



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 09:14 pm GMT

PDB ID : 3VTA  
Title : Crystal Structure of cucumisin, a subtilisin-like endoprotease from Cucumis melo L  
Authors : Murayama, K.; Kato-Murayama, M.; Hosaka, T.; Sotokawauchi, A.; Shirouzu, M.; Arima, K.; Yokoyama, S.  
Deposited on : 2012-05-23  
Resolution : 2.75 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467  
Mogul : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : trunk28620  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
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) : recalc28949

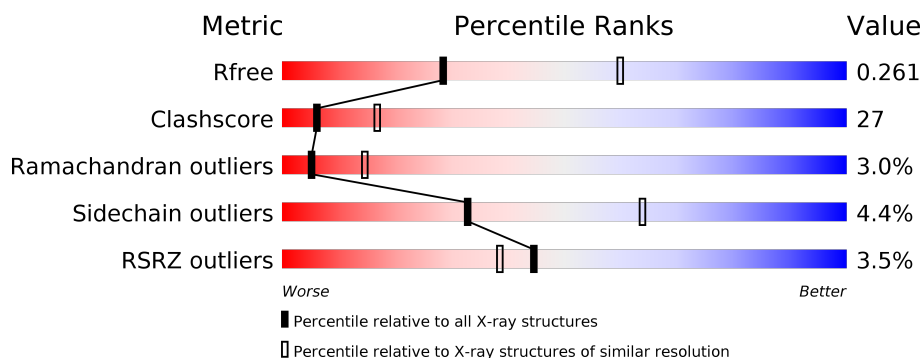
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.75 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))
$R_{free}$	100719	3666 (2.80-2.72)
Clashscore	112137	4174 (2.80-2.72)
Ramachandran outliers	110173	4103 (2.80-2.72)
Sidechain outliers	110143	4106 (2.80-2.72)
RSRZ outliers	101464	3697 (2.80-2.72)

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

Mol	Chain	Length	Quality of chain
1	A	621	<div> <div>%</div> <div> <div></div> <div>57%</div> <div>34%</div> <div>• 5%</div> </div> </div>
1	B	621	<div> <div>6%</div> <div> <div></div> <div>50%</div> <div>42%</div> <div>• •</div> </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	DFP	B	801	-	-	-	X
4	NAG	B	808	-	-	-	X

## 2 Entry composition [i](#)

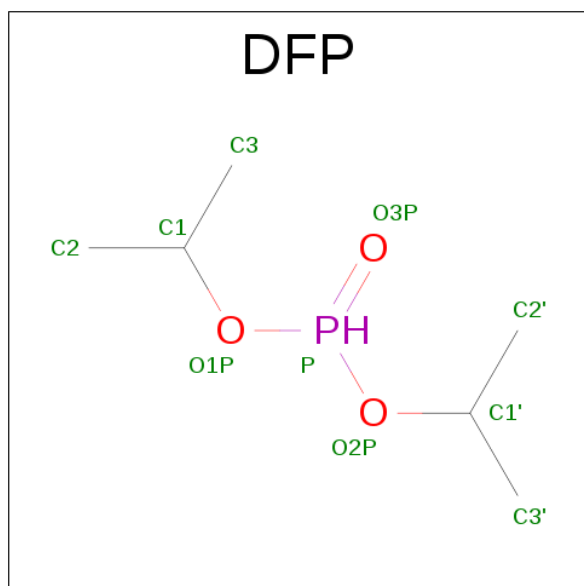
There are 5 unique types of molecules in this entry. The entry contains 9322 atoms, of which 0 are hydrogens and 0 are deuteriums.

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

- Molecule 1 is a protein called Cucumisn.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	591	Total	C	N	O	S	0	0	0
			4455	2818	773	851	13			
1	B	594	Total	C	N	O	S	0	0	0
			4483	2839	777	854	13			

- Molecule 2 is DIISOPROPYL PHOSPHONATE (three-letter code: DFP) (formula:  $C_6H_{15}O_3P$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	O	P	0	0
			10	6	3	1		
2	B	1	Total	C	O	P	0	0
			10	6	3	1		

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

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	5	Total	C	N	O	0	0
			60	34	2	24		
3	B	5	Total	C	N	O	0	0
			60	34	2	24		

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

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	3	Total	C	N	O	0	0
			38	22	2	14		

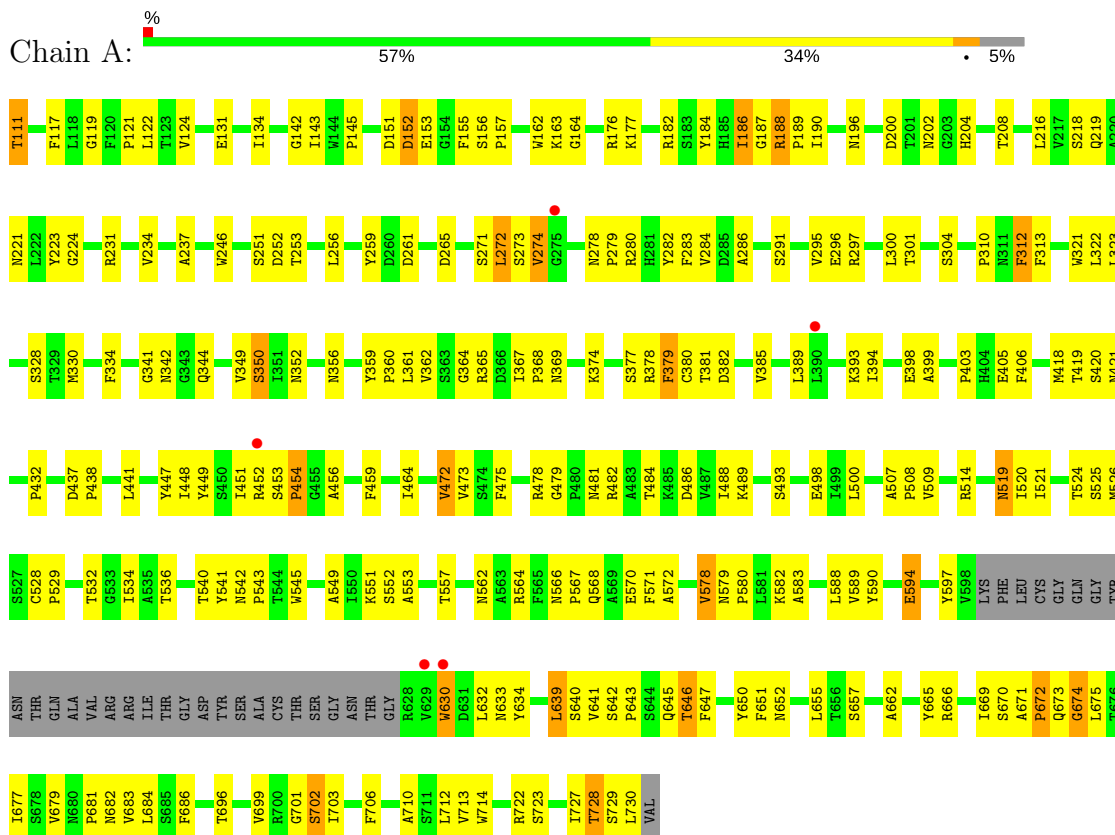
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	129	Total	O	0	0
			129	129		
5	B	77	Total	O	0	0
			77	77		

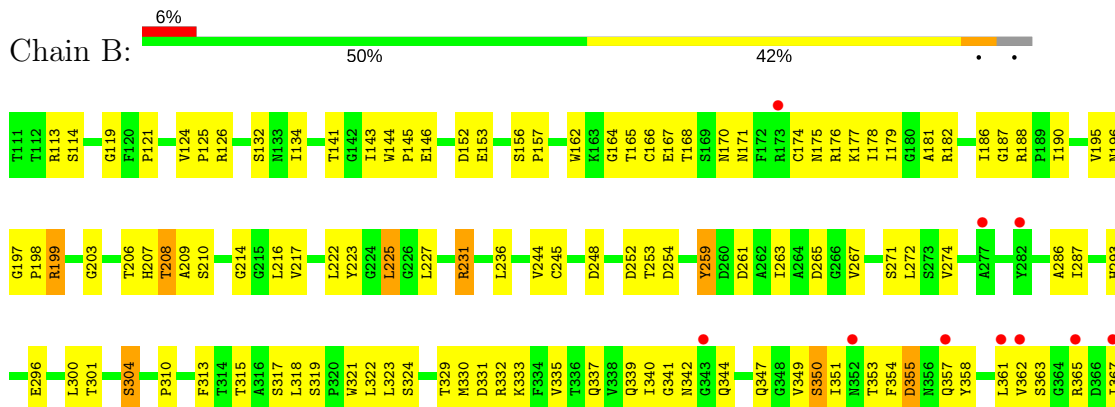
### 3 Residue-property plots

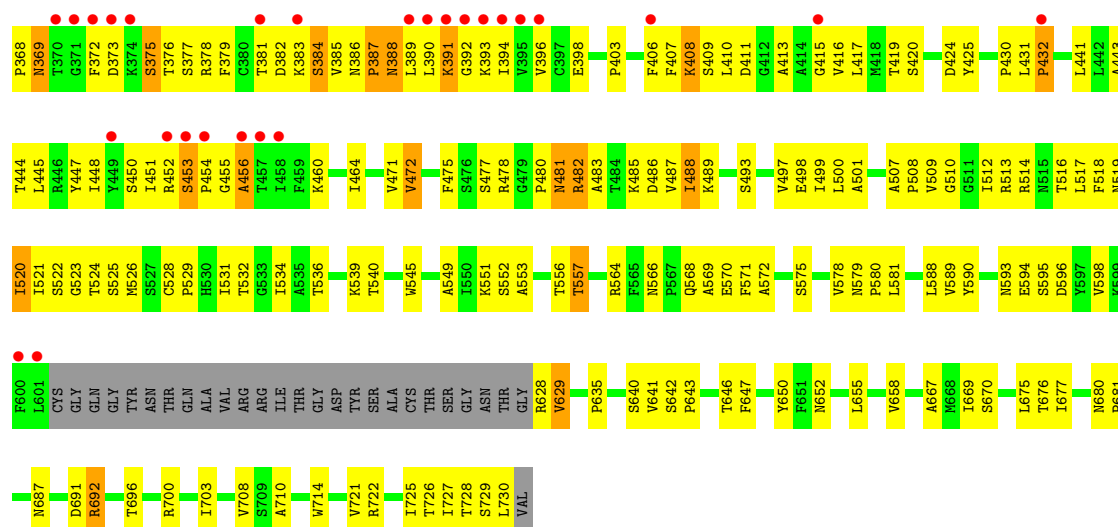
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

#### • Molecule 1: Cucumisins



#### • Molecule 1: Cucumisins





## 4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	149.48Å 149.48Å 218.03Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	44.64 – 2.75 44.64 – 2.76	Depositor EDS
% Data completeness (in resolution range)	99.9 (44.64-2.75) 100.0 (44.64-2.76)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.16	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.45 (at 2.77Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.214 , 0.262 0.220 , 0.261	Depositor DCC
$R_{free}$ test set	2331 reflections (4.97%)	DCC
Wilson B-factor (Å <sup>2</sup> )	42.7	Xtriage
Anisotropy	0.451	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 24.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.047 for h,-h-k,-l	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	9322	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.06% of the height of the origin peak. No significant pseudotranslation is detected.*

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

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 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: FUC, MAN, BMA, NAG, DFP

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	0.39	0/4570	0.65	0/6241
1	B	0.37	0/4599	0.63	0/6279
All	All	0.38	0/9169	0.64	0/12520

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4455	0	4358	224	0
1	B	4483	0	4389	266	0
2	A	10	0	14	0	0
2	B	10	0	14	0	0
3	A	60	0	52	3	0
3	B	60	0	52	3	0
4	B	38	0	34	3	0
5	A	129	0	0	6	0
5	B	77	0	0	11	0
All	All	9322	0	8913	491	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including

hydrogen atoms). The all-atom clashscore for this structure is 27.

All (491) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:361:LEU:HB3	1:B:454:PRO:HB2	1.30	1.08
1:A:221:ASN:HD22	1:A:224:GLY:H	1.02	0.99
1:A:488:ILE:HD11	1:A:570:GLU:HB3	1.40	0.98
1:A:342:ASN:HD22	1:A:344:GLN:HE21	1.14	0.96
1:B:363:SER:HB3	1:B:448:ILE:HD13	1.48	0.94
1:B:545:TRP:CZ2	1:B:658:VAL:HG11	2.03	0.94
1:A:296:GLU:HG2	1:A:594:GLU:HG2	1.50	0.92
1:A:278:ASN:HD21	1:A:280:ARG:HB2	1.37	0.90
1:A:374:LYS:H	1:A:374:LYS:HD2	1.37	0.88
1:A:221:ASN:ND2	1:A:224:GLY:H	1.73	0.87
1:B:545:TRP:HZ2	1:B:658:VAL:HG11	1.39	0.86
1:B:165:THR:HG22	1:B:182:ARG:HH21	1.41	0.85
1:A:478:ARG:HD3	1:A:568:GLN:NE2	1.92	0.85
1:A:364:GLY:O	1:A:377:SER:HB3	1.77	0.84
1:B:570:GLU:HG3	1:B:726:THR:OG1	1.77	0.84
1:A:342:ASN:ND2	1:A:344:GLN:HE21	1.76	0.83
1:A:300:LEU:HD12	1:A:301:THR:H	1.43	0.83
1:B:641:VAL:HG22	1:B:642:SER:H	1.46	0.81
1:A:557:THR:HG21	1:A:583:ALA:HA	1.62	0.81
1:B:225:LEU:H	1:B:225:LEU:HD12	1.43	0.81
1:A:280:ARG:HG3	1:A:482:ARG:HH11	1.46	0.80
1:B:696:THR:OG1	4:B:808:NAG:H82	1.81	0.79
1:A:464:ILE:HA	1:B:670:SER:HB2	1.65	0.78
1:A:278:ASN:ND2	1:A:280:ARG:HB2	1.98	0.76
1:A:365:ARG:HE	1:A:378:ARG:HG2	1.50	0.76
1:A:712:LEU:HD21	1:A:714:TRP:HE1	1.48	0.76
1:A:553:ALA:O	1:A:557:THR:HG22	1.85	0.76
1:A:323:LEU:HD13	1:A:551:LYS:HG3	1.68	0.76
1:A:478:ARG:HD3	1:A:568:GLN:HE22	1.51	0.75
1:A:472:VAL:H	1:A:566:ASN:HD21	1.33	0.75
1:B:430:PRO:O	1:B:512:ILE:HD11	1.88	0.74
1:B:340:ILE:HG23	1:B:342:ASN:H	1.52	0.74
1:A:202:ASN:ND2	1:A:204:HIS:HB2	2.03	0.74
1:B:464:ILE:HD11	3:B:803:NAG:H81	1.70	0.74
1:B:335:VAL:HG11	1:B:347:GLN:HG3	1.69	0.73
1:A:488:ILE:CD1	1:A:570:GLU:HB3	2.17	0.73
1:A:341:GLY:HA3	1:A:447:TYR:OH	1.89	0.73
1:A:380:CYS:HB2	1:A:385:VAL:HG21	1.71	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:407:PHE:O	1:B:411:ASP:HB2	1.87	0.73
1:B:362:VAL:HG21	1:B:393:LYS:HD3	1.70	0.73
1:B:313:PHE:CE2	1:B:480:PRO:HG2	2.25	0.72
1:B:199:ARG:HB2	1:B:199:ARG:HH11	1.53	0.72
1:A:671:ALA:HB2	1:A:677:ILE:HD12	1.72	0.72
1:B:669:ILE:HD13	1:B:677:ILE:HG22	1.72	0.72
1:B:367:ILE:HG22	1:B:369:ASN:ND2	2.05	0.72
1:A:703:ILE:HD13	1:A:727:ILE:HG22	1.71	0.71
1:B:121:PRO:O	1:B:124:VAL:HG23	1.90	0.71
1:B:271:SER:HB2	1:B:532:THR:HG21	1.71	0.71
1:A:313:PHE:HB3	1:A:482:ARG:HH21	1.55	0.71
1:A:552:SER:HB3	1:A:589:VAL:HG13	1.73	0.71
1:B:482:ARG:HA	1:B:482:ARG:HH11	1.55	0.71
1:B:358:TYR:HE1	1:B:456:ALA:H	1.38	0.70
1:A:231:ARG:HD2	5:A:918:HOH:O	1.90	0.70
1:A:478:ARG:HB2	1:A:571:PHE:O	1.92	0.70
1:B:369:ASN:HD22	1:B:386:ASN:H	1.39	0.70
1:B:481:ASN:HD22	1:B:481:ASN:C	1.95	0.69
1:A:421:ASN:O	3:A:805:BMA:H2	1.92	0.69
1:B:367:ILE:HG23	1:B:389:LEU:HD12	1.75	0.69
1:B:692:ARG:HH11	1:B:692:ARG:HB3	1.57	0.69
1:B:641:VAL:HG22	1:B:642:SER:N	2.07	0.69
1:A:452:ARG:C	1:A:454:PRO:HD3	2.13	0.69
1:B:486:ASP:HB2	1:B:629:VAL:HG11	1.73	0.69
1:B:552:SER:HB3	1:B:589:VAL:HG13	1.75	0.69
1:A:679:VAL:HG12	1:A:681:PRO:O	1.93	0.69
1:B:313:PHE:CD2	1:B:482:ARG:HD2	2.28	0.69
1:B:144:TRP:CZ3	1:B:146:GLU:HB2	2.28	0.68
1:B:333:LYS:HD3	3:B:803:NAG:H82	1.74	0.68
1:B:441:LEU:C	1:B:443:ALA:H	1.95	0.68
1:A:182:ARG:HG2	5:A:912:HOH:O	1.94	0.67
1:A:472:VAL:HG13	1:A:566:ASN:ND2	2.09	0.67
1:B:216:LEU:HD23	1:B:231:ARG:CB	2.24	0.67
1:A:500:LEU:HD23	1:A:500:LEU:C	2.14	0.67
1:A:669:ILE:HD13	1:A:677:ILE:HG22	1.76	0.67
1:B:403:PRO:O	1:B:406:PHE:HB2	1.93	0.67
1:B:497:VAL:HG12	1:B:498:GLU:HG3	1.75	0.67
1:A:641:VAL:HG12	1:A:642:SER:N	2.09	0.67
1:A:134:ILE:HD12	1:A:540:THR:HG22	1.76	0.67
1:A:478:ARG:CD	1:A:568:GLN:HE22	2.08	0.67
1:B:341:GLY:HA3	1:B:447:TYR:OH	1.94	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:509:VAL:HG13	1:A:519:ASN:ND2	2.10	0.67
1:B:394:ILE:HA	1:B:415:GLY:O	1.93	0.66
1:B:667:ALA:HB2	1:B:714:TRP:CZ3	2.30	0.66
1:A:221:ASN:HD22	1:A:224:GLY:N	1.86	0.66
1:A:399:ALA:O	1:A:418:MET:HB3	1.96	0.65
1:B:692:ARG:HH11	1:B:692:ARG:CB	2.08	0.65
1:B:369:ASN:HB3	1:B:386:ASN:HD22	1.61	0.65
1:B:481:ASN:ND2	1:B:483:ALA:H	1.94	0.65
1:B:628:ARG:HG2	1:B:629:VAL:H	1.61	0.65
1:A:509:VAL:HG13	1:A:519:ASN:HD21	1.60	0.65
1:B:199:ARG:NH1	1:B:199:ARG:HB2	2.11	0.65
1:A:188:ARG:HB3	1:A:189:PRO:CD	2.27	0.64
1:A:208:THR:HG23	1:A:529:PRO:HG3	1.78	0.64
1:A:156:SER:HB2	1:A:157:PRO:HD2	1.80	0.64
1:B:369:ASN:ND2	1:B:386:ASN:H	1.96	0.64
1:B:385:VAL:HG12	1:B:390:LEU:HD11	1.80	0.64
1:B:594:GLU:O	1:B:598:VAL:HG23	1.98	0.64
1:A:352:ASN:HD21	1:A:432:PRO:HA	1.62	0.63
1:A:528:CYS:HB3	1:A:529:PRO:HD3	1.80	0.63
1:B:444:THR:O	1:B:448:ILE:HG13	1.98	0.63
1:A:683:VAL:O	1:A:684:LEU:HD23	1.99	0.62
1:B:643:PRO:HG3	1:B:729:SER:HB2	1.80	0.62
1:B:406:PHE:CD2	1:B:430:PRO:HD2	2.34	0.62
1:B:223:TYR:H	1:B:225:LEU:HD12	1.62	0.62
1:A:310:PRO:O	1:A:478:ARG:HD2	2.00	0.61
1:B:408:LYS:HG2	1:B:409:SER:N	2.15	0.61
1:B:488:ILE:HG12	1:B:489:LYS:N	2.15	0.61
1:A:272:LEU:CD2	1:A:274:VAL:HG22	2.30	0.61
1:A:594:GLU:HA	1:A:597:TYR:HD1	1.65	0.61
1:A:328:SER:HA	1:A:472:VAL:HA	1.83	0.61
1:B:539:LYS:HE3	5:B:970:HOH:O	2.01	0.61
1:B:549:ALA:HA	1:B:589:VAL:HG11	1.83	0.61
1:A:361:LEU:HB2	1:A:454:PRO:HB2	1.83	0.61
1:B:478:ARG:HG3	1:B:568:GLN:OE1	2.00	0.61
1:B:368:PRO:HB3	1:B:372:PHE:O	2.02	0.60
1:A:419:THR:OG1	1:A:441:LEU:HD22	2.01	0.60
1:B:132:SER:HA	1:B:236:LEU:O	2.01	0.60
1:A:342:ASN:HD22	1:A:344:GLN:NE2	1.93	0.60
1:B:362:VAL:HG11	1:B:389:LEU:O	2.00	0.60
1:A:283:PHE:CD1	1:A:284:VAL:HG13	2.37	0.59
1:A:283:PHE:HD1	1:A:284:VAL:HG13	1.67	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:152:ASP:OD1	1:B:176:ARG:HB2	2.02	0.59
1:B:481:ASN:HD21	1:B:483:ALA:HB3	1.67	0.59
1:B:526:MET:O	1:B:529:PRO:HD2	2.01	0.59
1:A:280:ARG:HG3	1:A:482:ARG:HD3	1.84	0.59
1:B:134:ILE:HD12	1:B:540:THR:HG22	1.84	0.59
1:A:304:SER:HB3	1:A:524:THR:O	2.03	0.59
1:A:252:ASP:O	1:A:256:LEU:HG	2.02	0.59
1:B:143:ILE:HG21	1:B:178:ILE:HD13	1.84	0.59
1:B:386:ASN:HB2	1:B:388:ASN:ND2	2.18	0.59
1:B:464:ILE:HD11	3:B:803:NAG:C8	2.32	0.59
1:A:359:TYR:HB2	1:A:394:ILE:HG13	1.85	0.59
1:A:381:THR:HG22	1:A:382:ASP:N	2.17	0.59
1:A:557:THR:HG21	1:A:583:ALA:CA	2.32	0.58
1:B:526:MET:C	1:B:529:PRO:HD2	2.24	0.58
1:A:639:LEU:HD22	1:A:641:VAL:HG23	1.86	0.58
1:B:361:LEU:HA	1:B:394:ILE:CG2	2.34	0.58
1:B:367:ILE:HD13	1:B:389:LEU:HB2	1.85	0.58
1:B:387:PRO:O	1:B:391:LYS:HG3	2.03	0.58
1:B:217:VAL:HG11	1:B:518:PHE:CZ	2.38	0.58
1:B:383:LYS:HA	1:B:409:SER:HB3	1.86	0.58
1:B:369:ASN:CG	1:B:388:ASN:HD22	2.06	0.58
1:A:271:SER:HB2	1:A:532:THR:HG21	1.85	0.58
1:B:335:VAL:HG11	1:B:347:GLN:CG	2.33	0.58
1:A:549:ALA:HA	1:A:589:VAL:HG11	1.85	0.58
1:A:131:GLU:CG	1:A:234:VAL:HG13	2.34	0.58
1:B:362:VAL:HG21	1:B:393:LYS:CD	2.34	0.58
1:B:392:GLY:H	1:B:413:ALA:HA	1.68	0.58
1:A:300:LEU:HD12	1:A:301:THR:N	2.16	0.57
1:A:682:ASN:OD1	1:A:683:VAL:HG23	2.04	0.57
1:B:430:PRO:O	1:B:431:LEU:HD23	2.04	0.57
1:B:528:CYS:HB3	1:B:529:PRO:HD3	1.87	0.57
1:B:216:LEU:HD23	1:B:231:ARG:HB3	1.85	0.57
1:B:646:THR:HG22	1:B:647:PHE:N	2.20	0.57
1:B:680:ASN:HD22	1:B:681:PRO:HA	1.70	0.57
1:A:234:VAL:HG12	1:A:234:VAL:O	2.05	0.57
1:B:430:PRO:HG3	1:B:510:GLY:O	2.05	0.57
1:A:291:SER:O	1:A:295:VAL:HG23	2.05	0.57
1:B:222:LEU:HB3	1:B:225:LEU:HD13	1.86	0.57
1:B:377:SER:HA	1:B:384:SER:HB3	1.86	0.57
1:A:706:PHE:HA	1:A:729:SER:OG	2.05	0.56
1:B:658:VAL:O	1:B:658:VAL:HG12	2.03	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:493:SER:O	1:A:578:VAL:HG22	2.04	0.56
1:B:373:ASP:O	1:B:377:SER:HB2	2.06	0.56
1:A:398:GLU:HG2	1:A:441:LEU:HD21	1.88	0.56
1:B:304:SER:HB3	1:B:524:THR:O	2.05	0.56
1:B:519:ASN:ND2	1:B:520:ILE:H	2.04	0.56
1:B:475:PHE:CE1	1:B:523:GLY:HA2	2.41	0.56
1:B:337:GLN:HB3	5:B:966:HOH:O	2.05	0.56
1:A:684:LEU:HG	1:A:714:TRP:CZ3	2.41	0.55
1:B:340:ILE:HD11	1:B:444:THR:HA	1.88	0.55
1:B:304:SER:HB2	1:B:524:THR:OG1	2.07	0.55
1:B:598:VAL:O	1:B:598:VAL:HG12	2.06	0.55
1:B:480:PRO:HB3	1:B:571:PHE:CE1	2.41	0.55
1:B:593:ASN:ND2	1:B:595:SER:HB3	2.21	0.55
1:B:361:LEU:CB	1:B:454:PRO:HB2	2.20	0.55
1:A:448:ILE:HA	1:A:454:PRO:HG2	1.88	0.55
1:B:488:ILE:HG22	5:B:922:HOH:O	2.07	0.55
1:B:553:ALA:O	1:B:557:THR:HB	2.07	0.54
1:A:231:ARG:NH2	5:A:945:HOH:O	2.40	0.54
1:B:658:VAL:CG1	1:B:658:VAL:O	2.54	0.54
1:A:646:THR:HA	1:A:702:SER:HB3	1.88	0.54
1:B:509:VAL:HB	1:B:514:ARG:HG3	1.90	0.54
1:A:451:ILE:O	1:A:452:ARG:HB2	2.08	0.54
1:B:728:THR:HG22	1:B:730:LEU:H	1.72	0.54
1:A:218:SER:O	1:A:219:GLN:HB2	2.06	0.54
1:A:349:VAL:O	1:A:350:SER:HB3	2.08	0.54
1:B:165:THR:HG22	1:B:182:ARG:NH2	2.19	0.54
1:A:453:SER:N	1:A:454:PRO:HD3	2.23	0.54
1:A:478:ARG:CB	1:A:571:PHE:O	2.56	0.54
1:B:186:ILE:HD12	1:B:253:THR:HG22	1.88	0.54
1:B:569:ALA:HB1	1:B:570:GLU:OE2	2.08	0.54
1:B:319:SER:HB3	1:B:321:TRP:CE2	2.43	0.54
1:A:639:LEU:HD22	1:A:641:VAL:CG2	2.39	0.53
1:B:680:ASN:ND2	1:B:681:PRO:HA	2.23	0.53
1:A:712:LEU:HD21	1:A:714:TRP:NE1	2.21	0.53
1:B:481:ASN:HD22	1:B:483:ALA:H	1.56	0.53
1:A:632:LEU:O	1:A:634:TYR:N	2.38	0.53
1:A:119:GLY:C	1:A:121:PRO:HD3	2.29	0.53
1:B:341:GLY:HA3	1:B:447:TYR:CZ	2.43	0.53
1:B:329:THR:HG23	1:B:471:VAL:O	2.08	0.53
1:A:162:TRP:C	1:A:163:LYS:HD2	2.29	0.53
1:A:669:ILE:HG22	1:A:670:SER:N	2.24	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:142:GLY:C	1:A:200:ASP:HB2	2.28	0.53
1:B:188:ARG:NH2	1:B:254:ASP:OD2	2.42	0.53
1:A:322:LEU:O	1:A:489:LYS:NZ	2.42	0.53
1:B:556:THR:HA	1:B:722:ARG:O	2.08	0.53
1:A:557:THR:HG23	5:A:907:HOH:O	2.09	0.53
1:A:570:GLU:C	1:A:572:ALA:H	2.12	0.53
1:B:488:ILE:HG23	5:B:903:HOH:O	2.09	0.53
1:B:512:ILE:HG22	1:B:513:ARG:N	2.24	0.53
1:A:594:GLU:HA	1:A:597:TYR:CD1	2.43	0.52
1:A:641:VAL:HG12	1:A:642:SER:H	1.74	0.52
1:B:272:LEU:CD1	1:B:274:VAL:HG12	2.40	0.52
1:B:385:VAL:CG1	1:B:390:LEU:HD11	2.38	0.52
1:A:186:ILE:HD12	1:A:253:THR:HG22	1.90	0.52
1:A:448:ILE:HA	1:A:454:PRO:CG	2.38	0.52
1:A:256:LEU:HD21	1:A:286:ALA:HB1	1.92	0.52
1:A:481:ASN:HB3	1:A:484:THR:O	2.09	0.52
1:A:567:PRO:HG2	3:A:806:MAN:O4	2.09	0.52
1:B:170:ASN:HD22	1:B:197:GLY:HA3	1.73	0.52
1:A:374:LYS:HD2	1:A:374:LYS:N	2.15	0.52
1:A:117:PHE:CE2	1:A:330:MET:HA	2.45	0.52
1:B:354:PHE:CD1	1:B:432:PRO:HD3	2.45	0.52
1:B:441:LEU:C	1:B:443:ALA:N	2.62	0.52
1:B:323:LEU:HD13	1:B:551:LYS:HG3	1.91	0.52
1:B:119:GLY:C	1:B:121:PRO:HD3	2.30	0.52
1:B:223:TYR:H	1:B:225:LEU:CD1	2.23	0.52
1:A:273:SER:HB3	1:A:525:SER:HB2	1.92	0.52
1:A:486:ASP:OD1	1:A:639:LEU:HA	2.10	0.51
1:B:167:GLU:OE1	1:B:198:PRO:HD3	2.11	0.51
1:B:692:ARG:HH11	1:B:692:ARG:CG	2.23	0.51
1:A:451:ILE:O	1:A:451:ILE:HG22	2.10	0.51
1:A:374:LYS:CD	1:A:374:LYS:H	2.15	0.51
1:A:665:TYR:HD2	1:A:714:TRP:O	1.94	0.51
1:A:272:LEU:HD21	1:A:274:VAL:HG22	1.91	0.51
1:B:340:ILE:CD1	1:B:444:THR:HA	2.40	0.51
1:A:362:VAL:HG13	1:A:393:LYS:HD3	1.92	0.51
1:A:641:VAL:CG1	1:A:642:SER:N	2.74	0.51
1:B:575:SER:HA	5:B:918:HOH:O	2.09	0.51
1:B:182:ARG:HH11	1:B:261:ASP:HB3	1.74	0.51
1:B:355:ASP:O	1:B:357:GLN:HG3	2.11	0.51
1:B:519:ASN:HD22	1:B:520:ILE:H	1.59	0.51
1:A:646:THR:HA	1:A:701:GLY:O	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:485:LYS:HE3	1:B:640:SER:OG	2.11	0.50
1:B:549:ALA:HA	1:B:589:VAL:CG1	2.41	0.50
1:A:312:PHE:O	1:A:479:GLY:O	2.28	0.50
1:A:728:THR:HG22	1:A:730:LEU:H	1.76	0.50
1:B:141:THR:HG22	1:B:245:CYS:HB2	1.93	0.50
1:B:416:VAL:O	1:B:417:LEU:HD23	2.12	0.50
1:B:588:LEU:HD13	1:B:655:LEU:HD13	1.94	0.50
1:A:360:PRO:HB2	1:A:393:LYS:HG2	1.93	0.50
1:A:119:GLY:O	1:A:121:PRO:HD3	2.11	0.50
1:A:521:ILE:HG13	1:A:526:MET:HE3	1.94	0.50
1:A:472:VAL:N	1:A:566:ASN:HD21	2.07	0.50
1:A:641:VAL:HG11	1:A:647:PHE:CD2	2.47	0.50
1:B:497:VAL:CG1	1:B:498:GLU:HG3	2.41	0.50
1:A:184:TYR:OH	1:A:261:ASP:OD1	2.26	0.49
1:B:272:LEU:HG	1:B:274:VAL:HG12	1.94	0.49
1:B:340:ILE:HG22	1:B:344:GLN:HB2	1.94	0.49
1:B:646:THR:CG2	1:B:647:PHE:N	2.74	0.49
1:B:222:LEU:HB3	1:B:225:LEU:CD1	2.43	0.49
1:B:481:ASN:HD22	1:B:482:ARG:N	2.09	0.49
1:B:472:VAL:HG11	1:B:572:ALA:HB1	1.94	0.49
1:B:318:LEU:HD22	1:B:481:ASN:ND2	2.27	0.49
1:A:202:ASN:HD22	1:A:204:HIS:HB2	1.73	0.49
1:A:639:LEU:HD23	1:A:640:SER:N	2.28	0.49
1:A:706:PHE:O	1:A:728:THR:HA	2.12	0.49
1:B:195:VAL:HG12	1:B:196:ASN:O	2.13	0.49
1:B:383:LYS:C	1:B:385:VAL:H	2.17	0.49
1:A:379:PHE:CD1	1:A:379:PHE:N	2.81	0.48
1:B:165:THR:CG2	1:B:182:ARG:HH21	2.20	0.48
1:B:164:GLY:HA2	1:B:265:ASP:OD1	2.13	0.48
1:B:641:VAL:HG21	1:B:647:PHE:CD2	2.48	0.48
1:A:403:PRO:O	1:A:406:PHE:HB2	2.14	0.48
1:B:259:TYR:O	1:B:263:ILE:HG13	2.14	0.48
1:B:441:LEU:O	1:B:445:LEU:HG	2.13	0.48
1:B:676:THR:HB	1:B:700:ARG:HG2	1.95	0.48
1:B:453:SER:N	1:B:454:PRO:HD3	2.29	0.48
1:A:111:THR:HB	1:A:500:LEU:O	2.13	0.48
1:A:500:LEU:HB2	1:A:520:ILE:HG12	1.96	0.48
1:A:570:GLU:H	1:A:570:GLU:CD	2.17	0.48
1:B:162:TRP:NE1	1:B:179:ILE:O	2.44	0.48
1:A:367:ILE:HG22	1:A:367:ILE:O	2.13	0.48
1:B:710:ALA:HB3	1:B:725:ILE:HB	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:202:ASN:HD21	1:A:204:HIS:HB2	1.79	0.48
1:A:131:GLU:CD	1:A:234:VAL:HG13	2.34	0.48
1:A:304:SER:HB2	1:A:524:THR:OG1	2.14	0.47
1:B:447:TYR:OH	1:B:455:GLY:N	2.40	0.47
1:B:641:VAL:CG2	1:B:642:SER:N	2.77	0.47
1:B:650:TYR:HE1	1:B:652:ASN:OD1	1.97	0.47
1:B:317:SER:OG	1:B:322:LEU:HD23	2.14	0.47
1:B:389:LEU:O	1:B:390:LEU:HD23	2.14	0.47
1:A:473:VAL:HG11	1:A:475:PHE:CZ	2.50	0.47
1:B:225:LEU:HD11	1:B:353:THR:HG22	1.97	0.47
1:B:354:PHE:CB	1:B:432:PRO:HG3	2.44	0.47
1:B:551:LYS:HD3	5:B:950:HOH:O	2.13	0.47
1:A:729:SER:O	1:A:730:LEU:HB3	2.14	0.47
1:A:378:ARG:HD2	1:A:379:PHE:CE1	2.50	0.47
1:B:166:CYS:HA	1:B:181:ALA:O	2.15	0.47
1:B:641:VAL:CG2	1:B:642:SER:H	2.21	0.47
1:A:164:GLY:HA2	1:A:265:ASP:OD1	2.14	0.47
1:A:374:LYS:HA	1:A:377:SER:HB2	1.97	0.47
1:B:481:ASN:ND2	1:B:481:ASN:C	2.67	0.47
1:B:340:ILE:HD11	1:B:444:THR:HG23	1.97	0.47
1:A:367:ILE:HG21	1:A:389:LEU:HB2	1.96	0.46
1:B:369:ASN:HB3	1:B:386:ASN:ND2	2.29	0.46
1:B:493:SER:O	1:B:578:VAL:HG23	2.15	0.46
1:B:252:ASP:HB3	1:B:286:ALA:HB2	1.97	0.46
1:B:349:VAL:O	1:B:350:SER:HB3	2.15	0.46
1:A:472:VAL:H	1:A:566:ASN:ND2	2.06	0.46
1:B:188:ARG:NH1	1:B:248:ASP:OD2	2.48	0.46
1:B:367:ILE:N	1:B:368:PRO:HD3	2.30	0.46
1:A:367:ILE:CG2	1:A:389:LEU:HB2	2.46	0.46
1:A:202:ASN:HD22	1:A:204:HIS:CB	2.29	0.46
1:B:272:LEU:HD11	1:B:274:VAL:HG12	1.98	0.46
1:B:378:ARG:HD2	1:B:398:GLU:OE2	2.16	0.46
1:A:342:ASN:ND2	1:A:344:GLN:NE2	2.54	0.46
1:B:287:ILE:HG13	5:B:917:HOH:O	2.15	0.46
1:B:703:ILE:HD13	1:B:727:ILE:HG22	1.96	0.46
1:B:520:ILE:CG1	1:B:520:ILE:O	2.64	0.46
1:A:381:THR:HG22	1:A:382:ASP:H	1.78	0.46
1:A:662:ALA:HA	1:A:686:PHE:O	2.16	0.46
1:B:156:SER:HB2	1:B:157:PRO:CD	2.46	0.46
1:B:174:CYS:HB2	1:B:178:ILE:O	2.16	0.46
1:B:203:GLY:O	1:B:206:THR:N	2.48	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:481:ASN:ND2	1:B:483:ALA:N	2.63	0.46
1:B:181:ALA:HB1	1:B:198:PRO:HB3	1.97	0.45
1:B:507:ALA:HA	1:B:508:PRO:HD3	1.83	0.45
1:B:590:TYR:CZ	1:B:655:LEU:HD21	2.51	0.45
1:B:114:SER:HA	1:B:330:MET:HE2	1.97	0.45
1:A:362:VAL:CG1	1:A:393:LYS:HD3	2.47	0.45
1:A:280:ARG:CG	1:A:482:ARG:HH11	2.21	0.45
1:B:486:ASP:CB	1:B:629:VAL:HG11	2.45	0.45
1:B:409:SER:O	1:B:410:LEU:HD23	2.17	0.45
1:A:334:PHE:HD1	1:A:498:GLU:OE1	1.99	0.45
1:B:365:ARG:HB3	1:B:378:ARG:HB3	1.99	0.45
1:B:367:ILE:CG2	1:B:369:ASN:ND2	2.76	0.45
1:A:534:ILE:CD1	1:A:580:PRO:HG3	2.47	0.45
1:B:351:ILE:HD13	1:B:520:ILE:CG1	2.46	0.45
1:B:488:ILE:HA	5:B:913:HOH:O	2.17	0.45
1:B:536:THR:O	1:B:540:THR:HG23	2.16	0.45
1:B:628:ARG:CG	1:B:629:VAL:H	2.28	0.45
1:A:641:VAL:CG1	1:A:642:SER:H	2.30	0.45
1:A:500:LEU:HA	1:A:519:ASN:O	2.17	0.45
1:B:568:GLN:HG3	5:B:948:HOH:O	2.17	0.45
1:B:487:VAL:HG23	1:B:629:VAL:HG13	1.98	0.45
1:A:163:LYS:HD2	1:A:163:LYS:N	2.32	0.44
1:A:438:PRO:O	1:A:441:LEU:HB3	2.18	0.44
1:B:358:TYR:OH	1:B:455:GLY:HA2	2.17	0.44
1:B:198:PRO:O	1:B:199:ARG:C	2.56	0.44
1:B:488:ILE:HG21	5:B:928:HOH:O	2.17	0.44
1:A:542:ASN:N	1:A:543:PRO:HD3	2.32	0.44
1:A:647:PHE:HE1	1:A:699:VAL:HG12	1.81	0.44
1:A:729:SER:O	1:A:730:LEU:CB	2.66	0.44
1:B:315:THR:CA	1:B:477:SER:HB3	2.47	0.44
1:A:121:PRO:HD2	1:A:124:VAL:CG1	2.48	0.44
1:A:365:ARG:HH21	1:A:378:ARG:HD3	1.83	0.44
1:A:381:THR:HG23	1:A:405:GLU:CD	2.37	0.44
1:A:361:LEU:HD11	1:A:456:ALA:HB2	1.99	0.44
1:A:645:GLN:O	1:A:646:THR:C	2.56	0.44
1:B:168:THR:HG22	1:B:168:THR:O	2.17	0.44
1:B:315:THR:HA	1:B:477:SER:HB3	1.98	0.44
1:A:234:VAL:CG1	1:A:237:ALA:HB2	2.47	0.44
1:B:210:SER:O	1:B:214:GLY:N	2.46	0.44
1:B:375:SER:C	1:B:377:SER:H	2.20	0.44
1:B:628:ARG:HG2	1:B:629:VAL:HG23	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:387:PRO:HB2	1:B:391:LYS:HE3	1.99	0.44
1:B:647:PHE:CZ	1:B:727:ILE:HD13	2.53	0.44
1:B:650:TYR:HH	4:B:808:NAG:HO6	1.65	0.44
1:A:188:ARG:CB	1:A:189:PRO:CD	2.95	0.44
1:A:562:ASN:OD1	1:A:564:ARG:HB3	2.18	0.44
1:A:673:GLN:C	1:A:675:LEU:H	2.20	0.44
1:B:351:ILE:CD1	1:B:520:ILE:HD11	2.47	0.44
1:A:551:LYS:HD3	5:A:946:HOH:O	2.18	0.44
1:A:177:LYS:HE3	5:A:985:HOH:O	2.18	0.43
1:A:507:ALA:HA	1:A:508:PRO:HD3	1.90	0.43
1:A:672:PRO:HD3	1:A:710:ALA:HB2	1.99	0.43
1:B:339:GLN:O	1:B:456:ALA:HA	2.18	0.43
1:B:522:SER:O	1:B:526:MET:CE	2.66	0.43
1:A:341:GLY:HA3	1:A:447:TYR:CZ	2.52	0.43
1:A:666:ARG:NH1	1:A:666:ARG:HB2	2.32	0.43
1:B:190:ILE:HD11	1:B:196:ASN:OD1	2.18	0.43
1:A:419:THR:HG21	1:A:438:PRO:HA	2.01	0.43
1:A:669:ILE:CG2	1:A:670:SER:N	2.81	0.43
1:B:519:ASN:ND2	1:B:520:ILE:N	2.66	0.43
1:B:534:ILE:HD13	1:B:578:VAL:HG11	1.99	0.43
1:A:279:PRO:O	1:A:282:TYR:HD2	2.02	0.43
1:B:339:GLN:HG3	1:B:344:GLN:O	2.19	0.43
1:B:728:THR:HG22	1:B:730:LEU:N	2.33	0.43
4:B:808:NAG:O7	4:B:809:FUC:C1	2.67	0.43
1:A:155:PHE:CE2	1:A:177:LYS:HD3	2.53	0.43
1:A:699:VAL:HG11	1:A:727:ILE:HD11	2.00	0.43
1:B:208:THR:HG22	1:B:209:ALA:N	2.33	0.43
1:B:416:VAL:HG12	1:B:417:LEU:N	2.34	0.43
1:B:482:ARG:CA	1:B:482:ARG:HH11	2.28	0.43
1:A:234:VAL:HG11	1:A:237:ALA:HB2	2.01	0.43
1:A:190:ILE:HD12	1:A:246:TRP:CZ3	2.53	0.43
1:B:272:LEU:HD11	1:B:274:VAL:CG1	2.49	0.43
1:B:315:THR:HG22	1:B:478:ARG:O	2.18	0.43
1:A:121:PRO:O	1:A:124:VAL:HG13	2.18	0.43
1:A:500:LEU:C	1:A:500:LEU:CD2	2.87	0.43
1:A:552:SER:CB	1:A:589:VAL:HG13	2.45	0.43
1:A:650:TYR:HE1	1:A:652:ASN:ND2	2.17	0.43
1:B:381:THR:HG22	1:B:382:ASP:N	2.34	0.43
1:A:280:ARG:HG3	1:A:482:ARG:CD	2.49	0.43
1:B:175:ASN:OD1	1:B:177:LYS:N	2.49	0.43
1:B:497:VAL:O	1:B:499:ILE:HD12	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:121:PRO:HD2	1:A:124:VAL:HG12	2.01	0.43
1:A:280:ARG:HG3	1:A:482:ARG:NH1	2.25	0.43
1:A:541:TYR:C	1:A:543:PRO:HD3	2.39	0.43
1:B:217:VAL:HG11	1:B:518:PHE:HZ	1.82	0.42
1:A:295:VAL:HG21	1:A:321:TRP:HB2	2.00	0.42
1:A:223:TYR:CE1	1:A:514:ARG:HB3	2.54	0.42
1:B:453:SER:H	1:B:454:PRO:HD3	1.84	0.42
1:A:208:THR:HG23	1:A:529:PRO:CG	2.46	0.42
1:B:472:VAL:HG13	1:B:566:ASN:CG	2.39	0.42
1:A:545:TRP:HZ3	1:A:553:ALA:HB2	1.85	0.42
1:B:351:ILE:HD13	1:B:520:ILE:HG12	2.00	0.42
1:A:557:THR:OG1	1:A:582:LYS:HB3	2.19	0.42
1:B:390:LEU:O	1:B:391:LYS:C	2.57	0.42
1:A:728:THR:HG22	1:A:730:LEU:N	2.34	0.42
1:B:675:LEU:HD11	1:B:708:VAL:HG11	2.02	0.42
1:A:588:LEU:HD23	1:A:657:SER:HA	2.02	0.42
1:B:378:ARG:HG3	1:B:379:PHE:CD1	2.54	0.42
1:A:589:VAL:HG12	1:A:590:TYR:N	2.35	0.42
1:B:207:HIS:CE1	1:B:501:ALA:HB3	2.55	0.42
1:B:424:ASP:O	1:B:425:TYR:HB3	2.20	0.42
1:B:588:LEU:HB2	1:B:721:VAL:HG21	2.02	0.42
1:A:151:ASP:OD1	1:A:152:ASP:N	2.52	0.42
1:A:155:PHE:CZ	1:A:177:LYS:HD3	2.55	0.42
3:A:802:NAG:O5	3:A:804:FUC:H5	2.20	0.42
1:A:419:THR:HG22	1:A:420:SER:N	2.35	0.42
1:A:449:TYR:O	1:A:449:TYR:CG	2.73	0.42
1:A:356:ASN:OD1	1:A:459:PHE:HA	2.20	0.42
1:B:522:SER:C	1:B:526:MET:HE3	2.39	0.42
1:B:676:THR:HB	1:B:700:ARG:CG	2.50	0.42
1:A:509:VAL:CG1	1:A:519:ASN:ND2	2.81	0.41
1:A:674:GLY:O	1:A:675:LEU:HD23	2.20	0.41
1:A:713:VAL:HG22	1:A:722:ARG:CB	2.50	0.41
1:B:430:PRO:HA	1:B:512:ILE:CD1	2.50	0.41
1:A:152:ASP:O	1:A:153:GLU:C	2.58	0.41
1:A:369:ASN:HB2	1:A:385:VAL:C	2.40	0.41
1:A:650:TYR:HD1	1:A:696:THR:CG2	2.34	0.41
1:B:113:ARG:HA	1:B:113:ARG:HD2	1.90	0.41
1:B:375:SER:C	1:B:377:SER:N	2.74	0.41
1:B:361:LEU:HD11	1:B:396:VAL:CG2	2.51	0.41
1:B:500:LEU:HD23	1:B:500:LEU:C	2.41	0.41
1:A:367:ILE:N	1:A:368:PRO:HD3	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:478:ARG:CG	1:A:568:GLN:HE22	2.33	0.41
1:B:408:LYS:CG	1:B:409:SER:N	2.81	0.41
1:A:134:ILE:HG21	1:A:536:THR:HG23	2.03	0.41
1:A:536:THR:O	1:A:540:THR:HG23	2.20	0.41
1:B:293:HIS:O	1:B:296:GLU:HB3	2.20	0.41
1:B:386:ASN:O	1:B:388:ASN:N	2.53	0.41
1:B:455:GLY:O	1:B:456:ALA:HB2	2.21	0.41
1:A:381:THR:CG2	1:A:382:ASP:N	2.83	0.41
1:B:207:HIS:ND1	1:B:501:ALA:HB3	2.35	0.41
1:A:364:GLY:HA3	1:A:377:SER:O	2.21	0.41
1:A:630:TRP:HB3	1:A:651:PHE:CE2	2.55	0.41
1:B:125:PRO:O	1:B:126:ARG:HG3	2.20	0.41
1:B:318:LEU:HD22	1:B:481:ASN:CG	2.41	0.41
1:A:143:ILE:O	1:A:145:PRO:HD3	2.21	0.41
1:A:488:ILE:CD1	1:A:570:GLU:CB	2.96	0.41
1:B:324:SER:OG	1:B:489:LYS:HE3	2.21	0.41
1:B:516:THR:OG1	1:B:517:LEU:N	2.54	0.41
1:B:531:ILE:HD11	1:B:578:VAL:HG21	2.02	0.41
1:A:475:PHE:CD1	1:A:475:PHE:C	2.95	0.41
1:A:650:TYR:HE1	1:A:652:ASN:HD21	1.69	0.41
1:B:354:PHE:HD1	1:B:432:PRO:CD	2.34	0.41
1:B:520:ILE:HG13	1:B:520:ILE:O	2.18	0.41
1:B:596:ASP:HB3	5:B:909:HOH:O	2.20	0.41
1:B:687:ASN:HB2	1:B:691:ASP:OD2	2.21	0.41
1:A:304:SER:HB2	1:A:524:THR:CB	2.51	0.40
1:A:590:TYR:CZ	1:A:655:LEU:HD21	2.57	0.40
1:B:379:PHE:HD1	1:B:379:PHE:H	1.69	0.40
1:B:448:ILE:HG22	1:B:448:ILE:O	2.21	0.40
1:B:499:ILE:O	1:B:520:ILE:HA	2.21	0.40
1:A:156:SER:O	1:A:176:ARG:HD3	2.20	0.40
1:A:188:ARG:HB3	1:A:189:PRO:HD2	2.02	0.40
1:A:272:LEU:HD23	1:A:274:VAL:HG13	2.02	0.40
1:B:124:VAL:HG13	1:B:581:LEU:HD21	2.03	0.40
1:B:165:THR:OG1	1:B:166:CYS:N	2.53	0.40
1:B:187:GLY:O	1:B:188:ARG:HB3	2.21	0.40
1:B:419:THR:CG2	1:B:420:SER:N	2.84	0.40
1:A:643:PRO:HG3	1:A:729:SER:HB2	2.02	0.40
1:B:272:LEU:CG	1:B:274:VAL:HG12	2.51	0.40
1:B:300:LEU:HG	1:B:301:THR:N	2.37	0.40
1:B:354:PHE:CD1	1:B:432:PRO:CD	3.05	0.40
1:B:369:ASN:HD22	1:B:386:ASN:N	2.12	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:528:CYS:N	1:A:529:PRO:CD	2.85	0.40
1:B:419:THR:HG22	1:B:420:SER:N	2.35	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	587/621 (94%)	527 (90%)	50 (8%)	10 (2%)	11	30
1	B	590/621 (95%)	497 (84%)	68 (12%)	25 (4%)	3	9
All	All	1177/1242 (95%)	1024 (87%)	118 (10%)	35 (3%)	5	15

All (35) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	633	ASN
1	B	375	SER
1	B	388	ASN
1	B	391	LYS
1	B	408	LYS
1	B	450	SER
1	A	187	GLY
1	A	454	PRO
1	B	171	ASN
1	B	267	VAL
1	B	355	ASP
1	B	369	ASN
1	B	451	ILE
1	B	452	ARG
1	B	453	SER
1	B	629	VAL

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Mol	Chain	Res	Type
1	A	312	PHE
1	A	646	THR
1	A	702	SER
1	B	153	GLU
1	B	350	SER
1	B	456	ALA
1	A	350	SER
1	B	384	SER
1	B	145	PRO
1	B	199	ARG
1	B	376	THR
1	B	387	PRO
1	B	635	PRO
1	A	186	ILE
1	A	672	PRO
1	B	432	PRO
1	B	580	PRO
1	A	674	GLY
1	B	244	VAL

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	489/512 (96%)	467 (96%)	22 (4%)	32	63
1	B	492/512 (96%)	471 (96%)	21 (4%)	33	64
All	All	981/1024 (96%)	938 (96%)	43 (4%)	33	64

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	111	THR
1	A	122	LEU
1	A	152	ASP
1	A	188	ARG

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Mol	Chain	Res	Type
1	A	196	ASN
1	A	216	LEU
1	A	251	SER
1	A	259	TYR
1	A	272	LEU
1	A	274	VAL
1	A	297	ARG
1	A	379	PHE
1	A	437	ASP
1	A	472	VAL
1	A	519	ASN
1	A	578	VAL
1	A	579	ASN
1	A	594	GLU
1	A	630	TRP
1	A	639	LEU
1	A	723	SER
1	A	728	THR
1	B	208	THR
1	B	225	LEU
1	B	227	LEU
1	B	231	ARG
1	B	259	TYR
1	B	304	SER
1	B	310	PRO
1	B	331	ASP
1	B	332	ARG
1	B	460	LYS
1	B	472	VAL
1	B	481	ASN
1	B	482	ARG
1	B	488	ILE
1	B	520	ILE
1	B	521	ILE
1	B	525	SER
1	B	557	THR
1	B	564	ARG
1	B	579	ASN
1	B	692	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:



Mol	Chain	Res	Type
1	A	202	ASN
1	A	221	ASN
1	A	278	ASN
1	A	303	ASN
1	A	339	GLN
1	A	342	ASN
1	A	352	ASN
1	A	386	ASN
1	A	519	ASN
1	A	566	ASN
1	A	568	GLN
1	A	652	ASN
1	A	719	HIS
1	B	311	ASN
1	B	369	ASN
1	B	386	ASN
1	B	388	ASN
1	B	481	ASN
1	B	519	ASN
1	B	577	HIS
1	B	579	ASN
1	B	680	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 ⓘ

13 carbohydrates are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	NAG	A	802	3	14,14,15	0.51	0	15,19,21	0.95	1 (6%)
3	NAG	A	803	1,3	14,14,15	0.48	0	15,19,21	0.71	0
3	FUC	A	804	3	9,10,11	0.51	0	13,14,16	0.40	0
3	BMA	A	805	3	11,11,12	0.42	0	13,15,17	0.52	0
3	MAN	A	806	3	11,11,12	0.45	0	13,15,17	0.90	1 (7%)
3	NAG	B	802	3	14,14,15	0.59	0	15,19,21	0.72	0
3	NAG	B	803	1,3	14,14,15	0.47	0	15,19,21	0.97	1 (6%)
3	FUC	B	804	3	9,10,11	0.45	0	13,14,16	0.42	0
3	BMA	B	805	3	11,11,12	0.60	0	13,15,17	0.68	0
3	MAN	B	806	3	11,11,12	0.56	0	13,15,17	0.88	1 (7%)
4	NAG	B	807	4	14,14,15	0.98	0	15,19,21	0.68	0
4	NAG	B	808	1,4	14,14,15	1.06	0	15,19,21	1.10	1 (6%)
4	FUC	B	809	4	9,10,11	0.89	0	13,14,16	1.15	1 (7%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	A	802	3	-	0/6/23/26	0/1/1/1
3	NAG	A	803	1,3	-	0/6/23/26	0/1/1/1
3	FUC	A	804	3	-	0/0/17/20	0/1/1/1
3	BMA	A	805	3	-	0/2/19/22	0/1/1/1
3	MAN	A	806	3	-	0/2/19/22	0/1/1/1
3	NAG	B	802	3	-	0/6/23/26	0/1/1/1
3	NAG	B	803	1,3	-	1/6/23/26	0/1/1/1
3	FUC	B	804	3	-	0/0/17/20	0/1/1/1
3	BMA	B	805	3	-	0/2/19/22	0/1/1/1
3	MAN	B	806	3	-	0/2/19/22	0/1/1/1
4	NAG	B	807	4	-	0/6/23/26	0/1/1/1
4	NAG	B	808	1,4	-	0/6/23/26	0/1/1/1
4	FUC	B	809	4	-	0/0/17/20	0/1/1/1

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	803	NAG	C2-N2-C7	-2.40	119.44	122.94
3	A	802	NAG	C2-N2-C7	-2.30	119.59	122.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	808	NAG	O4-C4-C3	-2.02	105.96	110.36
3	A	806	MAN	C1-O5-C5	2.44	115.54	112.17
3	B	806	MAN	C1-O5-C5	2.69	115.88	112.17
4	B	809	FUC	C1-C2-C3	3.23	113.75	109.65

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	803	NAG	O7-C7-N2-C2

There are no ring outliers.

7 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	802	NAG	1	0
3	A	804	FUC	1	0
3	A	805	BMA	1	0
3	A	806	MAN	1	0
3	B	803	NAG	3	0
4	B	808	NAG	3	0
4	B	809	FUC	1	0

## 5.6 Ligand geometry [i](#)

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
2	DFP	A	801	1	6,9,9	1.35	1 (16%)	6,11,11	0.48	0
2	DFP	B	801	1	6,9,9	1.10	1 (16%)	6,11,11	0.41	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	DFP	A	801	1	-	0/4/8/8	0/0/0/0
2	DFP	B	801	1	-	0/4/8/8	0/0/0/0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	801	DFP	O1P-C1	-2.26	1.43	1.46
2	B	801	DFP	O1P-C1	-2.18	1.43	1.46

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.7 Other polymers [i](#)

There are no such residues in this entry.

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

There are no chain breaks in this entry.

## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	591/621 (95%)	0.19	5 (0%) 86 83	13, 38, 78, 97	0
1	B	594/621 (95%)	0.45	37 (6%) 21 16	22, 45, 116, 140	0
All	All	1185/1242 (95%)	0.32	42 (3%) 44 38	13, 41, 98, 140	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	372	PHE	6.8
1	B	600	PHE	5.0
1	B	370	THR	5.0
1	B	601	LEU	4.6
1	B	452	ARG	4.4
1	B	371	GLY	4.0
1	B	389	LEU	3.8
1	B	391	LYS	3.3
1	B	373	ASP	3.3
1	B	458	ILE	3.3
1	B	394	ILE	3.2
1	B	453	SER	3.1
1	B	457	THR	3.0
1	B	396	VAL	3.0
1	B	456	ALA	2.9
1	B	365	ARG	2.9
1	B	343	GLY	2.8
1	B	392	GLY	2.8
1	B	415	GLY	2.8
1	B	362	VAL	2.7
1	B	374	LYS	2.7
1	B	454	PRO	2.6
1	B	406	PHE	2.5
1	B	357	GLN	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	361	LEU	2.5
1	B	277	ALA	2.4
1	A	275	GLY	2.3
1	A	452	ARG	2.3
1	B	282	TYR	2.3
1	B	449	TYR	2.3
1	A	629	VAL	2.3
1	B	395	VAL	2.3
1	B	432	PRO	2.3
1	A	630	TRP	2.3
1	B	390	LEU	2.2
1	B	173	ARG	2.2
1	B	352	ASN	2.1
1	B	393	LYS	2.1
1	B	383	LYS	2.0
1	A	390	LEU	2.0
1	B	367	ILE	2.0
1	B	381	THR	2.0

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

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

## 6.3 Carbohydrates [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
4	NAG	B	808	14/15	0.81	0.21	2.78	70,75,80,80	0
3	NAG	A	802	14/15	0.96	0.17	-0.58	40,49,53,55	0
3	NAG	B	802	14/15	0.95	0.17	-0.74	54,57,62,62	0
3	NAG	A	803	14/15	0.96	0.14	-1.18	34,39,43,44	0
3	NAG	B	803	14/15	0.96	0.15	-1.26	47,51,56,57	0
3	MAN	B	806	11/12	0.77	0.22	-	80,83,84,85	0
3	FUC	B	804	10/11	0.95	0.19	-	57,59,61,62	0
3	BMA	B	805	11/12	0.93	0.15	-	65,68,73,77	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
4	FUC	B	809	10/11	0.76	0.25	-	78,80,81,81	0
3	FUC	A	804	10/11	0.96	0.16	-	42,47,48,49	0
4	NAG	B	807	14/15	0.82	0.23	-	81,84,85,86	0
3	BMA	A	805	11/12	0.92	0.15	-	60,65,68,72	0
3	MAN	A	806	11/12	0.84	0.22	-	77,79,82,83	0

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	DFP	B	801	10/10	0.94	0.29	3.07	53,57,63,64	0
2	DFP	A	801	10/10	0.95	0.27	1.58	34,38,44,46	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.