



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 06:47 PM UTC

PDB ID : 1IPJ / pdb_00001ipj
Title : CRYSTAL STRUCTURES OF RECOMBINANT AND NATIVE SOYBEAN BETA-CONGLYCININ BETA HOMOTRIMERS COMPLEXES WITH N-A CETYL-D-GLUCOSAMINE
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Deposited on : 2001-05-16
Resolution : 2.70 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

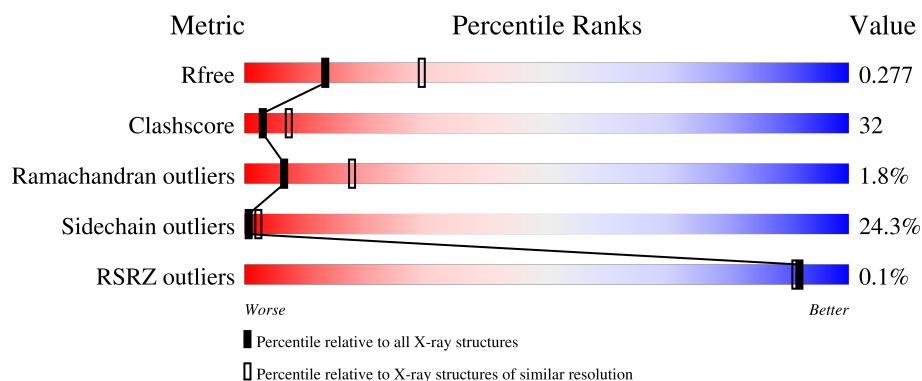
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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}	180053	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	416	 35% 40% 10% 14%
1	B	416	 29% 44% 14% 13%
1	C	416	 38% 38% 12% • 12%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8881 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.

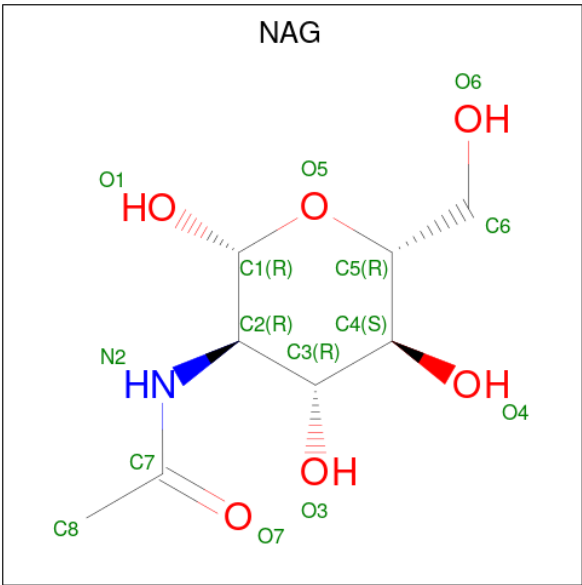
- Molecule 1 is a protein called BETA-CONGLYCININ, BETA CHAIN.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	357	Total	C	N	O	0	0	0
			2892	1832	511	549			
1	B	363	Total	C	N	O	0	0	0
			2952	1869	527	556			
1	C	368	Total	C	N	O	0	0	0
			2995	1890	537	568			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	14	LEU	PHE	conflict	UNP P25974
A	29	GLY	VAL	conflict	UNP P25974
A	175	LEU	PHE	conflict	UNP P25974
B	14	LEU	PHE	conflict	UNP P25974
B	29	GLY	VAL	conflict	UNP P25974
B	175	LEU	PHE	conflict	UNP P25974
C	14	LEU	PHE	conflict	UNP P25974
C	29	GLY	VAL	conflict	UNP P25974
C	175	LEU	PHE	conflict	UNP P25974

- Molecule 2 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).

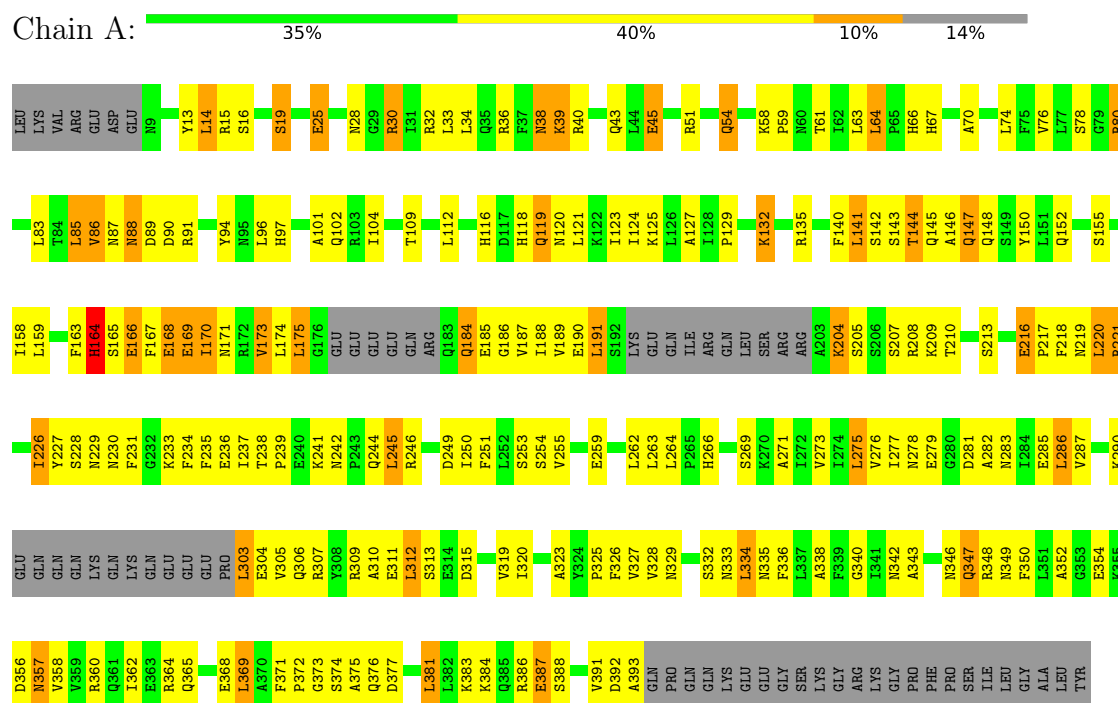


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			14	8	1	5		
2	B	1	Total	C	N	O	0	0
			14	8	1	5		
2	C	1	Total	C	N	O	0	0
			14	8	1	5		

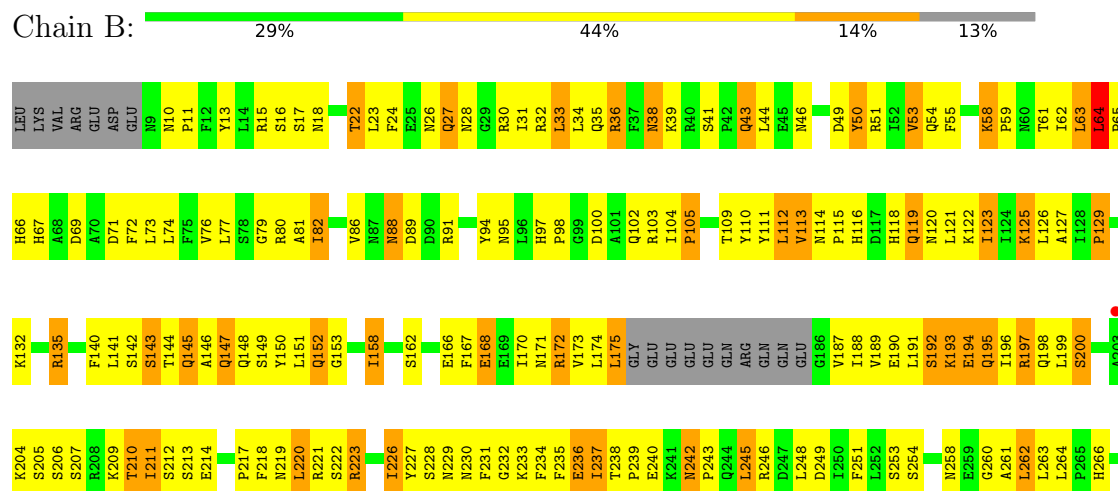
3 Residue-property plots

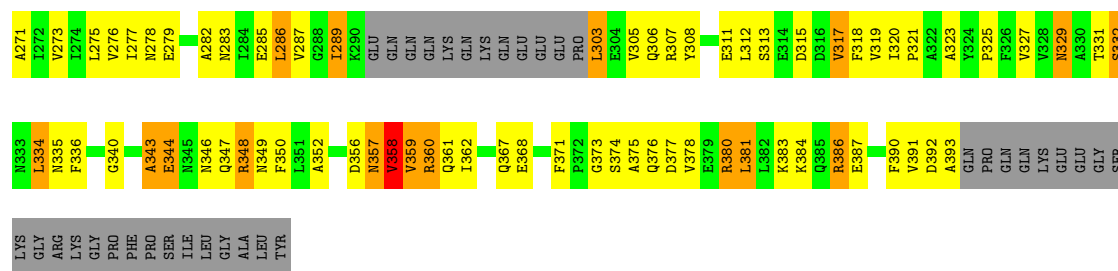
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: BETA-CONGLYCININ, BETA CHAIN



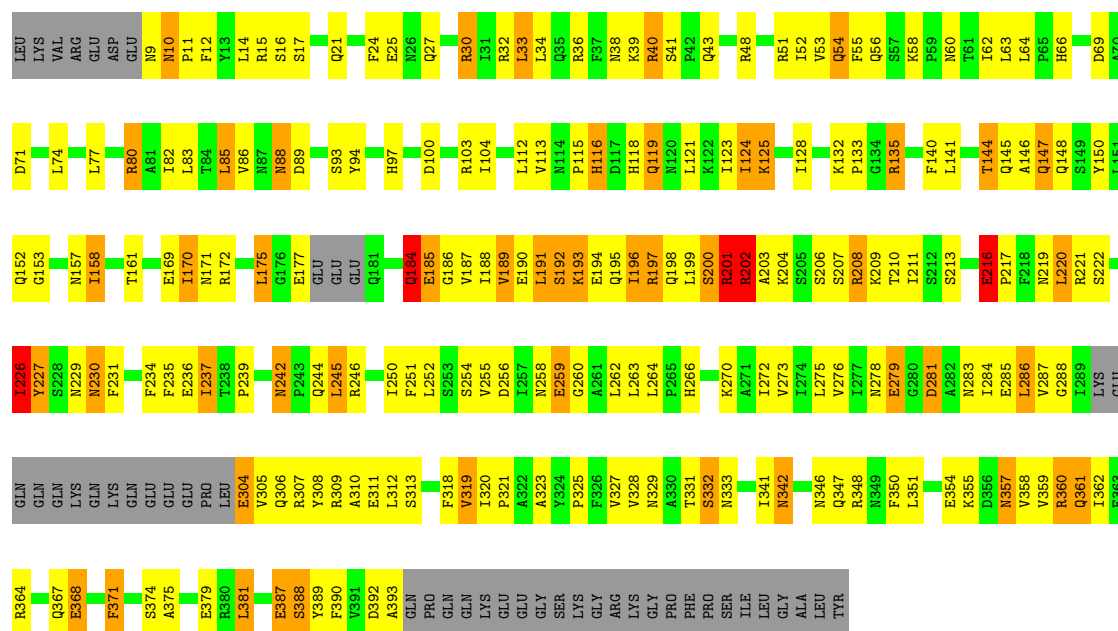
• Molecule 1: BETA-CONGLYCININ, BETA CHAIN





• Molecule 1: BETA-CONGLYCININ, BETA CHAIN

Chain C: 38% 38% 12% • 12%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	82.78Å 69.47Å 125.33Å 90.00° 97.22° 90.00°	Depositor
Resolution (Å)	7.00 – 2.70 7.00 – 2.80	Depositor EDS
% Data completeness (in resolution range)	(Not available) (7.00-2.70) 72.3 (7.00-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.50 (at 2.81Å)	Xtriage
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.202 , 0.284 0.215 , 0.277	Depositor DCC
R_{free} test set	2577 reflections (10.14%)	wwPDB-VP
Wilson B-factor (Å ²)	48.9	Xtriage
Anisotropy	0.347	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 84.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	8881	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.67% 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, 0.333 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: NAG

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	0.40	0/2949	0.91	1/3991 (0.0%)
1	B	0.38	0/3010	0.90	9/4072 (0.2%)
1	C	0.39	0/3053	0.88	4/4129 (0.1%)
All	All	0.39	0/9012	0.90	14/12192 (0.1%)

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	23	LEU	N-CA-C	-8.32	103.12	113.28
1	C	320	ILE	N-CA-C	7.64	115.12	107.55
1	C	201	ARG	N-CA-C	7.44	119.02	108.54
1	C	367	GLN	N-CA-C	-7.29	103.33	111.28
1	B	343	ALA	N-CA-C	6.76	121.54	113.16
1	B	64	LEU	CA-C-N	6.59	128.07	119.84
1	B	64	LEU	C-N-CA	6.59	128.07	119.84
1	B	273	VAL	N-CA-C	6.15	116.78	108.17
1	C	226	ILE	N-CA-C	-6.08	105.14	111.58
1	B	50	TYR	N-CA-C	5.85	119.11	110.46
1	B	358	VAL	N-CA-C	5.69	115.82	110.30
1	B	367	GLN	N-CA-C	-5.58	105.29	111.71
1	A	39	LYS	N-CA-C	-5.44	105.90	112.54
1	B	143	SER	N-CA-C	-5.16	101.31	109.76

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	2892	0	2835	187	0
1	B	2952	0	2911	222	0
1	C	2995	0	2939	195	0
2	A	14	0	13	2	0
2	B	14	0	13	2	0
2	C	14	0	13	3	0
All	All	8881	0	8724	559	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 32.

All (559) 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:95:ASN:HB3	1:B:205:SER:HB3	1.21	1.09
1:A:144:THR:HG22	1:A:146:ALA:H	1.17	1.04
1:C:25:GLU:HB2	1:C:30:ARG:HB2	1.44	1.00
1:A:25:GLU:HB2	1:A:30:ARG:HB2	1.45	0.94
1:C:364:ARG:HG2	1:C:375:ALA:HB1	1.50	0.94
1:A:242:ASN:HD21	1:A:244:GLN:HB2	1.34	0.93
1:B:220:LEU:HD22	1:B:237:ILE:HD11	1.52	0.92
1:C:329:ASN:OD1	2:C:904:NAG:H62	1.70	0.90
1:B:170:ILE:HA	1:C:381:LEU:HD11	1.54	0.89
1:A:365:GLN:HB3	1:C:195:GLN:HE21	1.38	0.88
1:C:144:THR:HG22	1:C:146:ALA:H	1.39	0.87
1:A:286:LEU:HD21	1:A:320:ILE:HG23	1.55	0.87
1:C:14:LEU:HD21	1:C:40:ARG:HD2	1.60	0.83
1:B:219:ASN:HD22	1:B:222:SER:H	1.26	0.83
1:C:220:LEU:HD21	1:C:252:LEU:HB3	1.62	0.82
1:A:305:VAL:HG21	1:C:150:TYR:HA	1.63	0.80
1:B:102:GLN:HB2	1:B:217:PRO:HB3	1.64	0.80
1:C:38:ASN:ND2	1:C:48:ARG:HA	1.97	0.80
1:A:15:ARG:O	1:A:19:SER:HB2	1.80	0.80
1:A:88:ASN:C	1:A:88:ASN:HD22	1.91	0.79
1:A:368:GLU:HB2	1:A:375:ALA:HB2	1.62	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:194:GLU:HG2	1:C:198:GLN:HE21	1.47	0.79
1:A:271:ALA:H	1:A:346:ASN:ND2	1.82	0.78
1:A:387:GLU:HB3	1:A:391:VAL:HG12	1.65	0.78
1:A:76:VAL:HG21	1:A:96:LEU:HB3	1.67	0.77
1:C:103:ARG:HB3	1:C:244:GLN:HE21	1.49	0.76
1:B:144:THR:HG22	1:B:146:ALA:H	1.50	0.76
1:B:58:LYS:HB3	1:B:59:PRO:HD2	1.66	0.76
1:B:357:ASN:HD21	1:B:359:VAL:HG13	1.49	0.76
1:B:170:ILE:HG13	1:B:171:ASN:N	2.01	0.76
1:A:283:ASN:OD1	1:A:311:GLU:HG2	1.87	0.75
1:A:38:ASN:H	1:A:38:ASN:HD22	1.32	0.75
1:B:58:LYS:HB3	1:B:59:PRO:CD	2.17	0.75
1:C:32:ARG:HH21	1:C:54:GLN:NE2	1.84	0.75
1:B:147:GLN:NE2	1:B:187:VAL:HG12	2.02	0.75
1:A:251:PHE:HD1	1:A:343:ALA:HB2	1.51	0.74
1:B:211:ILE:HG21	1:B:223:ARG:NH1	2.02	0.74
1:C:283:ASN:OD1	1:C:311:GLU:HG2	1.88	0.74
1:A:290:LYS:HG3	1:A:306:GLN:HB2	1.69	0.74
1:C:219:ASN:HB3	1:C:222:SER:HB3	1.68	0.73
1:B:144:THR:HG22	1:B:145:GLN:N	2.03	0.73
1:A:169:GLU:O	1:A:173:VAL:HG23	1.89	0.73
1:C:239:PRO:HB3	1:C:246:ARG:HA	1.70	0.73
1:C:97:HIS:O	1:C:100:ASP:HB2	1.89	0.72
1:B:153:GLY:O	1:C:307:ARG:HD3	1.88	0.72
1:B:173:VAL:HB	1:C:381:LEU:HD12	1.72	0.72
1:B:381:LEU:HD23	1:B:384:LYS:HG3	1.71	0.71
1:B:174:LEU:O	1:C:371:PHE:HB3	1.90	0.71
1:A:116:HIS:CD2	1:A:119:GLN:HB2	2.26	0.70
1:B:150:TYR:HA	1:C:305:VAL:HG21	1.72	0.70
1:A:204:LYS:HD3	1:A:205:SER:H	1.56	0.70
1:A:365:GLN:HB3	1:C:195:GLN:NE2	2.06	0.70
1:A:251:PHE:HD1	1:A:343:ALA:CB	2.05	0.70
1:A:307:ARG:HD3	1:C:153:GLY:O	1.92	0.70
1:C:38:ASN:HD22	1:C:48:ARG:HA	1.57	0.70
1:A:167:PHE:CE2	1:A:171:ASN:HB2	2.26	0.70
1:C:144:THR:CG2	1:C:146:ALA:H	2.03	0.69
1:B:144:THR:CG2	1:B:145:GLN:N	2.54	0.69
1:C:53:VAL:HB	1:C:125:LYS:HG3	1.73	0.69
1:A:251:PHE:CE1	1:A:340:GLY:HA3	2.28	0.69
1:B:38:ASN:HD22	1:B:38:ASN:H	1.41	0.69
1:B:383:LYS:HZ3	1:B:386:ARG:HH21	1.39	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:221:ARG:C	1:B:223:ARG:H	2.01	0.69
1:C:74:LEU:HD23	1:C:104:ILE:HD11	1.75	0.69
1:B:277:ILE:HD12	1:B:334:LEU:HD11	1.75	0.69
1:B:51:ARG:O	1:B:126:LEU:HD12	1.93	0.68
1:C:266:HIS:HB3	1:C:350:PHE:CD1	2.29	0.68
1:C:144:THR:HG22	1:C:146:ALA:N	2.09	0.68
1:A:278:ASN:HB3	1:A:335:ASN:ND2	2.09	0.68
1:B:147:GLN:HE21	1:B:187:VAL:HG12	1.59	0.68
1:C:116:HIS:CD2	1:C:119:GLN:HB2	2.29	0.67
1:A:220:LEU:HD13	1:A:254:SER:HB3	1.76	0.67
1:B:210:THR:HG23	1:B:210:THR:O	1.94	0.67
1:A:88:ASN:HD22	1:A:89:ASP:N	1.92	0.67
1:A:184:GLN:NE2	1:A:189:VAL:HG11	2.09	0.67
1:B:231:PHE:CE1	1:B:392:ASP:HB2	2.29	0.67
1:B:116:HIS:HD2	1:B:118:HIS:H	1.44	0.66
1:B:22:THR:HG23	1:B:32:ARG:HG2	1.76	0.66
1:B:285:GLU:HB2	1:B:327:VAL:HG12	1.76	0.66
1:C:231:PHE:CZ	1:C:392:ASP:HB2	2.31	0.66
1:A:303:LEU:HB2	1:C:147:GLN:HB2	1.78	0.65
1:B:167:PHE:CD2	1:B:170:ILE:HD11	2.31	0.65
1:B:50:TYR:O	1:B:51:ARG:HG2	1.96	0.65
1:B:357:ASN:ND2	1:B:359:VAL:HG13	2.10	0.65
1:A:85:LEU:HD12	1:A:94:TYR:HE2	1.61	0.65
1:A:287:VAL:CG2	1:C:153:GLY:HA3	2.25	0.65
1:B:116:HIS:CD2	1:B:119:GLN:H	2.15	0.65
1:C:275:LEU:HD11	1:C:318:PHE:HB3	1.79	0.65
1:A:83:LEU:HA	1:A:112:LEU:HD12	1.80	0.64
1:B:220:LEU:CD2	1:B:237:ILE:HD11	2.27	0.64
1:B:242:ASN:C	1:B:242:ASN:HD22	2.05	0.64
1:B:357:ASN:O	1:B:360:ARG:HB2	1.97	0.64
1:B:53:VAL:HB	1:B:125:LYS:HG3	1.78	0.64
1:C:192:SER:HB2	1:C:195:GLN:HG3	1.78	0.64
1:C:112:LEU:HD21	1:C:123:ILE:HD12	1.80	0.64
1:C:144:THR:HB	1:C:147:GLN:OE1	1.96	0.64
1:B:350:PHE:O	1:B:357:ASN:HA	1.97	0.64
1:C:196:ILE:O	1:C:200:SER:HB2	1.97	0.64
1:A:327:VAL:HG11	1:C:158:ILE:HG12	1.80	0.63
1:C:36:ARG:NH1	1:C:39:LYS:HE2	2.12	0.63
1:A:278:ASN:HB3	1:A:335:ASN:HD22	1.64	0.63
1:C:10:ASN:HD22	1:C:12:PHE:H	1.45	0.63
1:B:144:THR:CG2	1:B:145:GLN:H	2.11	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:82:ILE:HD11	1:C:200:SER:OG	1.99	0.63
1:A:226:ILE:HG13	1:A:234:PHE:O	1.98	0.63
1:B:62:ILE:HA	1:B:112:LEU:O	1.99	0.63
1:C:14:LEU:CD2	1:C:40:ARG:HD2	2.28	0.63
1:C:341:ILE:HG22	1:C:342:ASN:ND2	2.14	0.63
1:A:242:ASN:ND2	1:A:244:GLN:HB2	2.09	0.63
1:B:231:PHE:CZ	1:B:392:ASP:HB2	2.34	0.63
1:C:32:ARG:HH21	1:C:54:GLN:HE22	1.46	0.63
1:A:155:SER:HG	1:A:158:ILE:HD12	1.63	0.62
1:A:242:ASN:HD22	1:A:245:LEU:H	1.46	0.62
1:C:60:ASN:OD1	1:C:196:ILE:HD12	1.99	0.62
2:C:904:NAG:O7	2:C:904:NAG:H3	1.98	0.62
1:C:85:LEU:HD12	1:C:94:TYR:HE2	1.65	0.62
1:A:263:LEU:HB3	1:A:328:VAL:HB	1.80	0.62
1:B:36:ARG:NH1	1:B:135:ARG:HD3	2.14	0.62
1:B:194:GLU:HG2	1:B:197:ARG:NH2	2.14	0.62
1:C:342:ASN:N	1:C:342:ASN:HD22	1.97	0.62
1:A:85:LEU:CD2	1:A:104:ILE:HG23	2.29	0.62
1:A:287:VAL:HG21	1:C:153:GLY:HA3	1.81	0.62
1:B:88:ASN:HD22	1:B:89:ASP:N	1.97	0.62
1:C:219:ASN:HD22	1:C:222:SER:H	1.47	0.61
1:C:285:GLU:HB2	1:C:327:VAL:HG12	1.81	0.61
1:B:264:LEU:HD13	1:B:352:ALA:O	1.98	0.61
1:B:368:GLU:HB2	1:B:375:ALA:HB2	1.81	0.61
1:C:77:LEU:HD13	1:C:278:ASN:ND2	2.15	0.61
1:A:80:ARG:O	1:A:121:LEU:HD13	2.00	0.61
1:A:266:HIS:HB2	1:A:349:ASN:O	2.01	0.61
1:B:31:ILE:HG12	1:B:55:PHE:HD1	1.65	0.61
1:A:207:SER:OG	1:A:208:ARG:N	2.34	0.61
1:A:286:LEU:CD2	1:A:320:ILE:HG23	2.31	0.61
1:B:170:ILE:HA	1:C:381:LEU:CD1	2.30	0.61
1:B:73:LEU:HD22	1:B:248:LEU:HD12	1.83	0.60
1:B:13:TYR:OH	1:B:15:ARG:HG3	2.01	0.60
1:A:213:SER:OG	1:A:216:GLU:HG3	2.01	0.60
1:C:36:ARG:HB2	1:C:39:LYS:HG3	1.81	0.60
1:A:144:THR:HG22	1:A:146:ALA:N	2.03	0.60
1:B:193:LYS:HE2	1:B:194:GLU:N	2.16	0.60
1:B:28:ASN:OD1	1:B:61:THR:HG21	2.01	0.60
1:A:184:GLN:HE21	1:A:189:VAL:HG11	1.66	0.60
1:C:25:GLU:CB	1:C:30:ARG:HB2	2.25	0.60
1:C:147:GLN:NE2	1:C:187:VAL:HG12	2.17	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:28:ASN:HB3	1:A:61:THR:HG21	1.84	0.59
1:A:220:LEU:HD13	1:A:254:SER:CB	2.33	0.59
1:A:285:GLU:HB2	1:A:327:VAL:HG12	1.85	0.59
1:A:383:LYS:NZ	1:A:386:ARG:HH22	2.01	0.59
1:B:263:LEU:O	1:B:327:VAL:HG23	2.01	0.59
1:B:26:ASN:C	1:B:28:ASN:H	2.11	0.59
1:B:66:HIS:HB3	1:B:140:PHE:CD2	2.36	0.59
1:C:364:ARG:CG	1:C:375:ALA:HB1	2.30	0.59
1:C:10:ASN:HD22	1:C:10:ASN:C	2.11	0.59
1:B:148:GLN:OE1	1:C:304:GLU:HA	2.03	0.59
1:A:34:LEU:O	1:A:51:ARG:NH2	2.36	0.59
1:C:194:GLU:HG2	1:C:198:GLN:NE2	2.17	0.58
1:A:266:HIS:O	1:A:325:PRO:HA	2.02	0.58
1:B:36:ARG:HB2	1:B:39:LYS:HG3	1.86	0.58
1:C:220:LEU:CD2	1:C:252:LEU:HB3	2.33	0.58
1:B:219:ASN:HD22	1:B:222:SER:N	1.99	0.58
1:B:196:ILE:O	1:B:200:SER:OG	2.21	0.58
1:B:239:PRO:O	1:B:246:ARG:HD2	2.04	0.58
1:B:278:ASN:HB3	1:B:335:ASN:HD22	1.68	0.58
1:B:275:LEU:HD23	1:B:320:ILE:HD11	1.86	0.57
1:B:94:TYR:CD1	1:B:217:PRO:HD3	2.38	0.57
1:B:173:VAL:HG12	1:B:173:VAL:O	2.03	0.57
1:A:147:GLN:NE2	1:A:187:VAL:HG12	2.19	0.57
1:C:387:GLU:HG3	1:C:388:SER:H	1.69	0.57
1:B:195:GLN:O	1:B:199:LEU:HG	2.04	0.57
1:B:229:ASN:C	1:B:231:PHE:H	2.13	0.57
1:C:237:ILE:HG22	1:C:245:LEU:HD23	1.86	0.57
1:A:165:SER:HB3	1:A:169:GLU:OE2	2.05	0.57
1:C:361:GLN:HA	1:C:361:GLN:HE21	1.70	0.57
1:A:144:THR:CG2	1:A:145:GLN:N	2.68	0.56
1:A:58:LYS:HB3	1:A:59:PRO:HD2	1.87	0.56
1:B:34:LEU:HG	1:B:35:GLN:H	1.70	0.56
1:C:284:ILE:O	1:C:309:ARG:HA	2.05	0.56
1:B:95:ASN:HB3	1:B:205:SER:CB	2.14	0.56
1:C:305:VAL:HG12	1:C:306:GLN:N	2.21	0.56
1:B:278:ASN:HB3	1:B:335:ASN:ND2	2.20	0.56
1:C:201:ARG:O	1:C:202:ARG:HB2	2.04	0.56
1:B:251:PHE:CE1	1:B:340:GLY:HA3	2.41	0.56
1:A:185:GLU:HG3	1:A:185:GLU:O	2.04	0.55
1:C:10:ASN:HD21	1:C:12:PHE:HD2	1.54	0.55
1:B:167:PHE:CE2	1:B:170:ILE:HD11	2.41	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:211:ILE:HG21	1:B:223:ARG:HH12	1.70	0.55
1:A:38:ASN:H	1:A:38:ASN:ND2	2.02	0.55
1:A:303:LEU:CB	1:C:147:GLN:HB2	2.35	0.55
1:C:193:LYS:O	1:C:197:ARG:HB2	2.07	0.55
1:C:286:LEU:HB3	1:C:308:TYR:HB2	1.89	0.55
1:A:190:GLU:C	1:A:191:LEU:HD23	2.31	0.55
1:A:381:LEU:HD11	1:C:170:ILE:HG22	1.89	0.55
1:B:245:LEU:O	1:B:249:ASP:N	2.39	0.55
1:B:383:LYS:NZ	1:B:386:ARG:NH2	2.55	0.55
1:C:113:VAL:HG12	1:C:115:PRO:HD3	1.89	0.55
1:A:148:GLN:NE2	1:A:152:GLN:HE22	2.04	0.55
1:C:34:LEU:O	1:C:51:ARG:NH2	2.39	0.55
1:A:333:ASN:H	1:A:333:ASN:HD22	1.55	0.55
1:B:24:PHE:CD2	1:B:187:VAL:HG22	2.42	0.54
1:A:25:GLU:HB2	1:A:30:ARG:CB	2.28	0.54
1:B:383:LYS:HZ3	1:B:386:ARG:NH2	2.06	0.54
1:C:33:LEU:HD21	1:C:51:ARG:NH2	2.22	0.54
1:C:219:ASN:HD22	1:C:222:SER:N	2.05	0.54
1:A:290:LYS:HB2	1:A:304:GLU:O	2.06	0.54
1:C:231:PHE:CE1	1:C:392:ASP:HB2	2.43	0.54
1:B:387:GLU:HB2	1:B:391:VAL:HG12	1.90	0.54
1:B:210:THR:O	1:B:210:THR:CG2	2.56	0.54
1:C:357:ASN:C	1:C:357:ASN:HD22	2.15	0.54
1:C:191:LEU:HD23	1:C:196:ILE:HG12	1.90	0.54
1:C:263:LEU:HD13	1:C:390:PHE:CE1	2.43	0.54
1:C:112:LEU:HD21	1:C:123:ILE:CD1	2.37	0.54
1:A:88:ASN:HB3	1:B:349:ASN:HD22	1.73	0.54
1:C:145:GLN:HB2	1:C:185:GLU:O	2.08	0.54
1:C:359:VAL:HG12	1:C:359:VAL:O	2.08	0.54
1:A:144:THR:HG23	1:A:184:GLN:O	2.08	0.54
1:A:220:LEU:HD22	1:A:235:PHE:HB3	1.89	0.54
1:A:357:ASN:HD22	1:A:360:ARG:H	1.56	0.54
1:B:236:GLU:HG3	1:B:236:GLU:O	2.08	0.54
1:B:266:HIS:HB3	1:B:350:PHE:CD1	2.43	0.54
1:A:102:GLN:HB2	1:A:217:PRO:HB3	1.90	0.53
1:B:86:VAL:HG23	1:B:109:THR:HB	1.90	0.53
1:B:192:SER:HB2	1:B:195:GLN:HG3	1.90	0.53
1:B:289:ILE:HG21	1:B:303:LEU:HD13	1.90	0.53
1:C:55:PHE:CE2	1:C:112:LEU:HD23	2.43	0.53
1:A:173:VAL:CG1	1:B:377:ASP:HB3	2.37	0.53
1:B:167:PHE:O	1:B:170:ILE:HG12	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:263:LEU:HD23	1:A:328:VAL:HG21	1.90	0.53
1:A:369:LEU:HD23	1:C:184:GLN:OE1	2.07	0.53
1:B:275:LEU:HD23	1:B:320:ILE:CD1	2.38	0.53
1:B:271:ALA:H	1:B:346:ASN:ND2	2.07	0.53
1:C:144:THR:CG2	1:C:145:GLN:N	2.71	0.53
1:C:226:ILE:HD11	1:C:236:GLU:HB3	1.91	0.53
1:A:33:LEU:HD21	1:A:51:ARG:NH1	2.23	0.53
1:B:111:TYR:CE1	1:C:362:ILE:HG23	2.44	0.53
1:A:372:PRO:HG3	1:C:175:LEU:O	2.08	0.53
1:B:32:ARG:HH21	1:B:54:GLN:CD	2.17	0.53
2:C:904:NAG:O7	2:C:904:NAG:C3	2.56	0.53
1:A:287:VAL:HG12	1:A:325:PRO:O	2.09	0.53
1:C:10:ASN:ND2	1:C:12:PHE:H	2.05	0.53
1:C:116:HIS:HD2	1:C:119:GLN:H	1.56	0.53
1:C:287:VAL:CG2	1:C:325:PRO:HD2	2.38	0.53
1:B:103:ARG:O	1:B:105:PRO:HD3	2.09	0.53
1:B:266:HIS:O	1:B:325:PRO:HA	2.09	0.53
1:B:227:TYR:HB2	1:B:234:PHE:CB	2.39	0.52
1:A:346:ASN:O	1:A:347:GLN:HG2	2.09	0.52
1:B:235:PHE:HB2	1:B:254:SER:O	2.10	0.52
1:C:33:LEU:CD2	1:C:51:ARG:NH2	2.72	0.52
1:A:58:LYS:HA	1:A:120:ASN:OD1	2.10	0.52
1:A:279:GLU:HG2	1:A:335:ASN:HB3	1.91	0.52
1:C:270:LYS:H	1:C:346:ASN:HD22	1.57	0.52
1:A:170:ILE:HD13	1:A:171:ASN:N	2.24	0.52
1:A:374:SER:HB3	1:A:377:ASP:CG	2.35	0.52
1:B:334:LEU:HG	1:B:335:ASN:N	2.25	0.52
1:A:86:VAL:HG22	1:A:109:THR:HB	1.90	0.52
1:A:173:VAL:HG11	1:B:377:ASP:HB3	1.91	0.52
1:B:88:ASN:HD22	1:B:89:ASP:H	1.57	0.52
1:B:194:GLU:HG2	1:B:197:ARG:HH21	1.72	0.52
1:C:33:LEU:CD2	1:C:51:ARG:HH22	2.23	0.52
1:B:212:SER:HA	1:B:243:PRO:HD2	1.91	0.52
1:B:262:LEU:HB2	1:B:329:ASN:HA	1.92	0.52
1:B:162:SER:HB3	1:C:264:LEU:HD22	1.92	0.51
1:A:144:THR:CG2	1:A:186:GLY:H	2.22	0.51
1:A:271:ALA:N	1:A:346:ASN:ND2	2.56	0.51
1:A:287:VAL:CG1	1:A:325:PRO:HD2	2.41	0.51
1:C:237:ILE:CG2	1:C:245:LEU:HD23	2.40	0.51
1:A:210:THR:HG23	1:A:218:PHE:HA	1.91	0.51
1:A:271:ALA:H	1:A:346:ASN:HD22	1.57	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:360:ARG:HH21	1:A:383:LYS:HD3	1.74	0.51
1:A:155:SER:OG	1:A:158:ILE:HD12	2.10	0.51
1:A:204:LYS:HD3	1:A:205:SER:N	2.24	0.51
1:B:229:ASN:O	1:B:231:PHE:N	2.44	0.51
1:C:30:ARG:NH2	1:C:32:ARG:CZ	2.73	0.51
1:C:242:ASN:HD22	1:C:244:GLN:H	1.59	0.51
1:B:97:HIS:O	1:B:100:ASP:HB2	2.11	0.51
1:B:251:PHE:C	1:B:251:PHE:CD2	2.88	0.51
1:C:236:GLU:OE1	1:C:348:ARG:NH2	2.39	0.51
1:A:309:ARG:CG	1:A:310:ALA:N	2.74	0.51
1:A:227:TYR:OH	1:A:348:ARG:HD2	2.10	0.51
1:A:235:PHE:O	1:A:253:SER:HA	2.11	0.51
1:C:94:TYR:HA	1:C:204:LYS:O	2.12	0.50
1:B:383:LYS:NZ	1:B:386:ARG:HH21	2.09	0.50
1:B:228:SER:HB2	1:B:233:LYS:HG2	1.92	0.50
1:A:383:LYS:HZ1	1:A:386:ARG:HH22	1.59	0.50
1:C:281:ASP:HB2	1:C:331:THR:OG1	2.11	0.50
1:C:24:PHE:HZ	1:C:186:GLY:HA3	1.77	0.50
1:C:189:VAL:CG2	1:C:190:GLU:N	2.75	0.50
1:C:251:PHE:C	1:C:251:PHE:CD2	2.89	0.50
1:C:103:ARG:H	1:C:244:GLN:HE22	1.60	0.50
1:A:332:SER:O	1:A:333:ASN:C	2.54	0.50
1:B:113:VAL:HG12	1:B:115:PRO:HD3	1.94	0.50
1:A:219:ASN:ND2	1:A:221:ARG:H	2.10	0.49
1:A:237:ILE:HG22	1:A:245:LEU:HD23	1.94	0.49
1:B:63:LEU:HD23	1:B:187:VAL:O	2.11	0.49
1:B:312:LEU:HD21	1:B:318:PHE:CD2	2.47	0.49
1:C:226:ILE:HG13	1:C:234:PHE:O	2.12	0.49
1:C:275:LEU:CD1	1:C:318:PHE:HB3	2.42	0.49
1:A:70:ALA:HB2	1:A:129:PRO:HA	1.92	0.49
1:A:238:THR:HB	1:A:239:PRO:HD2	1.93	0.49
1:C:144:THR:HG23	1:C:145:GLN:N	2.28	0.49
1:C:236:GLU:CD	1:C:348:ARG:HH12	2.19	0.49
1:A:80:ARG:HD3	1:A:97:HIS:CE1	2.48	0.49
1:A:87:ASN:O	1:B:361:GLN:HG3	2.12	0.49
1:A:245:LEU:O	1:A:249:ASP:N	2.44	0.49
1:B:82:ILE:O	1:B:82:ILE:HG22	2.11	0.49
1:B:151:LEU:CD2	1:C:351:LEU:HD13	2.42	0.49
1:B:221:ARG:C	1:B:223:ARG:N	2.69	0.49
1:B:260:GLY:O	1:B:393:ALA:HB2	2.12	0.49
1:C:10:ASN:HD22	1:C:12:PHE:N	2.09	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:74:LEU:HD23	1:A:104:ILE:HD11	1.93	0.49
1:A:309:ARG:HG3	1:A:310:ALA:N	2.28	0.49
1:C:89:ASP:OD1	1:C:89:ASP:C	2.56	0.49
1:C:227:TYR:OH	1:C:348:ARG:HD2	2.12	0.49
1:A:271:ALA:N	1:A:346:ASN:HD22	2.10	0.49
1:C:242:ASN:HB3	1:C:245:LEU:HB2	1.94	0.49
1:A:245:LEU:HG	1:A:250:ILE:O	2.13	0.49
1:B:175:LEU:N	1:B:175:LEU:CD1	2.75	0.49
1:B:276:VAL:HG22	1:B:317:VAL:HG22	1.94	0.49
1:A:66:HIS:HB3	1:A:140:PHE:CD2	2.47	0.48
1:A:132:LYS:CD	1:A:135:ARG:HD3	2.43	0.48
1:A:287:VAL:HG23	1:C:153:GLY:HA3	1.93	0.48
1:B:102:GLN:OE1	1:B:217:PRO:HB3	2.13	0.48
1:C:24:PHE:CE2	1:C:188:ILE:HG13	2.48	0.48
1:C:88:ASN:HD22	1:C:89:ASP:N	2.11	0.48
1:C:305:VAL:HG12	1:C:306:GLN:H	1.77	0.48
1:B:77:LEU:HD12	1:B:122:LYS:HE3	1.95	0.48
1:C:74:LEU:HD12	1:C:74:LEU:O	2.13	0.48
1:C:342:ASN:ND2	1:C:342:ASN:N	2.59	0.48
1:A:251:PHE:CD1	1:A:343:ALA:HB2	2.40	0.48
1:B:343:ALA:O	1:B:344:GLU:C	2.56	0.48
1:B:381:LEU:C	1:B:383:LYS:H	2.21	0.48
1:C:216:GLU:HB3	1:C:217:PRO:HD2	1.96	0.48
1:C:288:GLY:O	1:C:306:GLN:N	2.44	0.48
1:C:332:SER:O	1:C:333:ASN:C	2.56	0.48
1:A:264:LEU:HD22	1:A:352:ALA:HB3	1.96	0.48
1:B:16:SER:HB2	1:B:315:ASP:O	2.13	0.48
1:B:149:SER:O	1:B:152:GLN:HB2	2.14	0.48
1:B:283:ASN:OD1	1:B:311:GLU:HG2	2.14	0.48
1:C:32:ARG:NH2	1:C:54:GLN:HE22	2.10	0.48
1:C:62:ILE:HA	1:C:112:LEU:O	2.14	0.48
1:C:85:LEU:CD2	1:C:104:ILE:HG23	2.43	0.48
1:A:242:ASN:ND2	1:A:244:GLN:H	2.12	0.48
1:C:15:ARG:HD3	1:C:17:SER:OG	2.13	0.48
1:C:62:ILE:O	1:C:62:ILE:HG23	2.14	0.48
1:C:10:ASN:OD1	1:C:308:TYR:HB3	2.13	0.47
1:C:279:GLU:O	1:C:279:GLU:HG2	2.14	0.47
1:B:28:ASN:OD1	1:B:61:THR:CG2	2.63	0.47
1:B:51:ARG:O	1:B:126:LEU:HA	2.14	0.47
1:A:88:ASN:C	1:A:88:ASN:ND2	2.63	0.47
1:B:95:ASN:CB	1:B:205:SER:HB3	2.16	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:220:LEU:HD22	1:A:235:PHE:CB	2.45	0.47
1:B:260:GLY:HA2	2:B:903:NAG:H83	1.95	0.47
1:C:33:LEU:HD23	1:C:51:ARG:HH22	1.78	0.47
1:A:16:SER:HB3	1:A:315:ASP:O	2.15	0.47
1:A:231:PHE:CE1	1:A:392:ASP:HB2	2.49	0.47
1:B:226:ILE:HD11	1:B:234:PHE:CD1	2.49	0.47
1:C:36:ARG:HH21	1:C:135:ARG:HG3	1.79	0.47
1:A:132:LYS:NZ	1:A:135:ARG:HD3	2.29	0.47
1:A:144:THR:HG22	1:A:145:GLN:N	2.29	0.47
1:A:239:PRO:O	1:A:246:ARG:HB2	2.14	0.47
1:C:309:ARG:HG2	1:C:310:ALA:N	2.28	0.47
1:C:358:VAL:C	1:C:360:ARG:H	2.22	0.47
1:A:253:SER:OG	1:A:338:ALA:HB3	2.15	0.47
1:B:229:ASN:C	1:B:231:PHE:N	2.72	0.47
1:B:49:ASP:O	1:B:129:PRO:HD2	2.15	0.47
1:B:76:VAL:HG22	1:B:123:ILE:HD11	1.97	0.47
1:B:227:TYR:HB2	1:B:234:PHE:HB3	1.97	0.47
1:B:58:LYS:O	1:B:114:ASN:ND2	2.48	0.46
1:B:144:THR:HG23	1:B:145:GLN:H	1.78	0.46
1:A:277:ILE:HD11	1:A:312:LEU:CD1	2.46	0.46
1:A:286:LEU:HD23	1:A:326:PHE:HB3	1.97	0.46
1:C:66:HIS:HB3	1:C:140:PHE:CD2	2.50	0.46
1:C:88:ASN:ND2	1:C:88:ASN:H	2.13	0.46
1:C:103:ARG:H	1:C:244:GLN:NE2	2.14	0.46
1:A:32:ARG:HH21	1:A:54:GLN:NE2	2.13	0.46
1:B:251:PHE:HD1	1:B:343:ALA:CB	2.28	0.46
1:C:194:GLU:O	1:C:198:GLN:HG3	2.15	0.46
1:C:210:THR:HG21	1:C:217:PRO:O	2.16	0.46
1:A:32:ARG:HH21	1:A:54:GLN:CD	2.23	0.46
1:A:163:PHE:O	1:A:164:HIS:CB	2.64	0.46
1:A:231:PHE:CZ	1:A:392:ASP:HB2	2.50	0.46
1:B:79:GLY:O	1:B:98:PRO:HG3	2.15	0.46
1:B:50:TYR:C	1:B:51:ARG:HG2	2.40	0.46
1:B:89:ASP:C	1:C:361:GLN:OE1	2.58	0.46
1:C:171:ASN:OD1	1:C:177:GLU:HG3	2.16	0.46
1:C:124:ILE:HD12	1:C:276:VAL:HG11	1.98	0.46
1:C:103:ARG:CB	1:C:244:GLN:HE21	2.24	0.46
1:A:13:TYR:CE1	1:A:15:ARG:HB2	2.50	0.46
1:A:116:HIS:HD2	1:A:119:GLN:H	1.63	0.46
1:A:287:VAL:HG13	1:A:325:PRO:HD2	1.98	0.46
1:A:334:LEU:HG	1:A:335:ASN:N	2.30	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:41:SER:OG	1:B:43:GLN:HB2	2.16	0.46
1:B:218:PHE:CD1	1:B:218:PHE:N	2.84	0.46
1:B:381:LEU:C	1:B:383:LYS:N	2.74	0.46
1:C:74:LEU:HD12	1:C:74:LEU:C	2.41	0.46
1:C:286:LEU:HD11	1:C:321:PRO:HD2	1.97	0.46
1:B:38:ASN:H	1:B:38:ASN:ND2	2.11	0.45
1:B:227:TYR:HB2	1:B:234:PHE:HB2	1.98	0.45
1:C:34:LEU:HB2	1:C:52:ILE:HB	1.98	0.45
1:B:86:VAL:CG2	1:B:109:THR:HB	2.47	0.45
1:B:346:ASN:C	1:B:347:GLN:HG2	2.42	0.45
1:B:350:PHE:HB2	1:B:356:ASP:O	2.17	0.45
1:C:252:LEU:HD23	1:C:252:LEU:HA	1.82	0.45
1:A:28:ASN:HB2	1:A:188:ILE:HG22	1.99	0.45
1:B:80:ARG:HD2	1:B:97:HIS:CE1	2.52	0.45
1:A:70:ALA:CB	1:A:129:PRO:HA	2.46	0.45
1:B:229:ASN:ND2	1:B:390:PHE:O	2.48	0.45
1:A:88:ASN:HB3	1:B:349:ASN:ND2	2.31	0.45
1:A:281:ASP:O	1:A:282:ALA:HB2	2.17	0.45
1:A:271:ALA:H	1:A:346:ASN:HD21	1.58	0.45
1:A:323:ALA:HB3	1:C:69:ASP:HB3	1.99	0.45
1:B:228:SER:HA	1:B:232:GLY:O	2.17	0.45
1:A:158:ILE:HD12	1:B:285:GLU:OE1	2.17	0.45
1:B:226:ILE:HD11	1:B:234:PHE:HD1	1.80	0.45
1:B:226:ILE:CD1	1:B:236:GLU:HG2	2.47	0.45
1:C:207:SER:OG	1:C:208:ARG:N	2.49	0.45
1:A:393:ALA:HB1	2:A:902:NAG:H83	1.98	0.45
1:B:282:ALA:HB2	1:B:334:LEU:HD22	1.99	0.45
1:A:371:PHE:C	1:A:373:GLY:H	2.24	0.44
1:C:24:PHE:CZ	1:C:186:GLY:HA3	2.52	0.44
1:C:80:ARG:HD2	1:C:97:HIS:CE1	2.53	0.44
1:A:309:ARG:HG3	1:A:310:ALA:H	1.80	0.44
2:A:902:NAG:C8	1:C:161:THR:OG1	2.65	0.44
1:B:69:ASP:HB3	1:C:323:ALA:HB3	1.97	0.44
1:B:88:ASN:H	1:B:88:ASN:ND2	2.14	0.44
1:A:269:SER:OG	1:A:347:GLN:N	2.46	0.44
1:C:38:ASN:ND2	1:C:48:ARG:CA	2.77	0.44
1:A:175:LEU:N	1:A:175:LEU:CD1	2.80	0.44
1:A:251:PHE:C	1:A:251:PHE:CD2	2.95	0.44
1:A:383:LYS:NZ	1:A:386:ARG:NH2	2.65	0.44
1:B:72:PHE:HE2	1:B:127:ALA:HB2	1.82	0.44
1:B:286:LEU:HB3	1:B:308:TYR:HB2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:33:LEU:HD21	1:C:51:ARG:CZ	2.48	0.44
1:C:229:ASN:C	1:C:231:PHE:H	2.25	0.44
1:C:235:PHE:HB2	1:C:254:SER:O	2.17	0.44
1:A:226:ILE:HD11	1:A:236:GLU:HB3	1.98	0.44
1:B:86:VAL:HG23	1:B:86:VAL:O	2.18	0.44
1:B:238:THR:C	1:B:240:GLU:N	2.75	0.44
1:B:74:LEU:HB3	1:B:125:LYS:HB3	1.99	0.44
1:A:36:ARG:HD2	1:A:39:LYS:HD3	2.00	0.44
1:A:51:ARG:HB2	1:A:127:ALA:HB3	2.00	0.44
1:B:88:ASN:ND2	1:B:88:ASN:N	2.65	0.44
1:C:116:HIS:CD2	1:C:119:GLN:CB	3.00	0.44
1:C:145:GLN:HG3	1:C:184:GLN:O	2.17	0.44
1:B:238:THR:C	1:B:240:GLU:H	2.25	0.44
1:A:45:GLU:OE2	1:C:133:PRO:HG3	2.18	0.44
1:A:158:ILE:HG13	2:B:903:NAG:H61	1.99	0.44
1:B:26:ASN:C	1:B:28:ASN:N	2.76	0.44
1:B:175:LEU:N	1:B:175:LEU:HD12	2.33	0.44
1:C:287:VAL:HA	1:C:306:GLN:O	2.18	0.44
1:C:262:LEU:HB2	1:C:328:VAL:O	2.18	0.43
1:A:64:LEU:HB3	1:A:142:SER:HB3	1.99	0.43
1:C:260:GLY:O	1:C:393:ALA:HB3	2.19	0.43
1:B:66:HIS:NE2	1:B:110:TYR:CE1	2.86	0.43
1:C:145:GLN:HB2	1:C:185:GLU:HA	2.01	0.43
1:A:173:VAL:CG1	1:B:373:GLY:HA3	2.48	0.43
1:C:226:ILE:HD11	1:C:236:GLU:CB	2.47	0.43
1:B:235:PHE:O	1:B:253:SER:HA	2.18	0.43
1:A:13:TYR:CE2	1:A:312:LEU:HD23	2.53	0.43
1:B:76:VAL:HA	1:B:123:ILE:HG12	2.00	0.43
1:C:275:LEU:HD12	1:C:275:LEU:O	2.19	0.43
1:C:259:GLU:HB2	1:C:332:SER:HA	2.01	0.43
1:C:357:ASN:C	1:C:357:ASN:ND2	2.77	0.43
1:A:350:PHE:O	1:A:357:ASN:HA	2.19	0.43
1:B:226:ILE:HD11	1:B:236:GLU:HG2	2.01	0.43
1:B:278:ASN:C	1:B:278:ASN:HD22	2.26	0.43
1:A:357:ASN:ND2	1:A:360:ARG:H	2.17	0.43
1:B:381:LEU:HD23	1:B:384:LYS:CG	2.44	0.43
1:C:83:LEU:O	1:C:93:SER:HA	2.19	0.43
1:C:93:SER:O	1:C:203:ALA:HA	2.19	0.43
1:B:10:ASN:HA	1:B:11:PRO:HD2	1.83	0.43
1:A:14:LEU:CD1	1:A:40:ARG:HD3	2.48	0.42
1:A:14:LEU:HD13	1:A:40:ARG:HD3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:67:HIS:CD2	1:B:67:HIS:C	2.96	0.42
1:B:219:ASN:ND2	1:B:222:SER:N	2.66	0.42
1:B:278:ASN:C	1:B:278:ASN:ND2	2.77	0.42
1:C:147:GLN:HG2	1:C:148:GLN:O	2.19	0.42
1:B:94:TYR:HA	1:B:204:LYS:O	2.19	0.42
1:C:12:PHE:HB3	1:C:319:VAL:O	2.19	0.42
1:C:250:ILE:HG22	1:C:251:PHE:N	2.33	0.42
1:C:258:ASN:O	1:C:259:GLU:C	2.63	0.42
1:B:81:ALA:HB2	1:B:121:LEU:HD11	2.00	0.42
1:B:170:ILE:CA	1:C:381:LEU:HD11	2.37	0.42
1:A:356:ASP:HA	1:C:88:ASN:HB2	2.01	0.42
1:B:170:ILE:HG22	1:C:381:LEU:HD21	2.01	0.42
1:A:150:TYR:HA	1:B:305:VAL:HG21	2.01	0.42
1:A:221:ARG:NH1	1:A:235:PHE:CE1	2.87	0.42
1:B:74:LEU:HD21	1:B:104:ILE:HD11	2.00	0.42
1:B:191:LEU:HD12	1:B:191:LEU:HA	1.91	0.42
1:C:286:LEU:CB	1:C:308:TYR:HB2	2.48	0.42
1:A:67:HIS:NE2	1:B:323:ALA:HB1	2.34	0.42
1:A:67:HIS:CD2	1:B:323:ALA:HB1	2.55	0.42
1:A:90:ASP:CG	1:A:91:ARG:H	2.26	0.42
1:A:239:PRO:CB	1:A:246:ARG:HA	2.49	0.42
1:B:168:GLU:O	1:B:172:ARG:HB2	2.20	0.42
1:B:74:LEU:HD12	1:B:74:LEU:O	2.20	0.42
1:B:62:ILE:HG13	1:B:112:LEU:O	2.19	0.42
1:C:287:VAL:HG22	1:C:325:PRO:HD2	2.01	0.42
1:A:286:LEU:HD21	1:A:320:ILE:CG2	2.39	0.42
1:A:372:PRO:HG3	1:C:175:LEU:C	2.44	0.42
1:B:170:ILE:CG1	1:B:171:ASN:N	2.76	0.42
1:B:358:VAL:C	1:B:360:ARG:N	2.78	0.42
1:A:173:VAL:HG12	1:B:373:GLY:HA3	2.02	0.41
1:A:218:PHE:CD1	1:A:218:PHE:N	2.87	0.41
1:B:33:LEU:HD21	1:B:51:ARG:NH2	2.34	0.41
1:B:151:LEU:HD21	1:C:351:LEU:HD13	2.02	0.41
1:B:220:LEU:HD13	1:B:237:ILE:CG1	2.50	0.41
1:B:236:GLU:OE2	1:B:348:ARG:NH1	2.53	0.41
1:C:229:ASN:O	1:C:231:PHE:N	2.53	0.41
1:A:251:PHE:CD1	1:A:343:ALA:CB	2.95	0.41
1:B:374:SER:O	1:B:378:VAL:HG23	2.21	0.41
1:B:392:ASP:OD2	1:B:393:ALA:N	2.53	0.41
1:C:24:PHE:CD2	1:C:187:VAL:HG22	2.55	0.41
1:A:132:LYS:HZ2	1:A:135:ARG:HD3	1.84	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:147:GLN:HE21	1:A:187:VAL:HG12	1.86	0.41
1:A:276:VAL:O	1:A:336:PHE:HB2	2.21	0.41
1:B:239:PRO:HB3	1:B:246:ARG:HA	2.02	0.41
1:B:374:SER:O	1:B:375:ALA:C	2.63	0.41
1:C:10:ASN:ND2	1:C:12:PHE:HD2	2.17	0.41
1:A:166:GLU:H	1:A:166:GLU:HG3	1.63	0.41
1:B:28:ASN:HB3	1:B:188:ILE:O	2.20	0.41
1:B:158:ILE:HD13	1:C:285:GLU:CD	2.45	0.41
1:B:167:PHE:CE1	1:B:171:ASN:HB2	2.56	0.41
1:B:170:ILE:HG13	1:B:171:ASN:H	1.80	0.41
1:B:192:SER:CB	1:B:195:GLN:HG3	2.50	0.41
1:A:229:ASN:C	1:A:231:PHE:H	2.27	0.41
1:B:242:ASN:O	1:B:243:PRO:C	2.64	0.41
1:B:357:ASN:CG	1:B:360:ARG:HG3	2.45	0.41
1:C:36:ARG:HH11	1:C:39:LYS:HE2	1.84	0.41
1:C:41:SER:C	1:C:43:GLN:N	2.79	0.41
1:A:229:ASN:O	1:A:231:PHE:N	2.53	0.41
1:A:275:LEU:HD23	1:A:320:ILE:HD11	2.03	0.41
1:B:26:ASN:O	1:B:28:ASN:N	2.54	0.41
1:B:81:ALA:HB2	1:B:121:LEU:CD1	2.50	0.41
1:B:251:PHE:CD1	1:B:340:GLY:HA3	2.56	0.41
1:B:278:ASN:HB3	1:B:335:ASN:O	2.20	0.41
1:C:34:LEU:CB	1:C:52:ILE:HB	2.51	0.41
1:A:88:ASN:O	1:B:356:ASP:HA	2.21	0.41
1:A:141:LEU:HD11	1:B:359:VAL:HB	2.02	0.41
1:B:147:GLN:HE21	1:B:187:VAL:CG1	2.28	0.41
1:B:220:LEU:HD13	1:B:237:ILE:HD11	2.03	0.41
1:B:251:PHE:HD1	1:B:343:ALA:HB1	1.85	0.41
1:B:331:THR:O	1:B:332:SER:CB	2.69	0.41
1:B:380:ARG:HE	1:B:380:ARG:HB3	1.64	0.41
1:C:11:PRO:HA	1:C:40:ARG:HH21	1.85	0.41
1:C:379:GLU:O	1:C:379:GLU:HG2	2.18	0.41
1:A:132:LYS:H	1:A:132:LYS:HG3	1.60	0.41
1:A:251:PHE:CD1	1:A:340:GLY:HA3	2.56	0.41
1:A:271:ALA:CB	1:A:346:ASN:ND2	2.84	0.41
1:B:65:PRO:HG3	1:B:111:TYR:CD1	2.56	0.41
1:C:229:ASN:C	1:C:229:ASN:OD1	2.63	0.41
1:C:368:GLU:OE1	1:C:375:ALA:N	2.50	0.41
1:B:64:LEU:HD23	1:B:142:SER:OG	2.21	0.40
1:A:174:LEU:HD22	1:B:371:PHE:CE2	2.56	0.40
1:A:74:LEU:O	1:A:101:ALA:HA	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:163:PHE:O	1:A:164:HIS:HB2	2.20	0.40
1:A:190:GLU:O	1:A:191:LEU:HD23	2.20	0.40
1:B:276:VAL:O	1:B:336:PHE:HB2	2.22	0.40
1:C:56:GLN:HA	1:C:121:LEU:O	2.21	0.40
1:C:124:ILE:HD11	1:C:276:VAL:HG21	2.03	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	349/416 (84%)	316 (90%)	30 (9%)	3 (1%)	14	35
1	B	357/416 (86%)	301 (84%)	45 (13%)	11 (3%)	3	8
1	C	362/416 (87%)	309 (85%)	48 (13%)	5 (1%)	9	23
All	All	1068/1248 (86%)	926 (87%)	123 (12%)	19 (2%)	6	18

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	332	SER
1	B	344	GLU
1	C	202	ARG
1	A	164	HIS
1	A	230	ASN
1	B	27	GLN
1	B	46	ASN
1	B	206	SER
1	B	230	ASN
1	B	321	PRO
1	C	230	ASN
1	A	168	GLU

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Mol	Chain	Res	Type
1	B	129	PRO
1	C	184	GLN
1	B	58	LYS
1	B	261	ALA
1	C	389	TYR
1	B	105	PRO
1	C	216	GLU

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	321/375 (86%)	251 (78%)	70 (22%)	1	3
1	B	328/375 (88%)	247 (75%)	81 (25%)	1	2
1	C	332/375 (88%)	245 (74%)	87 (26%)	0	2
All	All	981/1125 (87%)	743 (76%)	238 (24%)	1	2

All (238) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	LEU
1	A	19	SER
1	A	25	GLU
1	A	30	ARG
1	A	38	ASN
1	A	43	GLN
1	A	45	GLU
1	A	54	GLN
1	A	63	LEU
1	A	64	LEU
1	A	78	SER
1	A	80	ARG
1	A	85	LEU
1	A	86	VAL
1	A	88	ASN

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Mol	Chain	Res	Type
1	A	118	HIS
1	A	119	GLN
1	A	123	ILE
1	A	124	ILE
1	A	125	LYS
1	A	132	LYS
1	A	141	LEU
1	A	143	SER
1	A	144	THR
1	A	147	GLN
1	A	159	LEU
1	A	164	HIS
1	A	166	GLU
1	A	168	GLU
1	A	169	GLU
1	A	170	ILE
1	A	173	VAL
1	A	175	LEU
1	A	184	GLN
1	A	191	LEU
1	A	204	LYS
1	A	209	LYS
1	A	216	GLU
1	A	220	LEU
1	A	221	ARG
1	A	226	ILE
1	A	228	SER
1	A	233	LYS
1	A	241	LYS
1	A	245	LEU
1	A	255	VAL
1	A	259	GLU
1	A	262	LEU
1	A	273	VAL
1	A	275	LEU
1	A	286	LEU
1	A	303	LEU
1	A	312	LEU
1	A	313	SER
1	A	319	VAL
1	A	329	ASN
1	A	334	LEU

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Mol	Chain	Res	Type
1	A	342	ASN
1	A	347	GLN
1	A	354	GLU
1	A	357	ASN
1	A	358	VAL
1	A	362	ILE
1	A	364	ARG
1	A	369	LEU
1	A	376	GLN
1	A	381	LEU
1	A	384	LYS
1	A	387	GLU
1	A	388	SER
1	B	17	SER
1	B	18	ASN
1	B	22	THR
1	B	27	GLN
1	B	30	ARG
1	B	33	LEU
1	B	36	ARG
1	B	38	ASN
1	B	43	GLN
1	B	44	LEU
1	B	53	VAL
1	B	63	LEU
1	B	64	LEU
1	B	71	ASP
1	B	82	ILE
1	B	88	ASN
1	B	91	ARG
1	B	112	LEU
1	B	113	VAL
1	B	119	GLN
1	B	120	ASN
1	B	123	ILE
1	B	125	LYS
1	B	132	LYS
1	B	135	ARG
1	B	141	LEU
1	B	143	SER
1	B	145	GLN
1	B	147	GLN

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Mol	Chain	Res	Type
1	B	152	GLN
1	B	158	ILE
1	B	166	GLU
1	B	168	GLU
1	B	172	ARG
1	B	175	LEU
1	B	189	VAL
1	B	190	GLU
1	B	192	SER
1	B	193	LYS
1	B	194	GLU
1	B	195	GLN
1	B	197	ARG
1	B	198	GLN
1	B	200	SER
1	B	207	SER
1	B	209	LYS
1	B	210	THR
1	B	211	ILE
1	B	213	SER
1	B	214	GLU
1	B	220	LEU
1	B	223	ARG
1	B	226	ILE
1	B	236	GLU
1	B	237	ILE
1	B	242	ASN
1	B	245	LEU
1	B	258	ASN
1	B	262	LEU
1	B	279	GLU
1	B	286	LEU
1	B	287	VAL
1	B	289	ILE
1	B	303	LEU
1	B	306	GLN
1	B	307	ARG
1	B	313	SER
1	B	317	VAL
1	B	319	VAL
1	B	329	ASN
1	B	334	LEU

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Mol	Chain	Res	Type
1	B	348	ARG
1	B	357	ASN
1	B	358	VAL
1	B	359	VAL
1	B	360	ARG
1	B	362	ILE
1	B	376	GLN
1	B	380	ARG
1	B	381	LEU
1	B	386	ARG
1	C	9	ASN
1	C	10	ASN
1	C	16	SER
1	C	21	GLN
1	C	27	GLN
1	C	30	ARG
1	C	33	LEU
1	C	40	ARG
1	C	54	GLN
1	C	58	LYS
1	C	63	LEU
1	C	64	LEU
1	C	71	ASP
1	C	80	ARG
1	C	85	LEU
1	C	86	VAL
1	C	88	ASN
1	C	116	HIS
1	C	118	HIS
1	C	119	GLN
1	C	124	ILE
1	C	125	LYS
1	C	128	ILE
1	C	132	LYS
1	C	135	ARG
1	C	141	LEU
1	C	144	THR
1	C	147	GLN
1	C	152	GLN
1	C	157	ASN
1	C	158	ILE
1	C	169	GLU

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Mol	Chain	Res	Type
1	C	170	ILE
1	C	172	ARG
1	C	175	LEU
1	C	184	GLN
1	C	185	GLU
1	C	189	VAL
1	C	191	LEU
1	C	192	SER
1	C	193	LYS
1	C	196	ILE
1	C	197	ARG
1	C	199	LEU
1	C	200	SER
1	C	201	ARG
1	C	202	ARG
1	C	206	SER
1	C	208	ARG
1	C	209	LYS
1	C	211	ILE
1	C	213	SER
1	C	216	GLU
1	C	220	LEU
1	C	221	ARG
1	C	226	ILE
1	C	227	TYR
1	C	230	ASN
1	C	237	ILE
1	C	242	ASN
1	C	245	LEU
1	C	255	VAL
1	C	256	ASP
1	C	259	GLU
1	C	272	ILE
1	C	273	VAL
1	C	279	GLU
1	C	281	ASP
1	C	286	LEU
1	C	304	GLU
1	C	312	LEU
1	C	313	SER
1	C	319	VAL
1	C	332	SER

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Mol	Chain	Res	Type
1	C	342	ASN
1	C	347	GLN
1	C	354	GLU
1	C	355	LYS
1	C	357	ASN
1	C	360	ARG
1	C	361	GLN
1	C	368	GLU
1	C	371	PHE
1	C	374	SER
1	C	381	LEU
1	C	387	GLU
1	C	388	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (79) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	18	ASN
1	A	38	ASN
1	A	43	GLN
1	A	88	ASN
1	A	97	HIS
1	A	116	HIS
1	A	119	GLN
1	A	147	GLN
1	A	152	GLN
1	A	157	ASN
1	A	171	ASN
1	A	183	GLN
1	A	184	GLN
1	A	219	ASN
1	A	242	ASN
1	A	244	GLN
1	A	258	ASN
1	A	278	ASN
1	A	306	GLN
1	A	333	ASN
1	A	335	ASN
1	A	342	ASN
1	A	345	ASN
1	A	346	ASN
1	A	347	GLN

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Mol	Chain	Res	Type
1	A	349	ASN
1	A	357	ASN
1	A	361	GLN
1	B	9	ASN
1	B	38	ASN
1	B	43	GLN
1	B	88	ASN
1	B	95	ASN
1	B	97	HIS
1	B	116	HIS
1	B	119	GLN
1	B	145	GLN
1	B	147	GLN
1	B	156	HIS
1	B	157	ASN
1	B	195	GLN
1	B	219	ASN
1	B	242	ASN
1	B	258	ASN
1	B	266	HIS
1	B	278	ASN
1	B	306	GLN
1	B	335	ASN
1	B	345	ASN
1	B	346	ASN
1	B	349	ASN
1	B	357	ASN
1	B	361	GLN
1	C	10	ASN
1	C	21	GLN
1	C	38	ASN
1	C	54	GLN
1	C	88	ASN
1	C	116	HIS
1	C	119	GLN
1	C	120	ASN
1	C	145	GLN
1	C	157	ASN
1	C	195	GLN
1	C	198	GLN
1	C	219	ASN
1	C	242	ASN

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Mol	Chain	Res	Type
1	C	244	GLN
1	C	258	ASN
1	C	266	HIS
1	C	278	ASN
1	C	335	ASN
1	C	342	ASN
1	C	346	ASN
1	C	347	GLN
1	C	349	ASN
1	C	357	ASN
1	C	361	GLN
1	C	365	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

3 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	NAG	C	904	1	14,14,15	0.86	1 (7%)	17,19,21	1.47	3 (17%)
2	NAG	B	903	1	14,14,15	0.53	0	17,19,21	0.57	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	A	902	1	14,14,15	0.59	0	17,19,21	0.81	1 (5%)

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	C	904	1	-	3/6/23/26	0/1/1/1
2	NAG	B	903	1	-	4/6/23/26	0/1/1/1
2	NAG	A	902	1	-	2/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	904	NAG	O5-C5	2.09	1.47	1.43

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	904	NAG	C1-O5-C5	4.38	118.06	112.19
2	C	904	NAG	C4-C3-C2	-2.89	106.78	111.02
2	A	902	NAG	C2-N2-C7	-2.27	119.85	122.90
2	C	904	NAG	O5-C1-C2	2.08	114.51	111.29

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	903	NAG	C1-C2-N2-C7
2	C	904	NAG	C3-C2-N2-C7
2	B	903	NAG	O5-C5-C6-O6
2	B	903	NAG	C4-C5-C6-O6
2	C	904	NAG	O5-C5-C6-O6
2	A	902	NAG	C4-C5-C6-O6
2	A	902	NAG	O5-C5-C6-O6
2	C	904	NAG	C1-C2-N2-C7
2	B	903	NAG	C3-C2-N2-C7

There are no ring outliers.

3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	904	NAG	3	0
2	B	903	NAG	2	0
2	A	902	NAG	2	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

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

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	357/416 (85%)	-0.53	0 100 100	3, 20, 36, 50	0
1	B	363/416 (87%)	-0.29	1 (0%) 90 89	6, 30, 48, 64	0
1	C	368/416 (88%)	-0.45	0 100 100	3, 23, 42, 53	0
All	All	1088/1248 (87%)	-0.42	1 (0%) 92 91	3, 24, 44, 64	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	203	ALA	2.1

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

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. 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	B-factors(Å ²)	Q<0.9
2	NAG	C	904	14/15	0.70	0.17	45,51,55,55	0
2	NAG	A	902	14/15	0.73	0.16	41,45,48,49	0
2	NAG	B	903	14/15	0.76	0.18	56,60,64,66	0

6.5 Other polymers [i](#)

There are no such residues in this entry.