



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:14 AM GMT

PDB ID : 2C6F  
Title : STRUCTURE OF HUMAN SOMATIC ANGIOTENSIN-I CONVERTING  
ENZYME N DOMAIN  
Authors : Corradi, H.R.; Schwager, S.L.U.; Nichinda, A.; Sturrock, E.D.; Acharya, K.R.  
Deposited on : 2005-11-09  
Resolution : 3.01 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 (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

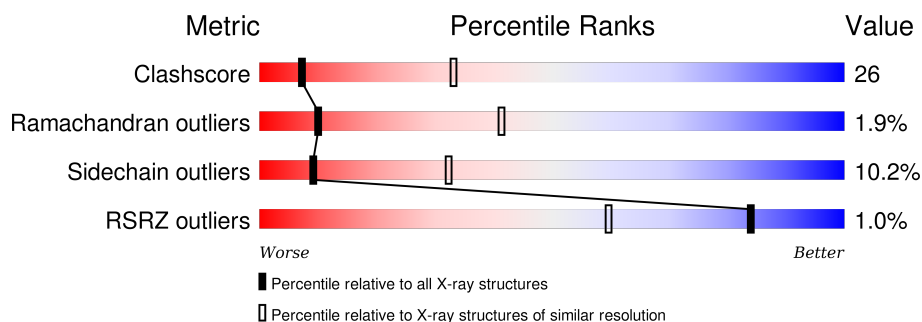
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.01 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	2117 (3.04-3.00)
Ramachandran outliers	100387	2050 (3.04-3.00)
Sidechain outliers	100360	2053 (3.04-3.00)
RSRZ outliers	91569	1788 (3.04-3.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	612	
1	B	612	

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	A	691	-	-	-	X
2	NAG	B	691	-	-	-	X

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	NAG	A	695	-	-	-	X
3	NAG	A	697	-	-	-	X
3	NAG	B	692	-	-	-	X
3	NAG	B	695	-	-	-	X
6	ACT	A	801	-	-	-	X
6	ACT	B	801	-	-	-	X
7	GOL	A	2434	-	X	-	X
7	GOL	B	2433	-	X	-	X

## 2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 9970 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 ANGIOTENSIN-CONVERTING ENZYME, SOMATIC ISOFORM.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	612	Total	C	N	O	S	6	0	0
			4884	3145	840	880	19			
1	B	611	Total	C	N	O	S	14	0	0
			4852	3124	830	879	19			

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

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

- Molecule 3 is SUGAR (N-ACETYL-D-GLUCOSAMINE) (three-letter code: NAG) (formula: C<sub>8</sub>H<sub>15</sub>NO<sub>6</sub>).



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

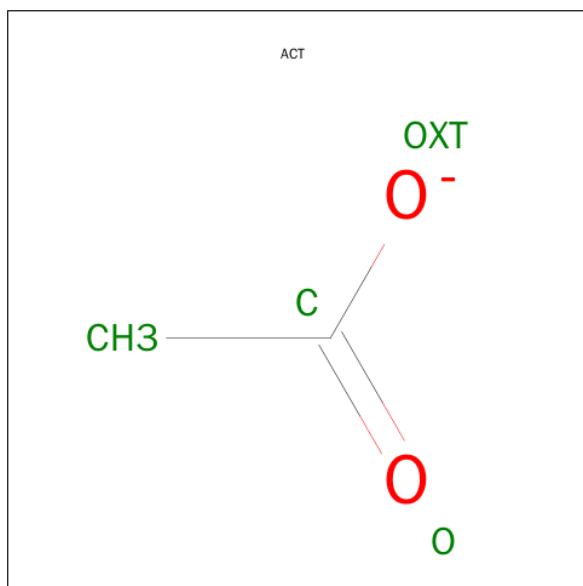
- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Zn	0	0
			1	1		
4	A	1	Total	Zn	0	0
			1	1		

- Molecule 5 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

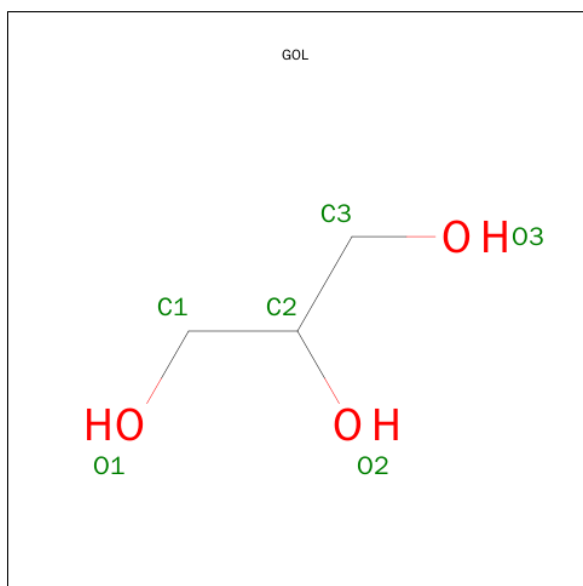
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total	Cl	0	0
			1	1		
5	A	1	Total	Cl	0	0
			1	1		

- Molecule 6 is ACETATE ION (three-letter code: ACT) (formula:  $\text{C}_2\text{H}_3\text{O}_2$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			4	2	2		
6	A	1	Total	C	O	0	0
			4	2	2		
6	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 7 is GLYCEROL (three-letter code: GOL) (formula:  $\text{C}_3\text{H}_8\text{O}_3$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			6	3	3		
7	B	1	Total	C	O	0	0
			6	3	3		

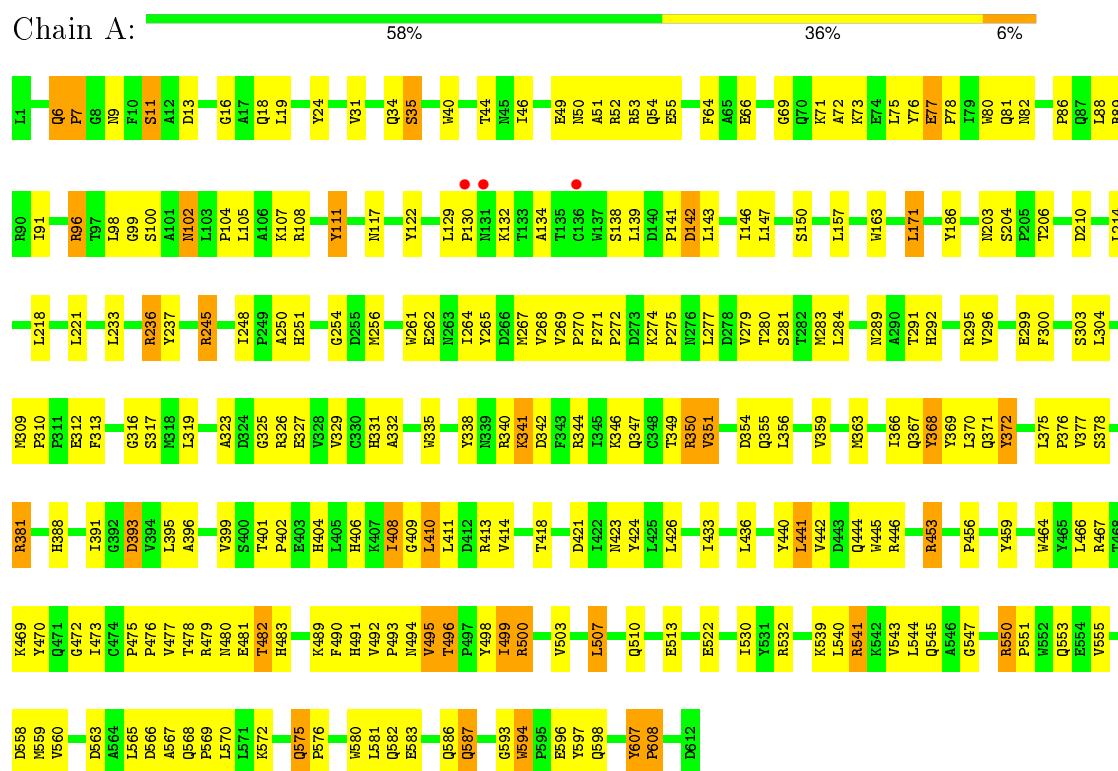
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	11	Total	O	0	0
			11	11		
8	B	13	Total	O	0	0
			13	13		

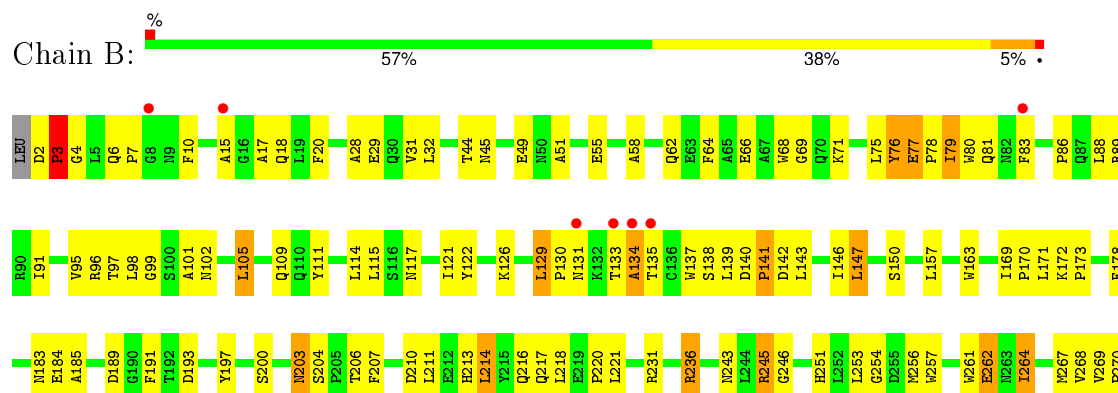
### 3 Residue-property plots

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

#### • Molecule 1: ANGIOTENSIN-CONVERTING ENZYME, SOMATIC ISOFORM



#### • Molecule 1: ANGIOTENSIN-CONVERTING ENZYME, SOMATIC ISOFORM





F271	P272	D273	K274	V275	V279	T282	M283	L284	Q285	N288	N289	H292	H293	F294	H296	V296	E299	F300	L304	M309	P310	P311	E312	F313	M314	E316	G317	E320	D324	G325	R326	E327	V328	V329	A332	V338	N339	R340	R341	D342	F343	R344	T345	R346	Q347	C348	T349	
R350	V351	T352	N353	D354	Q355	L356	E362	H365	I366	Q367	Y368	Y369	Y372	K373	D374	V377	R381	G382	A383	F387	R388	E389	A390	I391	H404	L405	H406	K407	T408	C409	L410	L411	V414	D417	T418	E419	S420	L426	K427	M428	A429	I433	A434	F435	L436	P437	F438	G439
Y440	L441	V442	D443	Q444	N445	R453	P456	Y459	N460	N464	P475	P476	V477	T478	R479	N480	E481	T482	A486	K489	F490	H491	V492	P493	H494	V495	S504	L507	Q508	F509	Q510	D529	T530	Y531	R532	A536	K539	L540	R541	K542	V543	L544	S548	P551	N552			
Q553	E554	K557	D558	M559	Y560	G561	L562	D563	A564	L565	D566	A567	Q568	P569	L570	L571	K572	Y573	F574	Q575	V580	Q584	H585	Q586	E590	G593	H594	P595	E596	Y597	Q598	M606	Y607	P608	D612													

## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	101.12Å 211.32Å 171.27Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.91 – 3.01 47.91 – 3.01	Depositor EDS
% Data completeness (in resolution range)	96.7 (47.91-3.01) 96.8 (47.91-3.01)	Depositor EDS
$R_{merge}$	0.14	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.33 (at 3.01Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.224 , 0.273 0.222 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	45.6	Xtriage
Anisotropy	0.335	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 50.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 35643 reflections	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	9970	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.90% 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, 0.375 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: GOL, ZN, ACT, NAG, CL

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	0.51	1/5042 (0.0%)	0.74	3/6885 (0.0%)
1	B	0.49	0/5010	0.86	6/6845 (0.1%)
All	All	0.50	1/10052 (0.0%)	0.80	9/13730 (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
1	A	0	1
1	B	3	2
All	All	3	3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	117	ASN	CB-CG	-5.47	1.38	1.51

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	3	PRO	CA-N-CD	-22.64	79.80	111.50
1	B	608	PRO	CA-N-CD	-18.89	85.05	111.50
1	B	607	TYR	CB-CG-CD1	-9.99	115.01	121.00
1	B	607	TYR	CB-CG-CD2	7.49	125.49	121.00
1	B	409	GLY	N-CA-C	5.92	127.91	113.10
1	B	3	PRO	CA-CB-CG	-5.55	93.45	104.00
1	A	102	ASN	CB-CA-C	-5.48	99.43	110.40
1	A	142	ASP	CB-CG-OD2	5.22	123.00	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	409	GLY	C-N-CA	-5.12	108.89	121.70

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	B	79	ILE	CA
1	B	410	LEU	CA
1	B	530	ILE	CB

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	607	TYR	Peptide
1	B	275	PRO	Peptide
1	B	409	GLY	Peptide

## 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	4884	0	4582	252	0
1	B	4852	0	4514	243	0
2	A	56	0	50	7	0
2	B	56	0	50	5	0
3	A	42	0	36	4	0
3	B	28	0	23	2	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	1	0	0	0	0
5	B	1	0	0	1	0
6	A	8	0	6	0	0
6	B	4	0	3	0	0
7	A	6	0	6	1	0
7	B	6	0	6	1	0
8	A	11	0	0	0	0
8	B	13	0	0	0	0
All	All	9970	0	9276	499	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 26.

All (499) 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:3:PRO:HD3	1:B:6:GLN:NE2	1.36	1.40
1:B:282:THR:HG21	1:B:410:LEU:CB	1.60	1.31
1:A:279:VAL:HG21	1:A:410:LEU:CD1	1.60	1.29
1:B:282:THR:CG2	1:B:410:LEU:CB	2.18	1.20
1:A:279:VAL:CG1	1:A:410:LEU:HD13	1.73	1.17
1:B:80:TRP:HA	1:B:83:PHE:CD2	1.78	1.17
1:A:279:VAL:HG21	1:A:410:LEU:HD12	1.17	1.13
1:B:80:TRP:O	1:B:89:ARG:HG2	1.48	1.12
1:A:206:THR:HG23	1:A:210:ASP:OD2	1.51	1.10
1:B:3:PRO:CD	1:B:6:GLN:NE2	2.13	1.10
1:A:279:VAL:HG11	1:A:410:LEU:HD13	1.09	1.09
3:B:692:NAG:H3	3:B:692:NAG:C8	1.48	1.09
1:A:139:LEU:HD22	1:A:163:TRP:CZ2	1.89	1.08
3:B:692:NAG:C3	3:B:692:NAG:C8	2.30	1.08
1:A:6:GLN:HG3	1:A:7:PRO:HD2	1.35	1.07
1:B:80:TRP:HA	1:B:83:PHE:CE2	1.88	1.07
1:A:6:GLN:HG3	1:A:7:PRO:CD	1.85	1.06
1:A:279:VAL:HG11	1:A:410:LEU:CD1	1.85	1.04
1:B:3:PRO:CD	1:B:6:GLN:HE22	1.70	1.03
1:A:340:ARG:NH2	2:A:694:NAG:HN2	1.56	1.03
1:A:279:VAL:CG2	1:A:410:LEU:CD1	2.37	1.01
1:A:279:VAL:CG2	1:A:410:LEU:HD12	1.91	1.00
1:A:279:VAL:HG13	1:A:410:LEU:HB2	1.49	0.95
1:A:340:ARG:HH22	2:A:694:NAG:HN2	1.09	0.95
1:B:482:THR:HG21	2:B:691:NAG:H82	1.49	0.93
1:A:72:ALA:O	1:A:76:TYR:HB2	1.71	0.91
1:B:129:LEU:HD23	1:B:612:ASP:CB	2.00	0.91
1:B:10:PHE:O	1:B:75:LEU:HD21	1.72	0.90
1:A:274:LYS:HB3	1:A:275:PRO:HD2	1.52	0.90
1:A:139:LEU:HD22	1:A:163:TRP:CH2	2.08	0.88
1:A:279:VAL:HG21	1:A:410:LEU:HD13	1.53	0.88
1:A:46:ILE:HG13	1:A:327:GLU:O	1.72	0.88
1:A:279:VAL:CB	1:A:410:LEU:HD13	2.04	0.87
1:B:29:GLU:OE1	2:B:693:NAG:H5	1.74	0.87
1:B:80:TRP:CE2	1:B:81:GLN:HG3	2.10	0.86
1:A:157:LEU:HG	1:A:607:TYR:OH	1.76	0.85
1:A:279:VAL:HG13	1:A:410:LEU:CB	2.07	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:408:ILE:O	1:A:408:ILE:HG12	1.77	0.84
1:A:279:VAL:CG2	1:A:410:LEU:HD13	2.04	0.84
1:A:270:PRO:HD3	1:A:426:LEU:HD22	1.60	0.84
1:B:353:MET:HE1	1:B:427:LYS:HE3	1.60	0.84
1:B:3:PRO:HD3	1:B:6:GLN:HE22	0.76	0.83
1:A:566:ASP:OD2	1:A:568:GLN:HB2	1.79	0.83
1:A:441:LEU:HD23	1:A:442:VAL:N	1.94	0.82
1:B:464:TRP:CE3	1:B:475:PRO:HD3	2.16	0.81
1:B:282:THR:HG21	1:B:410:LEU:CA	2.10	0.80
1:A:80:TRP:O	1:A:81:GLN:HB2	1.79	0.80
1:B:289:ASN:HB3	1:B:292:HIS:H	1.45	0.80
1:B:236:ARG:HE	1:B:267:MET:HE2	1.47	0.80
1:B:77:GLU:H	1:B:78:PRO:HD2	1.46	0.79
1:A:500:ARG:HG2	5:B:703:CL:CL	2.19	0.79
1:B:406:HIS:HA	1:B:411:LEU:O	1.82	0.79
1:B:91:ILE:O	1:B:95:VAL:HG23	1.81	0.79
1:B:2:ASP:CB	1:B:3:PRO:HD2	2.13	0.79
1:A:597:TYR:O	1:A:598:GLN:HB3	1.83	0.79
3:A:692:NAG:C8	3:A:692:NAG:O3	2.31	0.79
1:B:507:LEU:HA	1:B:510:GLN:HG3	1.66	0.78
1:B:129:LEU:CD2	1:B:612:ASP:CB	2.61	0.78
1:A:480:ASN:OD1	1:A:482:THR:HG23	1.84	0.78
1:A:329:VAL:O	1:A:346:LYS:HE3	1.83	0.77
1:B:441:LEU:HD23	1:B:442:VAL:N	1.99	0.77
1:A:129:LEU:HD12	1:A:129:LEU:N	1.99	0.76
1:A:372:TYR:HB2	1:A:375:LEU:HD12	1.68	0.76
1:B:295:ARG:HG2	1:B:314:TRP:CH2	2.20	0.76
1:A:221:LEU:HD23	1:A:433:ILE:HD12	1.67	0.76
1:B:147:LEU:HD22	1:B:256:MET:HA	1.68	0.76
1:B:78:PRO:O	1:B:79:ILE:CB	2.32	0.75
1:A:441:LEU:O	1:A:444:GLN:HB2	1.87	0.74
1:B:80:TRP:CA	1:B:83:PHE:CE2	2.70	0.74
1:B:77:GLU:H	1:B:78:PRO:CD	2.01	0.74
1:B:530:ILE:O	1:B:530:ILE:HG23	1.86	0.74
1:A:96:ARG:HG3	1:A:96:ARG:HH21	1.54	0.73
1:B:3:PRO:HA	1:B:6:GLN:NE2	2.03	0.73
1:A:369:TYR:HA	1:A:372:TYR:CE1	2.24	0.73
1:B:405:LEU:O	1:B:408:ILE:HG12	1.87	0.73
1:A:283:MET:HE1	1:A:356:LEU:HB2	1.70	0.72
1:B:129:LEU:O	1:B:131:ASN:N	2.22	0.72
1:A:279:VAL:CG1	1:A:410:LEU:CD1	2.54	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:540:LEU:O	1:B:543:VAL:HG22	1.89	0.72
1:A:69:GLY:HA3	1:A:98:LEU:HD21	1.70	0.72
1:B:332:ALA:HB2	1:B:347:GLN:HG3	1.71	0.72
1:A:279:VAL:CG1	1:A:410:LEU:CB	2.68	0.71
1:A:279:VAL:CG1	1:A:410:LEU:HB2	2.19	0.71
1:B:79:ILE:O	1:B:80:TRP:CD1	2.43	0.71
1:B:236:ARG:NE	1:B:267:MET:CE	2.53	0.71
1:B:129:LEU:N	1:B:129:LEU:HD12	2.05	0.71
1:A:279:VAL:HG12	1:A:283:MET:HG3	1.71	0.71
1:B:482:THR:CG2	2:B:691:NAG:H82	2.20	0.71
1:A:132:LYS:C	1:A:134:ALA:H	1.93	0.70
1:A:6:GLN:HG3	1:A:7:PRO:HD3	1.74	0.70
1:B:329:VAL:O	1:B:346:LYS:HE3	1.90	0.70
1:A:292:HIS:O	1:A:296:VAL:HG23	1.92	0.70
1:A:456:PRO:HA	1:A:459:TYR:CD2	2.26	0.69
1:B:282:THR:HG22	1:B:410:LEU:CB	2.19	0.69
1:A:489:LYS:HE3	1:A:491:HIS:HD2	1.56	0.69
1:B:284:LEU:HD23	1:B:351:VAL:HG11	1.73	0.69
1:B:129:LEU:C	1:B:131:ASN:H	1.96	0.69
1:B:464:TRP:CZ3	1:B:475:PRO:HD3	2.28	0.69
1:B:105:LEU:O	1:B:109:GLN:HG3	1.93	0.69
1:A:291:THR:HB	3:A:697:NAG:H2	1.73	0.69
1:A:157:LEU:HD11	1:A:477:VAL:HG13	1.73	0.68
1:B:489:LYS:O	1:B:493:PRO:HD2	1.93	0.68
1:A:147:LEU:HD22	1:A:256:MET:HA	1.75	0.68
1:A:279:VAL:HG11	1:A:410:LEU:CG	2.23	0.68
1:B:279:VAL:HG12	1:B:279:VAL:O	1.91	0.68
1:A:86:PRO:HA	1:A:89:ARG:NH1	2.09	0.68
1:B:129:LEU:HD21	1:B:137:TRP:CZ2	2.30	0.67
1:A:489:LYS:O	1:A:493:PRO:HD2	1.94	0.67
1:A:464:TRP:CE3	1:A:475:PRO:HD3	2.30	0.67
1:A:6:GLN:CG	1:A:7:PRO:HD2	2.19	0.67
1:B:203:ASN:N	1:B:203:ASN:HD22	1.93	0.66
1:B:264:ILE:HG23	1:B:264:ILE:O	1.94	0.66
1:B:171:LEU:HD21	1:B:493:PRO:HB3	1.77	0.66
1:B:58:ALA:O	1:B:62:GLN:HG3	1.95	0.66
1:B:157:LEU:HD13	1:B:476:PRO:HB2	1.77	0.66
1:A:206:THR:CG2	1:A:210:ASP:OD2	2.36	0.65
1:B:29:GLU:HG2	1:B:338:TYR:O	1.97	0.65
1:A:66:GLU:HA	1:A:98:LEU:HD22	1.78	0.65
1:B:353:MET:HE1	1:B:427:LYS:CE	2.27	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:440:TYR:O	1:A:444:GLN:HG2	1.97	0.65
1:A:395:LEU:O	1:A:399:VAL:HG23	1.98	0.64
1:B:316:GLY:HA3	1:B:344:ARG:HD3	1.79	0.64
1:B:264:ILE:CG2	1:B:264:ILE:O	2.46	0.64
1:B:221:LEU:HD12	1:B:433:ILE:HD13	1.80	0.64
1:A:340:ARG:NH2	2:A:694:NAG:N2	2.37	0.64
1:A:146:ILE:O	1:A:150:SER:HB3	1.98	0.64
1:B:236:ARG:NE	1:B:267:MET:HE2	2.13	0.64
1:A:73:LYS:HB2	1:A:77:GLU:HG3	1.78	0.64
1:B:441:LEU:HD23	1:B:441:LEU:C	2.17	0.64
1:B:183:ASN:HD21	1:B:193:ASP:HB2	1.64	0.63
1:A:332:ALA:HB2	1:A:347:GLN:HG3	1.80	0.63
1:B:607:TYR:CG	1:B:608:PRO:HD2	2.34	0.63
1:A:347:GLN:NE2	1:A:349:THR:OG1	2.31	0.63
1:A:31:VAL:HG21	1:A:64:PHE:CD1	2.33	0.63
1:A:541:ARG:HG3	1:A:541:ARG:O	1.99	0.63
1:A:251:HIS:CE1	1:A:476:PRO:HB3	2.34	0.62
1:B:203:ASN:HD22	1:B:203:ASN:H	1.45	0.62
1:A:340:ARG:NH1	2:A:694:NAG:H2	2.14	0.62
1:B:271:PHE:CD1	1:B:419:GLU:HB3	2.34	0.62
1:B:329:VAL:O	1:B:346:LYS:CE	2.48	0.62
1:B:3:PRO:CD	1:B:6:GLN:HE21	2.10	0.62
1:B:10:PHE:O	1:B:75:LEU:CD2	2.47	0.62
1:B:129:LEU:C	1:B:131:ASN:N	2.53	0.62
1:B:312:GLU:HG3	1:B:341:LYS:O	2.00	0.62
1:A:157:LEU:HD13	1:A:476:PRO:HB2	1.81	0.61
1:B:324:ASP:OD1	1:B:326:ARG:HG2	2.00	0.61
1:B:3:PRO:CA	1:B:6:GLN:NE2	2.64	0.61
1:A:96:ARG:HG3	1:A:96:ARG:NH2	2.14	0.61
1:B:279:VAL:CG1	1:B:279:VAL:O	2.48	0.61
1:A:270:PRO:HD3	1:A:426:LEU:CD2	2.30	0.61
1:B:404:HIS:HB2	1:B:529:ASP:OD2	2.01	0.61
1:B:279:VAL:HG13	1:B:282:THR:HB	1.83	0.60
1:A:73:LYS:CB	1:A:77:GLU:HG3	2.31	0.60
1:B:86:PRO:HA	1:B:89:ARG:HE	1.66	0.60
1:A:11:SER:C	1:A:13:ASP:H	2.05	0.60
1:A:147:LEU:HD22	1:A:256:MET:CA	2.32	0.59
1:B:316:GLY:C	1:B:344:ARG:HD3	2.22	0.59
1:A:377:VAL:HG13	1:A:378:SER:N	2.18	0.59
1:B:146:ILE:O	1:B:150:SER:HB3	2.02	0.59
1:B:441:LEU:CD2	1:B:442:VAL:N	2.64	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:513:GLU:OE2	1:A:572:LYS:HE3	2.02	0.59
1:B:530:ILE:CG2	1:B:530:ILE:O	2.51	0.59
1:B:142:ASP:O	1:B:146:ILE:HG13	2.02	0.59
1:A:300:PHE:HE2	1:A:304:LEU:HD11	1.68	0.59
1:B:98:LEU:HB2	1:B:102:ASN:ND2	2.18	0.59
1:B:504:SER:O	1:B:508:GLN:HB3	2.03	0.59
1:B:7:PRO:HB2	1:B:71:LYS:HG2	1.85	0.58
1:A:248:ILE:O	1:A:473:ILE:HA	2.03	0.58
1:B:77:GLU:N	1:B:78:PRO:CD	2.66	0.58
1:B:2:ASP:CB	1:B:3:PRO:CD	2.78	0.58
1:A:236:ARG:HG2	1:A:267:MET:CE	2.33	0.58
1:B:139:LEU:HD22	1:B:163:TRP:CZ2	2.39	0.58
1:B:580:TRP:O	1:B:584:GLN:HG2	2.04	0.58
1:A:147:LEU:HD22	1:A:256:MET:N	2.19	0.57
1:A:261:TRP:N	1:A:261:TRP:CD1	2.71	0.57
1:B:218:LEU:CD2	1:B:570:LEU:HD23	2.34	0.57
1:A:489:LYS:O	1:A:491:HIS:N	2.38	0.57
1:A:75:LEU:HD23	1:A:75:LEU:H	1.69	0.57
1:A:338:TYR:HD2	1:A:377:VAL:HG23	1.68	0.57
1:A:206:THR:HG23	1:A:210:ASP:CG	2.23	0.57
1:A:597:TYR:O	1:A:598:GLN:CB	2.53	0.57
1:A:142:ASP:O	1:A:146:ILE:HG13	2.05	0.57
1:B:541:ARG:O	1:B:541:ARG:HG2	2.04	0.57
1:A:51:ALA:O	1:A:55:GLU:HG3	2.04	0.57
1:B:97:THR:O	1:B:97:THR:HG22	2.04	0.56
1:B:80:TRP:CG	1:B:81:GLN:N	2.73	0.56
1:A:406:HIS:HD2	1:A:413:ARG:O	1.89	0.56
1:A:446:ARG:NH2	1:A:496:THR:O	2.35	0.56
1:A:404:HIS:O	1:A:408:ILE:HG23	2.05	0.56
1:A:75:LEU:N	1:A:75:LEU:HD23	2.21	0.56
1:A:129:LEU:HD12	1:A:129:LEU:H	1.71	0.55
1:B:283:MET:HE3	1:B:356:LEU:HD22	1.89	0.55
1:A:98:LEU:HB2	1:A:102:ASN:HD21	1.71	0.55
1:A:77:GLU:N	1:A:78:PRO:HD2	2.21	0.55
1:B:310:PRO:O	1:B:313:PHE:HB3	2.06	0.55
1:B:141:PRO:HG2	1:B:142:ASP:N	2.22	0.55
1:A:261:TRP:O	1:A:264:ILE:HG12	2.07	0.55
1:A:367:GLN:O	1:A:371:GLN:HG2	2.05	0.55
1:B:366:ILE:O	1:B:369:TYR:N	2.39	0.55
1:B:213:HIS:O	1:B:216:GLN:HB2	2.07	0.55
1:A:129:LEU:N	1:A:129:LEU:CD1	2.67	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:469:LYS:HE2	1:A:470:TYR:CZ	2.42	0.55
1:B:279:VAL:HG21	1:B:405:LEU:HD13	1.89	0.54
1:B:270:PRO:O	1:B:272:PRO:HD3	2.08	0.54
1:A:277:LEU:CD1	1:A:424:TYR:HA	2.37	0.54
1:B:270:PRO:HD3	1:B:426:LEU:HD22	1.89	0.54
1:A:91:ILE:HG23	1:A:378:SER:OG	2.06	0.54
1:B:316:GLY:CA	1:B:344:ARG:HD3	2.38	0.54
1:A:104:PRO:HG2	1:A:107:LYS:HD2	1.88	0.54
1:B:326:ARG:O	1:B:328:VAL:HG13	2.09	0.53
1:B:17:ALA:O	1:B:20:PHE:HB3	2.08	0.53
1:B:64:PHE:CE1	1:B:68:TRP:NE1	2.75	0.53
1:B:478:THR:O	1:B:479:ARG:HD2	2.08	0.53
1:A:492:VAL:N	1:A:493:PRO:HD2	2.23	0.53
1:B:607:TYR:C	1:B:608:PRO:O	2.43	0.53
1:B:251:HIS:CE1	1:B:476:PRO:HB3	2.43	0.53
1:A:147:LEU:O	1:A:254:GLY:HA2	2.09	0.53
1:B:99:GLY:O	1:B:102:ASN:ND2	2.36	0.53
1:B:147:LEU:O	1:B:254:GLY:HA2	2.08	0.53
1:A:54:GLN:HG3	1:A:55:GLU:N	2.24	0.53
1:A:607:TYR:CD1	1:A:608:PRO:HA	2.44	0.52
1:A:326:ARG:HD2	3:A:692:NAG:HN2	1.72	0.52
1:B:218:LEU:O	1:B:221:LEU:HB2	2.09	0.52
1:A:350:ARG:HD2	1:A:351:VAL:H	1.75	0.52
1:B:29:GLU:OE1	2:B:693:NAG:C5	2.54	0.52
1:B:29:GLU:CG	1:B:338:TYR:O	2.57	0.52
1:B:326:ARG:O	1:B:327:GLU:C	2.48	0.52
1:A:214:LEU:HD11	1:A:565:LEU:HB3	1.90	0.52
1:A:510:GLN:OE1	1:A:566:ASP:N	2.40	0.52
1:A:480:ASN:OD1	1:A:482:THR:CG2	2.57	0.52
1:B:426:LEU:O	1:B:429:ALA:HB3	2.10	0.52
1:A:441:LEU:C	1:A:441:LEU:HD23	2.29	0.52
1:A:98:LEU:HB2	1:A:102:ASN:ND2	2.24	0.52
1:B:292:HIS:O	1:B:296:VAL:HG23	2.09	0.52
1:A:69:GLY:HA3	1:A:98:LEU:CD2	2.38	0.52
1:A:147:LEU:CD2	1:A:256:MET:HA	2.41	0.51
1:B:408:ILE:CG1	1:B:409:GLY:N	2.70	0.51
1:A:66:GLU:HA	1:A:98:LEU:CD2	2.41	0.51
1:A:312:GLU:HG2	1:A:341:LYS:HE2	1.93	0.51
1:A:269:VAL:HG13	1:A:269:VAL:O	2.09	0.51
1:B:71:LYS:HE3	1:B:75:LEU:HD11	1.91	0.51
1:A:204:SER:OG	1:A:206:THR:HG22	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:294:PHE:HE1	1:B:349:THR:HG1	1.59	0.51
1:A:495:VAL:O	1:A:495:VAL:HG12	2.10	0.51
1:A:323:ALA:C	1:A:325:GLY:H	2.14	0.50
1:A:445:TRP:HB2	1:A:466:LEU:HD12	1.94	0.50
1:B:3:PRO:N	1:B:6:GLN:HE21	2.10	0.50
1:B:200:SER:O	1:B:203:ASN:ND2	2.44	0.50
1:A:583:GLU:O	1:A:587:GLN:HG3	2.12	0.50
1:B:510:GLN:OE1	1:B:566:ASP:N	2.44	0.49
1:B:563:ASP:OD1	1:B:563:ASP:O	2.30	0.49
1:A:139:LEU:CD2	1:A:163:TRP:CZ2	2.79	0.49
1:A:40:TRP:CE2	1:A:44:THR:HG21	2.46	0.49
1:A:478:THR:O	1:A:479:ARG:HD2	2.13	0.49
1:B:138:SER:O	1:B:143:LEU:HG	2.11	0.49
1:B:362:GLU:O	1:B:365:HIS:HB2	2.13	0.49
1:A:265:TYR:CZ	1:A:269:VAL:HG23	2.47	0.49
1:A:268:VAL:O	1:A:269:VAL:C	2.51	0.49
1:A:583:GLU:O	1:A:587:GLN:CG	2.61	0.49
1:A:218:LEU:HD13	1:A:436:LEU:HD13	1.95	0.49
1:B:2:ASP:O	1:B:3:PRO:HB2	2.12	0.49
1:B:129:LEU:CD1	1:B:129:LEU:N	2.76	0.49
1:B:332:ALA:CB	1:B:347:GLN:HG3	2.39	0.49
1:B:256:MET:HB3	1:B:257:TRP:CE3	2.48	0.48
1:B:3:PRO:CA	1:B:6:GLN:HE21	2.26	0.48
1:A:499:ILE:HG13	1:A:499:ILE:O	2.14	0.48
1:B:593:GLY:HA3	7:B:2433:GOL:O1	2.13	0.48
1:B:79:ILE:O	1:B:81:GLN:N	2.41	0.48
1:A:393:ASP:O	1:A:396:ALA:HB3	2.13	0.48
1:A:271:PHE:HB2	1:A:423:ASN:HD21	1.78	0.48
1:A:300:PHE:CE2	1:A:304:LEU:HD11	2.48	0.48
1:A:270:PRO:O	1:A:272:PRO:HD3	2.13	0.48
1:A:49:GLU:HG3	1:A:53:ARG:HE	1.79	0.48
1:A:99:GLY:HA2	1:A:186:TYR:CE2	2.48	0.48
1:B:169:ILE:N	1:B:170:PRO:HD2	2.29	0.48
1:A:399:VAL:HA	1:A:404:HIS:CD2	2.48	0.48
1:A:551:PRO:O	1:A:555:VAL:HG23	2.14	0.48
1:B:51:ALA:O	1:B:55:GLU:HG3	2.14	0.48
1:A:507:LEU:HB3	1:A:565:LEU:CD2	2.44	0.48
1:A:441:LEU:HD23	1:A:442:VAL:CA	2.43	0.48
1:A:171:LEU:HD21	1:A:493:PRO:HB3	1.95	0.48
1:B:482:THR:HG21	2:B:691:NAG:C8	2.34	0.48
1:B:295:ARG:HG2	1:B:314:TRP:HH2	1.72	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:139:LEU:HA	1:A:143:LEU:HB2	1.95	0.47
1:B:44:THR:O	1:B:328:VAL:HG12	2.14	0.47
1:A:453:ARG:O	1:A:453:ARG:HG2	2.14	0.47
1:A:539:LYS:HE3	1:A:559:MET:O	2.13	0.47
1:A:274:LYS:HB3	1:A:275:PRO:CD	2.34	0.47
1:B:606:ASN:O	1:B:608:PRO:O	2.32	0.47
1:B:80:TRP:CZ2	1:B:81:GLN:HG3	2.48	0.47
1:A:98:LEU:HD12	1:A:102:ASN:OD1	2.13	0.47
1:A:236:ARG:HG2	1:A:267:MET:HE3	1.96	0.47
1:B:77:GLU:O	1:B:77:GLU:CG	2.62	0.47
1:B:141:PRO:O	1:B:142:ASP:C	2.51	0.47
1:A:482:THR:HG21	2:A:691:NAG:HN2	1.80	0.47
1:B:299:GLU:HG2	1:B:300:PHE:N	2.30	0.47
1:B:243:ASN:OD1	1:B:245:ARG:N	2.41	0.47
1:A:355:GLN:O	1:A:359:VAL:HG23	2.15	0.47
1:A:274:LYS:CB	1:A:275:PRO:HD2	2.35	0.47
1:B:98:LEU:HB2	1:B:102:ASN:HD21	1.78	0.47
1:B:218:LEU:HD22	1:B:570:LEU:HD23	1.95	0.47
1:B:236:ARG:HE	1:B:267:MET:CE	2.11	0.47
1:A:122:TYR:CD1	1:A:493:PRO:HG3	2.49	0.47
1:A:277:LEU:HD11	1:A:424:TYR:HB2	1.97	0.47
1:A:597:TYR:CD1	1:A:597:TYR:C	2.88	0.47
1:A:490:PHE:O	1:A:493:PRO:HG2	2.15	0.47
1:B:551:PRO:HB2	1:B:554:GLU:HG3	1.96	0.47
1:B:329:VAL:HB	1:B:346:LYS:HE3	1.97	0.46
1:A:489:LYS:O	1:A:490:PHE:C	2.53	0.46
1:A:34:GLN:HG3	1:A:35:SER:H	1.80	0.46
1:B:418:THR:HG22	1:B:419:GLU:N	2.29	0.46
1:B:270:PRO:HD3	1:B:426:LEU:CD2	2.44	0.46
1:A:284:LEU:HD23	1:A:351:VAL:HG11	1.97	0.46
1:B:495:VAL:O	1:B:495:VAL:HG12	2.14	0.46
1:B:262:GLU:N	1:B:262:GLU:OE1	2.48	0.46
1:A:129:LEU:H	1:A:129:LEU:CD1	2.27	0.46
1:B:486:ALA:O	1:B:492:VAL:HG21	2.16	0.46
1:A:456:PRO:HA	1:A:459:TYR:CE2	2.50	0.46
1:A:77:GLU:N	1:A:78:PRO:CD	2.78	0.46
1:B:597:TYR:C	1:B:597:TYR:CD1	2.89	0.46
1:B:133:THR:O	1:B:134:ALA:C	2.53	0.46
1:B:217:GLN:O	1:B:220:PRO:HD2	2.15	0.46
1:A:507:LEU:HG	1:A:507:LEU:O	2.16	0.46
1:B:75:LEU:HD23	1:B:76:TYR:CZ	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:80:TRP:CE2	1:A:81:GLN:HG3	2.50	0.46
1:B:492:VAL:HB	1:B:493:PRO:CD	2.46	0.46
1:B:408:ILE:HG12	1:B:409:GLY:N	2.30	0.46
1:A:279:VAL:HG11	1:A:410:LEU:HD22	1.98	0.46
1:A:80:TRP:O	1:A:81:GLN:CB	2.57	0.46
1:B:437:PRO:O	1:B:441:LEU:HB3	2.15	0.46
1:B:529:ASP:OD1	1:B:531:TYR:HB2	2.16	0.46
1:A:143:LEU:O	1:A:147:LEU:HD12	2.15	0.46
1:A:338:TYR:HD2	1:A:377:VAL:CG2	2.29	0.46
1:A:481:GLU:OE2	1:A:481:GLU:HA	2.16	0.46
1:B:140:ASP:OD1	1:B:140:ASP:C	2.53	0.46
1:A:418:THR:O	1:A:421:ASP:N	2.49	0.46
1:A:510:GLN:HE21	1:A:560:VAL:HG11	1.81	0.46
1:A:340:ARG:CZ	2:A:694:NAG:HN2	2.21	0.46
1:B:147:LEU:HD22	1:B:256:MET:CA	2.44	0.46
1:B:607:TYR:CD2	1:B:608:PRO:HD2	2.50	0.46
1:A:204:SER:C	1:A:206:THR:H	2.18	0.45
1:B:45:ASN:OD1	1:B:326:ARG:HB2	2.17	0.45
1:B:44:THR:O	1:B:44:THR:HG22	2.15	0.45
1:B:435:PHE:CE2	1:B:439:GLY:HA3	2.51	0.45
1:B:507:LEU:O	1:B:507:LEU:HG	2.16	0.45
1:B:141:PRO:CG	1:B:142:ASP:N	2.79	0.45
1:B:369:TYR:HD2	1:B:372:TYR:HH	1.64	0.45
1:B:129:LEU:H	1:B:129:LEU:HD12	1.78	0.45
1:B:536:ALA:O	1:B:539:LYS:N	2.50	0.45
1:B:406:HIS:O	1:B:408:ILE:O	2.34	0.45
1:B:309:MET:HB3	1:B:313:PHE:CD2	2.52	0.45
1:B:438:PHE:O	1:B:442:VAL:HG23	2.17	0.45
1:B:570:LEU:CD1	1:B:574:PHE:HE1	2.30	0.45
1:B:607:TYR:CD2	1:B:608:PRO:CD	3.00	0.45
1:B:98:LEU:HB3	1:B:101:ALA:HB3	1.99	0.45
1:B:191:PHE:CZ	1:B:197:TYR:HD1	2.35	0.45
2:A:693:NAG:H61	2:A:694:NAG:C7	2.47	0.45
1:A:369:TYR:HA	1:A:372:TYR:HE1	1.79	0.45
1:A:469:LYS:HE2	1:A:470:TYR:CE1	2.51	0.45
1:B:539:LYS:HE2	1:B:559:MET:O	2.16	0.45
1:A:570:LEU:HD12	1:A:570:LEU:O	2.17	0.45
1:B:32:LEU:CD1	1:B:377:VAL:HG11	2.46	0.45
1:B:568:GLN:HA	1:B:571:LEU:HD12	1.98	0.45
1:B:71:LYS:HE3	1:B:75:LEU:CD1	2.47	0.45
1:B:289:ASN:HB2	1:B:292:HIS:HB2	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:289:ASN:HB3	1:A:292:HIS:H	1.82	0.45
1:A:31:VAL:O	1:A:34:GLN:HG3	2.17	0.45
1:B:3:PRO:O	1:B:4:GLY:C	2.52	0.45
1:A:377:VAL:CG1	1:A:378:SER:N	2.80	0.45
1:B:77:GLU:N	1:B:78:PRO:HD2	2.23	0.44
1:A:279:VAL:O	1:A:281:SER:N	2.50	0.44
1:B:218:LEU:HD21	1:B:570:LEU:HD23	1.99	0.44
1:A:575:GLN:HB3	1:A:576:PRO:HD3	1.99	0.44
1:B:111:TYR:CE2	1:B:115:LEU:HD11	2.53	0.44
1:A:480:ASN:OD1	1:A:483:HIS:CE1	2.71	0.44
1:B:417:ASP:HB3	1:B:420:SER:OG	2.16	0.44
1:B:86:PRO:HB3	1:B:89:ARG:HH21	1.81	0.44
1:A:300:PHE:CD2	1:A:395:LEU:HD23	2.53	0.44
1:B:172:LYS:N	1:B:173:PRO:HD2	2.32	0.44
1:A:18:GLN:HG2	1:A:88:LEU:HD22	2.00	0.44
1:B:80:TRP:NE1	1:B:81:GLN:HG3	2.32	0.44
1:A:77:GLU:H	1:A:78:PRO:HD2	1.82	0.44
1:B:117:ASN:O	1:B:121:ILE:HG13	2.18	0.44
1:A:218:LEU:O	1:A:221:LEU:HB2	2.18	0.44
1:A:279:VAL:HG11	1:A:410:LEU:CD2	2.47	0.44
1:A:283:MET:CE	1:A:356:LEU:HD22	2.47	0.44
1:B:288:TRP:CH2	1:B:296:VAL:HG21	2.51	0.44
1:A:44:THR:O	1:A:326:ARG:NH2	2.47	0.44
1:A:221:LEU:HD23	1:A:433:ILE:CD1	2.42	0.44
1:B:171:LEU:HD21	1:B:493:PRO:CB	2.47	0.44
1:B:183:ASN:HD21	1:B:193:ASP:CB	2.30	0.44
1:B:69:GLY:HA3	1:B:98:LEU:CD2	2.47	0.44
1:B:383:ALA:HB3	1:B:387:PHE:CG	2.53	0.44
1:B:75:LEU:HD23	1:B:76:TYR:OH	2.17	0.43
1:A:269:VAL:CG1	1:A:269:VAL:O	2.65	0.43
1:A:593:GLY:HA3	7:A:2434:GOL:O1	2.18	0.43
1:A:233:LEU:HD22	1:A:237:TYR:HE1	1.83	0.43
1:B:507:LEU:HB3	1:B:565:LEU:CD2	2.49	0.43
1:A:24:TYR:CE1	1:A:64:PHE:HE2	2.35	0.43
1:A:494:ASN:HD22	1:A:494:ASN:N	2.15	0.43
1:B:207:PHE:CZ	1:B:211:LEU:HD11	2.53	0.43
1:B:15:ALA:HA	1:B:18:GLN:HB2	2.00	0.43
1:A:441:LEU:CD1	1:A:467:ARG:HA	2.48	0.43
1:B:122:TYR:CE1	1:B:163:TRP:HZ2	2.36	0.43
1:A:71:LYS:O	1:A:75:LEU:HD23	2.17	0.43
1:B:352:THR:HG22	1:B:353:MET:N	2.32	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:489:LYS:O	1:B:490:PHE:C	2.56	0.43
1:A:464:TRP:CZ3	1:A:475:PRO:HD3	2.52	0.43
1:B:442:VAL:O	1:B:445:TRP:HB3	2.19	0.43
1:B:317:SER:HA	1:B:344:ARG:HB3	2.01	0.43
1:B:245:ARG:O	1:B:595:PRO:HD2	2.19	0.43
1:A:580:TRP:CZ3	1:A:581:LEU:HD23	2.54	0.43
1:B:83:PHE:CD2	1:B:88:LEU:HD12	2.53	0.43
1:A:503:VAL:HG13	1:A:565:LEU:HD11	2.00	0.43
1:A:366:ILE:O	1:A:370:LEU:HG	2.18	0.43
1:A:498:TYR:C	1:A:500:ARG:H	2.22	0.43
1:A:478:THR:HG22	1:A:479:ARG:N	2.34	0.43
1:A:510:GLN:HG2	1:A:569:PRO:HG3	2.00	0.43
1:B:567:ALA:O	1:B:568:GLN:C	2.57	0.43
1:A:245:ARG:O	1:A:596:GLU:HB2	2.18	0.43
1:A:472:GLY:CA	1:A:594:TRP:CE2	3.02	0.42
1:A:594:TRP:N	1:A:594:TRP:CE3	2.87	0.42
1:A:141:PRO:O	1:A:142:ASP:C	2.57	0.42
1:A:381:ARG:O	1:A:547:GLY:HA2	2.20	0.42
1:B:304:LEU:HD21	1:B:530:ILE:O	2.19	0.42
1:A:132:LYS:C	1:A:134:ALA:N	2.63	0.42
1:B:140:ASP:HA	1:B:141:PRO:HA	1.86	0.42
1:B:115:LEU:HD21	1:B:178:PHE:CE1	2.54	0.42
1:A:368:TYR:CE2	1:A:544:LEU:HA	2.54	0.42
1:A:366:ILE:HA	1:A:369:TYR:CD1	2.54	0.42
1:A:490:PHE:C	1:A:490:PHE:CD2	2.92	0.42
1:A:31:VAL:HG21	1:A:64:PHE:HD1	1.79	0.42
1:A:472:GLY:HA2	1:A:594:TRP:CE2	2.53	0.42
1:A:316:GLY:C	1:A:344:ARG:HD3	2.39	0.42
1:B:390:ALA:O	1:B:391:ILE:C	2.57	0.42
1:B:405:LEU:O	1:B:408:ILE:CG1	2.63	0.42
1:B:460:ASN:N	1:B:481:GLU:OE2	2.34	0.42
1:A:143:LEU:HD13	1:A:163:TRP:HB2	2.01	0.42
1:A:607:TYR:HA	1:A:608:PRO:HA	1.76	0.42
1:B:570:LEU:CD1	1:B:574:PHE:CE1	3.03	0.42
1:A:510:GLN:HB3	1:A:510:GLN:HE21	1.63	0.42
1:A:335:TRP:CD1	1:A:335:TRP:N	2.87	0.42
1:B:83:PHE:HB2	1:B:89:ARG:HG3	2.02	0.42
1:B:436:LEU:N	1:B:437:PRO:HD2	2.35	0.42
1:A:210:ASP:O	1:A:214:LEU:HB2	2.19	0.42
1:B:362:GLU:HA	1:B:362:GLU:OE1	2.19	0.42
1:B:283:MET:HE3	1:B:356:LEU:CD2	2.49	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:312:GLU:N	1:A:312:GLU:OE1	2.46	0.41
1:A:108:ARG:O	1:A:111:TYR:HB3	2.20	0.41
1:B:406:HIS:CA	1:B:411:LEU:O	2.61	0.41
1:B:531:TYR:O	1:B:532:ARG:HB2	2.20	0.41
1:A:567:ALA:O	1:A:570:LEU:HB3	2.20	0.41
1:A:309:MET:HA	1:A:310:PRO:HD3	1.93	0.41
1:B:541:ARG:CG	1:B:541:ARG:O	2.66	0.41
1:B:279:VAL:CG2	1:B:405:LEU:HD13	2.49	0.41
1:A:498:TYR:O	1:A:500:ARG:N	2.54	0.41
1:B:382:GLY:O	1:B:383:ALA:C	2.58	0.41
1:A:303:SER:HB2	1:A:532:ARG:HG2	2.03	0.41
1:A:157:LEU:HD23	1:A:607:TYR:CZ	2.56	0.41
1:A:326:ARG:HD2	3:A:692:NAG:N2	2.36	0.41
1:B:114:LEU:HA	1:B:114:LEU:HD23	1.93	0.41
1:B:339:ASN:OD1	1:B:342:ASP:N	2.50	0.41
1:A:350:ARG:HD2	1:A:351:VAL:HG23	2.01	0.41
1:B:268:VAL:O	1:B:269:VAL:C	2.57	0.41
1:B:557:LYS:O	1:B:561:GLY:HA2	2.19	0.41
1:A:317:SER:HB3	1:A:319:LEU:HD21	2.01	0.41
1:A:331:HIS:CB	1:A:490:PHE:CE1	3.04	0.41
1:A:441:LEU:HD23	1:A:442:VAL:HA	2.03	0.41
1:A:375:LEU:O	1:A:376:PRO:C	2.58	0.41
1:A:104:PRO:CG	1:A:107:LYS:HD2	2.51	0.41
1:B:79:ILE:O	1:B:80:TRP:CG	2.74	0.41
1:A:138:SER:O	1:A:139:LEU:C	2.59	0.41
1:B:147:LEU:HD13	1:B:256:MET:CE	2.51	0.41
1:A:11:SER:C	1:A:13:ASP:N	2.73	0.41
1:B:243:ASN:OD1	1:B:246:GLY:N	2.54	0.41
1:A:401:THR:HA	1:A:402:PRO:HD3	1.97	0.41
1:B:210:ASP:O	1:B:214:LEU:HB2	2.21	0.41
1:A:351:VAL:HG12	1:A:351:VAL:O	2.21	0.41
1:A:586:GLN:O	1:A:587:GLN:C	2.58	0.41
1:A:309:MET:HB3	1:A:313:PHE:CD2	2.56	0.41
1:B:368:TYR:CE2	1:B:544:LEU:HA	2.56	0.41
1:A:550:ARG:HG2	1:A:550:ARG:HH11	1.85	0.41
1:B:261:TRP:CD1	1:B:261:TRP:N	2.88	0.41
1:A:279:VAL:HG13	1:A:410:LEU:HB3	1.98	0.40
1:A:157:LEU:CD2	1:A:607:TYR:CZ	3.03	0.40
1:A:500:ARG:H	1:A:500:ARG:HG3	1.54	0.40
1:B:441:LEU:CD2	1:B:441:LEU:C	2.89	0.40
1:B:28:ALA:HA	1:B:31:VAL:HG12	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:550:ARG:HG2	1:A:550:ARG:NH1	2.36	0.40
1:A:279:VAL:O	1:A:279:VAL:HG12	2.22	0.40
1:A:6:GLN:CG	1:A:7:PRO:CD	2.77	0.40
1:B:253:LEU:HG	1:B:261:TRP:CE2	2.56	0.40
1:A:16:GLY:O	1:A:19:LEU:HB2	2.21	0.40
1:A:6:GLN:HE21	1:A:6:GLN:HB2	1.60	0.40
1:B:31:VAL:HG11	1:B:64:PHE:CD2	2.56	0.40
1:A:50:ASN:HA	1:A:53:ARG:HD2	2.02	0.40
1:B:204:SER:C	1:B:206:THR:H	2.23	0.40
1:B:381:ARG:HB2	1:B:548:SER:HB2	2.02	0.40
1:B:456:PRO:HA	1:B:459:TYR:CD2	2.57	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	610/612 (100%)	537 (88%)	60 (10%)	13 (2%)	9	38
1	B	609/612 (100%)	526 (86%)	73 (12%)	10 (2%)	12	46
All	All	1219/1224 (100%)	1063 (87%)	133 (11%)	23 (2%)	10	41

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	608	PRO
1	B	3	PRO
1	B	608	PRO
1	A	7	PRO
1	A	250	ALA
1	A	280	THR
1	A	545	GLN

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Mol	Chain	Res	Type
1	B	130	PRO
1	B	134	ALA
1	B	135	THR
1	B	391	ILE
1	A	351	VAL
1	A	111	TYR
1	A	414	VAL
1	B	77	GLU
1	A	391	ILE
1	A	495	VAL
1	B	185	ALA
1	B	408	ILE
1	A	130	PRO
1	B	79	ILE
1	A	499	ILE
1	A	530	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	496/526 (94%)	448 (90%)	48 (10%)	10	36
1	B	489/526 (93%)	437 (89%)	52 (11%)	8	30
All	All	985/1052 (94%)	885 (90%)	100 (10%)	9	32

All (100) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	GLN
1	A	9	ASN
1	A	11	SER
1	A	35	SER
1	A	52	ARG
1	A	77	GLU
1	A	82	ASN

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Mol	Chain	Res	Type
1	A	96	ARG
1	A	100	SER
1	A	105	LEU
1	A	171	LEU
1	A	203	ASN
1	A	236	ARG
1	A	245	ARG
1	A	262	GLU
1	A	295	ARG
1	A	299	GLU
1	A	341	LYS
1	A	342	ASP
1	A	350	ARG
1	A	354	ASP
1	A	363	MET
1	A	368	TYR
1	A	372	TYR
1	A	381	ARG
1	A	388	HIS
1	A	393	ASP
1	A	408	ILE
1	A	410	LEU
1	A	411	LEU
1	A	441	LEU
1	A	453	ARG
1	A	482	THR
1	A	496	THR
1	A	500	ARG
1	A	507	LEU
1	A	522	GLU
1	A	540	LEU
1	A	541	ARG
1	A	543	VAL
1	A	550	ARG
1	A	553	GLN
1	A	558	ASP
1	A	563	ASP
1	A	575	GLN
1	A	582	GLN
1	A	587	GLN
1	A	594	TRP
1	B	3	PRO

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Mol	Chain	Res	Type
1	B	49	GLU
1	B	66	GLU
1	B	76	TYR
1	B	96	ARG
1	B	105	LEU
1	B	126	LYS
1	B	129	LEU
1	B	141	PRO
1	B	147	LEU
1	B	184	GLU
1	B	189	ASP
1	B	203	ASN
1	B	214	LEU
1	B	231	ARG
1	B	236	ARG
1	B	245	ARG
1	B	262	GLU
1	B	264	ILE
1	B	274	LYS
1	B	285	GLN
1	B	289	ASN
1	B	299	GLU
1	B	312	GLU
1	B	320	GLU
1	B	342	ASP
1	B	354	ASP
1	B	368	TYR
1	B	372	TYR
1	B	374	ASP
1	B	388	HIS
1	B	408	ILE
1	B	411	LEU
1	B	414	VAL
1	B	418	THR
1	B	441	LEU
1	B	443	ASP
1	B	453	ARG
1	B	480	ASN
1	B	489	LYS
1	B	507	LEU
1	B	508	GLN
1	B	510	GLN

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Mol	Chain	Res	Type
1	B	542	LYS
1	B	553	GLN
1	B	572	LYS
1	B	575	GLN
1	B	586	GLN
1	B	590	GLU
1	B	598	GLN
1	B	607	TYR
1	B	608	PRO

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	GLN
1	A	50	ASN
1	A	81	GLN
1	A	347	GLN
1	A	406	HIS
1	A	491	HIS
1	A	494	ASN
1	A	545	GLN
1	A	575	GLN
1	B	6	GLN
1	B	18	GLN
1	B	50	ASN
1	B	183	ASN
1	B	203	ASN
1	B	416	ASN
1	B	545	GLN
1	B	575	GLN
1	B	579	GLN
1	B	582	GLN
1	B	587	GLN

### 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$
2	NAG	A	691	1,2	14,14,15	0.56	0	15,19,21	0.76	1 (6%)
2	NAG	A	693	1,2	14,14,15	0.56	0	15,19,21	0.76	1 (6%)
2	NAG	A	694	2	14,14,15	0.49	0	15,19,21	0.72	1 (6%)
2	NAG	A	696	2	14,14,15	0.50	0	15,19,21	0.72	1 (6%)
2	NAG	B	691	1,2	14,14,15	0.55	0	15,19,21	0.76	1 (6%)
2	NAG	B	693	1,2	14,14,15	2.43	1 (7%)	15,19,21	0.88	1 (6%)
2	NAG	B	694	2	14,14,15	2.42	1 (7%)	15,19,21	0.88	1 (6%)
2	NAG	B	696	2	14,14,15	0.50	0	15,19,21	0.72	1 (6%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	A	691	1,2	-	0/6/23/26	0/1/1/1
2	NAG	A	693	1,2	-	0/6/23/26	0/1/1/1
2	NAG	A	694	2	-	0/6/23/26	0/1/1/1
2	NAG	A	696	2	-	0/6/23/26	0/1/1/1
2	NAG	B	691	1,2	-	0/6/23/26	0/1/1/1
2	NAG	B	693	1,2	-	0/6/23/26	0/1/1/1
2	NAG	B	694	2	-	0/6/23/26	0/1/1/1
2	NAG	B	696	2	-	0/6/23/26	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	693	NAG	C8-C7	-9.05	1.32	1.50
2	B	694	NAG	C8-C7	-9.00	1.32	1.50

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	691	NAG	C2-N2-C7	-2.37	120.00	123.04
2	A	693	NAG	C2-N2-C7	-2.34	120.04	123.04
2	B	691	NAG	C2-N2-C7	-2.33	120.05	123.04
2	B	693	NAG	C3-C2-N2	-2.32	105.00	110.56
2	B	694	NAG	C3-C2-N2	-2.32	105.01	110.56
2	A	694	NAG	C2-N2-C7	-2.18	120.24	123.04
2	A	696	NAG	C2-N2-C7	-2.17	120.25	123.04
2	B	696	NAG	C2-N2-C7	-2.17	120.25	123.04

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	691	NAG	1	0
2	A	693	NAG	1	0
2	A	694	NAG	6	0
2	B	691	NAG	3	0
2	B	693	NAG	2	0

## 5.6 Ligand geometry

Of 14 ligands modelled in this entry, 4 are monoatomic - leaving 10 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	GOL	A	2434	-	5,5,5	3.46	3 (60%)	5,5,5	1.84	3 (60%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	NAG	A	692	1	14,14,15	4.05	3 (21%)	15,19,21	1.89	4 (26%)
3	NAG	A	695	1	14,14,15	0.56	0	15,19,21	0.75	1 (6%)
3	NAG	A	697	1	14,14,15	0.56	0	15,19,21	0.76	1 (6%)
6	ACT	A	800	-	1,3,3	3.08	1 (100%)	0,3,3	0.00	-
6	ACT	A	801	-	1,3,3	3.56	1 (100%)	0,3,3	0.00	-
7	GOL	B	2433	-	5,5,5	3.49	3 (60%)	5,5,5	2.01	4 (80%)
3	NAG	B	692	1	14,14,15	3.89	2 (14%)	15,19,21	1.43	3 (20%)
3	NAG	B	695	1	14,14,15	0.56	0	15,19,21	0.76	1 (6%)
6	ACT	B	801	-	1,3,3	3.09	1 (100%)	0,3,3	0.00	-

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	GOL	A	2434	-	-	0/4/4/4	0/0/0/0
3	NAG	A	692	1	-	0/6/23/26	0/1/1/1
3	NAG	A	695	1	-	0/6/23/26	0/1/1/1
3	NAG	A	697	1	-	0/6/23/26	0/1/1/1
6	ACT	A	800	-	-	0/0/0/0	0/0/0/0
6	ACT	A	801	-	-	0/0/0/0	0/0/0/0
7	GOL	B	2433	-	-	0/4/4/4	0/0/0/0
3	NAG	B	692	1	-	0/6/23/26	0/1/1/1
3	NAG	B	695	1	-	0/6/23/26	0/1/1/1
6	ACT	B	801	-	-	0/0/0/0	0/0/0/0

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	692	NAG	C8-C7	-14.01	1.22	1.50
3	B	692	NAG	C8-C7	-13.96	1.22	1.50
7	B	2433	GOL	C3-C2	-6.71	1.26	1.52
7	A	2434	GOL	C3-C2	-6.65	1.26	1.52
7	B	2433	GOL	C1-C2	2.31	1.61	1.52
7	A	2434	GOL	O2-C2	2.51	1.51	1.43
7	A	2434	GOL	C1-C2	2.55	1.62	1.52
7	B	2433	GOL	O2-C2	2.68	1.51	1.43
6	A	800	ACT	CH3-C	3.08	1.53	1.48
6	B	801	ACT	CH3-C	3.09	1.53	1.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	801	ACT	CH3-C	3.56	1.53	1.48
3	B	692	NAG	O7-C7	3.97	1.32	1.23
3	A	692	NAG	O7-C7	4.00	1.32	1.23
3	A	692	NAG	O5-C5	4.02	1.52	1.43

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	692	NAG	C1-O5-C5	-4.75	106.22	112.25
3	A	692	NAG	O7-C7-N2	-2.64	116.49	121.86
3	B	692	NAG	O7-C7-N2	-2.61	116.55	121.86
7	B	2433	GOL	O2-C2-C3	-2.38	97.75	108.65
3	A	697	NAG	C2-N2-C7	-2.35	120.02	123.04
3	B	695	NAG	C2-N2-C7	-2.34	120.03	123.04
3	A	692	NAG	C3-C2-N2	-2.33	104.97	110.56
3	A	695	NAG	C2-N2-C7	-2.32	120.05	123.04
3	B	692	NAG	C3-C2-N2	-2.31	105.03	110.56
7	B	2433	GOL	O3-C3-C2	-2.20	99.51	110.18
7	A	2434	GOL	O2-C2-C3	-2.16	98.73	108.65
7	B	2433	GOL	O2-C2-C1	-2.15	98.79	108.65
7	A	2434	GOL	O2-C2-C1	-2.00	99.46	108.65
7	A	2434	GOL	C3-C2-C1	2.08	119.26	111.12
7	B	2433	GOL	C3-C2-C1	2.24	119.90	111.12
3	B	692	NAG	C8-C7-N2	3.58	122.96	116.11
3	A	692	NAG	C8-C7-N2	3.59	122.98	116.11

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	2434	GOL	1	0
3	A	692	NAG	3	0
3	A	697	NAG	1	0
7	B	2433	GOL	1	0
3	B	692	NAG	2	0

## 5.7 Other polymers

There are no such residues in this entry.

## 5.8 Polymer linkage issues

There are no chain breaks in this entry.

## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	612/612 (100%)	-0.39	3 (0%) 91 76	12, 36, 58, 83	3 (0%)
1	B	611/612 (99%)	-0.26	9 (1%) 76 48	12, 38, 68, 84	6 (0%)
All	All	1223/1224 (99%)	-0.33	12 (0%) 84 60	12, 37, 64, 84	9 (0%)

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	131	ASN	3.2
1	B	612	ASP	3.2
1	B	134	ALA	3.1
1	B	135	THR	2.9
1	B	410	LEU	2.9
1	B	131	ASN	2.6
1	B	133	THR	2.5
1	A	136	CYS	2.4
1	B	8	GLY	2.3
1	B	83	PHE	2.2
1	A	130	PRO	2.1
1	B	15	ALA	2.1

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

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

### 6.3 Carbohydrates

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

of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	NAG	B	691	14/15	0.80	0.33	10.89	55,57,62,69	0
2	NAG	A	691	14/15	0.74	0.34	6.66	56,60,68,72	0
2	NAG	B	693	14/15	0.86	0.24	0.74	61,65,67,68	0
2	NAG	A	693	14/15	0.90	0.23	-0.07	56,58,60,60	0
2	NAG	B	696	14/15	0.63	0.41	-	86,90,91,91	0
2	NAG	A	696	14/15	0.72	0.37	-	95,96,97,97	0
2	NAG	B	694	14/15	0.78	0.34	-	79,80,81,81	0
2	NAG	A	694	14/15	0.70	0.29	-	69,72,73,75	0

## 6.4 Ligands ⓘ

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
3	NAG	A	697	14/15	0.51	0.87	12.38	116,117,118,119	0
6	ACT	A	801	4/4	0.92	0.39	11.07	32,33,34,35	0
6	ACT	B	801	4/4	0.86	0.41	9.53	52,52,53,53	0
7	GOL	A	2434	6/6	0.91	0.26	5.92	29,31,32,33	0
7	GOL	B	2433	6/6	0.93	0.25	4.59	31,34,35,39	0
3	NAG	A	695	14/15	0.86	0.32	3.89	75,76,79,79	0
3	NAG	B	695	14/15	0.84	0.36	3.81	76,78,80,83	0
3	NAG	B	692	14/15	0.70	0.35	2.29	83,86,86,87	0
3	NAG	A	692	14/15	0.72	0.34	1.16	66,68,70,71	0
5	CL	B	703	1/1	0.95	0.13	-0.64	29,29,29,29	0
5	CL	A	702	1/1	0.95	0.12	-1.86	37,37,37,37	0
4	ZN	B	701	1/1	0.99	0.17	-	29,29,29,29	0
4	ZN	A	701	1/1	1.00	0.12	-	31,31,31,31	0
6	ACT	A	800	4/4	0.76	0.46	-	49,51,52,53	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.