



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:17 PM GMT

PDB ID : 1OAN
Title : CRYSTAL STRUCTURE OF THE DENGUE 2 VIRUS ENVELOPE PROTEIN
Authors : Modis, Y.; Harrison, S.C.
Deposited on : 2003-01-16
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
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

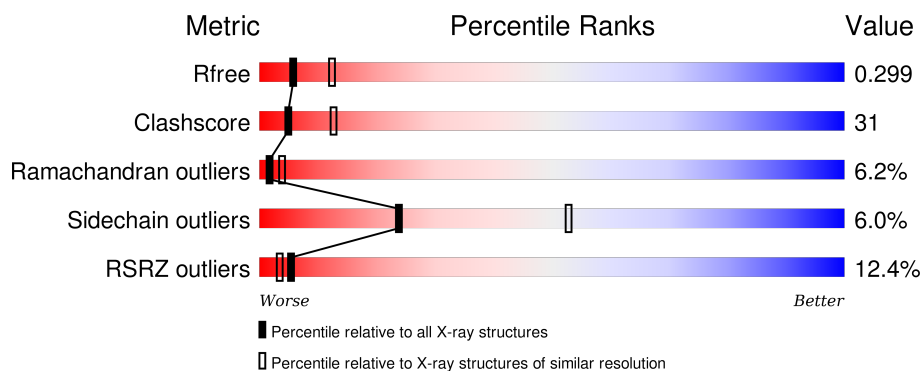
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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}	91344	3340 (2.80-2.72)
Clashscore	102246	3829 (2.80-2.72)
Ramachandran outliers	100387	3767 (2.80-2.72)
Sidechain outliers	100360	3770 (2.80-2.72)
RSRZ outliers	91569	3352 (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 ≥ 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	394	<div> <div>16%</div> <div>52%</div> <div>42%</div> <div>5%</div> </div>
1	B	394	<div> <div>9%</div> <div>47%</div> <div>47%</div> <div>5%</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	NA	A	1395	-	-	-	X
4	FUC	A	1399	X	-	-	-
4	FUC	B	1398	X	-	-	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6296 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 ENVELOPE GLYCOPROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	394	Total	C	N	O	S	0	0	0
			3062	1931	526	579	26			
1	B	394	Total	C	N	O	S	0	0	0
			3062	1931	526	579	26			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	71	GLU	ASP	CONFLICT	UNP P12823
A	390	ASN	ASP	CONFLICT	UNP P12823
B	71	GLU	ASP	CONFLICT	UNP P12823
B	390	ASN	ASP	CONFLICT	UNP P12823

- Molecule 2 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Na	0	0
			1	1		

- Molecule 3 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: C₈H₁₅NO₆).



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

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

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	4	Total	C	N	O	0	0
			49	28	2	19		
4	B	4	Total	C	N	O	0	0
			49	28	2	19		

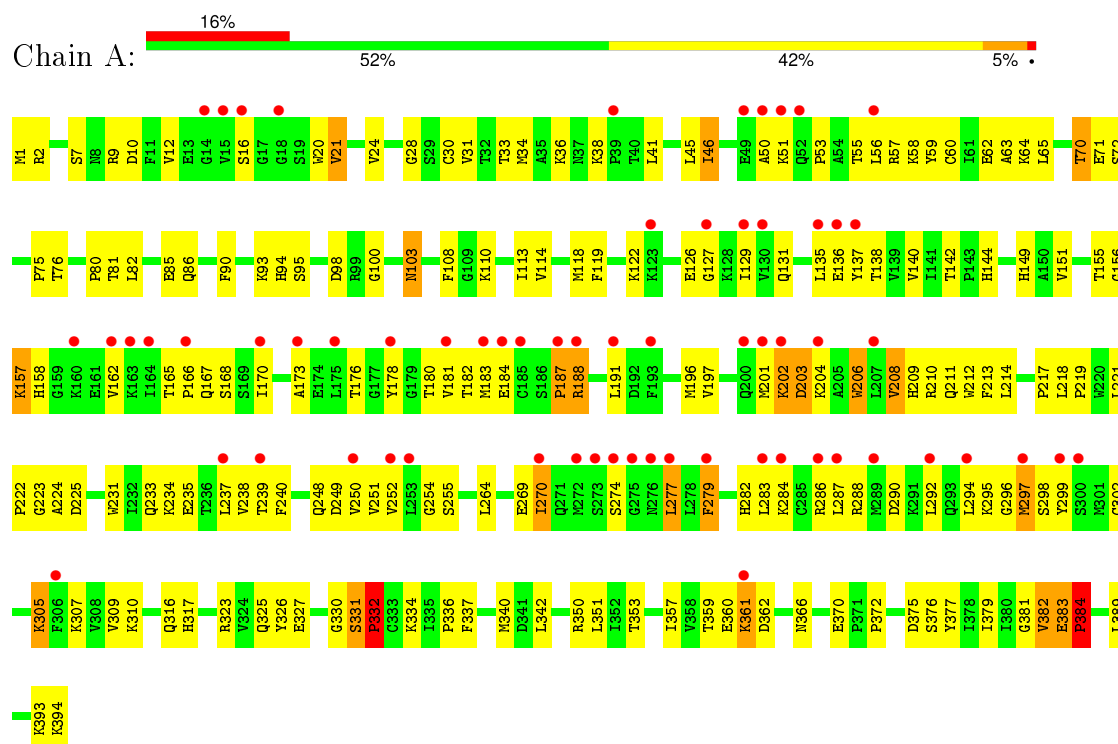
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	20	Total	O	0	0
			20	20		
5	B	25	Total	O	0	0
			25	25		

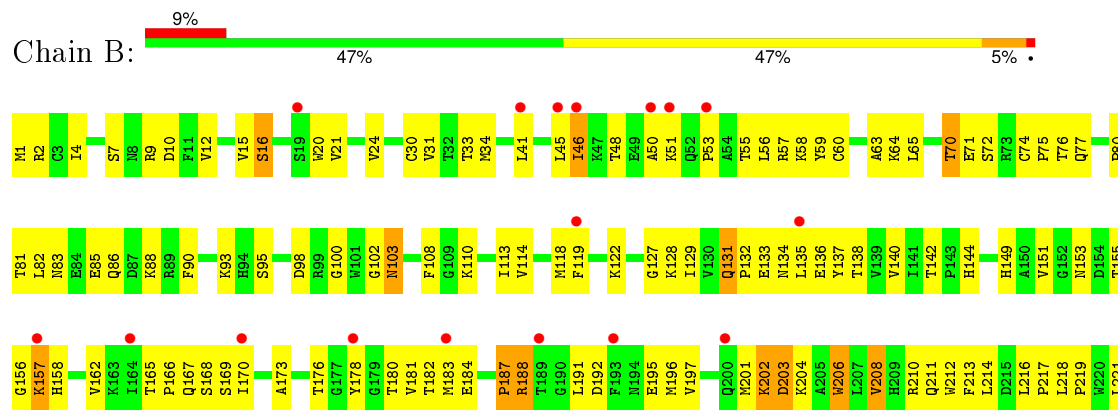
3 Residue-property plots

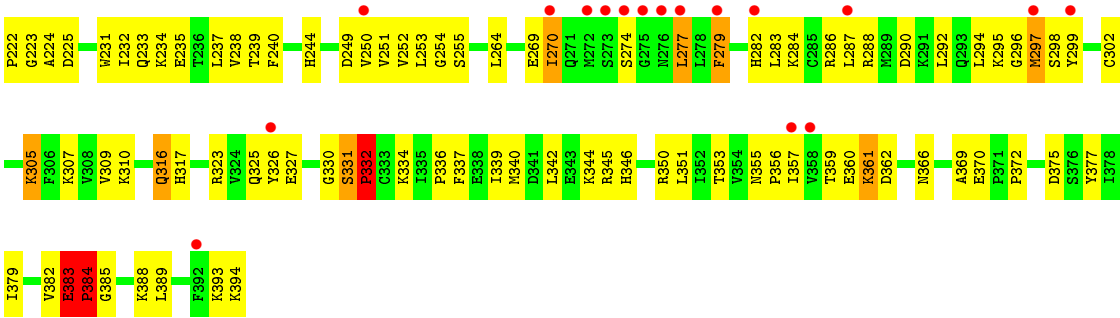
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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: ENVELOPE GLYCOPROTEIN



• Molecule 1: ENVELOPE GLYCOPROTEIN





4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	81.54Å 81.54Å 288.62Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.10 – 2.75 48.10 – 2.75	Depositor EDS
% Data completeness (in resolution range)	83.2 (48.10-2.75) 83.2 (48.10-2.75)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.15 (at 2.77Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.261 , 0.296 0.265 , 0.299	Depositor DCC
R_{free} test set	1281 reflections (5.15%)	DCC
Wilson B-factor (Å ²)	59.7	Xtriage
Anisotropy	0.262	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 83.4	EDS
Estimated twinning fraction	0.034 for -h,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 29971 reflections	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	6296	wwPDB-VP
Average B, all atoms (Å ²)	76.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.62% 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: NA, BMA, NAG, FUC

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	0.44	0/3124	0.68	1/4219 (0.0%)
1	B	0.45	0/3124	0.68	1/4219 (0.0%)
All	All	0.44	0/6248	0.68	2/8438 (0.0%)

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
4	A	1	0
4	B	1	0
All	All	2	0

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	384	PRO	CA-N-CD	-6.91	101.83	111.50
1	B	384	PRO	CA-N-CD	-5.65	103.59	111.50

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
4	A	1399	FUC	C1
4	B	1398	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	3062	0	3065	193	0
1	B	3062	0	3065	203	0
2	A	1	0	0	0	0
3	A	14	0	13	0	0
3	B	14	0	13	0	0
4	A	49	0	43	2	0
4	B	49	0	43	2	0
5	A	20	0	0	4	0
5	B	25	0	0	5	1
All	All	6296	0	6242	390	1

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 31.

All (390) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:383:GLU:HB3	1:A:384:PRO:HD2	1.43	1.01
1:B:323:ARG:HE	1:B:366:ASN:HD21	1.13	0.96
1:A:323:ARG:HE	1:A:366:ASN:HD21	1.09	0.95
1:B:182:THR:HG22	1:B:288:ARG:HB2	1.56	0.87
1:B:344:LYS:HG3	5:B:2017:HOH:O	1.77	0.85
1:A:182:THR:HG22	1:A:288:ARG:HB2	1.57	0.85
1:A:71:GLU:HG2	1:A:81:THR:H	1.40	0.84
1:B:71:GLU:HG2	1:B:81:THR:H	1.41	0.84
1:A:323:ARG:NE	1:A:366:ASN:HD21	1.75	0.83
1:B:345:ARG:HG2	1:B:346:HIS:CD2	2.17	0.79
1:B:323:ARG:NE	1:B:366:ASN:HD21	1.79	0.79
1:A:383:GLU:HB3	1:A:384:PRO:CD	2.13	0.79
1:A:342:LEU:HD23	1:A:377:TYR:CE1	2.17	0.79
1:A:342:LEU:HA	1:A:377:TYR:CE1	2.18	0.79
1:A:269:GLU:O	1:A:270:ILE:HG13	1.85	0.77
1:A:56:LEU:HD22	1:A:213:PHE:HE2	1.49	0.77
1:A:56:LEU:HD21	1:A:214:LEU:HD21	1.68	0.75
1:A:277:LEU:HB2	1:A:279:PHE:CE1	2.22	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:202:LYS:O	1:B:203:ASP:CG	2.25	0.75
1:B:394:LYS:HD3	1:B:394:LYS:OXT	1.87	0.74
1:A:394:LYS:OXT	1:A:394:LYS:HD3	1.86	0.74
1:B:277:LEU:HB2	1:B:279:PHE:CE1	2.23	0.74
1:B:383:GLU:HB3	1:B:384:PRO:HD2	1.67	0.74
1:B:269:GLU:O	1:B:270:ILE:HG13	1.88	0.74
1:B:56:LEU:HD21	1:B:214:LEU:HD21	1.70	0.73
1:B:56:LEU:HD22	1:B:213:PHE:HE2	1.53	0.73
1:A:305:LYS:NZ	1:A:305:LYS:H	1.86	0.72
1:B:383:GLU:OE1	5:B:2022:HOH:O	2.07	0.72
1:A:305:LYS:HZ2	1:A:305:LYS:H	1.36	0.72
1:A:165:THR:HB	1:A:168:SER:OG	1.89	0.72
1:B:217:PRO:O	1:B:218:LEU:HD23	1.90	0.71
1:B:192:ASP:OD2	1:B:195:GLU:HG2	1.90	0.71
1:B:336:PRO:HG2	1:B:382:VAL:HG23	1.70	0.71
1:B:305:LYS:CE	1:B:305:LYS:H	2.03	0.71
1:B:296:GLY:C	1:B:298:SER:H	1.94	0.71
1:B:100:GLY:H	1:B:103:ASN:HD21	1.37	0.71
1:A:359:THR:O	1:A:360:GLU:HG3	1.91	0.71
1:A:342:LEU:HD23	1:A:377:TYR:HE1	1.54	0.70
1:B:165:THR:HB	1:B:168:SER:OG	1.90	0.70
1:B:155:THR:HG22	1:B:157:LYS:H	1.57	0.70
1:A:20:TRP:HA	1:A:287:LEU:O	1.91	0.70
1:B:240:PHE:CE1	1:B:250:VAL:HG22	2.27	0.69
1:A:98:ASP:OD2	1:B:7:SER:HB2	1.93	0.69
1:B:240:PHE:HE1	1:B:250:VAL:HG22	1.57	0.69
1:A:305:LYS:H	1:A:305:LYS:CE	2.05	0.69
1:B:305:LYS:H	1:B:305:LYS:NZ	1.90	0.69
1:A:155:THR:HG22	1:A:157:LYS:H	1.58	0.69
1:B:20:TRP:HA	1:B:287:LEU:O	1.92	0.68
1:A:204:LYS:HB3	1:A:206:TRP:CH2	2.28	0.68
1:A:100:GLY:H	1:A:103:ASN:HD21	1.39	0.68
1:B:56:LEU:HD23	1:B:56:LEU:C	2.14	0.68
1:B:305:LYS:HE3	1:B:305:LYS:N	2.08	0.67
1:B:206:TRP:H	1:B:206:TRP:HE3	1.43	0.67
1:B:359:THR:O	1:B:360:GLU:HG3	1.94	0.67
1:A:173:ALA:HB3	1:A:181:VAL:HG13	1.77	0.67
1:A:71:GLU:HG3	1:A:80:PRO:HB3	1.75	0.67
1:B:20:TRP:CE3	1:B:286:ARG:HD2	2.30	0.67
1:B:344:LYS:NZ	1:B:344:LYS:HB3	2.10	0.66
1:A:20:TRP:CE3	1:A:286:ARG:HD2	2.31	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:344:LYS:HE3	1:B:388:LYS:HE3	1.77	0.66
1:B:71:GLU:HG3	1:B:80:PRO:HB3	1.76	0.66
1:B:206:TRP:CE3	1:B:206:TRP:N	2.64	0.66
1:B:337:PHE:CD2	1:B:351:LEU:HD21	2.31	0.66
1:B:170:ILE:HG13	1:B:184:GLU:HG3	1.77	0.66
1:A:305:LYS:HE3	1:A:305:LYS:N	2.11	0.65
1:B:305:LYS:H	1:B:305:LYS:HZ2	1.42	0.65
1:A:316:GLN:HE22	1:B:110:LYS:HE2	1.60	0.65
1:A:56:LEU:C	1:A:56:LEU:HD23	2.17	0.65
1:A:296:GLY:C	1:A:298:SER:H	1.99	0.65
1:B:170:ILE:CG1	1:B:184:GLU:HG3	2.27	0.65
1:A:382:VAL:HG22	5:A:2018:HOH:O	1.96	0.64
1:B:82:LEU:O	1:B:85:GLU:HB2	1.97	0.64
1:B:383:GLU:HB3	1:B:384:PRO:CD	2.27	0.64
1:A:240:PHE:CE1	1:A:250:VAL:HG22	2.32	0.64
1:B:286:ARG:HE	1:B:288:ARG:HH21	1.44	0.63
1:A:206:TRP:CE3	1:A:206:TRP:N	2.67	0.63
1:B:202:LYS:O	1:B:203:ASP:OD2	2.17	0.63
1:A:162:VAL:HG11	1:A:183:MET:CE	2.28	0.63
1:B:173:ALA:HB3	1:B:181:VAL:HG13	1.80	0.63
1:B:162:VAL:HG11	1:B:183:MET:CE	2.29	0.63
1:A:217:PRO:O	1:A:218:LEU:HD23	1.99	0.63
1:B:240:PHE:CE1	1:B:250:VAL:HG13	2.34	0.63
1:B:137:TYR:CD2	1:B:283:LEU:HD22	2.34	0.63
1:A:286:ARG:HE	1:A:288:ARG:HH21	1.47	0.63
1:A:206:TRP:H	1:A:206:TRP:HE3	1.47	0.63
4:A:1398:NAG:H4	4:A:1400:BMA:O2	1.98	0.62
1:A:383:GLU:CB	1:A:384:PRO:CD	2.77	0.62
1:A:376:SER:C	1:A:377:TYR:CD1	2.73	0.62
1:B:296:GLY:HA3	1:B:299:TYR:CE2	2.34	0.62
1:B:100:GLY:H	1:B:103:ASN:ND2	1.96	0.62
1:B:166:PRO:CG	1:B:187:PRO:HG3	2.30	0.62
1:A:240:PHE:HE1	1:A:250:VAL:HG22	1.65	0.62
1:B:297:MET:HA	1:B:334:LYS:NZ	2.15	0.61
1:A:100:GLY:H	1:A:103:ASN:ND2	1.98	0.61
1:A:24:VAL:HG22	1:A:284:LYS:HG2	1.81	0.61
1:A:127:GLY:HA3	1:A:213:PHE:CZ	2.35	0.61
1:B:127:GLY:HA3	1:B:213:PHE:CZ	2.35	0.61
1:B:24:VAL:HG22	1:B:284:LYS:HG2	1.81	0.61
1:A:170:ILE:CG1	1:A:184:GLU:HG3	2.31	0.61
4:B:1397:NAG:H4	4:B:1399:BMA:O2	1.99	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:166:PRO:CG	1:A:187:PRO:HG3	2.31	0.61
1:A:221:LEU:HD13	1:A:225:ASP:OD2	2.01	0.60
1:A:137:TYR:CD2	1:A:283:LEU:HD22	2.36	0.60
1:A:270:ILE:HG22	1:A:270:ILE:O	2.02	0.60
1:B:206:TRP:HE3	1:B:206:TRP:N	1.97	0.60
1:A:206:TRP:N	1:A:206:TRP:HE3	1.98	0.60
1:A:82:LEU:O	1:A:85:GLU:HB2	2.01	0.60
1:A:340:MET:HG3	1:A:379:ILE:HG13	1.82	0.60
1:B:182:THR:CG2	1:B:288:ARG:HB2	2.31	0.60
1:B:204:LYS:HB3	1:B:206:TRP:CH2	2.36	0.60
1:B:340:MET:HG3	1:B:379:ILE:HG13	1.84	0.60
1:A:20:TRP:HE3	1:A:286:ARG:HD2	1.67	0.60
1:B:56:LEU:O	1:B:56:LEU:HD23	2.02	0.60
1:A:71:GLU:HG2	1:A:81:THR:N	2.16	0.59
1:B:270:ILE:HG22	1:B:270:ILE:O	2.02	0.59
1:B:191:LEU:HD13	1:B:196:MET:CE	2.33	0.58
1:A:383:GLU:HA	1:A:383:GLU:OE1	2.03	0.58
1:A:323:ARG:HE	1:A:366:ASN:ND2	1.91	0.58
1:A:182:THR:CG2	1:A:288:ARG:HB2	2.30	0.58
1:A:129:ILE:CD1	1:A:210:ARG:HH22	2.16	0.58
1:A:206:TRP:NE1	1:A:264:LEU:HB3	2.18	0.58
1:B:383:GLU:HA	1:B:383:GLU:OE1	2.03	0.57
1:B:221:LEU:HD13	1:B:225:ASP:OD2	2.04	0.57
1:B:20:TRP:HE3	1:B:286:ARG:HD2	1.67	0.57
1:A:55:THR:O	1:A:223:GLY:HA3	2.04	0.57
1:A:297:MET:HA	1:A:334:LYS:NZ	2.19	0.57
1:A:50:ALA:HA	1:A:135:LEU:HD23	1.87	0.57
1:B:296:GLY:HA3	1:B:299:TYR:CD2	2.40	0.57
1:A:34:MET:SD	1:A:350:ARG:NH2	2.78	0.57
1:B:50:ALA:HA	1:B:135:LEU:HD23	1.85	0.57
1:A:56:LEU:HD12	1:A:129:ILE:HD11	1.86	0.57
1:B:221:LEU:CD2	1:B:231:TRP:HA	2.35	0.57
1:A:53:PRO:HB3	1:A:129:ILE:O	2.04	0.56
1:A:65:LEU:HG	1:A:252:VAL:HG22	1.87	0.56
1:A:251:VAL:HG21	1:B:204:LYS:HE3	1.86	0.56
1:A:162:VAL:HG11	1:A:183:MET:HE1	1.87	0.56
1:A:350:ARG:HD3	1:A:370:GLU:OE1	2.05	0.56
1:B:55:THR:O	1:B:223:GLY:HA3	2.05	0.56
1:A:7:SER:HB2	1:B:98:ASP:OD2	2.06	0.56
1:B:162:VAL:HG11	1:B:183:MET:HE3	1.86	0.56
1:B:202:LYS:HE2	1:B:202:LYS:HA	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:191:LEU:HD13	1:B:196:MET:HE3	1.88	0.56
1:B:307:LYS:HB3	1:B:307:LYS:NZ	2.21	0.56
1:A:383:GLU:HG3	1:A:384:PRO:HD3	1.88	0.56
1:B:286:ARG:NE	1:B:288:ARG:HH21	2.04	0.56
1:A:296:GLY:HA3	1:A:299:TYR:CE2	2.41	0.56
1:B:350:ARG:HD3	1:B:370:GLU:OE1	2.06	0.56
1:B:305:LYS:CE	1:B:305:LYS:N	2.68	0.55
1:A:221:LEU:CD2	1:A:231:TRP:HA	2.37	0.55
1:B:129:ILE:CD1	1:B:210:ARG:HH22	2.18	0.55
1:B:9:ARG:HB3	1:B:317:HIS:CE1	2.41	0.55
1:B:206:TRP:NE1	1:B:264:LEU:HB3	2.22	0.55
1:A:151:VAL:HB	5:A:2009:HOH:O	2.07	0.55
1:B:342:LEU:HD21	1:B:375:ASP:CB	2.37	0.55
1:A:305:LYS:CE	1:A:305:LYS:N	2.68	0.55
1:A:337:PHE:CD2	1:A:351:LEU:HD21	2.41	0.55
1:A:197:VAL:HG23	1:A:210:ARG:HA	1.89	0.54
1:B:197:VAL:HG23	1:B:210:ARG:HA	1.88	0.54
1:A:307:LYS:NZ	1:A:307:LYS:HB3	2.22	0.54
1:A:234:LYS:O	1:A:238:VAL:HG23	2.07	0.54
1:A:240:PHE:CE1	1:A:250:VAL:HG13	2.43	0.54
1:B:53:PRO:HB3	1:B:129:ILE:O	2.07	0.54
1:A:165:THR:HB	1:A:168:SER:CB	2.38	0.54
1:A:70:THR:HA	1:A:82:LEU:HD11	1.90	0.54
1:B:344:LYS:HE3	1:B:388:LYS:CE	2.38	0.54
1:A:1:MET:HG3	1:A:144:HIS:HA	1.90	0.53
1:B:221:LEU:HD21	1:B:231:TRP:HA	1.90	0.53
1:A:286:ARG:NE	1:A:288:ARG:HH21	2.06	0.53
1:A:342:LEU:HD23	1:A:377:TYR:OH	2.09	0.53
1:B:235:GLU:CD	1:B:235:GLU:H	2.12	0.53
1:B:80:PRO:HG2	1:B:113:ILE:CA	2.39	0.53
1:B:65:LEU:HG	1:B:252:VAL:HG22	1.91	0.53
1:B:34:MET:SD	1:B:350:ARG:NH2	2.82	0.53
1:A:155:THR:CG2	1:A:157:LYS:HG3	2.39	0.53
1:B:1:MET:HG3	1:B:144:HIS:HA	1.90	0.53
1:B:56:LEU:HD12	1:B:129:ILE:HD11	1.90	0.52
1:A:235:GLU:H	1:A:235:GLU:CD	2.13	0.52
1:B:165:THR:HB	1:B:168:SER:CB	2.39	0.52
1:A:165:THR:HG22	1:A:167:GLN:H	1.75	0.52
1:B:336:PRO:HG2	1:B:382:VAL:CG2	2.37	0.52
1:A:204:LYS:HE3	1:B:251:VAL:HG21	1.92	0.52
1:B:64:LYS:HB2	1:B:122:LYS:HD2	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:202:LYS:O	1:A:203:ASP:OD2	2.27	0.52
1:A:75:PRO:O	1:A:76:THR:OG1	2.23	0.52
1:B:90:PHE:CZ	1:B:118:MET:HB2	2.44	0.52
1:B:114:VAL:O	1:B:114:VAL:HG13	2.10	0.52
1:B:296:GLY:C	1:B:298:SER:N	2.63	0.51
1:B:155:THR:CG2	1:B:157:LYS:HG3	2.39	0.51
1:A:64:LYS:HB2	1:A:122:LYS:HD2	1.92	0.51
1:A:80:PRO:HG2	1:A:113:ILE:CA	2.41	0.51
1:B:165:THR:HG22	1:B:167:GLN:H	1.75	0.51
1:A:296:GLY:HA3	1:A:299:TYR:CD2	2.46	0.51
1:A:202:LYS:O	1:A:203:ASP:CG	2.48	0.51
1:B:57:ARG:HG2	1:B:58:LYS:N	2.26	0.51
1:B:331:SER:CB	1:B:332:PRO:CD	2.88	0.51
1:A:381:GLY:HA2	5:A:2018:HOH:O	2.10	0.51
1:B:70:THR:HA	1:B:82:LEU:HD11	1.93	0.51
1:B:166:PRO:HG2	1:B:187:PRO:HG3	1.93	0.51
1:A:221:LEU:HD21	1:A:231:TRP:HA	1.92	0.51
1:B:45:LEU:HD12	1:B:138:THR:O	2.11	0.51
1:A:56:LEU:O	1:A:56:LEU:HD23	2.11	0.51
1:A:170:ILE:HG13	1:A:184:GLU:HG3	1.93	0.51
1:A:166:PRO:HB3	1:A:187:PRO:HD3	1.92	0.51
1:B:234:LYS:O	1:B:238:VAL:HG23	2.11	0.50
1:A:98:ASP:CG	1:B:7:SER:HB2	2.31	0.50
1:A:166:PRO:HG2	1:A:187:PRO:HG3	1.93	0.50
1:A:28:GLY:HA3	1:B:244:HIS:CD2	2.46	0.50
1:A:342:LEU:HD23	1:A:377:TYR:CZ	2.45	0.50
1:A:191:LEU:HD13	1:A:196:MET:CE	2.41	0.50
1:A:90:PHE:CZ	1:A:118:MET:HB2	2.47	0.50
1:A:80:PRO:HB2	1:A:114:VAL:HG12	1.94	0.50
1:A:342:LEU:HD21	1:A:375:ASP:CB	2.41	0.50
1:A:239:THR:HB	1:A:251:VAL:CG1	2.42	0.50
1:A:9:ARG:HB3	1:A:317:HIS:CE1	2.47	0.50
1:B:71:GLU:HG2	1:B:81:THR:N	2.18	0.50
1:A:45:LEU:HD12	1:A:138:THR:O	2.11	0.50
1:A:114:VAL:HG13	1:A:114:VAL:O	2.12	0.50
1:B:100:GLY:HA3	1:B:108:PHE:CD1	2.46	0.50
1:B:239:THR:HB	1:B:251:VAL:CG1	2.42	0.50
1:A:100:GLY:HA3	1:A:108:PHE:CD1	2.47	0.50
1:A:63:ALA:O	1:A:252:VAL:HG11	2.12	0.50
1:B:170:ILE:N	1:B:170:ILE:HD12	2.28	0.49
1:A:219:PRO:HG2	1:A:237:LEU:HD12	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:46:ILE:CG1	1:B:140:VAL:HG23	2.43	0.49
1:B:173:ALA:O	1:B:180:THR:HG23	2.12	0.48
1:B:166:PRO:HB3	1:B:187:PRO:HD3	1.94	0.48
1:A:202:LYS:HA	1:A:202:LYS:HE2	1.95	0.48
1:A:277:LEU:HD12	1:A:277:LEU:H	1.79	0.48
1:B:95:SER:HB3	1:B:113:ILE:CG2	2.43	0.48
1:A:307:LYS:HZ3	1:A:307:LYS:HB3	1.77	0.48
1:A:46:ILE:CG1	1:A:140:VAL:HG23	2.44	0.48
1:B:202:LYS:CA	1:B:202:LYS:HE2	2.44	0.48
1:B:197:VAL:CG2	1:B:210:ARG:HA	2.44	0.48
1:A:57:ARG:HG2	1:A:58:LYS:N	2.29	0.48
1:B:131:GLN:HG3	1:B:134:ASN:HD22	1.78	0.48
1:B:223:GLY:C	1:B:225:ASP:H	2.17	0.48
1:B:46:ILE:HG12	1:B:140:VAL:HG23	1.96	0.48
1:A:110:LYS:HE2	1:B:316:GLN:HE22	1.79	0.48
1:A:331:SER:CB	1:A:332:PRO:CD	2.92	0.48
1:B:277:LEU:H	1:B:277:LEU:HD12	1.79	0.47
1:B:131:GLN:HB2	1:B:133:GLU:OE1	2.14	0.47
1:A:173:ALA:HB3	1:A:181:VAL:CG1	2.43	0.47
1:B:233:GLN:HA	1:B:235:GLU:OE2	2.14	0.47
1:A:309:VAL:CG2	1:A:325:GLN:HB2	2.44	0.47
1:A:95:SER:HB3	1:A:113:ILE:CG2	2.44	0.47
1:B:345:ARG:HB2	1:B:345:ARG:NH1	2.29	0.47
1:A:197:VAL:CG2	1:A:210:ARG:HA	2.45	0.47
1:B:201:MET:O	1:B:203:ASP:N	2.47	0.47
1:B:342:LEU:HD21	1:B:375:ASP:HB2	1.95	0.47
1:B:251:VAL:HG22	1:B:252:VAL:N	2.30	0.47
1:A:56:LEU:HB2	1:A:129:ILE:HG13	1.97	0.47
1:A:223:GLY:C	1:A:225:ASP:H	2.17	0.47
1:B:383:GLU:O	1:B:385:GLY:N	2.48	0.47
1:B:309:VAL:CG2	1:B:325:GLN:HB2	2.44	0.47
1:A:51:LYS:HE3	1:A:136:GLU:HB2	1.96	0.47
1:A:204:LYS:HB3	1:A:206:TRP:CZ3	2.49	0.47
1:A:233:GLN:HA	1:A:235:GLU:OE2	2.14	0.47
1:B:51:LYS:HE3	1:B:136:GLU:HB2	1.96	0.47
1:A:309:VAL:HG21	1:A:325:GLN:HB2	1.97	0.47
1:A:342:LEU:HD21	1:A:375:ASP:HB2	1.97	0.46
1:A:296:GLY:C	1:A:298:SER:N	2.68	0.46
1:A:208:VAL:HG21	1:A:212:TRP:CZ3	2.50	0.46
1:B:72:SER:HB3	1:B:113:ILE:HG13	1.96	0.46
1:A:377:TYR:CD1	1:A:377:TYR:N	2.83	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:56:LEU:CD2	1:B:56:LEU:C	2.82	0.46
1:B:372:PRO:O	5:B:2020:HOH:O	2.21	0.46
1:B:155:THR:HG22	1:B:157:LYS:HG3	1.98	0.46
1:A:155:THR:HG22	1:A:157:LYS:HG3	1.96	0.46
1:B:65:LEU:HD11	1:B:238:VAL:HG13	1.98	0.46
1:A:119:PHE:CB	1:A:234:LYS:HD2	2.45	0.46
1:B:57:ARG:CG	1:B:58:LYS:N	2.79	0.46
1:B:83:ASN:ND2	5:B:2003:HOH:O	2.47	0.46
1:B:309:VAL:HG21	1:B:325:GLN:HB2	1.98	0.46
1:A:208:VAL:HG21	1:A:212:TRP:HZ3	1.81	0.46
1:A:360:GLU:O	1:A:362:ASP:N	2.49	0.46
1:A:239:THR:HB	1:A:251:VAL:HG12	1.98	0.46
1:B:60:CYS:HB2	1:B:231:TRP:CZ3	2.51	0.46
1:A:60:CYS:HB2	1:A:231:TRP:CZ3	2.51	0.45
1:B:221:LEU:HG	1:B:231:TRP:CE3	2.51	0.45
1:B:119:PHE:CB	1:B:234:LYS:HD2	2.45	0.45
1:A:170:ILE:HG12	1:A:184:GLU:HG3	1.98	0.45
1:A:206:TRP:CD1	1:A:264:LEU:HB3	2.50	0.45
1:A:173:ALA:O	1:A:180:THR:HG23	2.15	0.45
1:A:296:GLY:O	1:A:299:TYR:HD2	2.00	0.45
1:B:173:ALA:HB3	1:B:181:VAL:CG1	2.46	0.45
1:A:310:LYS:HB3	1:A:323:ARG:HB3	1.97	0.45
1:B:63:ALA:O	1:B:252:VAL:HG11	2.17	0.45
1:B:323:ARG:HE	1:B:366:ASN:ND2	1.96	0.45
1:A:201:MET:O	1:A:203:ASP:N	2.50	0.45
1:B:339:ILE:HA	1:B:377:TYR:O	2.17	0.45
1:A:251:VAL:HG22	1:A:252:VAL:N	2.32	0.45
1:A:65:LEU:HD11	1:A:238:VAL:HG13	1.99	0.45
1:B:208:VAL:HG21	1:B:212:TRP:HZ3	1.82	0.45
1:A:36:LYS:O	1:A:38:LYS:HG2	2.17	0.45
1:A:95:SER:HB2	1:A:248:GLN:NE2	2.32	0.44
1:A:302:CYS:HB3	1:A:326:TYR:CZ	2.52	0.44
1:B:41:LEU:HD11	1:B:292:LEU:HD11	1.99	0.44
1:B:208:VAL:HG21	1:B:212:TRP:CZ3	2.52	0.44
1:A:296:GLY:O	1:A:299:TYR:CD2	2.70	0.44
1:A:336:PRO:HG2	5:A:2018:HOH:O	2.17	0.44
1:B:7:SER:O	1:B:317:HIS:CE1	2.71	0.44
1:B:372:PRO:O	1:B:393:LYS:HB3	2.18	0.44
1:B:240:PHE:HA	1:B:249:ASP:O	2.18	0.44
1:B:219:PRO:HG2	1:B:237:LEU:HD12	2.00	0.44
1:A:302:CYS:SG	1:A:334:LYS:O	2.75	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:178:TYR:CE1	1:B:295:LYS:HE3	2.52	0.44
1:B:239:THR:HB	1:B:251:VAL:HG12	1.99	0.43
1:B:377:TYR:HB3	1:B:379:ILE:HD11	2.00	0.43
1:A:56:LEU:C	1:A:56:LEU:CD2	2.85	0.43
1:A:62:GLU:HG2	1:A:122:LYS:HB2	2.01	0.43
1:A:334:LYS:HG2	1:A:357:ILE:HG22	2.00	0.43
1:A:46:ILE:HG12	1:A:140:VAL:HG23	2.01	0.43
1:A:210:ARG:O	1:A:213:PHE:HB3	2.19	0.43
1:A:72:SER:HB3	1:A:113:ILE:HG13	2.01	0.43
1:A:353:THR:O	1:A:353:THR:HG22	2.19	0.43
1:A:221:LEU:HG	1:A:231:TRP:CE3	2.54	0.43
1:A:12:VAL:O	1:A:33:THR:HA	2.18	0.43
1:B:80:PRO:HB2	1:B:114:VAL:HG12	2.00	0.43
1:B:151:VAL:HB	5:B:2010:HOH:O	2.18	0.43
1:B:12:VAL:O	1:B:33:THR:HA	2.18	0.43
1:B:59:TYR:CD1	1:B:218:LEU:HB2	2.53	0.43
1:B:344:LYS:HB3	1:B:344:LYS:HZ3	1.83	0.42
1:B:383:GLU:CB	1:B:384:PRO:CD	2.96	0.42
1:A:41:LEU:HD11	1:A:292:LEU:HD11	2.00	0.42
1:A:327:GLU:O	1:A:361:LYS:HE3	2.19	0.42
1:A:10:ASP:O	1:A:31:VAL:HA	2.18	0.42
1:B:10:ASP:O	1:B:31:VAL:HA	2.19	0.42
1:B:310:LYS:HB3	1:B:323:ARG:HB3	2.00	0.42
1:B:331:SER:HB2	1:B:332:PRO:CD	2.49	0.42
1:B:302:CYS:HB3	1:B:326:TYR:CZ	2.54	0.42
1:B:360:GLU:O	1:B:362:ASP:N	2.52	0.42
1:B:74:CYS:HB2	1:B:77:GLN:HG3	2.01	0.42
1:A:222:PRO:O	1:A:225:ASP:HB2	2.20	0.42
1:B:75:PRO:O	1:B:76:THR:OG1	2.28	0.42
1:B:355:ASN:N	1:B:356:PRO:HD3	2.34	0.42
1:A:206:TRP:CD2	1:A:264:LEU:HD13	2.55	0.42
1:A:269:GLU:C	1:A:270:ILE:HG13	2.40	0.42
1:A:85:GLU:OE1	1:A:94:HIS:NE2	2.31	0.42
1:A:57:ARG:CG	1:A:58:LYS:N	2.83	0.42
1:A:178:TYR:CE1	1:A:295:LYS:HE3	2.54	0.42
1:B:296:GLY:O	1:B:299:TYR:HD2	2.03	0.42
1:A:188:ARG:CZ	1:A:284:LYS:HD2	2.50	0.42
1:B:339:ILE:HD11	1:B:369:ALA:HB1	2.02	0.42
1:B:56:LEU:HB2	1:B:129:ILE:HG13	2.02	0.42
1:B:57:ARG:CG	1:B:58:LYS:H	2.32	0.42
1:A:59:TYR:CD1	1:A:218:LEU:HB2	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:331:SER:CB	1:B:332:PRO:HD2	2.49	0.42
1:B:327:GLU:O	1:B:361:LYS:HE3	2.20	0.42
1:A:119:PHE:HB3	1:A:234:LYS:HD2	2.02	0.41
1:A:12:VAL:HG12	1:A:21:VAL:HG11	2.01	0.41
1:B:149:HIS:O	1:B:153:ASN:HB2	2.20	0.41
1:B:188:ARG:CZ	1:B:284:LYS:HD2	2.50	0.41
1:B:156:GLY:C	1:B:158:HIS:H	2.24	0.41
1:B:119:PHE:HB3	1:B:234:LYS:HD2	2.02	0.41
1:B:334:LYS:HG2	1:B:357:ILE:HG22	2.03	0.41
1:A:170:ILE:HD12	1:A:170:ILE:N	2.36	0.41
1:B:307:LYS:HB3	1:B:307:LYS:HZ3	1.84	0.41
1:B:149:HIS:HB3	4:B:1396:NAG:H2	2.01	0.41
1:A:156:GLY:C	1:A:158:HIS:H	2.23	0.41
1:B:169:SER:C	1:B:170:ILE:HD12	2.41	0.41
1:B:48:THR:O	1:B:279:PHE:CE1	2.73	0.41
1:B:206:TRP:CD1	1:B:264:LEU:HB3	2.56	0.41
1:B:232:ILE:HG22	1:B:233:GLN:HG3	2.02	0.41
1:B:210:ARG:O	1:B:213:PHE:HB3	2.20	0.41
1:B:296:GLY:O	1:B:299:TYR:CD2	2.74	0.41
1:A:359:THR:O	1:A:360:GLU:CG	2.65	0.41
1:B:238:VAL:CG1	1:B:250:VAL:HG12	2.51	0.41
1:B:206:TRP:CD2	1:B:264:LEU:HD13	2.56	0.41
1:A:240:PHE:HA	1:A:249:ASP:O	2.21	0.41
1:A:202:LYS:CA	1:A:202:LYS:HE2	2.51	0.41
1:A:331:SER:CB	1:A:332:PRO:HD2	2.51	0.41
1:B:184:GLU:HB3	1:B:286:ARG:HB3	2.02	0.41
1:B:213:PHE:O	1:B:216:LEU:HG	2.20	0.41
1:A:173:ALA:O	1:A:180:THR:HA	2.21	0.41
1:A:372:PRO:O	1:A:393:LYS:HB3	2.21	0.41
1:B:253:LEU:HD12	1:B:253:LEU:HA	1.96	0.41
1:B:15:VAL:O	1:B:16:SER:C	2.59	0.40
1:B:128:LYS:O	1:B:197:VAL:HG13	2.21	0.40
1:B:191:LEU:HD22	1:B:196:MET:HE2	2.03	0.40
1:A:191:LEU:HD13	1:A:196:MET:HE3	2.03	0.40
1:B:131:GLN:O	1:B:132:PRO:C	2.60	0.40
1:A:51:LYS:HE3	1:A:136:GLU:CB	2.51	0.40
1:A:129:ILE:HD11	1:A:210:ARG:HH22	1.84	0.40
1:B:269:GLU:C	1:B:270:ILE:HG13	2.42	0.40
1:A:100:GLY:N	1:A:103:ASN:HD21	2.14	0.40
1:A:188:ARG:HH11	1:A:188:ARG:HB2	1.86	0.40
1:A:221:LEU:HD23	1:A:231:TRP:HA	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:222:PRO:O	1:B:225:ASP:HB2	2.22	0.40
1:B:353:THR:O	1:B:353:THR:HG22	2.21	0.40
1:A:149:HIS:HB3	4:A:1397:NAG:H2	2.03	0.40

All (1) 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)
5:B:2003:HOH:O	5:B:2003:HOH:O[6_555]	1.99	0.21

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	392/394 (100%)	321 (82%)	48 (12%)	23 (6%)	2	4
1	B	392/394 (100%)	318 (81%)	48 (12%)	26 (7%)	1	3
All	All	784/788 (100%)	639 (82%)	96 (12%)	49 (6%)	2	4

All (49) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	176	THR
1	A	187	PRO
1	A	202	LYS
1	A	383	GLU
1	B	176	THR
1	B	187	PRO
1	B	202	LYS
1	B	383	GLU
1	A	16	SER
1	A	203	ASP

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Mol	Chain	Res	Type
1	A	224	ALA
1	A	254	GLY
1	A	270	ILE
1	A	279	PHE
1	A	330	GLY
1	A	331	SER
1	A	361	LYS
1	B	16	SER
1	B	203	ASP
1	B	224	ALA
1	B	254	GLY
1	B	270	ILE
1	B	279	PHE
1	B	330	GLY
1	B	331	SER
1	A	274	SER
1	A	282	HIS
1	A	332	PRO
1	A	384	PRO
1	B	274	SER
1	B	282	HIS
1	B	332	PRO
1	B	361	LYS
1	B	384	PRO
1	A	157	LYS
1	A	255	SER
1	A	294	LEU
1	B	46	ILE
1	B	157	LYS
1	B	294	LEU
1	B	297	MET
1	A	46	ILE
1	A	297	MET
1	B	88	LYS
1	A	126	GLU
1	B	255	SER
1	B	316	GLN
1	B	102	GLY
1	B	4	ILE

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	342/342 (100%)	321 (94%)	21 (6%)	23	52
1	B	342/342 (100%)	322 (94%)	20 (6%)	25	54
All	All	684/684 (100%)	643 (94%)	41 (6%)	24	53

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	ARG
1	A	21	VAL
1	A	30	CYS
1	A	70	THR
1	A	86	GLN
1	A	93	LYS
1	A	103	ASN
1	A	131	GLN
1	A	142	THR
1	A	188	ARG
1	A	206	TRP
1	A	208	VAL
1	A	209	HIS
1	A	211	GLN
1	A	277	LEU
1	A	290	ASP
1	A	305	LYS
1	A	332	PRO
1	A	382	VAL
1	A	384	PRO
1	A	389	LEU
1	B	2	ARG
1	B	21	VAL
1	B	30	CYS
1	B	70	THR
1	B	86	GLN
1	B	93	LYS

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Mol	Chain	Res	Type
1	B	103	ASN
1	B	131	GLN
1	B	142	THR
1	B	188	ARG
1	B	206	TRP
1	B	208	VAL
1	B	211	GLN
1	B	277	LEU
1	B	290	ASP
1	B	305	LYS
1	B	332	PRO
1	B	383	GLU
1	B	384	PRO
1	B	389	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	77	GLN
1	A	86	GLN
1	A	103	ASN
1	A	131	GLN
1	A	149	HIS
1	A	200	GLN
1	A	244	HIS
1	A	248	GLN
1	A	256	GLN
1	A	316	GLN
1	A	366	ASN
1	B	77	GLN
1	B	83	ASN
1	B	86	GLN
1	B	103	ASN
1	B	131	GLN
1	B	134	ASN
1	B	149	HIS
1	B	200	GLN
1	B	244	HIS
1	B	248	GLN
1	B	316	GLN
1	B	317	HIS
1	B	366	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 ⓘ

8 carbohydrates are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	NAG	A	1397	1,4	14,14,15	0.67	0	15,19,21	0.74	0
4	NAG	A	1398	4	14,14,15	0.72	0	15,19,21	0.98	1 (6%)
4	FUC	A	1399	4	10,10,11	0.40	0	14,14,16	0.57	0
4	BMA	A	1400	4	11,11,12	0.57	0	14,15,17	0.38	0
4	NAG	B	1396	1,4	14,14,15	0.60	0	15,19,21	0.68	0
4	NAG	B	1397	4	14,14,15	0.74	0	15,19,21	0.99	1 (6%)
4	FUC	B	1398	4	10,10,11	0.43	0	14,14,16	0.61	0
4	BMA	B	1399	4	11,11,12	0.63	0	14,15,17	0.35	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
4	NAG	A	1397	1,4	-	0/6/23/26	0/1/1/1
4	NAG	A	1398	4	-	0/6/23/26	0/1/1/1
4	FUC	A	1399	4	1/1/4/5	0/0/17/20	0/1/1/1
4	BMA	A	1400	4	-	0/2/19/22	0/1/1/1
4	NAG	B	1396	1,4	-	0/6/23/26	0/1/1/1
4	NAG	B	1397	4	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	FUC	B	1398	4	1/1/4/5	0/0/17/20	0/1/1/1
4	BMA	B	1399	4	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1397	NAG	C4-C3-C2	2.17	114.61	111.23
4	A	1398	NAG	C4-C3-C2	2.25	114.73	111.23

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
4	A	1399	FUC	C1
4	B	1398	FUC	C1

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1397	NAG	1	0
4	A	1398	NAG	1	0
4	A	1400	BMA	1	0
4	B	1396	NAG	1	0
4	B	1397	NAG	1	0
4	B	1399	BMA	1	0

5.6 Ligand geometry

Of 3 ligands modelled in this entry, 1 is 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
3	NAG	A	1396	1	14,14,15	0.81	1 (7%)	15,19,21	0.64	0
3	NAG	B	1395	1	14,14,15	0.67	0	15,19,21	0.65	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
3	NAG	A	1396	1	-	0/6/23/26	0/1/1/1
3	NAG	B	1395	1	-	0/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1396	NAG	C1-C2	2.07	1.55	1.52

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, 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	394/394 (100%)	0.86	64 (16%) 3 1	29, 73, 131, 149	0
1	B	394/394 (100%)	0.69	34 (8%) 13 8	24, 69, 116, 149	0
All	All	788/788 (100%)	0.78	98 (12%) 5 3	24, 71, 126, 149	0

All (98) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	272	MET	7.0
1	A	276	ASN	7.0
1	B	50	ALA	6.0
1	A	162	VAL	5.2
1	A	164	ILE	5.1
1	A	51	LYS	5.0
1	A	14	GLY	4.9
1	B	276	ASN	4.9
1	A	297	MET	4.3
1	A	50	ALA	4.3
1	A	275	GLY	4.2
1	B	275	GLY	4.2
1	A	200	GLN	4.2
1	A	137	TYR	4.1
1	B	297	MET	4.1
1	A	183	MET	4.0
1	A	279	PHE	3.9
1	B	193	PHE	3.7
1	B	157	LYS	3.6
1	B	183	MET	3.6
1	A	277	LEU	3.5
1	A	52	GLN	3.5
1	B	299	TYR	3.4
1	A	178	TYR	3.4

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Mol	Chain	Res	Type	RSRZ
1	B	282	HIS	3.4
1	A	135	LEU	3.3
1	A	173	ALA	3.3
1	A	289	MET	3.3
1	A	191	LEU	3.3
1	B	274	SER	3.3
1	A	253	LEU	3.3
1	A	16	SER	3.2
1	A	15	VAL	3.2
1	B	119	PHE	3.1
1	B	19	SER	3.1
1	B	326	TYR	3.0
1	A	127	GLY	3.0
1	A	272	MET	3.0
1	A	130	VAL	3.0
1	A	18	GLY	2.9
1	A	286	ARG	2.9
1	A	163	LYS	2.9
1	A	292	LEU	2.8
1	A	193	PHE	2.8
1	B	135	LEU	2.7
1	A	270	ILE	2.7
1	B	164	ILE	2.7
1	B	392	PHE	2.7
1	A	129	ILE	2.7
1	A	166	PRO	2.7
1	A	274	SER	2.6
1	A	204	LYS	2.5
1	A	273	SER	2.5
1	A	299	TYR	2.5
1	B	200	GLN	2.5
1	A	136	GLU	2.5
1	B	287	LEU	2.5
1	A	250	VAL	2.4
1	A	202	LYS	2.4
1	B	358	VAL	2.4
1	A	175	LEU	2.4
1	A	123	LYS	2.4
1	A	181	VAL	2.3
1	A	185	CYS	2.3
1	B	279	PHE	2.3
1	B	270	ILE	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	239	THR	2.3
1	B	170	ILE	2.3
1	B	41	LEU	2.2
1	A	56	LEU	2.2
1	A	237	LEU	2.2
1	B	45	LEU	2.2
1	A	188	ARG	2.2
1	B	53	PRO	2.2
1	B	189	THR	2.2
1	A	284	LYS	2.2
1	A	361	LYS	2.2
1	A	187	PRO	2.2
1	B	250	VAL	2.2
1	A	252	VAL	2.2
1	A	49	GLU	2.2
1	B	178	TYR	2.1
1	B	51	LYS	2.1
1	B	46	ILE	2.1
1	A	39	PRO	2.1
1	A	207	LEU	2.1
1	A	287	LEU	2.1
1	A	306	PHE	2.1
1	A	283	LEU	2.1
1	A	300	SER	2.1
1	A	294	LEU	2.1
1	B	273	SER	2.1
1	B	277	LEU	2.0
1	A	170	ILE	2.0
1	A	184	GLU	2.0
1	A	201	MET	2.0
1	B	357	ILE	2.0
1	A	160	LYS	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, 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	NAG	A	1397	14/15	0.88	0.25	0.80	75,81,89,90	0
4	NAG	B	1396	14/15	0.94	0.21	-0.09	45,67,72,74	0
4	FUC	A	1399	10/11	0.94	0.23	-	70,75,78,88	0
4	NAG	A	1398	14/15	0.91	0.18	-	82,94,105,119	0
4	BMA	B	1399	11/12	0.70	0.17	-	108,121,127,129	0
4	BMA	A	1400	11/12	0.44	0.21	-	124,127,129,131	0
4	NAG	B	1397	14/15	0.89	0.18	-	58,80,95,105	0
4	FUC	B	1398	10/11	0.94	0.21	-	63,71,77,81	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
2	NA	A	1395	1/1	0.97	0.34	6.01	12,12,12,12	0
3	NAG	A	1396	14/15	0.81	0.15	-	88,98,103,104	0
3	NAG	B	1395	14/15	0.82	0.15	-	74,91,101,105	0

6.5 Other polymers

There are no such residues in this entry.