



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 08:07 PM GMT

PDB ID : 1IVD
Title : STRUCTURES OF AROMATIC INHIBITORS OF INFLUENZA VIRUS
NEURAMINIDASE
Authors : Jedrzejewski, M.J.; Luo, M.
Deposited on : 1994-12-12
Resolution : 1.90 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

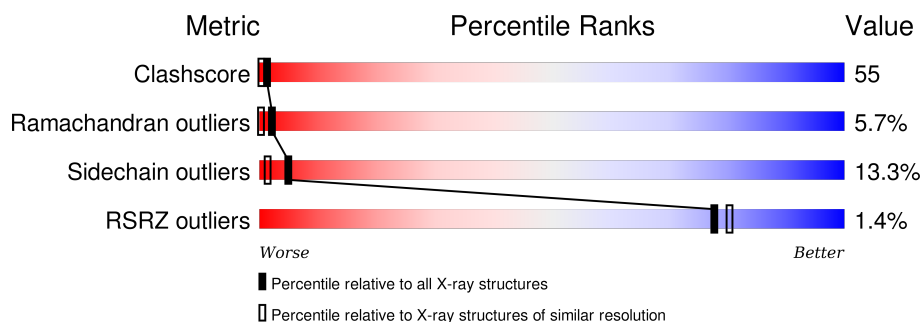
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	102246	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)
RSRZ outliers	91569	4766 (1.90-1.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	388	<div> <div>29%</div> <div>58%</div> <div>12%</div> </div>
1	B	388	<div> <div>28%</div> <div>60%</div> <div>11%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	NAG	B	484	-	-	-	X
4	NAG	A	478	-	-	X	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	MAN	A	483	-	-	X	X
4	NAG	B	478	-	-	-	X
4	MAN	B	483	-	-	-	X
5	FUC	B	477	X	-	-	-

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 8595 atoms, of which 2036 are hydrogens and 0 are deuteriums.

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

- Molecule 1 is a protein called INFLUENZA A SUBTYPE N2 NEURAMINIDASE.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	388	Total	C	H	N	O	S	0	0	0
			3745	1866	723	545	588	23			
1	B	388	Total	C	H	N	O	S	0	0	0
			3745	1866	723	545	588	23			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	339	ASP	ASN	CONFLICT	UNP P06820
B	339	ASP	ASN	CONFLICT	UNP P06820

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

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	2	Total	C	H	N	O		0	0
			55	16	27	2	10			
2	A	2	Total	C	H	N	O		0	0
			55	16	27	2	10			
2	B	2	Total	C	H	N	O		0	0
			55	16	27	2	10			
2	B	2	Total	C	H	N	O		0	0
			55	16	27	2	10			

- Molecule 3 is a polymer of unknown type called SUGAR (4-MER).

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	4	Total	C	H	N	O		0	0
			96	28	47	2	19			

- Molecule 4 is a polymer of unknown type called SUGAR (6-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	6	Total	C	H	N	O	0	0
			139	40	67	2	30		
4	B	6	Total	C	H	N	O	0	0
			139	40	67	2	30		

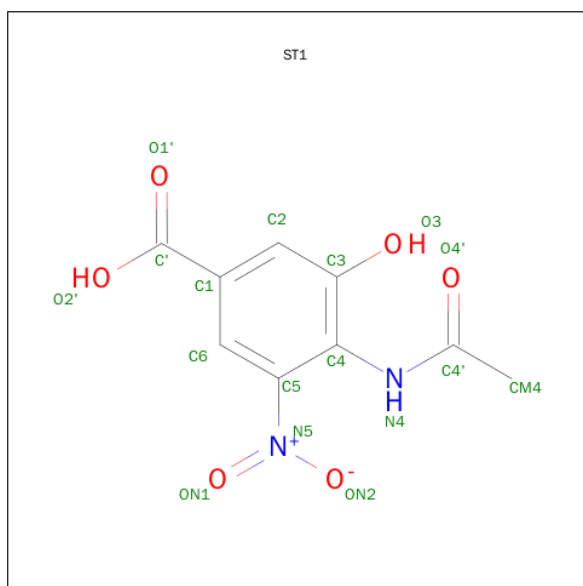
- Molecule 5 is a polymer of unknown type called SUGAR (4-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	B	4	Total	C	H	N	O	0	0
			96	28	47	2	19		

- Molecule 6 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	1	Total	Ca	0	0
			1	1		
6	A	1	Total	Ca	0	0
			1	1		

- Molecule 7 is 4-(ACETYLAMINO)-3-HYDROXY-5-NITROBENZOIC ACID (three-letter code: ST1) (formula: C₉H₈N₂O₆).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	A	1	Total	C	H	N	O	0	0
			19	9	2	2	6		
7	B	1	Total	C	H	N	O	0	0
			19	9	2	2	6		

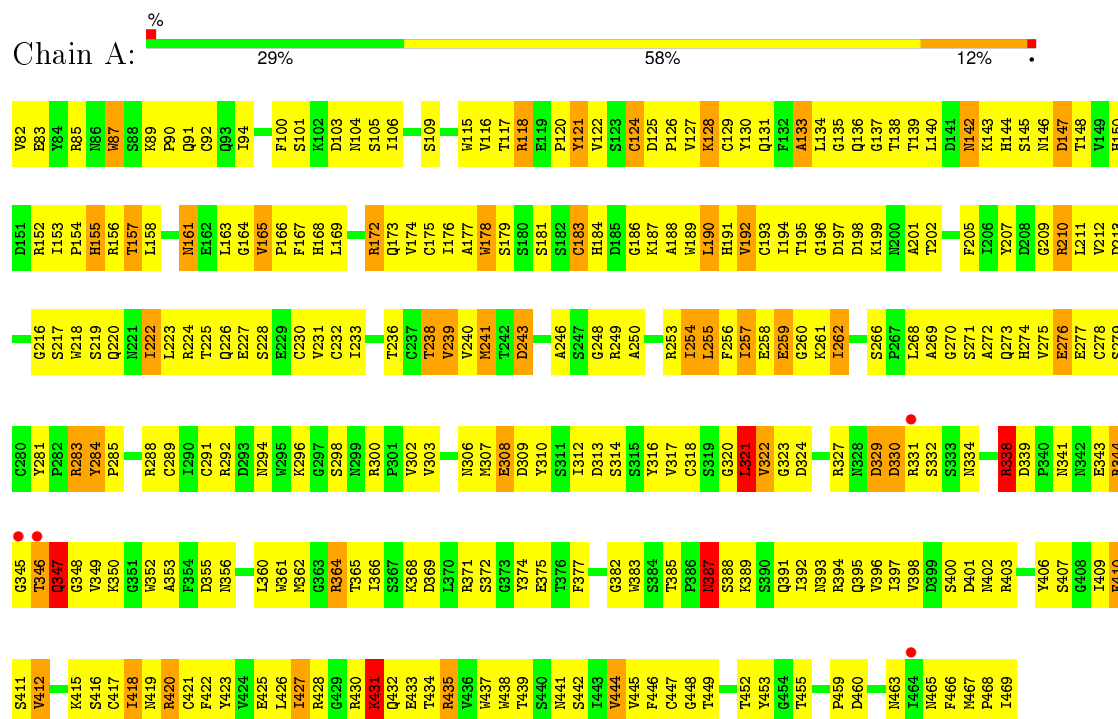
- Molecule 8 is water.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	121	Total	H	O	1	0
			363	242	121		
8	B	4	Total	H	O	0	0
			12	8	4		

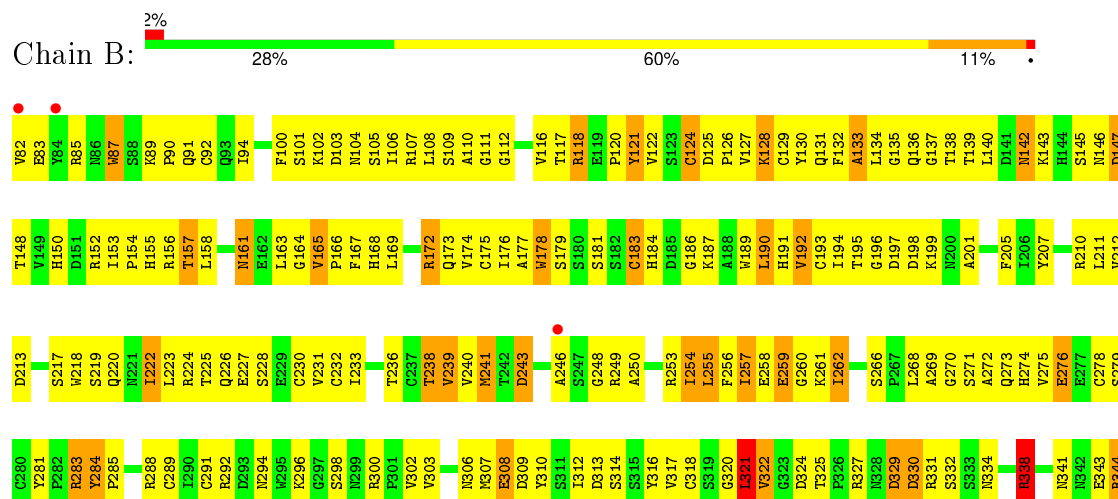
3 Residue-property plots

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: INFLUENZA A SUBTYPE N2 NEURAMINIDASE



• Molecule 1: INFLUENZA A SUBTYPE N2 NEURAMINIDASE





4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	120.42Å 139.83Å 140.07Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	5.00 – 1.90 25.93 – 1.76	Depositor EDS
% Data completeness (in resolution range)	(Not available) (5.00-1.90) 28.4 (25.93-1.76)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.71 (at 1.76Å)	Xtriage
Refinement program	X-PLOR	Depositor
R, R_{free}	0.225 , (Not available) 0.287 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	4.1	Xtriage
Anisotropy	3.656	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 65.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.24$	Xtriage
Outliers	0 of 33261 reflections	Xtriage
F_o, F_c correlation	0.72	EDS
Total number of atoms	8595	wwPDB-VP
Average B, all atoms (Å ²)	7.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.53% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: BMA, NAG, ST1, CA, FUC, FUL, MAN

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.64	0/3092	0.94	4/4194 (0.1%)
1	B	0.64	0/3092	0.94	4/4194 (0.1%)
All	All	0.64	0/6184	0.94	8/8388 (0.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
5	B	1	0

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	243	ASP	N-CA-C	-5.70	95.60	111.00
1	B	243	ASP	N-CA-C	-5.70	95.60	111.00
1	A	338	ARG	NE-CZ-NH1	5.12	122.86	120.30
1	B	338	ARG	NE-CZ-NH1	5.12	122.86	120.30
1	A	444	VAL	N-CA-C	-5.06	97.35	111.00
1	B	444	VAL	N-CA-C	-5.06	97.35	111.00
1	A	321	LEU	CA-CB-CG	-5.01	103.77	115.30
1	B	321	LEU	CA-CB-CG	-5.01	103.77	115.30

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
5	B	477	FUC	C1

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3022	723	2850	335	5
1	B	3022	723	2850	352	4
2	A	56	54	50	9	0
2	B	56	54	50	8	0
3	A	49	47	42	2	4
4	A	72	67	60	26	5
4	B	72	67	60	3	4
5	B	49	47	42	0	0
6	A	1	0	0	0	0
6	B	1	0	0	0	0
7	A	17	2	7	1	0
7	B	17	2	7	1	0
8	A	121	242	0	27	2
8	B	4	8	0	5	0
All	All	6559	2036	6018	678	12

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 55.

All (678) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:478:NAG:H61	1:B:455:THR:CG2	1.57	1.35
1:B:107:ARG:CZ	8:B:497:HOH:O	1.78	1.31
8:A:515:HOH:O	1:B:100:PHE:CZ	1.83	1.31
4:A:478:NAG:C6	1:B:455:THR:CG2	2.15	1.24
1:B:205:PHE:HD2	1:B:257:ILE:HD12	1.08	1.17
4:A:478:NAG:C6	1:B:455:THR:HG21	1.75	1.15
1:A:205:PHE:HD2	1:A:257:ILE:HD12	1.08	1.12
4:A:478:NAG:H61	1:B:455:THR:HG22	1.26	1.09
1:A:169:LEU:HD11	1:B:112:GLY:HA3	1.31	1.07
1:A:115:TRP:CZ2	1:B:108:LEU:HD11	1.90	1.05

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:205:PHE:CD2	1:A:257:ILE:HD12	1.95	1.01
1:A:142:ASN:ND2	1:B:110:ALA:HB3	1.74	1.01
1:B:205:PHE:CD2	1:B:257:ILE:HD12	1.95	1.00
1:A:177:ALA:HB2	1:A:193:CYS:HB3	1.42	0.99
1:B:177:ALA:HB2	1:B:193:CYS:HB3	1.42	0.99
1:A:136:GLN:HG3	1:A:148:THR:HG23	1.46	0.96
1:A:409:ILE:HD11	1:A:420:ARG:HB3	1.47	0.95
1:B:409:ILE:HD11	1:B:420:ARG:HB3	1.47	0.95
8:A:518:HOH:O	1:B:412:VAL:HG23	1.66	0.94
1:A:142:ASN:HD21	1:B:110:ALA:HB3	1.28	0.94
1:B:136:GLN:HG3	1:B:148:THR:HG23	1.46	0.94
1:A:281:TYR:HB3	8:A:525:HOH:O	1.66	0.93
4:A:478:NAG:O5	1:B:455:THR:HB	1.69	0.92
1:A:106:ILE:HG12	8:A:557:HOH:O	1.69	0.91
4:A:483:MAN:O2	1:B:396:VAL:CG1	2.18	0.90
1:A:202:THR:HB	1:B:454:GLY:H	1.35	0.90
1:A:136:GLN:O	1:A:138:THR:HG22	1.72	0.90
1:B:136:GLN:O	1:B:138:THR:HG22	1.72	0.89
4:A:478:NAG:H61	1:B:455:THR:CB	2.03	0.88
1:B:257:ILE:HD11	1:B:260:GLY:HA2	1.55	0.87
1:A:257:ILE:HD11	1:A:260:GLY:HA2	1.55	0.86
1:B:272:ALA:HA	1:B:316:TYR:CE1	2.11	0.86
1:A:272:ALA:HA	1:A:316:TYR:CE1	2.11	0.86
1:A:195:THR:HG21	1:B:458:TRP:HE1	1.43	0.83
1:A:169:LEU:CD1	1:B:112:GLY:HA3	2.09	0.83
4:A:483:MAN:O2	1:B:396:VAL:HG11	1.79	0.82
4:A:478:NAG:H5	1:B:455:THR:HG21	1.60	0.80
1:B:240:VAL:HG12	1:B:254:ILE:HG23	1.63	0.80
1:A:240:VAL:HG12	1:A:254:ILE:HG23	1.63	0.79
1:B:249:ARG:HG3	1:B:250:ALA:H	1.47	0.79
1:B:331:ARG:HA	1:B:331:ARG:NH1	1.99	0.78
1:A:331:ARG:HH11	1:A:331:ARG:HA	1.48	0.78
1:B:331:ARG:HA	1:B:331:ARG:HH11	1.48	0.77
4:A:478:NAG:C5	1:B:455:THR:HG21	2.13	0.77
1:A:331:ARG:NH1	1:A:331:ARG:HA	1.99	0.77
1:A:249:ARG:HG3	1:A:250:ALA:H	1.47	0.77
1:A:138:THR:HA	1:B:108:LEU:HD23	1.64	0.77
1:A:101:SER:HA	8:A:500:HOH:O	1.84	0.76
1:A:272:ALA:HA	1:A:316:TYR:HE1	1.48	0.76
1:B:452:THR:HG22	1:B:453:TYR:H	1.51	0.76
1:A:318:CYS:HB3	1:A:385:THR:O	1.86	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:452:THR:HG22	1:A:453:TYR:H	1.51	0.76
1:A:296:LYS:O	1:A:345:GLY:HA3	1.86	0.76
1:B:318:CYS:HB3	1:B:385:THR:O	1.86	0.75
1:B:272:ALA:HA	1:B:316:TYR:HE1	1.48	0.74
4:A:478:NAG:C6	1:B:455:THR:HG22	2.00	0.74
1:A:115:TRP:HZ2	1:B:108:LEU:HD11	1.51	0.73
1:B:438:TRP:HD1	1:B:469:ILE:HG13	1.53	0.73
1:A:158:LEU:O	1:A:174:VAL:HG12	1.88	0.73
1:B:296:LYS:O	1:B:345:GLY:HA3	1.86	0.73
1:B:271:SER:HB3	1:B:338:ARG:O	1.89	0.73
1:B:158:LEU:O	1:B:174:VAL:HG12	1.88	0.73
1:A:283:ARG:HE	1:A:288:ARG:HE	1.36	0.72
1:A:438:TRP:HD1	1:A:469:ILE:HG13	1.53	0.72
1:B:283:ARG:HE	1:B:288:ARG:HE	1.36	0.72
1:A:271:SER:HB3	1:A:338:ARG:O	1.89	0.72
4:A:478:NAG:C5	1:B:455:THR:CB	2.68	0.71
1:B:338:ARG:HB2	1:B:338:ARG:HH11	1.55	0.71
1:B:257:ILE:HD11	1:B:260:GLY:CA	2.19	0.71
1:A:257:ILE:HD11	1:A:260:GLY:CA	2.19	0.71
1:A:268:LEU:HD12	1:A:312:ILE:HG21	1.73	0.71
1:B:438:TRP:CD1	1:B:469:ILE:HG13	2.26	0.70
1:B:183:CYS:HB2	1:B:232:CYS:SG	2.31	0.70
1:A:183:CYS:HB2	1:A:232:CYS:SG	2.31	0.70
1:A:438:TRP:CD1	1:A:469:ILE:HG13	2.26	0.70
1:A:338:ARG:HB2	1:A:338:ARG:HH11	1.55	0.70
1:B:157:THR:HG22	1:B:176:ILE:HA	1.73	0.70
1:A:394:ARG:HG3	1:A:395:GLN:N	2.06	0.70
1:B:125:ASP:HB3	1:B:184:HIS:ND1	2.06	0.70
1:A:125:ASP:HB3	1:A:184:HIS:ND1	2.06	0.70
8:A:501:HOH:O	1:B:108:LEU:HD22	1.91	0.70
1:A:430:ARG:HD3	1:A:437:TRP:HA	1.72	0.70
1:B:394:ARG:HG3	1:B:395:GLN:N	2.06	0.70
1:B:268:LEU:HD12	1:B:312:ILE:HG21	1.73	0.70
1:A:374:TYR:HB3	1:A:398:VAL:HG23	1.74	0.69
1:A:289:CYS:HB2	1:A:303:VAL:HB	1.74	0.69
1:B:321:LEU:O	1:B:322:VAL:HB	1.93	0.69
1:B:430:ARG:HG3	1:B:431:LYS:H	1.57	0.69
1:B:430:ARG:HD3	1:B:437:TRP:HA	1.72	0.69
1:A:157:THR:HG22	1:A:176:ILE:HA	1.73	0.69
1:B:87:TRP:HD1	1:B:233:ILE:HG12	1.58	0.69
1:B:217:SER:HA	1:B:223:LEU:HD22	1.75	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:430:ARG:HG3	1:A:431:LYS:H	1.57	0.69
1:A:155:HIS:HB3	1:B:104:ASN:ND2	2.06	0.69
1:A:87:TRP:HD1	1:A:233:ILE:HG12	1.58	0.68
1:B:289:CYS:HB2	1:B:303:VAL:HB	1.74	0.68
4:A:478:NAG:C5	1:B:455:THR:HB	2.22	0.68
1:B:419:ASN:HB3	8:B:517:HOH:O	1.92	0.68
4:A:478:NAG:C1	1:B:455:THR:OG1	2.41	0.68
1:A:134:LEU:HB2	1:A:156:ARG:NH2	2.08	0.68
1:B:134:LEU:HB2	1:B:156:ARG:NH2	2.08	0.68
1:A:321:LEU:O	1:A:322:VAL:HB	1.93	0.68
1:A:216:GLY:H	1:B:452:THR:HB	1.59	0.68
1:A:116:VAL:HG11	1:A:148:THR:HG21	1.76	0.68
1:A:103:ASP:HB3	8:A:502:HOH:O	1.94	0.68
1:B:116:VAL:HG11	1:B:148:THR:HG21	1.76	0.67
1:B:374:TYR:HB3	1:B:398:VAL:HG23	1.74	0.67
1:A:321:LEU:HD13	1:A:377:PHE:CE1	2.30	0.67
1:B:396:VAL:O	1:B:396:VAL:HG13	1.95	0.67
1:A:116:VAL:CG1	1:A:148:THR:HG21	2.24	0.67
1:A:396:VAL:O	1:A:396:VAL:HG13	1.95	0.67
1:A:217:SER:HA	1:A:223:LEU:HD22	1.75	0.67
1:B:238:THR:OG1	1:B:307:MET:SD	2.52	0.67
4:A:478:NAG:O5	1:B:455:THR:CB	2.41	0.67
1:A:238:THR:OG1	1:A:307:MET:SD	2.52	0.67
1:B:116:VAL:CG1	1:B:148:THR:HG21	2.24	0.66
1:A:144:HIS:CD2	1:B:466:PHE:HD2	2.13	0.66
1:B:300:ARG:O	1:B:317:VAL:HG23	1.95	0.66
1:A:249:ARG:HG3	1:A:250:ALA:N	2.11	0.66
1:B:321:LEU:HD13	1:B:377:PHE:CE1	2.30	0.66
4:A:478:NAG:O6	1:B:455:THR:HG21	1.93	0.66
1:A:300:ARG:O	1:A:317:VAL:HG23	1.95	0.66
1:B:226:GLN:HE21	1:B:240:VAL:H	1.44	0.66
4:A:478:NAG:H61	1:B:455:THR:HB	1.77	0.65
1:A:412:VAL:HG11	1:A:421:CYS:SG	2.36	0.65
1:B:412:VAL:HG11	1:B:421:CYS:SG	2.36	0.65
1:B:249:ARG:HG3	1:B:250:ALA:N	2.11	0.65
1:B:312:ILE:HG22	1:B:313:ASP:H	1.62	0.65
1:A:419:ASN:HD21	1:A:448:GLY:HA3	1.61	0.65
1:B:419:ASN:HD21	1:B:448:GLY:HA3	1.61	0.65
1:A:312:ILE:HG22	1:A:313:ASP:H	1.62	0.64
1:A:134:LEU:HD11	8:A:521:HOH:O	1.98	0.64
1:B:228:SER:HB2	1:B:350:LYS:NZ	2.13	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:366:ILE:HG21	1:A:400:SER:HB3	1.80	0.64
1:B:366:ILE:HG21	1:B:400:SER:HB3	1.80	0.64
1:A:228:SER:HB2	1:A:350:LYS:NZ	2.13	0.63
1:A:226:GLN:HE21	1:A:240:VAL:H	1.44	0.63
1:B:334:ASN:HA	1:B:387:ASN:HD21	1.63	0.63
1:B:219:SER:O	1:B:220:GLN:HB2	1.99	0.63
1:A:115:TRP:CZ2	1:B:108:LEU:CD1	2.76	0.63
1:A:101:SER:HB2	1:A:445:VAL:HG22	1.80	0.63
1:B:212:VAL:HB	1:B:260:GLY:O	1.98	0.63
1:A:212:VAL:HB	1:A:260:GLY:O	1.98	0.63
1:B:101:SER:HB2	1:B:445:VAL:HG22	1.80	0.63
1:A:430:ARG:HD2	1:A:434:THR:HA	1.81	0.63
1:B:320:GLY:H	1:B:388:SER:H	1.47	0.63
1:A:177:ALA:CB	1:A:193:CYS:HB3	2.24	0.63
1:B:452:THR:HG22	1:B:453:TYR:N	2.14	0.63
1:A:146:ASN:OD1	1:A:437:TRP:HB3	1.99	0.63
4:A:478:NAG:C5	1:B:455:THR:CG2	2.74	0.63
1:B:372:SER:HA	1:B:403:ARG:HA	1.81	0.63
1:B:430:ARG:HD2	1:B:434:THR:HA	1.81	0.63
1:A:372:SER:HA	1:A:403:ARG:HA	1.81	0.63
1:B:146:ASN:OD1	1:B:437:TRP:HB3	1.99	0.62
1:A:409:ILE:HD12	1:A:421:CYS:O	1.99	0.62
1:A:320:GLY:N	1:A:388:SER:H	1.97	0.62
1:A:412:VAL:CG1	1:A:421:CYS:SG	2.87	0.62
1:A:190:LEU:HD11	1:A:257:ILE:HD13	1.82	0.62
1:B:412:VAL:CG1	1:B:421:CYS:SG	2.87	0.62
1:A:317:VAL:HB	8:A:543:HOH:O	1.98	0.62
4:A:478:NAG:C6	1:B:455:THR:CB	2.71	0.62
1:B:320:GLY:N	1:B:388:SER:H	1.97	0.62
1:B:85:ARG:HH12	1:B:125:ASP:HA	1.65	0.62
1:A:334:ASN:HA	1:A:387:ASN:HD21	1.63	0.62
1:A:117:THR:HG21	1:A:167:PHE:HD2	1.65	0.62
1:B:409:ILE:HD12	1:B:421:CYS:O	1.99	0.62
1:A:85:ARG:HH12	1:A:125:ASP:HA	1.65	0.62
1:A:219:SER:O	1:A:220:GLN:HB2	1.99	0.62
1:B:117:THR:HG21	1:B:167:PHE:HD2	1.65	0.62
1:A:118:ARG:HA	1:A:441:ASN:OD1	1.99	0.61
1:B:196:GLY:HA3	1:B:201:ALA:HA	1.82	0.61
1:B:118:ARG:HA	1:B:441:ASN:OD1	1.99	0.61
1:A:320:GLY:H	1:A:388:SER:H	1.47	0.61
1:A:452:THR:HG22	1:A:453:TYR:N	2.14	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:146:ASN:O	1:A:147:ASP:HB2	2.01	0.61
1:B:107:ARG:NH2	8:B:497:HOH:O	2.04	0.61
1:B:275:VAL:HG13	1:B:278:CYS:SG	2.41	0.61
1:A:283:ARG:CZ	1:A:288:ARG:HH21	2.14	0.61
1:A:196:GLY:HA3	1:A:201:ALA:HA	1.82	0.61
1:B:283:ARG:CZ	1:B:288:ARG:HH21	2.14	0.61
1:B:302:VAL:HG21	1:B:383:TRP:CZ3	2.36	0.61
1:A:302:VAL:HG21	1:A:383:TRP:CZ3	2.36	0.60
1:B:146:ASN:O	1:B:147:ASP:HB2	2.01	0.60
1:B:190:LEU:HD11	1:B:257:ILE:HD13	1.82	0.60
1:A:463:ASN:HB3	1:A:466:PHE:CG	2.36	0.60
1:B:246:ALA:O	1:B:274:HIS:NE2	2.34	0.60
1:A:275:VAL:HG13	1:A:278:CYS:SG	2.41	0.60
1:B:329:ASP:HA	1:B:368:LYS:HD3	1.84	0.60
1:A:138:THR:HA	1:B:108:LEU:CD2	2.32	0.60
1:A:246:ALA:O	1:A:274:HIS:NE2	2.34	0.59
1:B:463:ASN:HB3	1:B:466:PHE:CG	2.36	0.59
1:A:312:ILE:HG22	1:A:313:ASP:N	2.17	0.59
1:A:172:ARG:HB3	1:B:165:VAL:HG12	1.83	0.59
1:B:257:ILE:O	1:B:257:ILE:HG12	2.03	0.59
1:B:402:ASN:ND2	1:B:459:PRO:HB3	2.18	0.59
1:A:329:ASP:HA	1:A:368:LYS:HD3	1.84	0.59
1:B:352:TRP:HD1	1:B:407:SER:HG	1.50	0.58
1:A:402:ASN:ND2	1:A:459:PRO:HB3	2.18	0.58
1:B:312:ILE:HG22	1:B:313:ASP:N	2.17	0.58
1:B:291:CYS:O	1:B:300:ARG:HD3	2.03	0.58
1:A:291:CYS:O	1:A:300:ARG:HD3	2.03	0.58
4:A:483:MAN:O2	1:B:396:VAL:HG12	2.01	0.58
1:B:116:VAL:O	1:B:135:GLY:HA2	2.02	0.58
1:A:116:VAL:O	1:A:135:GLY:HA2	2.02	0.58
1:A:258:GLU:O	1:A:259:GLU:HB2	2.03	0.58
1:A:139:THR:O	1:A:142:ASN:HB2	2.04	0.58
1:B:139:THR:O	1:B:142:ASN:HB2	2.04	0.58
1:B:327:ARG:HG2	1:B:327:ARG:HH11	1.68	0.58
2:A:472:NAG:C3	2:A:473:NAG:O5	2.51	0.58
1:B:258:GLU:O	1:B:259:GLU:HB2	2.03	0.58
1:B:177:ALA:CB	1:B:193:CYS:HB3	2.24	0.58
1:A:431:LYS:H	1:A:434:THR:HG22	1.69	0.57
4:A:478:NAG:O6	1:B:455:THR:CG2	2.50	0.57
1:A:218:TRP:HH2	1:A:241:MET:HB2	1.69	0.57
1:B:218:TRP:HH2	1:B:241:MET:HB2	1.69	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:283:ARG:NE	1:B:288:ARG:HH21	2.03	0.57
1:B:131:GLN:NE2	1:B:164:GLY:H	2.02	0.57
1:B:431:LYS:H	1:B:434:THR:HG22	1.69	0.57
1:B:91:GLN:HG3	1:B:420:ARG:NH1	2.20	0.57
1:A:131:GLN:NE2	1:A:164:GLY:H	2.02	0.57
1:A:327:ARG:HG2	1:A:327:ARG:HH11	1.68	0.57
1:A:283:ARG:NE	1:A:288:ARG:HH21	2.03	0.57
1:A:257:ILE:HG12	1:A:257:ILE:O	2.03	0.57
1:A:142:ASN:ND2	1:B:110:ALA:CB	2.60	0.56
1:A:341:ASN:OD1	1:A:343:GLU:HG3	2.05	0.56
1:A:91:GLN:HG3	1:A:420:ARG:NH1	2.20	0.56
1:B:430:ARG:HG3	1:B:431:LYS:N	2.21	0.56
1:A:155:HIS:HE2	1:B:461:GLY:HA3	1.70	0.56
1:B:324:ASP:O	1:B:327:ARG:HD3	2.05	0.56
1:A:430:ARG:HG3	1:A:431:LYS:N	2.21	0.56
1:B:341:ASN:OD1	1:B:343:GLU:HG3	2.05	0.56
2:B:472:NAG:C3	2:B:473:NAG:O5	2.51	0.56
4:A:478:NAG:C6	1:B:455:THR:HB	2.34	0.56
1:A:279:SER:HB3	1:A:409:ILE:HG22	1.87	0.56
1:B:279:SER:HB3	1:B:409:ILE:HG22	1.87	0.56
1:B:136:GLN:HA	1:B:136:GLN:OE1	2.06	0.55
1:A:410:PHE:CD1	1:A:410:PHE:C	2.80	0.55
1:B:147:ASP:O	1:B:150:HIS:HD2	1.90	0.55
1:A:197:ASP:O	1:A:201:ALA:HB2	2.07	0.55
1:A:324:ASP:O	1:A:327:ARG:HD3	2.05	0.55
2:B:472:NAG:N2	2:B:485:NAG:H82	2.22	0.55
1:B:338:ARG:CB	1:B:338:ARG:HH11	2.20	0.55
1:A:150:HIS:HB3	8:A:585:HOH:O	2.05	0.55
1:B:318:CYS:HB3	1:B:385:THR:C	2.28	0.55
1:A:276:GLU:O	1:A:292:ARG:HB3	2.07	0.55
1:B:276:GLU:O	1:B:292:ARG:HB3	2.07	0.55
1:B:197:ASP:O	1:B:201:ALA:HB2	2.07	0.54
1:B:288:ARG:NH1	1:B:383:TRP:CZ2	2.75	0.54
1:B:395:GLN:HA	8:B:511:HOH:O	2.06	0.54
1:A:136:GLN:HA	1:A:136:GLN:OE1	2.06	0.54
1:B:410:PHE:C	1:B:410:PHE:CD1	2.80	0.54
1:A:155:HIS:HB3	1:B:104:ASN:HD21	1.70	0.54
1:A:168:HIS:CD2	1:A:168:HIS:H	2.25	0.54
1:A:318:CYS:HB3	1:A:385:THR:C	2.28	0.54
1:B:228:SER:HB2	1:B:350:LYS:CE	2.38	0.54
1:A:228:SER:HB2	1:A:350:LYS:CE	2.38	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:228:SER:HB2	1:A:350:LYS:HE2	1.90	0.54
1:A:147:ASP:O	1:A:150:HIS:HD2	1.90	0.54
1:B:346:THR:O	1:B:347:GLN:O	2.26	0.54
1:A:143:LYS:NZ	1:B:466:PHE:HA	2.23	0.53
2:A:472:NAG:N2	2:A:485:NAG:H82	2.22	0.53
1:B:168:HIS:CD2	1:B:168:HIS:H	2.25	0.53
1:A:213:ASP:OD2	1:A:262:ILE:HG13	2.08	0.53
1:A:288:ARG:NH1	1:A:383:TRP:CZ2	2.75	0.53
1:A:411:SER:HB3	1:A:418:ILE:HD11	1.90	0.53
1:B:468:PRO:O	1:B:469:ILE:HB	2.08	0.53
1:B:199:LYS:HA	1:B:220:GLN:O	2.09	0.53
1:A:199:LYS:HA	1:A:220:GLN:O	2.09	0.53
1:B:213:ASP:OD2	1:B:262:ILE:HG13	2.08	0.53
1:A:468:PRO:O	1:A:469:ILE:HB	2.08	0.53
1:B:87:TRP:CD1	1:B:233:ILE:HG12	2.43	0.53
1:A:198:ASP:HB3	1:A:222:ILE:HG12	1.91	0.53
1:A:152:ARG:HD3	8:A:578:HOH:O	2.07	0.53
1:A:177:ALA:HB2	1:A:193:CYS:CB	2.28	0.53
1:A:338:ARG:CB	1:A:338:ARG:HH11	2.20	0.53
1:A:409:ILE:HG13	1:A:410:PHE:N	2.23	0.53
1:B:409:ILE:HG13	1:B:410:PHE:N	2.23	0.53
1:A:463:ASN:OD1	1:A:465:ASN:HB2	2.09	0.53
1:B:411:SER:HB3	1:B:418:ILE:HD11	1.90	0.53
1:B:463:ASN:OD1	1:B:465:ASN:HB2	2.09	0.53
1:A:176:ILE:O	1:A:177:ALA:HB2	2.09	0.52
1:A:346:THR:O	1:A:347:GLN:O	2.26	0.52
1:A:212:VAL:HG12	1:A:261:LYS:HE2	1.91	0.52
1:B:106:ILE:O	1:B:109:SER:HB2	2.10	0.52
1:B:228:SER:HB2	1:B:350:LYS:HE2	1.90	0.52
1:A:155:HIS:NE2	1:B:461:GLY:HA3	2.24	0.52
1:A:82:VAL:O	1:A:187:LYS:HE2	2.10	0.52
1:B:107:ARG:NH1	8:B:497:HOH:O	2.15	0.52
1:A:106:ILE:O	1:A:109:SER:HB2	2.10	0.52
1:A:421:CYS:HA	1:A:447:CYS:HA	1.92	0.52
1:A:281:TYR:OH	1:A:288:ARG:HD2	2.10	0.52
1:A:402:ASN:HD22	1:A:459:PRO:HB3	1.75	0.52
1:A:209:GLY:O	1:B:100:PHE:HZ	1.93	0.51
1:A:329:ASP:OD1	1:A:330:ASP:N	2.43	0.51
1:B:177:ALA:HB2	1:B:193:CYS:CB	2.28	0.51
1:B:329:ASP:OD1	1:B:330:ASP:N	2.43	0.51
4:B:479:NAG:H5	4:B:480:BMA:O2	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:134:LEU:HB2	1:B:156:ARG:HH21	1.75	0.51
1:B:281:TYR:OH	1:B:288:ARG:HD2	2.10	0.51
1:B:128:LYS:HB3	1:B:130:TYR:CZ	2.45	0.51
1:B:388:SER:C	1:B:389:LYS:HG2	2.30	0.51
1:B:82:VAL:O	1:B:187:LYS:HE2	2.10	0.51
1:B:402:ASN:HD22	1:B:459:PRO:HB3	1.75	0.51
4:A:479:NAG:H5	4:A:480:BMA:O2	2.10	0.51
1:A:388:SER:C	1:A:389:LYS:HG2	2.30	0.51
1:B:198:ASP:HB3	1:B:222:ILE:HG12	1.91	0.51
1:B:212:VAL:HG12	1:B:261:LYS:HE2	1.91	0.51
1:B:120:PRO:HG2	1:B:423:TYR:CE2	2.46	0.51
1:A:134:LEU:HB2	1:A:156:ARG:HH21	1.75	0.51
1:A:128:LYS:HB3	1:A:130:TYR:CZ	2.45	0.51
1:A:87:TRP:CD1	1:A:233:ILE:HG12	2.43	0.51
8:A:515:HOH:O	1:B:100:PHE:CE2	2.36	0.50
1:B:176:ILE:O	1:B:177:ALA:HB2	2.09	0.50
1:A:129:CYS:O	1:A:163:LEU:N	2.45	0.50
1:A:320:GLY:CA	1:A:387:ASN:HD22	2.24	0.50
1:B:421:CYS:HA	1:B:447:CYS:HA	1.92	0.50
1:A:142:ASN:OD1	1:B:111:GLY:HA3	2.12	0.50
1:A:120:PRO:HG2	1:A:423:TYR:CE2	2.46	0.50
1:B:172:ARG:HG3	1:B:172:ARG:HH11	1.77	0.50
8:A:515:HOH:O	1:B:100:PHE:HZ	1.50	0.50
1:A:101:SER:HB2	1:A:445:VAL:CG2	2.41	0.50
1:B:352:TRP:NE1	1:B:374:TYR:OH	2.45	0.50
1:B:101:SER:HB2	1:B:445:VAL:CG2	2.41	0.50
1:B:227:GLU:O	1:B:350:LYS:HE2	2.12	0.50
1:A:176:ILE:HD11	1:B:102:LYS:HB2	1.93	0.50
1:B:320:GLY:CA	1:B:387:ASN:HD22	2.24	0.50
1:B:117:THR:HG21	1:B:167:PHE:CD2	2.46	0.50
1:A:199:LYS:O	1:A:220:GLN:HA	2.12	0.50
2:B:484:NAG:O6	2:B:485:NAG:N2	2.45	0.50
1:B:199:LYS:O	1:B:220:GLN:HA	2.12	0.50
1:A:403:ARG:HG2	8:A:550:HOH:O	2.12	0.50
1:B:270:GLY:HA3	1:B:314:SER:OG	2.12	0.50
1:B:423:TYR:HB2	1:B:445:VAL:HG12	1.94	0.49
1:B:419:ASN:ND2	1:B:448:GLY:HA3	2.26	0.49
1:A:125:ASP:HB2	1:A:126:PRO:CD	2.42	0.49
1:B:348:GLY:O	1:B:371:ARG:NH2	2.44	0.49
1:B:302:VAL:HG21	1:B:383:TRP:CH2	2.47	0.49
1:B:125:ASP:HB2	1:B:126:PRO:CD	2.42	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:227:GLU:O	1:A:350:LYS:HE2	2.12	0.49
1:B:419:ASN:ND2	1:B:420:ARG:N	2.61	0.49
1:B:134:LEU:CB	1:B:156:ARG:HH21	2.26	0.49
1:A:344:ARG:NH2	1:A:368:LYS:O	2.46	0.49
1:A:419:ASN:ND2	1:A:420:ARG:N	2.61	0.49
1:B:344:ARG:NH2	1:B:368:LYS:O	2.46	0.49
1:A:172:ARG:HG3	1:A:172:ARG:HH11	1.77	0.49
1:A:212:VAL:HB	1:A:260:GLY:C	2.33	0.49
1:A:423:TYR:HB2	1:A:445:VAL:HG12	1.94	0.49
1:A:302:VAL:HG21	1:A:383:TRP:CH2	2.47	0.49
1:A:419:ASN:ND2	1:A:448:GLY:HA3	2.26	0.49
1:A:202:THR:HB	1:B:454:GLY:N	2.16	0.49
1:A:270:GLY:HA3	1:A:314:SER:OG	2.12	0.49
1:B:425:GLU:HB2	1:B:441:ASN:HD22	1.78	0.49
1:A:406:TYR:HD2	8:A:488:HOH:O	1.95	0.49
1:B:320:GLY:HA3	1:B:387:ASN:HD22	1.78	0.49
1:A:348:GLY:O	1:A:371:ARG:NH2	2.44	0.49
1:A:320:GLY:HA3	1:A:387:ASN:HD22	1.78	0.49
2:A:484:NAG:O6	2:A:485:NAG:N2	2.45	0.49
1:B:212:VAL:HB	1:B:260:GLY:C	2.33	0.49
1:A:403:ARG:NH1	1:A:433:GLU:HG3	2.28	0.49
1:A:172:ARG:CB	1:B:165:VAL:HG12	2.42	0.49
1:B:419:ASN:ND2	1:B:420:ARG:O	2.45	0.49
1:A:419:ASN:ND2	1:A:420:ARG:O	2.45	0.49
1:A:425:GLU:HB2	1:A:441:ASN:HD22	1.78	0.48
1:B:327:ARG:HG2	1:B:327:ARG:NH1	2.28	0.48
1:B:422:PHE:CE1	1:B:446:PHE:HB2	2.47	0.48
1:A:140:LEU:HD23	1:A:468:PRO:HG2	1.95	0.48
1:A:327:ARG:NH1	1:A:327:ARG:HG2	2.28	0.48
1:A:124:CYS:HB3	1:A:129:CYS:HA	1.96	0.48
1:A:179:SER:HB3	1:A:194:ILE:HB	1.95	0.48
1:A:422:PHE:CE1	1:A:446:PHE:HB2	2.47	0.48
1:A:426:LEU:HD21	1:A:444:VAL:HG23	1.95	0.48
1:B:281:TYR:CZ	1:B:288:ARG:HD2	2.49	0.48
1:B:85:ARG:NH1	1:B:184:HIS:HB3	2.29	0.48
8:A:501:HOH:O	1:B:108:LEU:HD13	2.14	0.48
1:B:129:CYS:SG	1:B:410:PHE:HZ	2.37	0.48
1:A:430:ARG:O	1:A:431:LYS:HB2	2.14	0.48
1:A:134:LEU:CB	1:A:156:ARG:HH21	2.26	0.48
1:A:128:LYS:HB3	1:A:130:TYR:CE1	2.49	0.48
1:A:441:ASN:C	8:A:557:HOH:O	2.51	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:124:CYS:CB	1:A:129:CYS:HA	2.43	0.48
1:B:192:VAL:HG11	1:B:239:VAL:HG21	1.96	0.48
1:A:138:THR:OG1	1:A:139:THR:N	2.44	0.48
1:A:144:HIS:HE1	1:B:462:ALA:HB1	1.77	0.48
1:A:352:TRP:NE1	1:A:374:TYR:OH	2.45	0.48
1:B:124:CYS:CB	1:B:129:CYS:HA	2.43	0.48
1:B:139:THR:HG22	1:B:140:LEU:N	2.28	0.48
1:B:410:PHE:O	1:B:421:CYS:N	2.43	0.47
1:B:128:LYS:HB3	1:B:130:TYR:CE1	2.49	0.47
1:B:430:ARG:O	1:B:431:LYS:HB2	2.14	0.47
1:B:165:VAL:HG23	1:B:166:PRO:O	2.14	0.47
2:B:472:NAG:H3	2:B:473:NAG:O5	2.14	0.47
1:B:140:LEU:HD23	1:B:468:PRO:HG2	1.95	0.47
1:A:239:VAL:O	1:A:239:VAL:HG13	2.13	0.47
1:A:85:ARG:NH1	1:A:184:HIS:HB3	2.29	0.47
1:B:239:VAL:O	1:B:239:VAL:HG13	2.13	0.47
1:B:407:SER:HB2	1:B:423:TYR:O	2.13	0.47
1:B:129:CYS:O	1:B:163:LEU:N	2.45	0.47
1:A:100:PHE:HB3	1:A:445:VAL:HG23	1.97	0.47
1:B:179:SER:HB3	1:B:194:ILE:HB	1.95	0.47
1:A:407:SER:HB2	1:A:423:TYR:O	2.13	0.47
1:B:153:ILE:HG13	1:B:153:ILE:O	2.14	0.47
1:B:125:ASP:HB3	1:B:184:HIS:CE1	2.49	0.47
1:A:431:LYS:NZ	1:A:431:LYS:HA	2.29	0.47
1:B:431:LYS:NZ	1:B:431:LYS:HA	2.29	0.47
2:A:472:NAG:H3	2:A:473:NAG:O5	2.14	0.47
1:A:165:VAL:HG23	1:A:166:PRO:O	2.14	0.47
1:A:189:TRP:O	1:A:207:TYR:HA	2.14	0.47
1:A:281:TYR:CZ	1:A:288:ARG:HD2	2.49	0.47
1:B:249:ARG:CG	1:B:250:ALA:N	2.78	0.47
1:B:426:LEU:HD21	1:B:444:VAL:HG23	1.95	0.47
1:A:139:THR:HG22	1:A:140:LEU:N	2.28	0.47
1:A:115:TRP:CH2	1:B:108:LEU:HD11	2.47	0.47
1:A:177:ALA:HA	1:A:195:THR:OG1	2.15	0.47
1:A:129:CYS:SG	1:A:410:PHE:HZ	2.37	0.47
1:A:410:PHE:O	1:A:421:CYS:N	2.43	0.47
1:A:268:LEU:HG	1:A:269:ALA:N	2.30	0.47
1:B:268:LEU:HG	1:B:269:ALA:N	2.30	0.47
1:B:403:ARG:NH1	1:B:433:GLU:HG3	2.28	0.47
1:A:320:GLY:H	1:A:388:SER:N	2.11	0.47
1:A:255:LEU:HD23	1:A:255:LEU:N	2.30	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:120:PRO:HG3	1:B:425:GLU:HB3	1.96	0.47
1:A:120:PRO:HG3	1:A:425:GLU:HB3	1.96	0.47
1:B:426:LEU:HD11	1:B:444:VAL:CG2	2.45	0.47
1:A:188:ALA:HB1	8:A:510:HOH:O	2.14	0.47
1:A:153:ILE:HG13	1:A:153:ILE:O	2.14	0.47
1:B:138:THR:OG1	1:B:139:THR:N	2.44	0.47
1:A:218:TRP:CG	1:A:253:ARG:NH2	2.83	0.47
1:A:117:THR:HG21	1:A:167:PHE:CD2	2.46	0.47
1:B:218:TRP:CG	1:B:253:ARG:NH2	2.83	0.47
1:A:306:ASN:OD1	1:A:308:GLU:HB2	2.15	0.47
1:A:152:ARG:O	1:A:153:ILE:HG23	2.15	0.46
1:A:352:TRP:HD1	1:A:407:SER:HG	1.61	0.46
1:A:211:LEU:HD23	1:A:211:LEU:O	2.15	0.46
1:A:236:THR:HG23	1:A:258:GLU:HG2	1.98	0.46
1:A:284:TYR:OH	2:A:484:NAG:H3	2.15	0.46
1:B:211:LEU:HD23	1:B:211:LEU:O	2.15	0.46
1:B:100:PHE:HB3	1:B:445:VAL:HG23	1.97	0.46
4:A:483:MAN:O3	1:B:457:SER:HB3	2.15	0.46
1:A:283:ARG:NE	1:A:288:ARG:HE	2.11	0.46
1:B:124:CYS:HB3	1:B:129:CYS:HA	1.96	0.46
1:B:152:ARG:O	1:B:153:ILE:HG23	2.15	0.46
1:A:179:SER:OG	1:A:225:THR:HG22	2.15	0.46
1:B:189:TRP:O	1:B:207:TYR:HA	2.14	0.46
1:B:177:ALA:HA	1:B:195:THR:OG1	2.15	0.46
1:B:283:ARG:NE	1:B:288:ARG:HE	2.11	0.46
1:A:426:LEU:HD11	1:A:444:VAL:CG2	2.45	0.46
1:B:248:GLY:O	1:B:274:HIS:CE1	2.69	0.46
1:A:248:GLY:O	1:A:274:HIS:CE1	2.69	0.46
1:B:306:ASN:OD1	1:B:308:GLU:HB2	2.15	0.46
1:A:106:ILE:HG13	1:A:428:ARG:HG3	1.98	0.46
1:A:352:TRP:HE1	1:A:374:TYR:HH	1.63	0.46
1:A:91:GLN:HG2	1:A:92:CYS:N	2.31	0.46
1:A:125:ASP:HB3	1:A:184:HIS:CE1	2.49	0.46
1:B:255:LEU:HD23	1:B:255:LEU:N	2.30	0.46
1:B:120:PRO:CG	1:B:425:GLU:HB3	2.46	0.46
1:B:410:PHE:CD1	1:B:411:SER:O	2.69	0.46
1:B:284:TYR:OH	2:B:484:NAG:H3	2.15	0.46
1:A:198:ASP:HB3	1:A:222:ILE:CG1	2.46	0.46
1:B:426:LEU:HD13	1:B:460:ASP:N	2.31	0.46
1:A:120:PRO:CG	1:A:425:GLU:HB3	2.46	0.46
1:A:410:PHE:HD1	1:A:410:PHE:C	2.19	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:249:ARG:HA	8:A:526:HOH:O	2.15	0.46
1:A:183:CYS:CB	1:A:232:CYS:SG	3.00	0.46
1:B:133:ALA:HB3	1:B:167:PHE:HE2	1.80	0.46
1:B:91:GLN:HG2	1:B:92:CYS:N	2.31	0.46
1:B:432:GLN:O	1:B:434:THR:HG23	2.16	0.46
1:A:284:TYR:CE2	1:A:285:PRO:HG3	2.51	0.46
1:B:143:LYS:HE3	1:B:143:LYS:HB2	1.79	0.46
1:A:321:LEU:O	1:A:322:VAL:CB	2.64	0.45
1:B:284:TYR:CE2	1:B:285:PRO:HG3	2.51	0.45
1:A:426:LEU:HD13	1:A:460:ASP:N	2.31	0.45
1:B:183:CYS:CB	1:B:232:CYS:SG	3.00	0.45
1:A:192:VAL:HG11	1:A:239:VAL:HG21	1.96	0.45
1:A:432:GLN:O	1:A:434:THR:HG23	2.16	0.45
1:A:392:ILE:HG12	1:A:393:ASN:N	2.32	0.45
1:A:115:TRP:CH2	1:B:108:LEU:HD21	2.52	0.45
1:B:179:SER:OG	1:B:225:THR:HG22	2.15	0.45
1:B:392:ILE:HG12	1:B:393:ASN:N	2.32	0.45
1:B:106:ILE:HG13	1:B:428:ARG:HG3	1.98	0.45
1:B:288:ARG:HH11	1:B:288:ARG:HG3	1.82	0.45
1:A:133:ALA:HB3	1:A:167:PHE:HE2	1.80	0.45
2:A:473:NAG:C6	2:A:473:NAG:H2	2.46	0.45
2:B:473:NAG:H2	2:B:473:NAG:C6	2.46	0.45
1:B:198:ASP:HB3	1:B:222:ILE:CG1	2.46	0.45
1:A:410:PHE:CD1	1:A:411:SER:O	2.69	0.45
1:A:87:TRP:CZ3	1:A:418:ILE:HD13	2.52	0.45
1:B:410:PHE:C	1:B:410:PHE:HD1	2.19	0.45
1:B:87:TRP:CZ3	1:B:418:ILE:HD13	2.52	0.45
1:A:144:HIS:CD2	1:B:466:PHE:CD2	3.00	0.45
1:A:387:ASN:HD22	1:A:387:ASN:HA	1.61	0.45
1:B:236:THR:HG23	1:B:258:GLU:HG2	1.98	0.45
1:B:181:SER:HB3	1:B:192:VAL:HG13	1.98	0.45
1:B:321:LEU:O	1:B:322:VAL:CB	2.64	0.45
1:B:120:PRO:O	1:B:121:TYR:HB3	2.17	0.45
1:A:210:ARG:NH2	8:A:518:HOH:O	2.49	0.45
1:B:298:SER:O	1:B:322:VAL:HG13	2.17	0.45
1:A:298:SER:O	1:A:322:VAL:HG13	2.17	0.45
1:B:211:LEU:HD23	1:B:211:LEU:C	2.37	0.45
1:A:140:LEU:CD2	1:A:468:PRO:HG2	2.47	0.44
1:A:442:SER:HB2	1:A:460:ASP:OD1	2.17	0.44
1:A:178:TRP:CE2	1:A:195:THR:HA	2.53	0.44
1:A:278:CYS:HB3	1:A:289:CYS:HB3	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:320:GLY:H	1:B:388:SER:N	2.11	0.44
1:B:306:ASN:OD1	1:B:308:GLU:CB	2.65	0.44
1:A:190:LEU:CD1	1:A:257:ILE:HD13	2.46	0.44
1:B:139:THR:HB	1:B:142:ASN:OD1	2.17	0.44
1:B:184:HIS:HD2	1:B:186:GLY:H	1.64	0.44
2:B:472:NAG:HN2	2:B:485:NAG:H82	1.82	0.44
1:B:442:SER:HB2	1:B:460:ASP:OD1	2.17	0.44
1:A:139:THR:HB	1:A:142:ASN:OD1	2.17	0.44
1:A:240:VAL:HG21	1:A:278:CYS:SG	2.57	0.44
1:A:184:HIS:HD2	1:A:186:GLY:H	1.64	0.44
1:A:211:LEU:HD23	1:A:211:LEU:C	2.37	0.44
1:A:396:VAL:O	1:A:398:VAL:N	2.51	0.44
1:B:195:THR:HG22	1:B:196:GLY:N	2.33	0.44
1:B:181:SER:HB3	1:B:192:VAL:CG1	2.48	0.44
1:A:181:SER:HB3	1:A:192:VAL:HG13	1.98	0.44
1:A:90:PRO:O	1:A:417:CYS:HA	2.18	0.44
1:B:90:PRO:O	1:B:417:CYS:HA	2.18	0.44
1:A:288:ARG:HH11	1:A:288:ARG:HG3	1.82	0.44
1:B:278:CYS:HB3	1:B:289:CYS:HB3	1.99	0.44
1:A:181:SER:HB3	1:A:192:VAL:CG1	2.48	0.44
1:A:415:LYS:HG2	1:A:415:LYS:O	2.18	0.44
1:A:195:THR:HG22	1:A:196:GLY:N	2.33	0.44
1:B:178:TRP:CE2	1:B:195:THR:HA	2.53	0.44
1:B:240:VAL:HG21	1:B:278:CYS:SG	2.57	0.44
1:A:306:ASN:OD1	1:A:308:GLU:CB	2.65	0.44
1:B:415:LYS:HG2	1:B:415:LYS:O	2.18	0.44
1:B:199:LYS:HG2	1:B:220:GLN:O	2.18	0.44
1:B:218:TRP:NE1	1:B:243:ASP:HB3	2.33	0.44
1:A:143:LYS:HZ3	1:B:466:PHE:HA	1.82	0.43
1:B:140:LEU:CD2	1:B:468:PRO:HG2	2.47	0.43
1:A:249:ARG:CG	1:A:250:ALA:N	2.78	0.43
1:B:131:GLN:HE21	1:B:164:GLY:H	1.66	0.43
1:B:406:TYR:H	1:B:425:GLU:HG2	1.83	0.43
1:B:190:LEU:CG	1:B:257:ILE:HD13	2.48	0.43
1:A:406:TYR:H	1:A:425:GLU:HG2	1.83	0.43
1:B:106:ILE:HG23	1:B:467:MET:CE	2.48	0.43
1:A:106:ILE:HG23	1:A:467:MET:CE	2.48	0.43
1:A:120:PRO:O	1:A:121:TYR:HB3	2.17	0.43
2:A:472:NAG:HN2	2:A:485:NAG:H82	1.82	0.43
1:B:169:LEU:HA	1:B:169:LEU:HD23	1.79	0.43
1:B:153:ILE:HG12	1:B:156:ARG:HG3	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:183:CYS:N	1:B:230:CYS:SG	2.91	0.43
1:B:396:VAL:O	1:B:398:VAL:N	2.51	0.43
1:B:153:ILE:CG1	1:B:156:ARG:HG3	2.49	0.43
1:A:183:CYS:N	1:A:230:CYS:SG	2.91	0.43
3:A:475:NAG:O6	3:A:476:BMA:O2	2.33	0.43
1:B:350:LYS:HG2	1:B:407:SER:O	2.18	0.43
1:A:190:LEU:CG	1:A:257:ILE:HD13	2.48	0.43
1:A:105:SER:HB2	1:A:442:SER:HA	2.00	0.43
1:A:153:ILE:HG12	1:A:156:ARG:HG3	2.01	0.43
1:A:350:LYS:HG2	1:A:407:SER:O	2.18	0.43
1:B:127:VAL:HG23	1:B:128:LYS:N	2.34	0.43
1:B:324:ASP:H	1:B:327:ARG:HD3	1.84	0.43
1:A:324:ASP:H	1:A:327:ARG:HD3	1.84	0.43
1:A:365:THR:HG21	1:A:371:ARG:HA	2.01	0.43
1:A:153:ILE:CG1	1:A:156:ARG:HG3	2.49	0.43
1:A:131:GLN:HE21	1:A:164:GLY:H	1.66	0.43
1:B:300:ARG:HH22	1:B:349:VAL:HG13	1.84	0.43
1:B:463:ASN:O	1:B:467:MET:HG2	2.19	0.43
1:A:366:ILE:HG21	1:A:400:SER:CB	2.49	0.43
1:A:218:TRP:NE1	1:A:243:ASP:HB3	2.33	0.43
1:A:161:ASN:HA	1:A:161:ASN:HD22	1.62	0.43
1:B:309:ASP:O	1:B:310:TYR:HB2	2.19	0.43
1:B:109:SER:HB3	1:B:140:LEU:HD13	2.00	0.42
1:B:153:ILE:HB	1:B:154:PRO:HD2	2.00	0.42
1:B:327:ARG:NH1	1:B:368:LYS:HA	2.34	0.42
1:B:325:THR:HG1	1:B:365:THR:HG1	1.66	0.42
1:A:161:ASN:HB3	1:A:165:VAL:HG22	2.01	0.42
1:A:442:SER:N	8:A:557:HOH:O	2.53	0.42
1:A:332:SER:O	1:A:334:ASN:OD1	2.37	0.42
1:A:327:ARG:NH1	1:A:368:LYS:HA	2.34	0.42
1:A:292:ARG:CZ	1:A:294:ASN:HD21	2.32	0.42
1:B:292:ARG:CZ	1:B:294:ASN:HD21	2.32	0.42
1:B:365:THR:HG21	1:B:371:ARG:HA	2.01	0.42
1:A:153:ILE:HB	1:A:154:PRO:HD2	2.00	0.42
1:A:463:ASN:O	1:A:467:MET:HG2	2.19	0.42
1:A:144:HIS:CE1	1:B:463:ASN:H	2.37	0.42
1:A:199:LYS:HG2	1:A:220:GLN:O	2.18	0.42
1:A:409:ILE:HG12	8:A:525:HOH:O	2.19	0.42
1:B:128:LYS:HE2	1:B:128:LYS:HB2	1.82	0.42
1:B:427:ILE:HD11	1:B:439:THR:HG23	2.02	0.42
1:A:109:SER:HB3	1:A:140:LEU:HD13	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:302:VAL:HG21	1:A:383:TRP:HZ3	1.84	0.42
1:B:103:ASP:O	1:B:104:ASN:HB2	2.20	0.42
1:B:332:SER:O	1:B:334:ASN:OD1	2.37	0.42
1:B:161:ASN:HB3	1:B:165:VAL:HG22	2.01	0.42
1:A:103:ASP:O	1:A:104:ASN:HB2	2.20	0.42
4:B:478:NAG:O4	4:B:479:NAG:O6	2.33	0.42
1:B:105:SER:HB2	1:B:442:SER:HA	2.00	0.42
2:A:484:NAG:H5	2:A:485:NAG:HN2	1.85	0.42
1:A:309:ASP:O	1:A:310:TYR:HB2	2.19	0.42
1:A:467:MET:HE3	1:A:468:PRO:HD2	2.02	0.42
1:B:157:THR:OG1	1:B:173:GLN:HG2	2.20	0.42
1:A:127:VAL:HG23	1:A:128:LYS:N	2.34	0.42
1:A:402:ASN:ND2	1:A:459:PRO:CB	2.83	0.42
1:A:323:GLY:HA3	8:A:534:HOH:O	2.19	0.42
4:A:479:NAG:O4	4:A:483:MAN:C6	2.68	0.42
1:B:419:ASN:HD21	1:B:448:GLY:CA	2.32	0.42
1:A:431:LYS:HZ2	1:A:431:LYS:HA	1.85	0.42
1:A:427:ILE:HD11	1:A:439:THR:HG23	2.02	0.42
1:A:455:THR:HG23	8:A:555:HOH:O	2.20	0.42
1:A:375:GLU:HA	1:A:395:GLN:O	2.20	0.41
1:B:402:ASN:ND2	1:B:459:PRO:CB	2.83	0.41
1:A:157:THR:OG1	1:A:173:GLN:HG2	2.20	0.41
1:B:302:VAL:HG21	1:B:383:TRP:HZ3	1.84	0.41
1:B:161:ASN:HD22	1:B:161:ASN:HA	1.62	0.41
2:B:484:NAG:H5	2:B:485:NAG:HN2	1.85	0.41
4:B:479:NAG:O4	4:B:483:MAN:C6	2.68	0.41
1:B:353:ALA:HB1	1:B:360:LEU:HD11	2.02	0.41
1:B:190:LEU:CD1	1:B:257:ILE:HD13	2.46	0.41
1:A:118:ARG:CA	1:A:441:ASN:OD1	2.67	0.41
1:B:134:LEU:CB	1:B:156:ARG:NH2	2.80	0.41
1:A:158:LEU:HB3	1:A:175:CYS:SG	2.60	0.41
1:A:256:PHE:CD2	1:A:310:TYR:CD1	3.08	0.41
1:A:94:ILE:HD11	1:A:361:TRP:CD2	2.55	0.41
1:A:136:GLN:O	1:A:138:THR:N	2.54	0.41
1:A:425:GLU:HB2	1:A:441:ASN:ND2	2.35	0.41
1:A:266:SER:OG	1:A:310:TYR:HB3	2.21	0.41
1:A:463:ASN:HB3	1:A:466:PHE:CD2	2.56	0.41
1:A:435:ARG:HH21	1:A:465:ASN:HA	1.86	0.41
1:A:409:ILE:CG1	8:A:525:HOH:O	2.68	0.41
1:B:136:GLN:O	1:B:138:THR:N	2.54	0.41
1:B:158:LEU:HB3	1:B:175:CYS:SG	2.60	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:125:ASP:O	1:B:127:VAL:N	2.54	0.41
1:A:300:ARG:HH22	1:A:349:VAL:HG13	1.84	0.41
1:B:94:ILE:HD11	1:B:361:TRP:CD2	2.55	0.41
1:B:375:GLU:HA	1:B:395:GLN:O	2.20	0.41
1:B:366:ILE:HG21	1:B:400:SER:CB	2.49	0.41
1:B:266:SER:OG	1:B:310:TYR:HB3	2.21	0.41
1:A:432:GLN:O	1:A:434:THR:N	2.54	0.41
1:A:327:ARG:C	1:A:344:ARG:HE	2.24	0.41
1:B:228:SER:HB2	1:B:350:LYS:HZ3	1.85	0.41
1:B:425:GLU:HB2	1:B:441:ASN:ND2	2.35	0.41
1:A:191:HIS:O	1:A:193:CYS:SG	2.79	0.41
1:A:283:ARG:O	1:A:283:ARG:HD3	2.20	0.41
1:B:463:ASN:HB3	1:B:466:PHE:CD2	2.56	0.41
1:A:181:SER:OG	1:A:230:CYS:HB3	2.21	0.41
1:A:125:ASP:O	1:A:127:VAL:N	2.54	0.41
1:A:147:ASP:O	1:A:150:HIS:CD2	2.73	0.41
1:B:391:GLN:HE22	1:B:394:ARG:HB2	1.86	0.41
1:B:387:ASN:HA	1:B:387:ASN:HD22	1.61	0.41
1:A:218:TRP:CH2	1:A:241:MET:HB2	2.51	0.41
1:B:256:PHE:CD2	1:B:310:TYR:CD1	3.08	0.41
7:A:471:ST1:O3	7:A:471:ST1:HM43	2.21	0.41
1:B:191:HIS:O	1:B:193:CYS:SG	2.79	0.41
1:A:216:GLY:N	1:B:452:THR:HB	2.31	0.41
1:B:283:ARG:O	1:B:283:ARG:HD3	2.20	0.41
1:B:414:GLY:N	1:B:417:CYS:O	2.54	0.41
1:B:132:PHE:CD1	1:B:132:PHE:N	2.89	0.41
7:B:471:ST1:HM43	7:B:471:ST1:O3	2.21	0.41
1:B:120:PRO:HG2	1:B:423:TYR:HE2	1.86	0.40
1:A:338:ARG:NH1	1:A:339:ASP:OD1	2.55	0.40
1:A:127:VAL:HG23	1:A:128:LYS:H	1.86	0.40
1:B:327:ARG:C	1:B:344:ARG:HE	2.24	0.40
2:A:472:NAG:O3	2:A:485:NAG:C8	2.69	0.40
1:A:362:MET:CE	1:A:364:ARG:HG2	2.51	0.40
4:A:479:NAG:O4	4:A:483:MAN:H61	2.22	0.40
1:B:432:GLN:O	1:B:434:THR:N	2.54	0.40
1:B:320:GLY:N	1:B:388:SER:N	2.68	0.40
1:A:353:ALA:HB1	1:A:360:LEU:HD11	2.02	0.40
1:A:277:GLU:N	8:A:522:HOH:O	2.53	0.40
1:A:136:GLN:OE1	1:A:156:ARG:CG	2.69	0.40
1:A:144:HIS:HE1	1:B:462:ALA:CB	2.34	0.40
1:A:128:LYS:HE2	1:A:128:LYS:HB2	1.82	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:317:VAL:HG21	8:A:535:HOH:O	2.21	0.40
1:A:142:ASN:HD21	1:B:110:ALA:CB	2.14	0.40
1:B:91:GLN:OE1	1:B:420:ARG:NH1	2.51	0.40
1:B:368:LYS:NZ	1:B:369:ASP:OD2	2.54	0.40
1:A:284:TYR:CG	1:A:285:PRO:HA	2.57	0.40
1:B:256:PHE:N	1:B:256:PHE:CD1	2.90	0.40
1:B:349:VAL:HG22	1:B:350:LYS:N	2.37	0.40
1:B:136:GLN:OE1	1:B:156:ARG:CG	2.69	0.40
1:B:467:MET:HE3	1:B:468:PRO:HD2	2.02	0.40
1:A:368:LYS:NZ	1:A:369:ASP:OD2	2.54	0.40
3:A:475:NAG:H3	3:A:476:BMA:O5	2.22	0.40
1:B:256:PHE:CD2	1:B:310:TYR:CE1	3.10	0.40

All (12) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:463:ASN:HD22	4:A:482:MAN:H62[3_654]	1.02	0.58
3:A:477:FUL:H1	4:A:481:MAN:O3[3_654]	1.02	0.58
3:A:477:FUL:H3	4:A:481:MAN:HO3[3_654]	1.04	0.56
1:A:455:THR:OG1	4:B:478:NAG:H1[4_555]	1.42	0.18
1:A:455:THR:CG2	4:B:478:NAG:H61[4_555]	1.43	0.17
1:B:270:GLY:CA	8:A:489:HOH:O[7_544]	2.10	0.10
1:A:164:GLY:O	1:B:173:GLN:H[4_555]	1.54	0.06
1:A:455:THR:CG2	4:B:478:NAG:C6[4_555]	2.15	0.05
3:A:477:FUL:H3	4:A:481:MAN:O2[3_654]	1.56	0.04
3:A:477:FUL:H5	4:A:481:MAN:O5[3_654]	1.58	0.02
1:B:270:GLY:C	8:A:489:HOH:O[7_544]	2.18	0.02
1:A:396:VAL:CG1	4:B:483:MAN:O2[4_555]	2.18	0.02

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	386/388 (100%)	294 (76%)	70 (18%)	22 (6%)	2	0
1	B	386/388 (100%)	294 (76%)	70 (18%)	22 (6%)	2	0
All	All	772/776 (100%)	588 (76%)	140 (18%)	44 (6%)	2	0

All (44) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	145	SER
1	A	284	TYR
1	A	329	ASP
1	A	346	THR
1	A	347	GLN
1	A	397	ILE
1	A	412	VAL
1	B	145	SER
1	B	284	TYR
1	B	329	ASP
1	B	346	THR
1	B	347	GLN
1	B	397	ILE
1	B	412	VAL
1	A	87	TRP
1	A	118	ARG
1	A	259	GLU
1	A	322	VAL
1	A	416	SER
1	B	87	TRP
1	B	118	ARG
1	B	259	GLU
1	B	322	VAL
1	B	416	SER
1	A	133	ALA
1	A	137	GLY
1	A	147	ASP
1	A	431	LYS
1	B	133	ALA
1	B	137	GLY
1	B	147	ASP
1	B	431	LYS
1	A	356	ASN
1	B	356	ASN
1	A	121	TYR

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Mol	Chain	Res	Type
1	A	222	ILE
1	B	121	TYR
1	B	222	ILE
1	A	239	VAL
1	A	387	ASN
1	B	239	VAL
1	B	387	ASN
1	A	382	GLY
1	B	382	GLY

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	338/338 (100%)	293 (87%)	45 (13%)	5	1
1	B	338/338 (100%)	293 (87%)	45 (13%)	5	1
All	All	676/676 (100%)	586 (87%)	90 (13%)	5	1

All (90) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	83	GLU
1	A	89	LYS
1	A	122	VAL
1	A	124	CYS
1	A	128	LYS
1	A	142	ASN
1	A	155	HIS
1	A	157	THR
1	A	161	ASN
1	A	165	VAL
1	A	172	ARG
1	A	178	TRP
1	A	183	CYS
1	A	190	LEU

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Mol	Chain	Res	Type
1	A	192	VAL
1	A	210	ARG
1	A	224	ARG
1	A	231	VAL
1	A	238	THR
1	A	241	MET
1	A	254	ILE
1	A	255	LEU
1	A	257	ILE
1	A	262	ILE
1	A	273	GLN
1	A	276	GLU
1	A	283	ARG
1	A	308	GLU
1	A	321	LEU
1	A	330	ASP
1	A	338	ARG
1	A	344	ARG
1	A	347	GLN
1	A	355	ASP
1	A	364	ARG
1	A	387	ASN
1	A	391	GLN
1	A	401	ASP
1	A	410	PHE
1	A	418	ILE
1	A	420	ARG
1	A	427	ILE
1	A	431	LYS
1	A	435	ARG
1	A	449	THR
1	B	83	GLU
1	B	89	LYS
1	B	122	VAL
1	B	124	CYS
1	B	128	LYS
1	B	142	ASN
1	B	155	HIS
1	B	157	THR
1	B	161	ASN
1	B	165	VAL
1	B	172	ARG

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Mol	Chain	Res	Type
1	B	178	TRP
1	B	183	CYS
1	B	190	LEU
1	B	192	VAL
1	B	210	ARG
1	B	224	ARG
1	B	231	VAL
1	B	238	THR
1	B	241	MET
1	B	254	ILE
1	B	255	LEU
1	B	257	ILE
1	B	262	ILE
1	B	273	GLN
1	B	276	GLU
1	B	283	ARG
1	B	308	GLU
1	B	321	LEU
1	B	330	ASP
1	B	338	ARG
1	B	344	ARG
1	B	347	GLN
1	B	355	ASP
1	B	364	ARG
1	B	387	ASN
1	B	391	GLN
1	B	401	ASP
1	B	410	PHE
1	B	418	ILE
1	B	420	ARG
1	B	427	ILE
1	B	431	LYS
1	B	435	ARG
1	B	449	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	104	ASN
1	A	131	GLN
1	A	144	HIS
1	A	150	HIS

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Mol	Chain	Res	Type
1	A	161	ASN
1	A	173	GLN
1	A	220	GLN
1	A	221	ASN
1	A	226	GLN
1	A	334	ASN
1	A	387	ASN
1	A	393	ASN
1	A	402	ASN
1	A	419	ASN
1	B	104	ASN
1	B	131	GLN
1	B	150	HIS
1	B	161	ASN
1	B	168	HIS
1	B	173	GLN
1	B	220	GLN
1	B	226	GLN
1	B	334	ASN
1	B	387	ASN
1	B	393	ASN
1	B	402	ASN
1	B	419	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

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	$\# Z > 2$
2	NAG	A	472	1,2	14,14,15	1.67	3 (21%)	15,19,21	2.56	8 (53%)
2	NAG	A	473	2	14,14,15	1.99	4 (28%)	15,19,21	2.64	3 (20%)
3	NAG	A	474	1,3	14,14,15	1.96	3 (21%)	15,19,21	2.04	5 (33%)
3	NAG	A	475	3	14,14,15	1.77	1 (7%)	15,19,21	3.32	6 (40%)
3	BMA	A	476	3	11,11,12	2.18	5 (45%)	14,15,17	2.43	3 (21%)
3	FUL	A	477	3,4	10,10,11	2.56	4 (40%)	14,14,16	2.79	8 (57%)
4	NAG	A	478	1,4	14,14,15	4.63	9 (64%)	15,19,21	3.51	8 (53%)
4	NAG	A	479	4	14,14,15	1.86	4 (28%)	15,19,21	2.08	6 (40%)
4	BMA	A	480	4	11,11,12	2.42	2 (18%)	14,15,17	3.06	5 (35%)
4	MAN	A	481	3,4	11,11,12	2.89	5 (45%)	14,15,17	2.08	5 (35%)
4	MAN	A	482	4	11,11,12	2.37	3 (27%)	14,15,17	3.04	4 (28%)
4	MAN	A	483	4	11,11,12	3.17	5 (45%)	14,15,17	2.84	8 (57%)
2	NAG	A	484	1,2	14,14,15	2.08	5 (35%)	15,19,21	2.62	5 (33%)
2	NAG	A	485	2	14,14,15	1.44	3 (21%)	15,19,21	3.71	11 (73%)
2	NAG	B	472	1,2	14,14,15	1.67	3 (21%)	15,19,21	2.56	8 (53%)
2	NAG	B	473	2	14,14,15	1.99	4 (28%)	15,19,21	2.64	3 (20%)
5	NAG	B	474	1,5	14,14,15	1.96	3 (21%)	15,19,21	2.04	5 (33%)
5	NAG	B	475	5	14,14,15	1.77	1 (7%)	15,19,21	3.32	6 (40%)
5	BMA	B	476	5	11,11,12	2.18	5 (45%)	14,15,17	2.43	3 (21%)
5	FUC	B	477	5	10,10,11	2.56	4 (40%)	14,14,16	2.79	8 (57%)
4	NAG	B	478	1,4	14,14,15	4.63	9 (64%)	15,19,21	3.51	8 (53%)
4	NAG	B	479	4	14,14,15	1.86	4 (28%)	15,19,21	2.08	6 (40%)
4	BMA	B	480	4	11,11,12	2.42	2 (18%)	14,15,17	3.06	5 (35%)
4	MAN	B	481	4	11,11,12	2.89	5 (45%)	14,15,17	2.08	5 (35%)
4	MAN	B	482	4	11,11,12	2.37	3 (27%)	14,15,17	3.04	4 (28%)
4	MAN	B	483	4	11,11,12	3.17	5 (45%)	14,15,17	2.84	8 (57%)
2	NAG	B	484	1,2	14,14,15	2.08	5 (35%)	15,19,21	2.62	5 (33%)
2	NAG	B	485	2	14,14,15	1.44	3 (21%)	15,19,21	3.71	11 (73%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	A	472	1,2	-	0/6/23/26	0/1/1/1
2	NAG	A	473	2	-	0/6/23/26	0/1/1/1
3	NAG	A	474	1,3	-	0/6/23/26	0/1/1/1
3	NAG	A	475	3	-	0/6/23/26	0/1/1/1
3	BMA	A	476	3	-	0/2/19/22	1/1/1/1
3	FUL	A	477	3,4	-	0/0/17/20	0/1/1/1
4	NAG	A	478	1,4	-	0/6/23/26	0/1/1/1
4	NAG	A	479	4	-	0/6/23/26	0/1/1/1
4	BMA	A	480	4	-	0/2/19/22	0/1/1/1
4	MAN	A	481	3,4	-	0/2/19/22	0/1/1/1
4	MAN	A	482	4	-	0/2/19/22	0/1/1/1
4	MAN	A	483	4	-	0/2/19/22	1/1/1/1
2	NAG	A	484	1,2	-	0/6/23/26	0/1/1/1
2	NAG	A	485	2	-	0/6/23/26	0/1/1/1
2	NAG	B	472	1,2	-	0/6/23/26	0/1/1/1
2	NAG	B	473	2	-	0/6/23/26	0/1/1/1
5	NAG	B	474	1,5	-	0/6/23/26	0/1/1/1
5	NAG	B	475	5	-	0/6/23/26	0/1/1/1
5	BMA	B	476	5	-	0/2/19/22	1/1/1/1
5	FUC	B	477	5	1/1/5/5	0/0/17/20	0/1/1/1
4	NAG	B	478	1,4	-	0/6/23/26	0/1/1/1
4	NAG	B	479	4	-	0/6/23/26	0/1/1/1
4	BMA	B	480	4	-	0/2/19/22	0/1/1/1
4	MAN	B	481	4	-	0/2/19/22	0/1/1/1
4	MAN	B	482	4	-	0/2/19/22	0/1/1/1
4	MAN	B	483	4	-	0/2/19/22	1/1/1/1
2	NAG	B	484	1,2	-	0/6/23/26	0/1/1/1
2	NAG	B	485	2	-	0/6/23/26	0/1/1/1

All (112) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	477	FUC	O2-C2	-3.39	1.35	1.43
3	A	477	FUL	O2-C2	-3.39	1.35	1.43
4	B	481	MAN	O4-C4	-3.24	1.35	1.43
4	A	481	MAN	O4-C4	-3.24	1.35	1.43
4	A	479	NAG	O5-C1	-2.44	1.39	1.43
4	B	479	NAG	O5-C1	-2.44	1.39	1.43
4	A	483	MAN	O2-C2	-2.11	1.38	1.43
4	B	483	MAN	O2-C2	-2.11	1.38	1.43
4	A	479	NAG	C4-C3	2.04	1.57	1.52
4	B	479	NAG	C4-C3	2.04	1.57	1.52
4	B	481	MAN	O5-C1	2.05	1.47	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	481	MAN	O5-C1	2.05	1.47	1.43
2	B	472	NAG	C1-C2	2.14	1.55	1.52
2	A	472	NAG	C1-C2	2.14	1.55	1.52
3	A	474	NAG	C4-C5	2.16	1.57	1.53
5	B	474	NAG	C4-C5	2.16	1.57	1.53
2	B	485	NAG	C4-C3	2.18	1.58	1.52
2	A	485	NAG	C4-C3	2.18	1.58	1.52
3	A	474	NAG	O5-C5	2.25	1.48	1.43
5	B	474	NAG	O5-C5	2.25	1.48	1.43
2	B	473	NAG	O4-C4	2.49	1.48	1.43
2	A	473	NAG	O4-C4	2.49	1.48	1.43
4	B	478	NAG	O6-C6	2.49	1.53	1.42
4	A	478	NAG	O6-C6	2.49	1.53	1.42
2	B	484	NAG	O5-C5	2.50	1.49	1.43
2	A	484	NAG	O5-C5	2.50	1.49	1.43
2	B	485	NAG	C4-C5	2.51	1.58	1.53
2	A	485	NAG	C4-C5	2.51	1.58	1.53
5	B	476	BMA	O5-C1	2.52	1.47	1.43
3	A	476	BMA	O5-C1	2.52	1.47	1.43
2	B	473	NAG	C8-C7	2.53	1.55	1.50
2	A	473	NAG	C8-C7	2.53	1.55	1.50
4	A	483	MAN	C1-C2	2.68	1.58	1.52
4	B	483	MAN	C1-C2	2.68	1.58	1.52
5	B	476	BMA	C4-C5	2.69	1.58	1.53
3	A	476	BMA	C4-C5	2.69	1.58	1.53
2	B	484	NAG	C2-N2	2.72	1.51	1.46
2	A	484	NAG	C2-N2	2.72	1.51	1.46
5	B	477	FUC	C6-C5	2.75	1.58	1.51
3	A	477	FUL	C6-C5	2.75	1.58	1.51
4	A	483	MAN	O4-C4	2.78	1.49	1.43
4	B	483	MAN	O4-C4	2.78	1.49	1.43
4	A	479	NAG	C4-C5	2.82	1.59	1.53
4	B	479	NAG	C4-C5	2.82	1.59	1.53
2	B	484	NAG	C1-C2	2.82	1.56	1.52
2	A	484	NAG	C1-C2	2.82	1.56	1.52
2	B	484	NAG	O4-C4	3.12	1.50	1.43
2	A	484	NAG	O4-C4	3.12	1.50	1.43
2	B	473	NAG	C4-C5	3.14	1.59	1.53
2	A	473	NAG	C4-C5	3.14	1.59	1.53
4	B	481	MAN	C1-C2	3.20	1.59	1.52
4	A	481	MAN	C1-C2	3.20	1.59	1.52
2	B	472	NAG	C4-C3	3.24	1.60	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	472	NAG	C4-C3	3.24	1.60	1.52
5	B	476	BMA	C4-C3	3.26	1.61	1.52
3	A	476	BMA	C4-C3	3.26	1.61	1.52
2	B	485	NAG	C3-C2	3.27	1.60	1.52
2	A	485	NAG	C3-C2	3.27	1.60	1.52
5	B	476	BMA	C2-C3	3.32	1.57	1.52
3	A	476	BMA	C2-C3	3.32	1.57	1.52
5	B	476	BMA	C1-C2	3.34	1.60	1.52
3	A	476	BMA	C1-C2	3.34	1.60	1.52
4	B	482	MAN	O5-C1	3.73	1.50	1.43
4	A	482	MAN	O5-C1	3.73	1.50	1.43
4	B	482	MAN	O5-C5	3.73	1.51	1.43
4	A	482	MAN	O5-C5	3.73	1.51	1.43
4	A	479	NAG	C6-C5	3.75	1.65	1.51
4	B	479	NAG	C6-C5	3.75	1.65	1.51
4	B	481	MAN	O5-C5	3.98	1.52	1.43
4	A	481	MAN	O5-C5	3.98	1.52	1.43
4	B	478	NAG	C2-N2	4.00	1.53	1.46
4	A	478	NAG	C2-N2	4.00	1.53	1.46
2	B	472	NAG	O5-C5	4.01	1.52	1.43
2	A	472	NAG	O5-C5	4.01	1.52	1.43
5	B	477	FUC	C4-C3	4.10	1.63	1.52
3	A	477	FUL	C4-C3	4.10	1.63	1.52
4	B	480	BMA	C4-C5	4.16	1.61	1.53
4	A	480	BMA	C4-C5	4.16	1.61	1.53
4	B	478	NAG	C6-C5	4.18	1.66	1.51
4	A	478	NAG	C6-C5	4.18	1.66	1.51
4	B	482	MAN	C4-C5	4.48	1.62	1.53
4	A	482	MAN	C4-C5	4.48	1.62	1.53
4	B	478	NAG	O5-C5	4.57	1.53	1.43
4	A	478	NAG	O5-C5	4.57	1.53	1.43
2	B	484	NAG	O5-C1	4.71	1.51	1.43
2	A	484	NAG	O5-C1	4.71	1.51	1.43
4	B	478	NAG	C4-C3	4.75	1.64	1.52
4	A	478	NAG	C4-C3	4.75	1.64	1.52
2	B	473	NAG	C4-C3	4.76	1.64	1.52
2	A	473	NAG	C4-C3	4.76	1.64	1.52
5	B	477	FUC	C4-C5	5.05	1.63	1.52
3	A	477	FUL	C4-C5	5.05	1.63	1.52
4	A	483	MAN	O5-C5	5.16	1.54	1.43
4	B	483	MAN	O5-C5	5.16	1.54	1.43
5	B	475	NAG	C4-C3	5.46	1.66	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	475	NAG	C4-C3	5.46	1.66	1.52
4	B	478	NAG	C3-C2	5.48	1.65	1.52
4	A	478	NAG	C3-C2	5.48	1.65	1.52
3	A	474	NAG	C1-C2	5.65	1.60	1.52
5	B	474	NAG	C1-C2	5.65	1.60	1.52
4	B	478	NAG	C4-C5	6.40	1.66	1.53
4	A	478	NAG	C4-C5	6.40	1.66	1.53
4	B	480	BMA	C1-C2	6.45	1.67	1.52
4	A	480	BMA	C1-C2	6.45	1.67	1.52
4	B	481	MAN	C2-C3	6.79	1.61	1.52
4	A	481	MAN	C2-C3	6.79	1.61	1.52
4	B	478	NAG	O5-C1	7.28	1.55	1.43
4	A	478	NAG	O5-C1	7.28	1.55	1.43
4	A	483	MAN	C4-C5	7.68	1.69	1.53
4	B	483	MAN	C4-C5	7.68	1.69	1.53
4	B	478	NAG	C1-C2	9.30	1.65	1.52
4	A	478	NAG	C1-C2	9.30	1.65	1.52

All (170) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	478	NAG	O7-C7-C8	-7.90	107.56	122.06
4	A	478	NAG	O7-C7-C8	-7.90	107.56	122.06
2	B	485	NAG	C2-N2-C7	-7.27	113.70	123.04
2	A	485	NAG	C2-N2-C7	-7.27	113.70	123.04
4	A	483	MAN	C3-C4-C5	-6.01	99.72	110.20
4	B	483	MAN	C3-C4-C5	-6.01	99.72	110.20
5	B	475	NAG	C4-C3-C2	-5.67	102.41	111.23
3	A	475	NAG	C4-C3-C2	-5.67	102.41	111.23
2	B	485	NAG	C4-C3-C2	-5.67	102.41	111.23
2	A	485	NAG	C4-C3-C2	-5.67	102.41	111.23
5	B	475	NAG	O3-C3-C2	-5.24	98.73	109.11
3	A	475	NAG	O3-C3-C2	-5.24	98.73	109.11
2	B	485	NAG	C3-C4-C5	-5.19	101.15	110.20
2	A	485	NAG	C3-C4-C5	-5.19	101.15	110.20
5	B	475	NAG	C2-N2-C7	-4.78	116.90	123.04
3	A	475	NAG	C2-N2-C7	-4.78	116.90	123.04
2	B	472	NAG	C4-C3-C2	-4.70	103.92	111.23
2	A	472	NAG	C4-C3-C2	-4.70	103.92	111.23
5	B	477	FUC	O5-C1-C2	-4.64	103.34	110.86
3	A	477	FUL	O5-C1-C2	-4.64	103.34	110.86
4	B	481	MAN	O4-C4-C3	-4.49	100.22	110.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	481	MAN	O4-C4-C3	-4.49	100.22	110.34
5	B	477	FUC	C2-C3-C4	-4.38	103.60	111.04
3	A	477	FUL	C2-C3-C4	-4.38	103.60	111.04
3	A	474	NAG	C3-C2-N2	-4.20	100.50	110.56
5	B	474	NAG	C3-C2-N2	-4.20	100.50	110.56
5	B	475	NAG	C3-C4-C5	-4.09	103.07	110.20
3	A	475	NAG	C3-C4-C5	-4.09	103.07	110.20
5	B	477	FUC	C1-C2-C3	-3.67	105.20	109.54
3	A	477	FUL	C1-C2-C3	-3.67	105.20	109.54
4	A	479	NAG	C4-C3-C2	-3.44	105.88	111.23
4	B	479	NAG	C4-C3-C2	-3.44	105.88	111.23
4	A	483	MAN	O2-C2-C3	-3.42	103.24	110.12
4	B	483	MAN	O2-C2-C3	-3.42	103.24	110.12
4	B	478	NAG	O3-C3-C4	-3.42	102.65	110.34
4	A	478	NAG	O3-C3-C4	-3.42	102.65	110.34
4	A	479	NAG	C2-N2-C7	-3.36	118.73	123.04
4	B	479	NAG	C2-N2-C7	-3.36	118.73	123.04
5	B	477	FUC	O3-C3-C2	-3.31	104.02	110.00
3	A	477	FUL	O3-C3-C2	-3.31	104.02	110.00
2	B	484	NAG	O7-C7-C8	-3.11	116.35	122.06
2	A	484	NAG	O7-C7-C8	-3.11	116.35	122.06
5	B	476	BMA	C1-C2-C3	-3.02	105.97	109.54
3	A	476	BMA	C1-C2-C3	-3.02	105.97	109.54
2	B	484	NAG	O4-C4-C3	-2.97	103.64	110.34
2	A	484	NAG	O4-C4-C3	-2.97	103.64	110.34
4	B	482	MAN	O2-C2-C1	-2.77	103.65	109.21
4	A	482	MAN	O2-C2-C1	-2.77	103.65	109.21
4	B	482	MAN	O2-C2-C3	-2.72	104.66	110.12
4	A	482	MAN	O2-C2-C3	-2.72	104.66	110.12
2	B	472	NAG	C8-C7-N2	-2.70	110.94	116.11
2	A	472	NAG	C8-C7-N2	-2.70	110.94	116.11
2	B	472	NAG	C3-C4-C5	-2.59	105.69	110.20
2	A	472	NAG	C3-C4-C5	-2.59	105.69	110.20
4	B	480	BMA	O2-C2-C3	-2.38	105.34	110.12
4	A	480	BMA	O2-C2-C3	-2.38	105.34	110.12
2	B	485	NAG	O7-C7-C8	-2.30	117.84	122.06
2	A	485	NAG	O7-C7-C8	-2.30	117.84	122.06
4	B	480	BMA	C3-C4-C5	-2.22	106.32	110.20
4	A	480	BMA	C3-C4-C5	-2.22	106.32	110.20
2	B	473	NAG	C2-N2-C7	-2.19	120.22	123.04
2	A	473	NAG	C2-N2-C7	-2.19	120.22	123.04
4	A	483	MAN	O6-C6-C5	-2.13	104.31	111.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	483	MAN	O6-C6-C5	-2.13	104.31	111.33
3	A	474	NAG	O4-C4-C3	-2.12	105.56	110.34
5	B	474	NAG	O4-C4-C3	-2.12	105.56	110.34
2	B	485	NAG	O6-C6-C5	-2.06	104.52	111.33
2	A	485	NAG	O6-C6-C5	-2.06	104.52	111.33
4	B	481	MAN	O3-C3-C4	-2.03	105.77	110.34
4	A	481	MAN	O3-C3-C4	-2.03	105.77	110.34
5	B	477	FUC	O2-C2-C1	2.01	113.24	109.21
3	A	477	FUL	O2-C2-C1	2.01	113.24	109.21
2	B	485	NAG	O3-C3-C4	2.02	114.88	110.34
2	A	485	NAG	O3-C3-C4	2.02	114.88	110.34
2	B	485	NAG	C8-C7-N2	2.05	120.03	116.11
2	A	485	NAG	C8-C7-N2	2.05	120.03	116.11
4	B	481	MAN	C1-C2-C3	2.09	112.02	109.54
4	A	481	MAN	C1-C2-C3	2.09	112.02	109.54
4	A	479	NAG	O6-C6-C5	2.21	118.63	111.33
4	B	479	NAG	O6-C6-C5	2.21	118.63	111.33
5	B	477	FUC	C6-C5-C4	2.24	117.50	113.08
3	A	477	FUL	C6-C5-C4	2.24	117.50	113.08
4	A	483	MAN	O2-C2-C1	2.30	113.81	109.21
4	B	483	MAN	O2-C2-C1	2.30	113.81	109.21
4	A	483	MAN	C6-C5-C4	2.31	118.72	113.02
4	B	483	MAN	C6-C5-C4	2.31	118.72	113.02
4	A	479	NAG	O3-C3-C2	2.45	113.98	109.11
4	B	479	NAG	O3-C3-C2	2.45	113.98	109.11
4	B	478	NAG	O6-C6-C5	2.48	119.53	111.33
4	A	478	NAG	O6-C6-C5	2.48	119.53	111.33
2	B	485	NAG	C3-C2-N2	2.51	116.56	110.56
2	A	485	NAG	C3-C2-N2	2.51	116.56	110.56
4	B	478	NAG	O5-C5-C6	2.59	112.95	107.35
4	A	478	NAG	O5-C5-C6	2.59	112.95	107.35
4	A	483	MAN	O3-C3-C2	2.63	114.76	110.00
4	B	483	MAN	O3-C3-C2	2.63	114.76	110.00
2	B	472	NAG	O3-C3-C4	2.64	116.28	110.34
2	A	472	NAG	O3-C3-C4	2.64	116.28	110.34
2	B	472	NAG	O3-C3-C2	2.64	114.34	109.11
2	A	472	NAG	O3-C3-C2	2.64	114.34	109.11
4	B	478	NAG	O4-C4-C5	2.65	116.25	109.24
4	A	478	NAG	O4-C4-C5	2.65	116.25	109.24
4	B	481	MAN	O3-C3-C2	2.66	114.81	110.00
4	A	481	MAN	O3-C3-C2	2.66	114.81	110.00
2	B	484	NAG	C8-C7-N2	2.90	121.65	116.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	484	NAG	C8-C7-N2	2.90	121.65	116.11
3	A	474	NAG	C3-C4-C5	3.15	115.69	110.20
5	B	474	NAG	C3-C4-C5	3.15	115.69	110.20
3	A	474	NAG	O3-C3-C2	3.20	115.45	109.11
5	B	474	NAG	O3-C3-C2	3.20	115.45	109.11
2	B	485	NAG	O4-C4-C5	3.21	117.73	109.24
2	A	485	NAG	O4-C4-C5	3.21	117.73	109.24
4	A	479	NAG	C1-O5-C5	3.21	116.33	112.25
4	B	479	NAG	C1-O5-C5	3.21	116.33	112.25
3	A	474	NAG	C2-N2-C7	3.33	127.31	123.04
5	B	474	NAG	C2-N2-C7	3.33	127.31	123.04
2	B	472	NAG	C2-N2-C7	3.35	127.34	123.04
2	A	472	NAG	C2-N2-C7	3.35	127.34	123.04
5	B	477	FUC	C3-C4-C5	3.35	115.36	109.72
3	A	477	FUL	C3-C4-C5	3.35	115.36	109.72
4	B	478	NAG	C4-C3-C2	3.51	116.69	111.23
4	A	478	NAG	C4-C3-C2	3.51	116.69	111.23
4	B	481	MAN	C3-C4-C5	3.69	116.62	110.20
4	A	481	MAN	C3-C4-C5	3.69	116.62	110.20
2	B	484	NAG	O3-C3-C2	3.71	116.46	109.11
2	A	484	NAG	O3-C3-C2	3.71	116.46	109.11
2	B	472	NAG	C1-O5-C5	3.79	117.06	112.25
2	A	472	NAG	C1-O5-C5	3.79	117.06	112.25
4	A	479	NAG	O5-C5-C6	3.79	115.56	107.35
4	B	479	NAG	O5-C5-C6	3.79	115.56	107.35
2	B	472	NAG	O4-C4-C3	3.90	119.11	110.34
2	A	472	NAG	O4-C4-C3	3.90	119.11	110.34
5	B	477	FUC	O3-C3-C4	3.95	119.22	110.34
3	A	477	FUL	O3-C3-C4	3.95	119.22	110.34
4	A	483	MAN	C1-C2-C3	4.19	114.49	109.54
4	B	483	MAN	C1-C2-C3	4.19	114.49	109.54
2	B	473	NAG	C6-C5-C4	4.24	123.49	113.02
2	A	473	NAG	C6-C5-C4	4.24	123.49	113.02
5	B	475	NAG	O3-C3-C4	4.39	120.22	110.34
3	A	475	NAG	O3-C3-C4	4.39	120.22	110.34
2	B	485	NAG	C6-C5-C4	4.57	124.28	113.02
2	A	485	NAG	C6-C5-C4	4.57	124.28	113.02
4	A	483	MAN	O4-C4-C5	4.58	121.39	109.24
4	B	483	MAN	O4-C4-C5	4.58	121.39	109.24
4	B	482	MAN	C1-C2-C3	5.04	115.51	109.54
4	A	482	MAN	C1-C2-C3	5.04	115.51	109.54
5	B	476	BMA	O2-C2-C1	5.29	119.82	109.21

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	476	BMA	O2-C2-C1	5.29	119.82	109.21
4	B	478	NAG	C6-C5-C4	5.41	126.35	113.02
4	A	478	NAG	C6-C5-C4	5.41	126.35	113.02
5	B	476	BMA	C1-O5-C5	5.47	119.19	112.25
3	A	476	BMA	C1-O5-C5	5.47	119.19	112.25
4	B	480	BMA	C1-O5-C5	5.54	119.28	112.25
4	A	480	BMA	C1-O5-C5	5.54	119.28	112.25
4	B	480	BMA	O5-C1-C2	5.69	120.08	110.86
4	A	480	BMA	O5-C1-C2	5.69	120.08	110.86
5	B	475	NAG	O4-C4-C3	5.85	123.52	110.34
3	A	475	NAG	O4-C4-C3	5.85	123.52	110.34
2	B	485	NAG	O3-C3-C2	5.99	120.97	109.11
2	A	485	NAG	O3-C3-C2	5.99	120.97	109.11
4	B	480	BMA	O2-C2-C1	6.38	121.99	109.21
4	A	480	BMA	O2-C2-C1	6.38	121.99	109.21
4	B	478	NAG	C8-C7-N2	6.73	128.98	116.11
4	A	478	NAG	C8-C7-N2	6.73	128.98	116.11
2	B	484	NAG	C1-O5-C5	7.21	121.40	112.25
2	A	484	NAG	C1-O5-C5	7.21	121.40	112.25
2	B	473	NAG	C1-O5-C5	8.29	122.77	112.25
2	A	473	NAG	C1-O5-C5	8.29	122.77	112.25
4	B	482	MAN	C1-O5-C5	8.99	123.66	112.25
4	A	482	MAN	C1-O5-C5	8.99	123.66	112.25

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
5	B	477	FUC	C1

There are no torsion outliers.

All (4) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	483	MAN	C1-C2-C3-C4-C5-O5
4	B	483	MAN	C1-C2-C3-C4-C5-O5
5	B	476	BMA	C1-C2-C3-C4-C5-O5
3	A	476	BMA	C1-C2-C3-C4-C5-O5

21 monomers are involved in 57 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	472	NAG	5	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	473	NAG	3	0
3	A	475	NAG	2	0
3	A	476	BMA	2	0
3	A	477	FUL	0	4
4	A	478	NAG	19	0
4	A	479	NAG	3	0
4	A	480	BMA	1	0
4	A	481	MAN	0	4
4	A	482	MAN	0	1
4	A	483	MAN	6	0
2	A	484	NAG	3	0
2	A	485	NAG	5	0
2	B	472	NAG	4	0
2	B	473	NAG	3	0
4	B	478	NAG	1	3
4	B	479	NAG	3	0
4	B	480	BMA	1	0
4	B	483	MAN	1	1
2	B	484	NAG	3	0
2	B	485	NAG	4	0

5.6 Ligand geometry

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 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	$\# Z > 2$
7	ST1	A	471	-	12,17,17	1.05	1 (8%)	13,24,24	3.02	3 (23%)
7	ST1	B	471	-	12,17,17	1.05	1 (8%)	13,24,24	3.02	3 (23%)

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
7	ST1	A	471	-	-	1/7/12/12	0/1/1/1
7	ST1	B	471	-	-	1/7/12/12	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	B	471	ST1	C4-C5	-2.16	1.38	1.42
7	A	471	ST1	C4-C5	-2.16	1.38	1.42

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	B	471	ST1	C4-N4-C4'	-9.70	110.54	123.33
7	A	471	ST1	C4-N4-C4'	-9.70	110.54	123.33
7	B	471	ST1	C5-C4-N4	-2.10	119.07	123.28
7	A	471	ST1	C5-C4-N4	-2.10	119.07	123.28
7	B	471	ST1	CM4-C4'-N4	3.23	119.56	114.97
7	A	471	ST1	CM4-C4'-N4	3.23	119.56	114.97

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	B	471	ST1	CM4-C4'-N4-C4
7	A	471	ST1	CM4-C4'-N4-C4

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	471	ST1	1	0
7	B	471	ST1	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	388/388 (100%)	0.28	4 (1%) 84 86	2, 8, 19, 43	0
1	B	388/388 (100%)	0.28	7 (1%) 71 74	2, 8, 19, 43	0
All	All	776/776 (100%)	0.28	11 (1%) 78 80	2, 8, 20, 43	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	345	GLY	5.2
1	B	345	GLY	3.3
1	A	346	THR	2.9
1	A	464	ILE	2.8
1	B	84	TYR	2.8
1	B	82	VAL	2.6
1	B	246	ALA	2.6
1	B	434	THR	2.4
1	B	436	VAL	2.2
1	A	331	ARG	2.2
1	B	410	PHE	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

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	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
4	MAN	A	483	11/12	0.88	0.34	11.24	0,0,19,20	0
4	MAN	B	483	11/12	0.91	0.28	7.48	0,0,19,20	0
2	NAG	B	484	14/15	0.90	0.18	4.07	0,0,12,16	0
4	NAG	A	478	14/15	0.87	0.19	3.58	0,2,14,17	0
4	NAG	B	478	14/15	0.86	0.18	2.70	0,2,14,17	0
2	NAG	A	484	14/15	0.91	0.17	1.73	0,0,12,16	0
2	NAG	B	472	14/15	0.94	0.16	0.22	0,0,12,16	0
2	NAG	A	472	14/15	0.95	0.13	-0.25	0,0,12,16	0
5	FUC	B	477	10/11	0.85	0.30	-	0,0,17,21	0
5	BMA	B	476	11/12	0.87	0.30	-	0,0,18,25	0
4	BMA	B	480	11/12	0.68	0.29	-	0,0,27,29	0
3	FUL	A	477	10/11	0.90	0.24	-	0,0,17,21	0
4	MAN	B	481	11/12	0.87	0.29	-	0,0,22,25	0
4	NAG	A	479	14/15	0.87	0.33	-	0,0,22,23	0
2	NAG	B	473	14/15	0.91	0.24	-	0,0,21,22	0
5	NAG	B	475	14/15	0.84	0.22	-	0,0,35,38	0
3	NAG	A	474	14/15	0.87	0.16	-	0,6,19,22	0
4	MAN	B	482	11/12	0.79	0.27	-	0,0,26,30	0
3	NAG	A	475	14/15	0.87	0.17	-	0,0,35,38	0
2	NAG	B	485	14/15	0.95	0.16	-	0,0,9,14	0
4	MAN	A	481	11/12	0.90	0.34	-	0,0,22,25	0
3	BMA	A	476	11/12	0.88	0.29	-	0,0,18,25	0
4	BMA	A	480	11/12	0.80	0.33	-	0,0,27,29	0
2	NAG	A	485	14/15	0.94	0.20	-	0,0,9,14	0
4	MAN	A	482	11/12	0.78	0.33	-	0,0,26,30	0
2	NAG	A	473	14/15	0.86	0.17	-	0,0,21,22	0
5	NAG	B	474	14/15	0.85	0.15	-	0,6,19,22	0
4	NAG	B	479	14/15	0.91	0.31	-	0,0,22,23	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	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
7	ST1	A	471	17/17	0.89	0.14	0.85	0,10,19,22	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
7	ST1	B	471	17/17	0.92	0.12	-0.69	0,10,19,22	0
6	CA	A	470	1/1	1.00	0.03	-1.63	15,15,15,15	0
6	CA	B	470	1/1	0.99	0.05	-1.86	15,15,15,15	0

6.5 Other polymers [i](#)

There are no such residues in this entry.