56	0.46
1:A:437:HIS:O	1:A:500:TRP:HB3	2.16	0.46
1:C:207:HIS:HD2	1:C:210:THR:H	1.64	0.45
1:A:493:LYS:O	1:A:495:ASP:N	2.48	0.45
1:A:521:ARG:N	1:A:522:PRO:HD3	2.31	0.45
1:C:255:MET:HG2	1:C:273:ARG:HH21	1.81	0.45
1:D:497:PRO:O	1:D:525:LEU:HD13	2.17	0.45
1:B:36:ALA:HA	1:B:79:GLU:O	2.15	0.45
1:A:349:ASN:N	1:A:349:ASN:OD1	2.48	0.45
1:B:441:PHE:O	1:B:476:ARG:HB2	2.16	0.45
1:A:388:ILE:HD12	1:A:405:ILE:HD11	1.98	0.45
1:A:425:ASN:OD1	1:A:427:ILE:HG23	2.17	0.45
1:A:358:THR:O	6:A:760:NAG:H82	2.16	0.45
1:A:266:LEU:HD21	1:A:274:TYR:CG	2.51	0.45
1:A:497:PRO:O	1:A:525:LEU:HD13	2.17	0.45
1:B:207:HIS:HD2	1:B:210:THR:H	1.65	0.45
1:D:496:ASN:O	1:D:497:PRO:C	2.54	0.45
1:C:538:ARG:O	1:C:541:PHE:HB3	2.17	0.45
1:B:255:MET:HG2	1:B:273:ARG:HH21	1.82	0.45
1:A:388:ILE:CD1	1:A:405:ILE:HD11	2.47	0.45
1:D:466:LEU:N	1:D:467:PRO:CD	2.80	0.45
1:D:137:TRP:HB2	1:D:152:GLY:O	2.16	0.45
1:C:36:ALA:HA	1:C:79:GLU:O	2.16	0.45
1:A:7:THR:HB	1:A:8:PRO:HD2	1.99	0.45
1:B:12:ALA:O	1:B:159:PRO:HB3	2.17	0.45
8:C:711:NAG:O3	8:C:712:BMA:O5	2.30	0.45
1:A:137:TRP:CZ3	1:A:233:THR:HB	2.52	0.45
1:D:408:VAL:CG1	1:D:494:THR:HG21	2.46	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:349:ASN:OD1	1:D:349:ASN:N	2.50	0.44
1:B:266:LEU:HD21	1:B:274:TYR:CD1	2.52	0.44
1:D:300:SER:HB2	9:D:668:HOH:O	2.17	0.44
1:D:41:LEU:CD2	1:D:61:VAL:HG22	2.48	0.44
1:A:207:HIS:HD2	1:A:210:THR:N	2.07	0.44
1:D:135:THR:O	1:D:273:ARG:HG3	2.18	0.44
1:D:387:PRO:O	1:D:390:ASP:HB2	2.18	0.44
1:D:23:ASP:HB3	1:D:26:VAL:CG2	2.48	0.44
1:D:137:TRP:CZ3	1:D:233:THR:HB	2.53	0.44
1:D:558:PRO:O	8:D:700:NAG:H83	2.17	0.44
1:A:380:ILE:HA	1:A:402:SER:O	2.18	0.44
1:B:166:ILE:HG23	7:B:714:MAN:H61	2.00	0.44
1:D:380:ILE:HA	1:D:402:SER:O	2.17	0.44
1:A:555:ASN:HA	1:A:556:PRO:HD3	1.75	0.44
1:C:499:ALA:HB3	9:C:694:HOH:O	2.18	0.43
1:D:442:LEU:HB3	1:D:474:PRO:HG2	1.99	0.43
1:D:422:ASP:OD1	1:D:425:ASN:HB2	2.18	0.43
1:D:545:CYS:O	1:D:549:ASN:ND2	2.51	0.43
1:A:172:PRO:HD2	1:A:202:ASN:HA	2.00	0.43
1:B:65:ILE:HG23	1:B:66:ILE:HG12	2.00	0.43
1:A:443:VAL:HA	1:A:490:LEU:HD23	2.00	0.43
1:C:284:VAL:HG13	1:C:312:TYR:HB3	2.00	0.43
1:A:538:ARG:NE	9:A:653:HOH:O	2.31	0.43
1:A:178:TYR:CE1	1:A:197:ASP:HB3	2.53	0.43
1:C:227:ARG:HB3	9:C:654:HOH:O	2.18	0.43
1:B:368:THR:O	1:B:368:THR:HG23	2.18	0.43
1:A:382:VAL:CG1	1:A:564:LEU:HD11	2.49	0.43
1:D:297:LEU:HD23	1:D:297:LEU:HA	1.82	0.43
1:D:178:TYR:CE1	1:D:197:ASP:HB3	2.54	0.43
1:D:388:ILE:CD1	1:D:405:ILE:HD11	2.48	0.43
1:D:465:ASP:HB2	9:D:650:HOH:O	2.18	0.43
1:D:351:PHE:HA	9:D:670:HOH:O	2.18	0.43
1:D:521:ARG:N	1:D:522:PRO:HD3	2.34	0.43
1:B:26:VAL:HG23	1:B:27:SER:N	2.34	0.43
1:D:425:ASN:OD1	1:D:427:ILE:HG23	2.19	0.43
1:A:441:PHE:O	1:A:476:ARG:HA	2.18	0.43
1:C:387:PRO:HB2	1:C:541:PHE:HZ	1.83	0.43
1:A:282:SER:HB2	1:A:312:TYR:HH	1.83	0.43
1:B:348:VAL:CG1	1:B:444:LEU:HD22	2.49	0.43
1:B:50:PRO:HB2	1:B:205:ALA:HB3	1.99	0.43
1:C:225:ARG:HB2	1:C:279:ASP:OD2	2.19	0.43
1:A:496:ASN:O	1:A:497:PRO:C	2.57	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:189:GLN:HB3	1:B:369:PRO:HB3	2.00	0.42
1:A:193:ALA:HA	1:A:194:PRO:HD3	1.81	0.42
1:D:113:THR:HA	1:D:389:LEU:HG	2.01	0.42
1:A:115:CYS:CB	1:A:545:CYS:HA	2.49	0.42
1:C:55:LYS:HB3	1:C:55:LYS:HE3	1.75	0.42
1:A:35:VAL:HG21	1:A:72:ALA:HB2	2.00	0.42
1:C:422:ASP:CG	1:C:425:ASN:HB2	2.40	0.42
1:B:207:HIS:HA	1:B:208:PRO:HD3	1.91	0.42
1:B:423:PRO:HA	9:B:661:HOH:O	2.19	0.42
1:C:207:HIS:HA	1:C:208:PRO:HD3	1.90	0.42
1:C:355:PRO:HB2	4:C:761:NAG:O6	2.20	0.42
1:A:23:ASP:HB3	1:A:26:VAL:CG2	2.49	0.42
1:A:137:TRP:HB2	1:A:152:GLY:O	2.20	0.42
1:C:226:HIS:O	1:C:277:THR:HA	2.19	0.42
1:D:35:VAL:HG21	1:D:72:ALA:HB2	2.01	0.42
1:C:552:TRP:HB3	1:C:553:PRO:HD3	2.02	0.42
1:C:60:LEU:HD22	1:C:64:LYS:O	2.20	0.42
1:B:370:LEU:N	9:B:624:HOH:O	2.52	0.42
1:B:41:LEU:CD2	1:B:61:VAL:HG22	2.49	0.42
4:B:740:NAG:HO3	4:B:741:NAG:H62	1.83	0.42
1:D:382:VAL:HG11	1:D:564:LEU:HD21	2.02	0.42
1:A:545:CYS:O	1:A:549:ASN:ND2	2.53	0.42
1:D:7:THR:HB	1:D:8:PRO:HD2	2.02	0.42
1:D:207:HIS:HD2	1:D:210:THR:N	2.06	0.42
1:A:427:ILE:HD11	1:B:427:ILE:HG13	2.00	0.42
1:A:142:PHE:CZ	4:A:700:NAG:H83	2.54	0.42
1:D:11:ARG:HG2	1:D:163:PRO:HA	2.01	0.42
1:B:137:TRP:CE3	1:B:233:THR:HB	2.55	0.42
1:B:534:LYS:O	1:B:535:ARG:C	2.58	0.42
1:B:165:ASP:HB2	1:B:166:ILE:HD12	2.00	0.42
1:D:115:CYS:CB	1:D:545:CYS:HA	2.49	0.42
1:B:193:ALA:HA	1:B:194:PRO:HD3	1.66	0.42
1:B:288:TRP:CE2	1:B:320:PRO:HB2	2.55	0.42
1:D:441:PHE:O	1:D:476:ARG:HA	2.20	0.42
1:A:7:THR:O	1:A:11:ARG:HD3	2.20	0.42
1:A:137:TRP:CZ2	1:A:271:GLY:HA2	2.54	0.42
1:A:382:VAL:HG13	1:A:564:LEU:HD11	2.01	0.42
1:D:555:ASN:HA	1:D:556:PRO:HD3	1.71	0.42
1:C:180:SER:CB	4:C:701:NAG:H81	2.50	0.42
1:A:387:PRO:O	1:A:390:ASP:HB2	2.20	0.42
1:C:288:TRP:CE2	1:C:320:PRO:HB2	2.55	0.42
1:C:499:ALA:N	9:C:694:HOH:O	2.28	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:60:LEU:HD22	1:B:64:LYS:O	2.20	0.41
1:C:136:SER:HA	1:C:273:ARG:HD2	2.01	0.41
1:B:251:ILE:O	1:B:252:ALA:HB2	2.20	0.41
1:B:425:ASN:HA	1:B:426:PRO:HD3	1.88	0.41
1:D:543:ARG:NH2	9:D:681:HOH:O	2.37	0.41
1:C:447:SER:HB3	1:C:487:TRP:CD1	2.55	0.41
1:B:289:PHE:O	1:B:307:ALA:HA	2.20	0.41
1:C:305:PHE:CE1	4:C:741:NAG:H5	2.55	0.41
1:A:11:ARG:HG2	1:A:163:PRO:HA	2.02	0.41
1:C:137:TRP:CE3	1:C:233:THR:HB	2.55	0.41
1:C:357:ASN:HB2	1:C:416:TYR:O	2.21	0.41
1:D:534:LYS:O	1:D:535:ARG:C	2.59	0.41
1:B:94:HIS:CE1	1:B:437:HIS:ND1	2.88	0.41
1:A:174:MET:HB2	1:A:233:THR:OG1	2.21	0.41
1:C:193:ALA:HA	1:C:194:PRO:HD3	1.69	0.41
1:A:425:ASN:HA	1:A:426:PRO:HD3	1.91	0.41
1:A:238:HIS:CD2	1:A:269:ALA:HA	2.56	0.41
1:D:99:ARG:O	1:D:100:GLN:C	2.59	0.41
1:B:254:ASP:HA	1:B:273:ARG:HG2	2.02	0.41
1:A:153:THR:HG21	1:A:172:PRO:HG3	2.02	0.41
1:A:382:VAL:HG11	1:A:564:LEU:HD21	2.02	0.41
1:D:72:ALA:O	1:D:156:ILE:HA	2.21	0.41
1:B:342:VAL:HG12	1:B:344:ARG:HH12	1.85	0.41
1:C:249:THR:O	1:C:276:VAL:HA	2.21	0.41
1:D:7:THR:O	1:D:11:ARG:HD3	2.20	0.41
1:D:534:LYS:O	1:D:537:ASP:N	2.54	0.41
1:A:297:LEU:HA	1:A:297:LEU:HD23	1.87	0.41
1:B:447:SER:HB3	1:B:487:TRP:CD1	2.56	0.41
1:D:92:SER:OG	1:D:110:ASN:CB	2.68	0.41
1:B:136:SER:HA	1:B:273:ARG:HD2	2.02	0.41
1:C:551:TYR:O	1:C:552:TRP:C	2.59	0.41
1:D:8:PRO:HA	1:D:164:TYR:O	2.21	0.41
1:A:347:PRO:HD2	1:A:416:TYR:OH	2.20	0.41
1:C:342:VAL:HG12	1:C:344:ARG:HH12	1.85	0.41
1:A:381:ASN:HB2	9:A:736:HOH:O	2.21	0.41
1:A:435:HIS:CE1	1:A:478:ASP:HB2	2.56	0.41
1:A:427:ILE:HG12	1:B:427:ILE:HG12	2.03	0.41
1:B:387:PRO:HB2	1:B:541:PHE:HZ	1.85	0.41
1:A:72:ALA:O	1:A:156:ILE:HA	2.21	0.41
1:D:11:ARG:HB3	1:D:160:ALA:HB3	2.02	0.41
1:D:106:ASN:HD22	1:D:106:ASN:N	2.19	0.41
1:D:443:VAL:HA	1:D:490:LEU:HD23	2.01	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:107:ASP:HB3	1:B:113:THR:HG21	2.03	0.40
1:A:534:LYS:O	1:A:535:ARG:C	2.59	0.40
1:B:545:CYS:O	1:B:549:ASN:ND2	2.50	0.40
1:B:180:SER:OG	1:B:183:GLU:HG3	2.21	0.40
1:B:473:ASN:N	1:B:474:PRO:HD3	2.36	0.40
1:B:20:ILE:HG13	1:B:20:ILE:O	2.22	0.40
1:A:259:ASN:OD1	1:A:341:VAL:HG22	2.21	0.40
1:B:365:ILE:O	1:B:365:ILE:HG22	2.20	0.40
1:C:106:ASN:N	1:C:106:ASN:ND2	2.66	0.40
1:A:187:PHE:CE2	1:A:194:PRO:HG3	2.56	0.40
1:D:137:TRP:CZ2	1:D:271:GLY:HA2	2.57	0.40
1:D:259:ASN:OD1	1:D:341:VAL:HG22	2.21	0.40
1:A:157:ASP:HA	9:A:652:HOH:O	2.21	0.40
1:A:336:LEU:O	1:A:475:VAL:HG23	2.21	0.40
1:D:133:TYR:CE1	1:D:159:PRO:HD2	2.57	0.40
1:D:325:LEU:HB3	1:D:326:PRO:HD2	2.02	0.40
1:C:252:ALA:HB3	1:C:275:ASP:HB2	2.03	0.40
1:D:134:GLY:HA2	9:D:653:HOH:O	2.20	0.40
7:B:710:NAG:H62	7:B:711:NAG:C7	2.51	0.40
1:C:348:VAL:CG1	1:C:444:LEU:HD22	2.51	0.40
5:A:711:NAG:C1	5:A:711:NAG:H82	2.52	0.40
1:A:389:LEU:HD22	1:A:529:LEU:HD11	2.03	0.40
1:C:89:ASN:HB3	1:C:145:GLN:NE2	2.36	0.40
1:A:428:VAL:HG11	1:B:427:ILE:HD11	2.02	0.40
1:A:442:LEU:HB3	1:A:474:PRO:HG2	2.04	0.40
1:D:112:VAL:HG21	1:D:501:LEU:HD21	2.03	0.40
1:A:316:PRO:HB2	1:A:318:THR:HG23	2.04	0.40
1:B:106:ASN:N	1:B:106:ASN:ND2	2.66	0.40
1:C:151:VAL:HG22	1:C:152:GLY:N	2.36	0.40
1:A:249:THR:O	1:A:276:VAL:HA	2.21	0.40
1:B:367:GLY:O	1:B:368:THR:C	2.59	0.40
1:C:282:SER:HB3	1:C:283:PRO:CD	2.52	0.40
1:C:41:LEU:CD2	1:C:61:VAL:HG22	2.51	0.40
1:A:162:LEU:HD12	1:A:162:LEU:HA	1.98	0.40
1:B:177:TYR:N	1:B:177:TYR:CD1	2.90	0.40
1:D:206:ARG:NE	9:D:660:HOH:O	2.54	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	562/604 (93%)	517 (92%)	42 (8%)	3 (0%)	38	60
1	B	562/604 (93%)	502 (89%)	60 (11%)	0	100	100
1	C	562/604 (93%)	505 (90%)	57 (10%)	0	100	100
1	D	562/604 (93%)	515 (92%)	44 (8%)	3 (0%)	38	60
All	All	2248/2416 (93%)	2039 (91%)	203 (9%)	6 (0%)	50	73

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	494	THR
1	D	494	THR
1	A	284	VAL
1	D	222	PRO
1	D	497	PRO
1	A	222	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	484/513 (94%)	477 (99%)	7 (1%)	78	94
1	B	484/513 (94%)	476 (98%)	8 (2%)	73	92
1	C	484/513 (94%)	477 (99%)	7 (1%)	78	94
1	D	484/513 (94%)	476 (98%)	8 (2%)	73	92
All	All	1936/2052 (94%)	1906 (98%)	30 (2%)	76	93

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	110	ASN
1	A	141	HIS
1	A	273	ARG
1	A	311	ARG
1	A	349	ASN
1	A	368	THR
1	A	427	ILE
1	B	91	THR
1	B	106	ASN
1	B	110	ASN
1	B	141	HIS
1	B	273	ARG
1	B	275	ASP
1	B	349	ASN
1	B	466	LEU
1	C	106	ASN
1	C	110	ASN
1	C	141	HIS
1	C	273	ARG
1	C	275	ASP
1	C	349	ASN
1	C	466	LEU
1	D	110	ASN
1	D	141	HIS
1	D	273	ARG
1	D	311	ARG
1	D	338	LEU
1	D	349	ASN
1	D	368	THR
1	D	427	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	46	ASN
1	A	103	ASN
1	A	207	HIS
1	B	10	ASN
1	B	73	ASN
1	B	103	ASN
1	B	207	HIS
1	B	240	GLN
1	C	10	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	103	ASN
1	C	145	GLN
1	C	155	GLN
1	C	207	HIS
1	C	240	GLN
1	D	103	ASN
1	D	207	HIS
1	D	237	ASN

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 ⓘ

38 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	$\# Z > 2$
4	NAG	A	700	1,4	12,14,15	0.65	0	15,19,21	1.44	4 (26%)
4	NAG	A	701	4	12,14,15	0.56	0	15,19,21	1.11	1 (6%)
5	NAG	A	710	1,5	12,14,15	0.65	1 (8%)	15,19,21	1.02	0
5	NAG	A	711	5	12,14,15	0.66	0	15,19,21	1.06	1 (6%)
5	BMA	A	712	5	10,11,12	0.73	0	11,15,17	1.24	1 (9%)
5	MAN	A	713	5	10,11,12	0.70	0	11,15,17	0.78	0
5	MAN	A	714	5	10,11,12	0.87	1 (10%)	11,15,17	1.00	1 (9%)
5	MAN	A	715	5	10,11,12	0.73	0	11,15,17	1.73	2 (18%)
4	NAG	A	740	1,4	12,14,15	0.82	1 (8%)	15,19,21	1.62	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	A	741	4	12,14,15	0.67	0	15,19,21	1.28	1 (6%)
4	NAG	B	700	1,4	12,14,15	0.71	1 (8%)	15,19,21	1.51	2 (13%)
4	NAG	B	701	4	12,14,15	0.64	0	15,19,21	1.10	1 (6%)
7	NAG	B	710	1,7	12,14,15	0.78	1 (8%)	15,19,21	1.04	0
7	NAG	B	711	7	12,14,15	0.82	1 (8%)	15,19,21	1.13	1 (6%)
7	BMA	B	712	7	10,11,12	0.84	1 (10%)	11,15,17	1.38	1 (9%)
7	MAN	B	714	7	10,11,12	0.83	1 (10%)	11,15,17	1.32	1 (9%)
7	MAN	B	715	7	10,11,12	0.56	0	11,15,17	2.39	5 (45%)
4	NAG	B	740	1,4	12,14,15	0.63	0	15,19,21	1.11	1 (6%)
4	NAG	B	741	4	12,14,15	0.65	0	15,19,21	1.23	1 (6%)
4	NAG	C	700	1,4	12,14,15	0.69	1 (8%)	15,19,21	1.23	2 (13%)
4	NAG	C	701	4	12,14,15	0.72	1 (8%)	15,19,21	1.03	1 (6%)
8	NAG	C	710	1,8	12,14,15	0.64	0	15,19,21	1.00	1 (6%)
8	NAG	C	711	8	12,14,15	0.62	0	15,19,21	1.11	1 (6%)
8	BMA	C	712	8	10,11,12	0.92	1 (10%)	11,15,17	1.85	3 (27%)
4	NAG	C	740	1,4	12,14,15	0.73	1 (8%)	15,19,21	1.28	2 (13%)
4	NAG	C	741	4	12,14,15	0.51	0	15,19,21	1.65	3 (20%)
4	NAG	C	760	1,4	12,14,15	0.63	0	15,19,21	1.10	1 (6%)
4	NAG	C	761	4	12,14,15	0.76	1 (8%)	15,19,21	1.36	3 (20%)
8	NAG	D	700	1,8	12,14,15	0.48	0	15,19,21	1.27	2 (13%)
8	NAG	D	701	8	12,14,15	0.70	0	15,19,21	1.01	1 (6%)
8	BMA	D	702	8	10,11,12	0.78	0	11,15,17	0.84	1 (9%)
7	NAG	D	710	1,7	12,14,15	0.62	0	15,19,21	1.54	3 (20%)
7	NAG	D	711	7	12,14,15	0.82	1 (8%)	15,19,21	1.42	1 (6%)
7	BMA	D	712	7	10,11,12	0.77	0	11,15,17	1.08	1 (9%)
7	MAN	D	714	7	10,11,12	0.84	1 (10%)	11,15,17	1.67	2 (18%)
7	MAN	D	715	7	10,11,12	0.75	0	11,15,17	1.59	2 (18%)
4	NAG	D	740	1,4	12,14,15	0.66	0	15,19,21	0.98	0
4	NAG	D	741	4	12,14,15	0.59	0	15,19,21	1.09	1 (6%)

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
4	NAG	A	700	1,4	-	0/6/23/26	0/1/1/1
4	NAG	A	701	4	-	0/6/23/26	0/1/1/1
5	NAG	A	710	1,5	-	0/6/23/26	0/1/1/1
5	NAG	A	711	5	-	0/6/23/26	0/1/1/1
5	BMA	A	712	5	-	0/2/19/22	0/1/1/1
5	MAN	A	713	5	-	0/2/19/22	0/1/1/1
5	MAN	A	714	5	-	0/2/19/22	0/1/1/1
5	MAN	A	715	5	-	0/2/19/22	0/1/1/1
4	NAG	A	740	1,4	-	0/6/23/26	0/1/1/1
4	NAG	A	741	4	-	0/6/23/26	0/1/1/1
4	NAG	B	700	1,4	-	0/6/23/26	0/1/1/1
4	NAG	B	701	4	-	0/6/23/26	0/1/1/1
7	NAG	B	710	1,7	-	0/6/23/26	0/1/1/1
7	NAG	B	711	7	-	0/6/23/26	0/1/1/1
7	BMA	B	712	7	-	0/2/19/22	0/1/1/1
7	MAN	B	714	7	-	0/2/19/22	0/1/1/1
7	MAN	B	715	7	-	0/2/19/22	0/1/1/1
4	NAG	B	740	1,4	-	0/6/23/26	0/1/1/1
4	NAG	B	741	4	-	0/6/23/26	0/1/1/1
4	NAG	C	700	1,4	-	0/6/23/26	0/1/1/1
4	NAG	C	701	4	-	0/6/23/26	0/1/1/1
8	NAG	C	710	1,8	-	0/6/23/26	0/1/1/1
8	NAG	C	711	8	-	0/6/23/26	0/1/1/1
8	BMA	C	712	8	-	0/2/19/22	0/1/1/1
4	NAG	C	740	1,4	-	0/6/23/26	0/1/1/1
4	NAG	C	741	4	-	0/6/23/26	0/1/1/1
4	NAG	C	760	1,4	-	0/6/23/26	0/1/1/1
4	NAG	C	761	4	-	0/6/23/26	0/1/1/1
8	NAG	D	700	1,8	-	0/6/23/26	0/1/1/1
8	NAG	D	701	8	-	0/6/23/26	0/1/1/1
8	BMA	D	702	8	-	0/2/19/22	0/1/1/1
7	NAG	D	710	1,7	-	0/6/23/26	0/1/1/1
7	NAG	D	711	7	-	0/6/23/26	0/1/1/1
7	BMA	D	712	7	-	0/2/19/22	0/1/1/1
7	MAN	D	714	7	-	0/2/19/22	0/1/1/1
7	MAN	D	715	7	-	0/2/19/22	0/1/1/1
4	NAG	D	740	1,4	-	0/6/23/26	0/1/1/1
4	NAG	D	741	4	-	0/6/23/26	0/1/1/1

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	D	711	NAG	O5-C5	-2.39	1.41	1.45

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	C	712	BMA	O5-C5	-2.32	1.41	1.45
7	B	710	NAG	O5-C5	-2.30	1.41	1.45
4	A	740	NAG	O5-C5	-2.23	1.41	1.45
4	B	700	NAG	O5-C5	-2.20	1.41	1.45
7	B	711	NAG	O5-C5	-2.19	1.41	1.45
4	C	701	NAG	O5-C5	-2.17	1.41	1.45
7	D	714	MAN	O5-C5	-2.16	1.41	1.45
5	A	714	MAN	O5-C5	-2.16	1.41	1.45
4	C	740	NAG	O5-C5	-2.15	1.41	1.45
4	C	761	NAG	O5-C5	-2.11	1.41	1.45
7	B	714	MAN	O5-C5	-2.10	1.41	1.45
4	C	700	NAG	O5-C5	-2.03	1.41	1.45
7	B	712	BMA	O5-C5	-2.02	1.41	1.45
5	A	710	NAG	O5-C5	-2.02	1.41	1.45

All (56) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	B	715	MAN	O5-C5-C6	5.60	112.86	106.98
4	A	740	NAG	O5-C5-C6	4.95	112.18	106.98
5	A	715	MAN	C4-C3-C2	-4.86	103.99	110.50
4	C	741	NAG	O5-C5-C6	4.43	111.63	106.98
7	D	714	MAN	C4-C3-C2	4.04	115.92	110.50
8	C	712	BMA	O5-C5-C6	3.72	110.89	106.98
4	A	741	NAG	O5-C5-C6	3.67	110.84	106.98
4	B	700	NAG	O5-C5-C4	3.64	115.27	110.65
7	D	715	MAN	C4-C3-C2	3.49	115.19	110.50
8	C	711	NAG	C2-N2-C7	-3.46	117.28	123.09
4	B	741	NAG	O5-C5-C6	3.45	110.60	106.98
8	C	712	BMA	O5-C5-C4	-3.43	106.30	110.65
7	D	715	MAN	O5-C5-C6	3.27	110.41	106.98
7	D	714	MAN	C3-C4-C5	3.22	115.95	110.20
4	B	740	NAG	O5-C5-C6	3.21	110.35	106.98
7	B	715	MAN	C4-C3-C2	-3.17	106.25	110.50
4	C	740	NAG	O5-C5-C6	3.11	110.25	106.98
4	C	761	NAG	O5-C5-C6	3.11	110.24	106.98
7	B	712	BMA	O5-C5-C6	3.10	110.23	106.98
4	C	760	NAG	C2-N2-C7	-3.08	117.92	123.09
5	A	712	BMA	O5-C5-C6	3.07	110.21	106.98
8	C	712	BMA	C4-C3-C2	3.07	114.63	110.50
8	D	700	NAG	O5-C5-C4	3.02	114.49	110.65
7	D	710	NAG	C3-C4-C5	3.02	115.59	110.20

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	701	NAG	O5-C5-C6	3.00	110.13	106.98
4	C	761	NAG	O5-C5-C4	-2.98	106.87	110.65
7	D	710	NAG	C2-N2-C7	-2.87	118.28	123.09
4	B	701	NAG	C2-N2-C7	-2.82	118.36	123.09
4	C	741	NAG	C2-N2-C7	-2.68	118.58	123.09
7	D	711	NAG	C4-C3-C2	2.63	117.76	111.32
4	C	701	NAG	O5-C5-C6	2.57	109.68	106.98
7	B	715	MAN	O3-C3-C4	2.55	116.06	110.35
4	A	700	NAG	C6-C5-C4	-2.54	106.86	113.00
8	C	710	NAG	O5-C5-C6	2.53	109.64	106.98
7	D	712	BMA	O5-C5-C6	2.53	109.63	106.98
4	A	700	NAG	O5-C5-C4	2.52	113.86	110.65
5	A	711	NAG	C2-N2-C7	2.46	127.22	123.09
7	B	714	MAN	C3-C4-C5	-2.42	105.88	110.20
4	B	700	NAG	C6-C5-C4	-2.42	107.16	113.00
7	B	715	MAN	C3-C4-C5	-2.41	105.89	110.20
4	C	741	NAG	C4-C3-C2	-2.41	105.42	111.32
4	C	740	NAG	O5-C5-C4	-2.39	107.63	110.65
4	A	700	NAG	C2-N2-C7	-2.34	119.17	123.09
7	B	711	NAG	C3-C2-N2	-2.26	108.33	111.76
8	D	702	BMA	O5-C5-C6	2.25	109.34	106.98
4	C	700	NAG	O5-C5-C6	2.23	109.32	106.98
7	D	710	NAG	O5-C5-C6	2.21	109.30	106.98
7	B	715	MAN	O4-C4-C3	2.20	115.29	110.35
4	C	700	NAG	C2-N2-C7	-2.19	119.42	123.09
5	A	714	MAN	O3-C3-C2	-2.17	105.97	109.94
4	A	700	NAG	O6-C6-C5	-2.14	103.98	111.36
4	D	741	NAG	C2-N2-C7	-2.13	119.50	123.09
8	D	700	NAG	C3-C2-N2	2.13	115.01	111.76
5	A	715	MAN	C6-C5-C4	-2.09	107.95	113.00
4	C	761	NAG	O3-C3-C2	2.07	113.44	109.09
8	D	701	NAG	O5-C5-C6	2.03	109.11	106.98

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.6 Ligand geometry ⓘ

Of 25 ligands modelled in this entry, 16 are monoatomic - leaving 9 for Mogul analysis.

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	OXY	A	620	2	1,1,1	0.44	0	0,0,0	0.00	-
6	NAG	A	720	1	12,14,15	0.71	1 (8%)	15,19,21	1.09	1 (6%)
6	NAG	A	730	1	12,14,15	0.68	1 (8%)	15,19,21	1.05	0
6	NAG	A	760	1	12,14,15	0.71	1 (8%)	15,19,21	1.37	2 (13%)
6	NAG	B	720	1	12,14,15	0.71	1 (8%)	15,19,21	0.79	0
6	NAG	B	760	1	12,14,15	0.63	0	15,19,21	1.46	2 (13%)
6	NAG	B	770	1	12,14,15	0.78	1 (8%)	15,19,21	1.63	2 (13%)
6	NAG	C	730	1	12,14,15	0.63	0	15,19,21	1.26	2 (13%)
6	NAG	D	720	1	12,14,15	0.69	1 (8%)	15,19,21	1.41	3 (20%)

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	OXY	A	620	2	-	0/0/0/0	0/0/0/0
6	NAG	A	720	1	-	0/6/23/26	0/1/1/1
6	NAG	A	730	1	-	0/6/23/26	0/1/1/1
6	NAG	A	760	1	-	0/6/23/26	0/1/1/1
6	NAG	B	720	1	-	0/6/23/26	0/1/1/1
6	NAG	B	760	1	-	0/6/23/26	0/1/1/1
6	NAG	B	770	1	-	1/6/23/26	0/1/1/1
6	NAG	C	730	1	-	0/6/23/26	0/1/1/1
6	NAG	D	720	1	-	0/6/23/26	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	770	NAG	O5-C5	-2.29	1.41	1.45
6	A	720	NAG	O5-C5	-2.21	1.41	1.45
6	A	760	NAG	O5-C5	-2.11	1.41	1.45
6	D	720	NAG	O5-C5	-2.07	1.41	1.45

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	730	NAG	O5-C5	-2.06	1.41	1.45
6	B	720	NAG	O5-C5	-2.02	1.41	1.45

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	770	NAG	O5-C5-C6	5.25	112.49	106.98
6	A	760	NAG	C2-N2-C7	-3.52	117.17	123.09
6	B	760	NAG	C2-N2-C7	-3.52	117.18	123.09
6	D	720	NAG	O5-C5-C4	3.02	114.49	110.65
6	A	720	NAG	O5-C5-C6	2.57	109.67	106.98
6	D	720	NAG	C3-C4-C5	2.48	114.63	110.20
6	B	760	NAG	O3-C3-C2	2.48	114.29	109.09
6	B	770	NAG	O5-C5-C4	-2.28	107.76	110.65
6	C	730	NAG	C4-C3-C2	-2.17	106.01	111.32
6	C	730	NAG	O5-C5-C4	2.16	113.39	110.65
6	D	720	NAG	C3-C2-N2	-2.14	108.50	111.76
6	A	760	NAG	C3-C2-N2	-2.14	108.50	111.76

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	B	770	NAG	O7-C7-N2-C2

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, 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	564/604 (93%)	-0.18	1 (0%) 93 94	11, 19, 31, 50	0
1	B	564/604 (93%)	-0.22	2 (0%) 90 92	9, 19, 34, 47	0
1	C	564/604 (93%)	-0.20	1 (0%) 93 94	10, 20, 34, 48	0
1	D	564/604 (93%)	-0.01	6 (1%) 77 79	11, 20, 32, 51	0
All	All	2256/2416 (93%)	-0.15	10 (0%) 90 92	9, 19, 33, 51	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	368	THR	3.7
1	B	367	GLY	3.5
1	B	366	GLY	2.3
1	D	395	GLY	2.3
1	D	368	THR	2.2
1	D	32	GLY	2.2
1	D	351	PHE	2.1
1	D	223	GLY	2.1
1	D	413	GLN	2.0
1	C	16	ASN	2.0

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, 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
4	NAG	C	761	14/15	0.25	4.53	18,33,42,48	0
7	MAN	D	714	11/12	0.20	4.50	24,41,53,54	0
5	MAN	A	715	11/12	0.18	4.38	18,26,37,41	0
4	NAG	A	741	14/15	0.18	3.24	19,25,31,35	0
5	MAN	A	714	11/12	0.13	2.56	17,24,33,39	0
4	NAG	C	741	14/15	0.19	1.87	12,24,29,34	0
4	NAG	B	741	14/15	0.18	1.86	17,28,33,34	0
7	NAG	D	711	14/15	0.15	1.86	21,28,39,44	0
8	NAG	D	701	14/15	0.17	1.56	15,26,42,54	0
7	MAN	B	715	11/12	0.18	1.18	18,29,36,38	0
4	NAG	C	760	14/15	0.15	0.74	8,27,36,45	0
4	NAG	D	740	14/15	0.17	0.66	12,23,34,35	0
4	NAG	A	740	14/15	0.14	0.61	11,16,25,28	0
8	NAG	C	711	14/15	0.15	0.50	19,25,33,38	0
4	NAG	C	740	14/15	0.15	0.29	13,23,30,36	0
4	NAG	B	701	14/15	0.14	0.25	10,14,25,26	0
4	NAG	B	740	14/15	0.14	0.16	8,24,32,34	0
8	NAG	C	710	14/15	0.14	0.08	12,20,29,29	0
4	NAG	B	700	14/15	0.14	-0.18	8,16,23,26	0
7	NAG	B	711	14/15	0.13	-0.20	8,22,28,30	0
7	MAN	D	715	11/12	0.17	-0.33	32,50,60,63	0
4	NAG	A	700	14/15	0.14	-0.43	9,16,22,30	0
5	NAG	A	711	14/15	0.13	-0.57	7,18,25,28	0
7	NAG	B	710	14/15	0.12	-0.57	12,19,27,34	0
4	NAG	D	741	14/15	0.16	-0.60	16,36,42,44	0
4	NAG	C	701	14/15	0.13	-0.75	14,28,40,40	0
4	NAG	A	701	14/15	0.12	-0.79	12,19,29,30	0
5	NAG	A	710	14/15	0.12	-1.06	11,15,24,25	0
8	NAG	D	700	14/15	0.12	-1.33	13,18,25,25	0
4	NAG	C	700	14/15	0.12	-1.37	12,20,28,29	0
7	NAG	D	710	14/15	0.10	-1.97	17,28,32,33	0
7	MAN	B	714	11/12	0.12	-3.53	16,30,35,37	0
8	BMA	D	702	11/12	0.20	-	38,51,61,65	0
7	BMA	B	712	11/12	0.14	-	9,24,28,40	0
7	BMA	D	712	11/12	0.13	-	32,48,57,68	0
5	BMA	A	712	11/12	0.14	-	12,25,33,34	0
5	MAN	A	713	11/12	0.23	-	16,32,37,42	0
8	BMA	C	712	11/12	0.15	-	12,28,36,37	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, 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
6	NAG	A	730	14/15	0.18	5.10	13,31,41,56	0
6	NAG	C	730	14/15	0.20	4.67	19,29,44,58	0
3	OXY	A	620	2/2	0.17	1.93	5,5,5,8	0
6	NAG	A	760	14/15	0.16	0.79	12,21,25,30	0
6	NAG	B	720	14/15	0.17	0.21	13,23,35,36	0
6	NAG	B	760	14/15	0.15	0.09	6,15,25,28	0
6	NAG	B	770	14/15	0.15	0.04	17,27,34,34	0
6	NAG	D	720	14/15	0.16	-0.04	14,20,28,32	0
6	NAG	A	720	14/15	0.13	-0.21	15,20,29,44	0
2	CU	D	602	1/1	0.10	-1.10	26,26,26,26	0
2	CU	C	603	1/1	0.10	-2.00	18,18,18,18	0
2	CU	D	603	1/1	0.09	-2.06	20,20,20,20	0
2	CU	C	601	1/1	0.06	-2.77	21,21,21,21	0
2	CU	A	603	1/1	0.07	-3.20	14,14,14,14	0
2	CU	B	603	1/1	0.09	-3.29	13,13,13,13	0
2	CU	B	602	1/1	0.09	-3.50	17,17,17,17	0
2	CU	D	604	1/1	0.06	-3.54	33,33,33,33	0
2	CU	A	601	1/1	0.06	-4.20	31,31,31,31	0
2	CU	C	604	1/1	0.06	-4.28	30,30,30,30	0
2	CU	A	604	1/1	0.07	-4.36	27,27,27,27	0
2	CU	C	602	1/1	0.08	-4.78	21,21,21,21	0
2	CU	B	604	1/1	0.05	-5.42	26,26,26,26	0
2	CU	D	601	1/1	0.06	-5.45	19,19,19,19	0
2	CU	B	601	1/1	0.05	-7.49	21,21,21,21	0
2	CU	A	602	1/1	0.09	-8.76	11,11,11,11	0

6.5 Other polymers ⓘ

There are no such residues in this entry.