



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 06:14 PM EDT

PDB ID : 6VCA / pdb_00006vca
Title : TB38 complex
Authors : Zhou, Y.F.; Lord, D.M.
Deposited on : 2019-12-20
Resolution : 3.73 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:

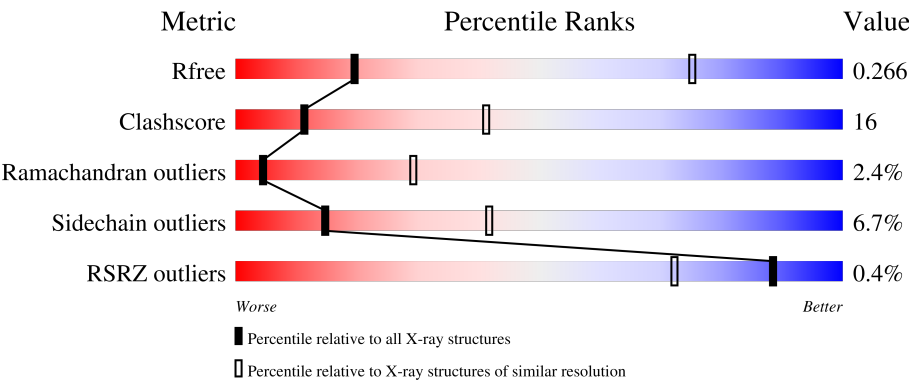
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry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

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.73 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.



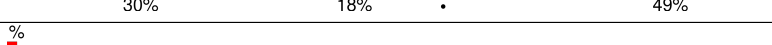
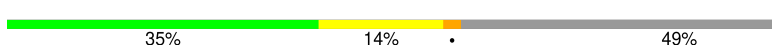
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R _{free}	180053	1234 (3.88-3.60)
Clashscore	190562	1274 (3.88-3.60)
Ramachandran outliers	187476	1226 (3.88-3.60)
Sidechain outliers	187428	1221 (3.88-3.60)
RSRZ outliers	180081	1233 (3.88-3.60)

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	529	71%24%...
1	B	529	70%26%..
1	C	529	69%25%5%.
1	D	529	% 69%27%..
2	E	233	58%30%5%6%

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Mol	Chain	Length	Quality of chain
2	H	233	 61% 29% 6%
2	K	233	 58% 31% 6%
2	P	233	 30% 18% 49%
3	F	216	 67% 27% 6%
3	J	216	 68% 26% 6%
3	L	216	 72% 21% 6%
3	O	216	 35% 14% 49%
4	G	2	 100%
5	I	3	 100%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 27834 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 5'-nucleotidase, ecto (CD73), isoform CRA_a.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	523	Total	C	N	O	S	0	1	0
			4087	2600	700	768	19			
1	B	523	Total	C	N	O	S	0	0	0
			4081	2595	699	768	19			
1	C	523	Total	C	N	O	S	0	0	0
			4071	2589	695	768	19			
1	D	523	Total	C	N	O	S	0	3	0
			4095	2605	701	770	19			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	550	HIS	-	expression tag	UNP Q53Z63
A	551	HIS	-	expression tag	UNP Q53Z63
A	552	HIS	-	expression tag	UNP Q53Z63
A	553	HIS	-	expression tag	UNP Q53Z63
A	554	HIS	-	expression tag	UNP Q53Z63
A	555	HIS	-	expression tag	UNP Q53Z63
B	550	HIS	-	expression tag	UNP Q53Z63
B	551	HIS	-	expression tag	UNP Q53Z63
B	552	HIS	-	expression tag	UNP Q53Z63
B	553	HIS	-	expression tag	UNP Q53Z63
B	554	HIS	-	expression tag	UNP Q53Z63
B	555	HIS	-	expression tag	UNP Q53Z63
C	550	HIS	-	expression tag	UNP Q53Z63
C	551	HIS	-	expression tag	UNP Q53Z63
C	552	HIS	-	expression tag	UNP Q53Z63
C	553	HIS	-	expression tag	UNP Q53Z63
C	554	HIS	-	expression tag	UNP Q53Z63
C	555	HIS	-	expression tag	UNP Q53Z63
D	550	HIS	-	expression tag	UNP Q53Z63
D	551	HIS	-	expression tag	UNP Q53Z63
D	552	HIS	-	expression tag	UNP Q53Z63

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Chain	Residue	Modelled	Actual	Comment	Reference
D	553	HIS	-	expression tag	UNP Q53Z63
D	554	HIS	-	expression tag	UNP Q53Z63
D	555	HIS	-	expression tag	UNP Q53Z63

- Molecule 2 is a protein called TB38 heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	220	Total	C	N	O	S	0	0	0
			1654	1047	278	322	7			
2	E	219	Total	C	N	O	S	0	1	0
			1649	1043	277	322	7			
2	P	119	Total	C	N	O	S	0	0	0
			928	587	159	177	5			
2	K	220	Total	C	N	O	S	0	0	0
			1654	1047	278	322	7			

- Molecule 3 is a protein called TB38 light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	212	Total	C	N	O	S	0	0	0
			1573	990	257	321	5			
3	F	212	Total	C	N	O	S	0	1	0
			1575	991	257	322	5			
3	O	110	Total	C	N	O	S	0	0	0
			813	512	132	166	3			
3	J	212	Total	C	N	O	S	0	0	0
			1573	990	257	321	5			

- Molecule 4 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



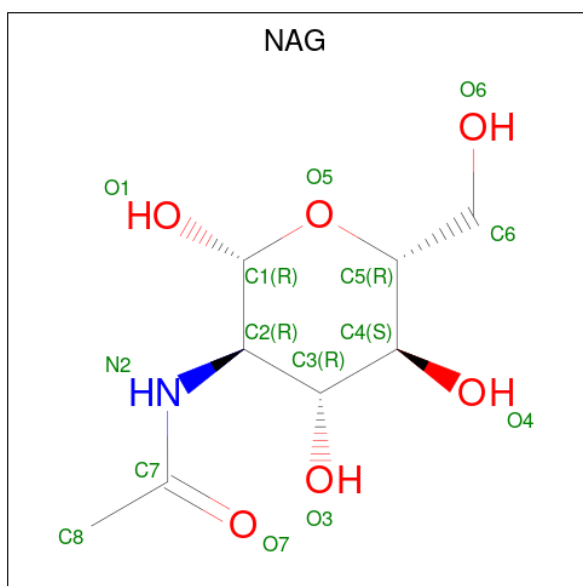
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	G	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 5 is an oligosaccharide called beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	I	3	Total	C	N	O	0	0	0
			39	22	2	15			

- Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).

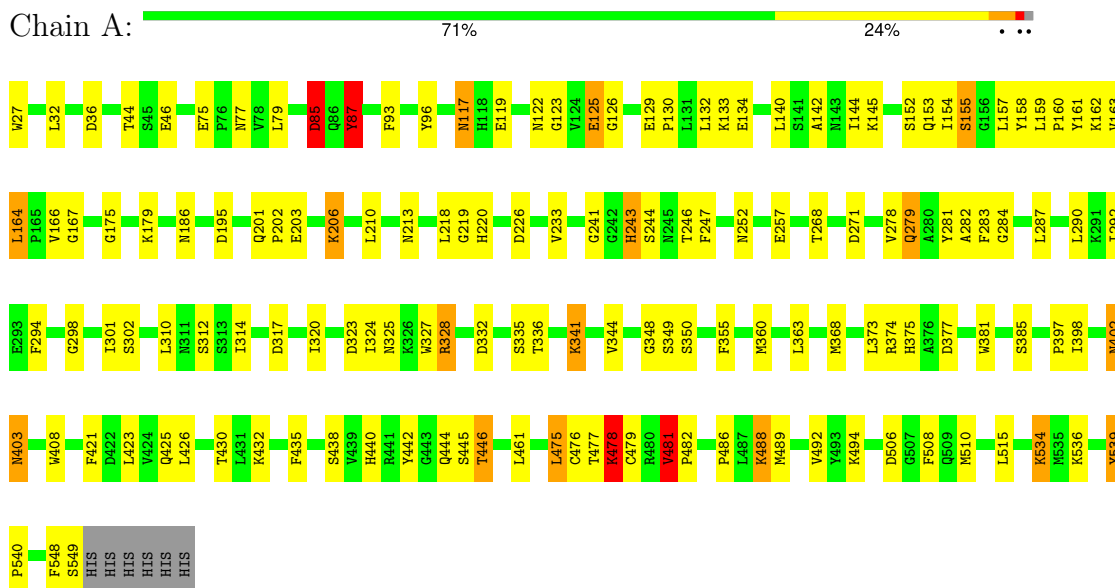


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	D	1	Total	C	N	O	0	0
			14	8	1	5		

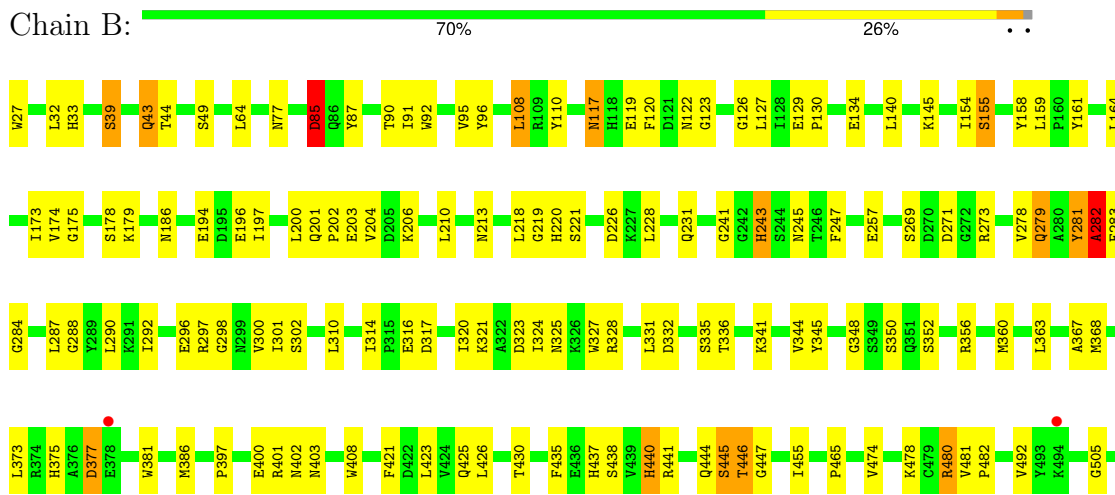
3 Residue-property plots

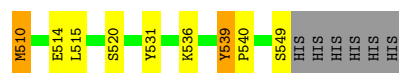
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: 5'-nucleotidase, ecto (CD73), isoform CRA_a



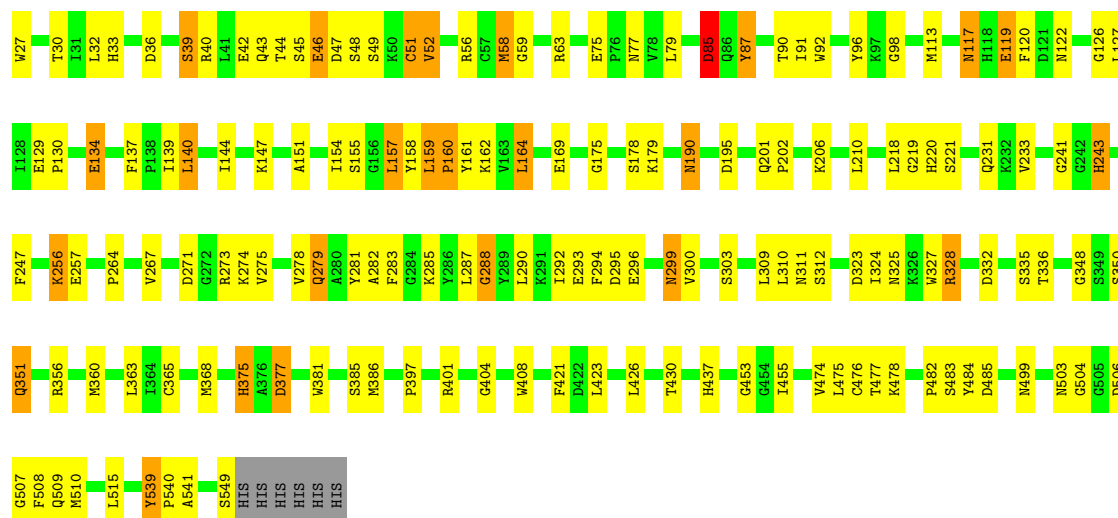
- Molecule 1: 5'-nucleotidase, ecto (CD73), isoform CRA_a





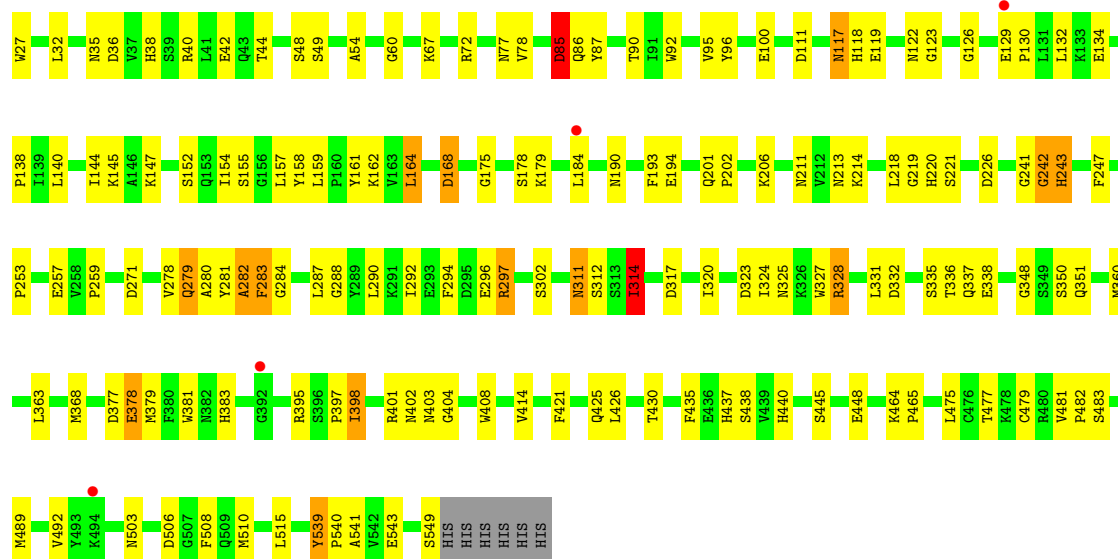
- Molecule 1: 5'-nucleotidase, ecto (CD73), isoform CRA_a

Chain C: 69% 25% 5% •



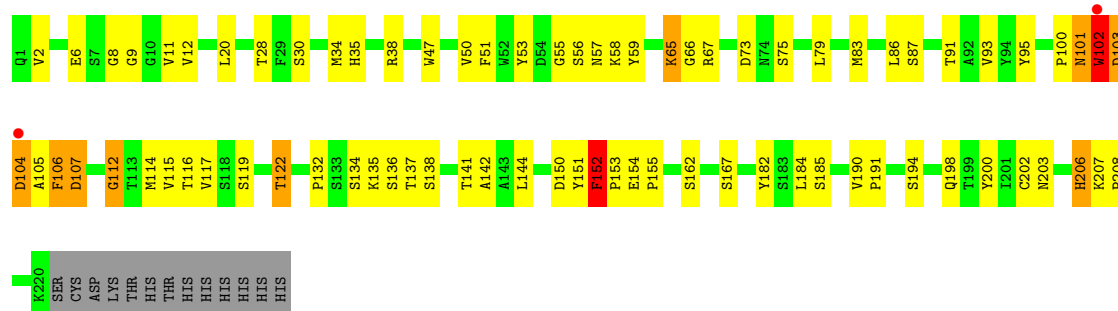
- Molecule 1: 5'-nucleotidase, ecto (CD73), isoform CRA_a

Chain D: 69% 27% • •



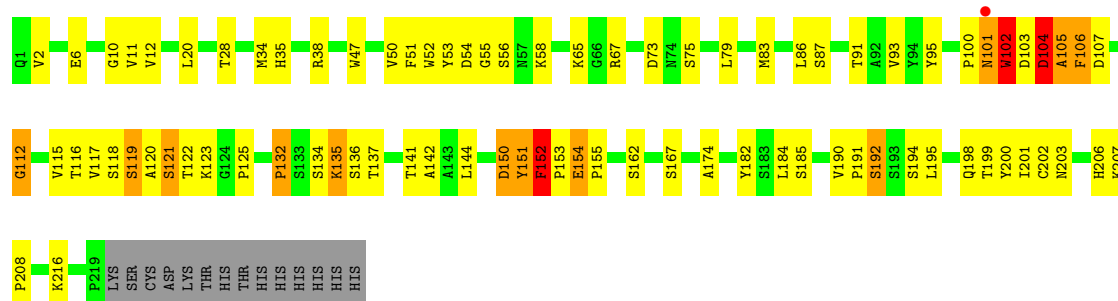
- Molecule 2: TB38 heavy chain

Chain H: 61% 29% • • 6%



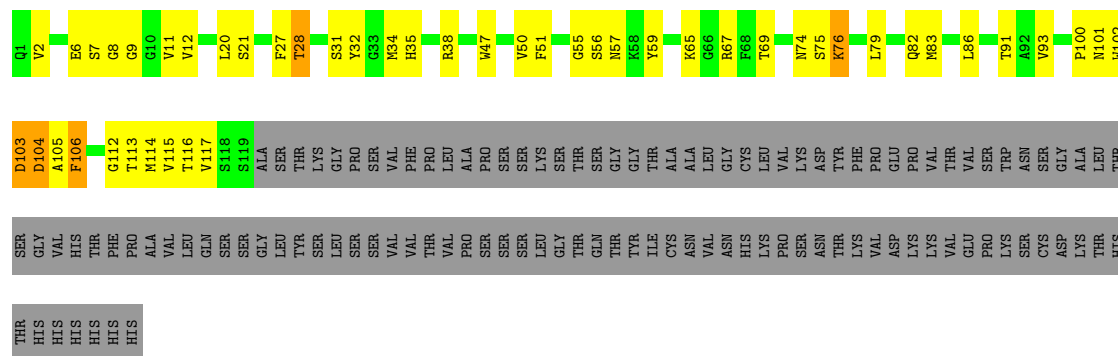
• Molecule 2: TB38 heavy chain

Chain E: 58% 30% 5% • 6%



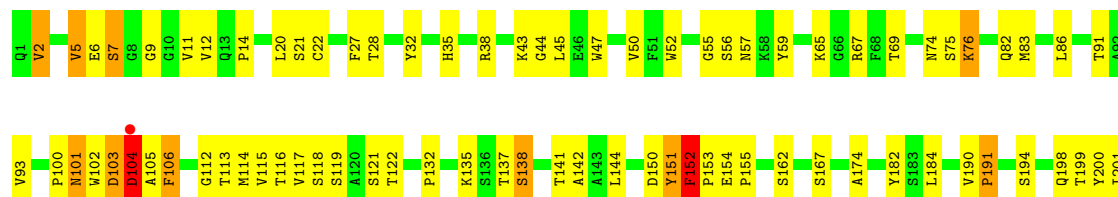
• Molecule 2: TB38 heavy chain

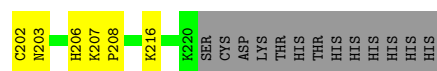
Chain P: 30% 18% • 49%



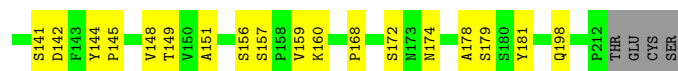
• Molecule 2: TB38 heavy chain

Chain K: 58% 31% • • 6%

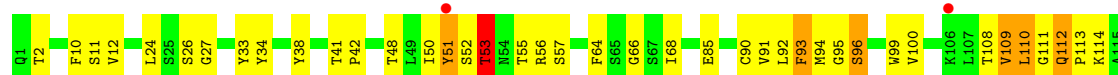




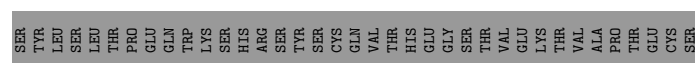
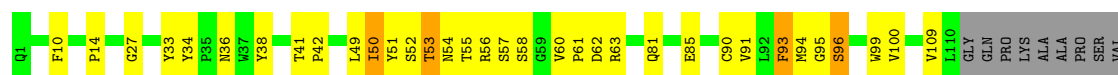
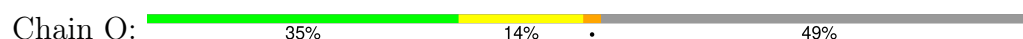
- Molecule 3: TB38 light chain



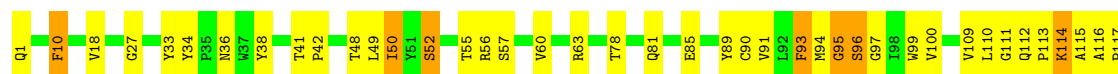
- Molecule 3: TB38 light chain



- Molecule 3: TB38 light chain



- Molecule 3: TB38 light chain



- Molecule 4: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain G:  100%

MAG1
MAG2

- Molecule 5: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain I:  100%

MAG1
MAG2
BMA3

4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	236.91Å 336.20Å 222.15Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 3.73 40.00 – 3.73	Depositor EDS
% Data completeness (in resolution range)	99.7 (40.00-3.73) 99.7 (40.00-3.73)	Depositor EDS
R_{merge}	0.26	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.12 (at 3.76Å)	Xtriage
Refinement program	REFMAC 5.8.0257	Depositor
R, R_{free}	0.224 , 0.264 0.229 , 0.266	Depositor DCC
R_{free} test set	4586 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	163.2	Xtriage
Anisotropy	0.035	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 182.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	27834	wwPDB-VP
Average B, all atoms (Å ²)	197.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.08% 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, BMA

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	1.08	1/4176 (0.0%)	1.52	15/5654 (0.3%)
1	B	1.09	0/4167	1.49	11/5643 (0.2%)
1	C	1.08	0/4157	1.48	15/5632 (0.3%)
1	D	1.07	1/4190 (0.0%)	1.51	15/5674 (0.3%)
2	E	1.16	6/1697 (0.4%)	1.48	14/2314 (0.6%)
2	H	1.17	0/1698	1.46	7/2315 (0.3%)
2	K	1.11	0/1698	1.47	9/2315 (0.4%)
2	P	1.12	0/953	1.41	4/1294 (0.3%)
3	F	1.16	0/1619	1.51	2/2217 (0.1%)
3	J	1.12	1/1613 (0.1%)	1.45	1/2209 (0.0%)
3	L	1.14	0/1613	1.51	7/2209 (0.3%)
3	O	1.16	1/832 (0.1%)	1.47	3/1137 (0.3%)
All	All	1.11	10/28413 (0.0%)	1.49	103/38613 (0.3%)

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	B	0	1
1	C	0	1
2	E	0	1
2	H	0	2
2	K	0	1
3	F	0	1
3	L	0	1
All	All	0	8

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	104	ASP	CG-OD2	6.67	1.38	1.25
2	E	10	GLY	C-O	6.36	1.30	1.23
2	E	104	ASP	CA-C	6.35	1.61	1.53
2	E	105	ALA	N-CA	6.25	1.54	1.46
1	D	314	ILE	N-CA	5.63	1.50	1.46
1	A	481	VAL	N-CA	5.62	1.50	1.46
2	E	104	ASP	C-N	5.35	1.41	1.33
3	O	34	TYR	N-CA	5.25	1.50	1.46
2	E	104	ASP	CA-CB	5.22	1.61	1.53
3	J	34	TYR	N-CA	5.19	1.50	1.46

All (103) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	104	ASP	CA-CB-CG	12.18	124.78	112.60
2	E	104	ASP	CB-CA-C	11.67	130.06	112.09
1	D	117	ASN	CA-CB-CG	9.72	122.32	112.60
2	H	104	ASP	CB-CA-C	8.82	125.68	112.09
1	A	117	ASN	CA-CB-CG	8.68	121.28	112.60
2	E	101	ASN	CA-CB-CG	8.46	121.06	112.60
1	C	117	ASN	CA-CB-CG	8.06	120.66	112.60
1	B	117	ASN	CA-CB-CG	8.04	120.64	112.60
2	E	104	ASP	N-CA-C	-7.98	100.22	111.24
2	H	101	ASN	CA-CB-CG	7.95	120.55	112.60
2	P	112	GLY	CA-C-O	-7.50	116.60	122.52
3	L	36	ASN	CA-CB-CG	7.31	119.91	112.60
1	D	168	ASP	CA-CB-CG	7.26	119.86	112.60
1	A	87	TYR	CA-C-O	-6.92	111.44	119.05
2	P	104	ASP	CB-CA-C	6.87	122.85	112.06
2	K	104	ASP	CA-CB-CG	6.66	119.26	112.60
3	O	27	GLY	CA-C-O	-6.64	117.67	122.45
1	C	231	GLN	CA-C-O	-6.62	113.86	120.82
1	D	60	GLY	CA-C-O	-6.62	116.72	121.88
1	D	85	ASP	CA-CB-CG	6.58	119.18	112.60
2	K	101	ASN	CA-CB-CG	6.56	119.16	112.60
2	E	151	TYR	CA-C-O	-6.52	114.03	121.40
2	K	112	GLY	CA-C-O	-6.48	116.61	122.57
3	J	27	GLY	CA-C-O	-6.41	117.84	122.45
2	K	104	ASP	CB-CA-C	6.39	122.10	112.06
2	E	104	ASP	CA-C-O	-6.36	114.26	122.14
2	E	106	PHE	CB-CA-C	-6.34	101.28	111.86
2	E	112	GLY	CA-C-O	-6.32	117.87	122.23
2	P	106	PHE	CB-CA-C	-6.30	99.86	110.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	27	GLY	CA-C-O	-6.17	118.01	122.45
1	A	403	ASN	CA-CB-CG	6.13	118.73	112.60
1	A	206	LYS	CB-CA-C	-6.07	100.72	110.79
1	D	243	HIS	N-CA-CB	-6.05	102.55	110.59
1	A	301	ILE	N-CA-C	-6.03	106.38	111.56
2	K	151	TYR	CA-C-O	-6.02	114.59	121.40
1	C	404	GLY	CA-C-N	6.02	129.26	120.71
1	C	404	GLY	C-N-CA	6.02	129.26	120.71
1	B	301	ILE	N-CA-C	-5.98	106.42	111.56
1	B	117	ASN	CB-CA-C	5.96	120.63	110.74
1	A	323	ASP	CA-CB-CG	5.95	118.55	112.60
1	B	281	TYR	CA-C-O	-5.94	114.61	121.15
1	D	403	ASN	CB-CA-C	5.94	122.25	110.42
1	B	85	ASP	CA-CB-CG	5.82	118.42	112.60
1	A	301	ILE	CB-CA-C	5.66	116.82	111.44
1	A	203	GLU	CB-CG-CD	5.64	122.19	112.60
1	B	226	ASP	CA-CB-CG	5.61	118.21	112.60
3	L	27	GLY	CA-C-O	-5.61	118.41	122.45
2	K	106	PHE	CB-CA-C	-5.58	101.05	110.19
2	H	65	LYS	CB-CA-C	5.55	118.16	110.16
3	L	51	TYR	CA-C-O	-5.55	114.67	121.78
1	A	117	ASN	CB-CA-C	5.55	119.95	110.74
3	L	33	TYR	CA-C-N	5.54	128.81	122.83
3	L	33	TYR	C-N-CA	5.54	128.81	122.83
2	E	152	PHE	CB-CA-C	-5.50	99.33	110.17
1	C	36	ASP	CA-CB-CG	5.50	118.10	112.60
1	D	288	GLY	CA-C-O	-5.48	116.59	121.41
1	A	271	ASP	CA-CB-CG	5.46	118.06	112.60
2	H	104	ASP	CA-C-N	5.46	130.78	121.29
2	H	104	ASP	C-N-CA	5.46	130.78	121.29
3	L	109	VAL	CA-C-O	-5.44	114.71	120.48
2	H	107	ASP	CB-CA-C	-5.41	99.66	110.42
1	C	58	MET	CA-C-O	-5.38	115.24	120.89
1	D	226	ASP	CA-CB-CG	5.38	117.98	112.60
1	A	36	ASP	CA-CB-CG	5.37	117.97	112.60
3	O	61	PRO	CA-C-N	5.34	128.42	120.31
3	O	61	PRO	C-N-CA	5.34	128.42	120.31
1	A	85	ASP	CA-CB-CG	5.29	117.89	112.60
1	C	117	ASN	CB-CA-C	5.28	119.51	110.74
1	B	445	SER	CA-C-O	-5.27	115.87	122.03
1	A	226	ASP	CA-CB-CG	5.25	117.85	112.60
2	E	101	ASN	OD1-CG-ND2	-5.25	117.35	122.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	42	GLU	CA-C-O	-5.24	115.44	121.47
1	D	54	ALA	CA-C-N	5.24	127.82	120.28
1	D	54	ALA	C-N-CA	5.24	127.82	120.28
1	D	323	ASP	CA-CB-CG	5.23	117.83	112.60
1	B	245	ASN	CA-C-N	5.23	129.10	121.31
1	B	245	ASN	C-N-CA	5.23	129.10	121.31
3	F	110	LEU	N-CA-C	-5.21	102.45	109.95
1	D	314	ILE	N-CA-CB	5.21	117.64	111.08
1	D	271	ASP	CA-CB-CG	5.20	117.80	112.60
1	C	85	ASP	CA-CB-CG	5.17	117.78	112.60
2	E	104	ASP	CA-C-N	5.17	131.42	121.54
2	E	104	ASP	C-N-CA	5.17	131.42	121.54
1	C	288	GLY	CA-C-O	-5.17	116.86	121.41
1	C	140	LEU	N-CA-C	5.17	116.66	108.96
1	A	506	ASP	CA-CB-CG	5.17	117.77	112.60
2	H	112	GLY	CA-C-O	-5.16	118.44	122.52
1	C	323	ASP	CA-CB-CG	5.15	117.75	112.60
2	K	191	PRO	CA-C-N	5.14	127.42	120.38
2	K	191	PRO	C-N-CA	5.14	127.42	120.38
1	C	42	GLU	CA-C-N	5.13	128.96	121.31
1	C	42	GLU	C-N-CA	5.13	128.96	121.31
1	A	284	GLY	O-C-N	5.13	126.66	122.51
1	C	509	GLN	N-CA-C	-5.11	107.43	113.97
1	D	117	ASN	OD1-CG-ND2	-5.10	117.50	122.60
1	C	134	GLU	CB-CG-CD	5.09	121.26	112.60
2	K	104	ASP	N-CA-C	-5.05	104.06	111.17
2	E	192	SER	CA-C-N	5.04	127.54	120.28
2	E	192	SER	C-N-CA	5.04	127.54	120.28
1	B	446	THR	CB-CA-C	5.04	120.44	110.42
3	L	50	ILE	CA-C-O	-5.03	115.19	120.27
2	P	106	PHE	N-CA-C	5.02	116.70	108.52
1	B	271	ASP	CA-CB-CG	5.02	117.62	112.60

There are no chirality outliers.

All (8) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	282	ALA	Peptide
1	C	75	GLU	Peptide
2	E	119	SER	Peptide
3	F	111	GLY	Peptide
2	H	152	PHE	Peptide

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Mol	Chain	Res	Type	Group
2	H	8	GLY	Peptide
2	K	152	PHE	Peptide
3	L	110	LEU	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	4087	0	4079	113	0
1	B	4081	0	4064	132	0
1	C	4071	0	4045	130	0
1	D	4095	0	4088	125	0
2	E	1649	0	1590	88	0
2	H	1654	0	1595	71	0
2	K	1654	0	1595	92	2
2	P	928	0	873	64	0
3	F	1575	0	1516	61	0
3	J	1573	0	1515	50	0
3	L	1573	0	1515	46	0
3	O	813	0	783	34	2
4	G	28	0	25	0	0
5	I	39	0	34	0	0
6	D	14	0	13	2	0
All	All	27834	0	27330	890	2

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 16.

All (890) 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)
2:K:105:ALA:HB1	3:J:36:ASN:ND2	1.28	1.42
2:K:105:ALA:CB	3:J:36:ASN:HD22	1.54	1.20
2:P:105:ALA:HB1	3:O:36:ASN:HD22	0.99	1.16
1:B:145:LYS:NZ	2:H:53:TYR:OH	1.87	1.07
2:H:153:PRO:HG3	2:H:206:HIS:CD2	1.93	1.04

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:292:ILE:HG23	1:C:300:VAL:HG21	1.40	1.04
1:B:321:LYS:O	1:B:324:ILE:HG22	1.58	1.03
1:A:117:ASN:OD1	1:A:220:HIS:NE2	1.91	1.03
1:B:161:TYR:HB2	2:H:102:TRP:CE3	1.94	1.02
1:D:336:THR:HG21	1:D:408:TRP:HB3	1.41	1.02
1:C:386:MET:CE	1:C:474:VAL:HB	1.90	1.01
1:B:117:ASN:OD1	1:B:220:HIS:NE2	1.92	1.01
1:D:117:ASN:OD1	1:D:220:HIS:NE2	1.93	1.01
2:P:105:ALA:HB1	3:O:36:ASN:ND2	1.76	1.00
1:C:117:ASN:OD1	1:C:220:HIS:NE2	1.92	1.00
3:F:110:LEU:HB2	3:F:144:TYR:OH	1.59	1.00
2:E:106:PHE:HB2	3:F:38:TYR:OH	1.61	0.99
2:K:153:PRO:HG3	2:K:206:HIS:CD2	1.95	0.99
1:C:44:THR:O	1:C:51:CYS:SG	2.21	0.98
3:F:110:LEU:HD12	3:F:144:TYR:HE2	1.26	0.98
1:C:47:ASP:OD2	1:C:499:ASN:ND2	1.96	0.98
1:C:58:MET:O	1:C:310:LEU:O	1.80	0.97
1:B:161:TYR:CE2	1:B:174:VAL:HG22	1.98	0.97
1:C:386:MET:HE2	1:C:474:VAL:HB	1.47	0.96
2:H:132:PRO:HB3	2:H:144:LEU:HB3	1.44	0.96
2:E:106:PHE:CE2	3:F:99:TRP:CH2	2.55	0.95
1:A:125:GLU:OE2	1:A:153:GLN:NE2	2.00	0.94
3:F:110:LEU:HD12	3:F:144:TYR:CE2	2.03	0.93
2:P:100:PRO:HD2	2:P:105:ALA:O	1.69	0.92
1:B:437:HIS:CE1	1:B:440:HIS:HE2	1.88	0.91
1:C:154:ILE:O	1:C:157:LEU:HG	1.69	0.91
3:J:56:ARG:NH2	3:J:60:VAL:O	2.05	0.90
1:C:44:THR:C	1:C:51:CYS:SG	2.56	0.89
3:L:109:VAL:O	3:L:110:LEU:O	1.91	0.89
2:E:153:PRO:HG2	2:E:206:HIS:NE2	1.88	0.88
2:K:103:ASP:OD2	2:K:105:ALA:HB3	1.75	0.87
3:O:56:ARG:NH2	3:O:60:VAL:O	2.07	0.87
1:C:91:ILE:CG2	1:C:408:TRP:CH2	2.58	0.87
2:P:105:ALA:CB	3:O:36:ASN:HD22	1.85	0.87
2:K:100:PRO:HD2	2:K:105:ALA:O	1.76	0.85
2:E:87:SER:O	2:E:117:VAL:CG1	2.24	0.84
2:E:154:GLU:OE1	2:E:155:PRO:HA	1.77	0.84
2:P:9:GLY:O	2:P:114:MET:O	1.96	0.84
1:C:45:SER:HB3	1:C:51:CYS:HB3	1.57	0.84
3:F:110:LEU:CB	3:F:144:TYR:OH	2.26	0.84
2:K:105:ALA:CB	3:J:36:ASN:ND2	2.25	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:445:SER:O	1:B:447:GLY:N	2.11	0.83
1:C:87:TYR:HD2	1:C:119:GLU:HA	1.44	0.82
3:F:110:LEU:HB2	3:F:144:TYR:CZ	2.13	0.82
1:C:59:GLY:HA2	1:C:63:ARG:HH11	1.44	0.82
1:C:91:ILE:HG22	1:C:408:TRP:CH2	2.15	0.81
2:E:87:SER:O	2:E:117:VAL:HG11	1.80	0.81
2:E:106:PHE:CE2	3:F:99:TRP:CZ3	2.69	0.81
1:D:36:ASP:HB2	1:D:282:ALA:HB1	1.61	0.81
2:H:87:SER:O	2:H:117:VAL:CG1	2.29	0.81
3:J:109:VAL:C	3:J:111:GLY:H	1.89	0.81
1:D:48[A]:SER:HB2	1:D:283:PHE:CE1	2.16	0.81
3:L:112:GLN:HE22	3:L:174:ASN:HB3	1.44	0.81
2:H:9:GLY:O	2:H:114:MET:O	1.98	0.80
2:H:87:SER:O	2:H:117:VAL:HG13	1.80	0.80
2:E:52:TRP:CH2	2:E:104:ASP:HB3	2.16	0.80
2:K:9:GLY:O	2:K:114:MET:O	2.00	0.80
3:J:94:MET:HE1	3:J:100:VAL:HG12	1.63	0.80
2:E:106:PHE:CZ	3:F:99:TRP:CZ3	2.70	0.79
2:K:132:PRO:HB3	2:K:144:LEU:HB3	1.62	0.79
3:F:52:SER:O	3:F:53:THR:OG1	2.00	0.79
3:L:14:PRO:HD3	3:L:110:LEU:HB2	1.64	0.78
3:L:52:SER:O	3:L:53:THR:OG1	2.02	0.78
2:E:106:PHE:CD2	3:F:99:TRP:CH2	2.71	0.78
2:E:132:PRO:HB3	2:E:144:LEU:HB3	1.66	0.77
3:F:110:LEU:CD1	3:F:144:TYR:HE2	1.98	0.77
3:L:93:PHE:CE2	3:L:99:TRP:CE3	2.73	0.77
2:P:35:HIS:CG	2:P:106:PHE:HZ	2.04	0.76
2:K:106:PHE:CE1	3:J:99:TRP:CH2	2.74	0.76
1:D:48[B]:SER:HB3	1:D:283:PHE:CE1	2.21	0.76
1:A:163:VAL:H	2:E:102:TRP:HE1	1.30	0.75
1:A:162:LYS:HD2	2:E:102:TRP:CZ2	2.21	0.75
3:O:94:MET:HE1	3:O:100:VAL:HG12	1.67	0.75
1:C:295:ASP:OD1	1:C:296:GLU:N	2.19	0.75
2:E:102:TRP:HE3	2:E:103:ASP:N	1.83	0.74
1:C:273:ARG:HG2	1:C:299:ASN:OD1	1.87	0.74
3:L:112:GLN:NE2	3:L:174:ASN:O	2.21	0.74
2:P:103:ASP:HB3	3:O:52:SER:HA	1.68	0.73
1:D:117:ASN:OD1	1:D:243:HIS:ND1	2.20	0.73
1:A:374:ARG:NH1	1:A:534:LYS:HE3	2.02	0.73
2:K:152:PHE:HD1	2:K:152:PHE:O	1.72	0.73
1:A:481:VAL:HG13	1:B:345:TYR:HB3	1.70	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:O:33:TYR:HE1	3:O:94:MET:HG3	1.53	0.73
2:K:83:MET:HB3	2:K:86:LEU:HD21	1.70	0.73
3:O:52:SER:OG	3:O:55:THR:HB	1.89	0.73
3:F:93:PHE:CD2	3:F:99:TRP:CD1	2.77	0.73
2:K:59:TYR:CD1	3:J:96:SER:O	2.41	0.73
3:L:109:VAL:O	3:L:110:LEU:C	2.32	0.72
3:F:94:MET:HE1	3:F:100:VAL:HG12	1.71	0.72
2:K:35:HIS:CG	2:K:106:PHE:HZ	2.07	0.72
2:P:101:ASN:O	2:P:104:ASP:OD1	2.07	0.72
3:J:33:TYR:HE1	3:J:94:MET:HG3	1.55	0.72
1:C:117:ASN:OD1	1:C:243:HIS:ND1	2.21	0.72
2:H:152:PHE:N	2:H:153:PRO:HD2	2.05	0.72
1:A:117:ASN:OD1	1:A:243:HIS:ND1	2.22	0.72
3:O:93:PHE:CD2	3:O:99:TRP:CD1	2.78	0.72
3:F:154:ALA:HB2	3:F:195:TYR:CE1	2.25	0.72
3:L:94:MET:HE1	3:L:100:VAL:HG12	1.72	0.71
1:C:91:ILE:HG21	1:C:408:TRP:CH2	2.25	0.71
1:D:118:HIS:NE2	1:D:395:ARG:NH1	2.38	0.71
3:J:93:PHE:CD2	3:J:99:TRP:CD1	2.78	0.71
1:B:400:GLU:C	1:B:402:ASN:H	1.97	0.71
1:A:179:LYS:HD2	1:A:195:ASP:HA	1.73	0.71
1:B:386:MET:HE2	1:B:455:ILE:HG12	1.71	0.71
1:C:122:ASN:HB3	1:C:126:GLY:HA3	1.73	0.70
2:E:106:PHE:CE2	3:F:99:TRP:CZ2	2.79	0.70
3:L:112:GLN:NE2	3:L:174:ASN:HB3	2.06	0.70
1:B:117:ASN:OD1	1:B:243:HIS:ND1	2.22	0.70
1:C:151:ALA:O	1:C:154:ILE:HG22	1.91	0.70
1:D:36:ASP:HB2	1:D:282:ALA:CB	2.21	0.70
2:P:106:PHE:HB2	3:O:38:TYR:OH	1.90	0.70
1:D:48[A]:SER:HB2	1:D:283:PHE:CZ	2.26	0.70
1:C:40:ARG:HE	1:C:48:SER:HB2	1.57	0.70
1:D:122:ASN:HB3	1:D:126:GLY:HA3	1.74	0.70
2:P:35:HIS:CG	2:P:106:PHE:CZ	2.79	0.70
2:K:152:PHE:O	2:K:152:PHE:CD1	2.45	0.70
2:H:153:PRO:CG	2:H:206:HIS:CD2	2.74	0.70
1:B:161:TYR:HB2	2:H:102:TRP:HE3	1.51	0.70
2:P:83:MET:HB3	2:P:86:LEU:HD21	1.73	0.69
6:D:601:NAG:O7	6:D:601:NAG:O3	2.09	0.69
1:C:45:SER:HB3	1:C:51:CYS:CB	2.23	0.69
2:E:87:SER:O	2:E:117:VAL:HG13	1.93	0.69
2:E:106:PHE:CD2	3:F:99:TRP:CZ2	2.80	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:K:153:PRO:CG	2:K:206:HIS:CD2	2.76	0.69
1:C:294:PHE:CD2	1:C:295:ASP:O	2.45	0.68
1:D:129:GLU:HB3	1:D:130:PRO:HD3	1.75	0.68
1:A:325:ASN:HA	1:A:328:ARG:HB2	1.74	0.68
1:C:179:LYS:HD2	1:C:195:ASP:HA	1.74	0.68
2:H:152:PHE:CD1	2:H:152:PHE:O	2.46	0.68
2:K:7:SER:OG	2:K:21:SER:O	2.09	0.68
2:K:35:HIS:CG	2:K:106:PHE:CZ	2.82	0.68
3:J:49:LEU:HB3	3:J:50:ILE:HD12	1.76	0.68
1:C:325:ASN:HA	1:C:328:ARG:HB2	1.76	0.68
2:E:52:TRP:HH2	2:E:104:ASP:CB	2.07	0.68
1:A:142:ALA:O	1:A:145:LYS:NZ	2.25	0.67
1:A:278:VAL:HG11	1:A:290:LEU:HD22	1.76	0.67
1:B:43:GLN:HG2	1:B:316:GLU:OE2	1.94	0.67
1:B:386:MET:CE	1:B:455:ILE:HG12	2.23	0.67
1:C:59:GLY:HA2	1:C:63:ARG:NH1	2.10	0.67
1:C:504:GLY:CA	1:C:508:PHE:O	2.42	0.67
1:A:432:LYS:HG2	1:A:489:MET:HE1	1.76	0.67
1:C:162:LYS:HD3	2:K:102:TRP:CE2	2.30	0.67
2:E:51:PHE:HE1	2:E:55:GLY:HA2	1.60	0.67
2:K:59:TYR:CG	3:J:96:SER:O	2.47	0.67
1:A:129:GLU:HB3	1:A:130:PRO:HD3	1.76	0.67
1:D:336:THR:HG21	1:D:408:TRP:CB	2.23	0.67
1:B:129:GLU:HB3	1:B:130:PRO:HD3	1.76	0.67
1:C:247:PHE:CD2	1:C:257:GLU:HB3	2.30	0.67
1:C:58:MET:HG2	1:C:285:LYS:HG3	1.76	0.66
1:D:206:LYS:HD3	2:P:32:TYR:CZ	2.30	0.66
1:B:161:TYR:CE1	1:B:203:GLU:HB3	2.31	0.66
1:B:247:PHE:CD2	1:B:257:GLU:HB3	2.30	0.66
1:B:444:GLN:N	1:B:444:GLN:OE1	2.29	0.66
1:B:122:ASN:HB3	1:B:126:GLY:HA3	1.77	0.66
1:C:129:GLU:HB3	1:C:130:PRO:HD3	1.77	0.66
1:D:325:ASN:HA	1:D:328:ARG:HB2	1.78	0.66
1:A:247:PHE:CD2	1:A:257:GLU:HB3	2.30	0.66
1:B:332:ASP:O	1:B:335:SER:OG	2.14	0.66
1:A:163:VAL:H	2:E:102:TRP:NE1	1.94	0.66
2:K:7:SER:CB	2:K:21:SER:O	2.43	0.66
2:K:106:PHE:CE1	3:J:99:TRP:CZ3	2.84	0.65
2:K:106:PHE:CD1	3:J:99:TRP:CH2	2.84	0.65
2:K:152:PHE:N	2:K:153:PRO:HD2	2.11	0.65
1:C:155:SER:OG	3:J:95:GLY:HA3	1.96	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:102:TRP:CE3	2:E:103:ASP:N	2.65	0.65
1:A:79:LEU:HD21	1:A:164:LEU:CD1	2.27	0.65
2:K:154:GLU:HG2	2:K:182:TYR:HE2	1.61	0.65
1:D:247:PHE:CD2	1:D:257:GLU:HB3	2.32	0.65
1:C:292:ILE:HG23	1:C:300:VAL:CG2	2.23	0.65
1:A:162:LYS:CD	2:E:102:TRP:CZ2	2.80	0.64
2:P:59:TYR:CD1	3:O:96:SER:O	2.49	0.64
1:B:145:LYS:HE3	1:B:194:GLU:CD	2.23	0.64
1:D:38:HIS:NE2	1:D:118:HIS:HD2	1.95	0.64
3:F:33:TYR:CE1	3:F:94:MET:HG3	2.33	0.64
2:E:52:TRP:CH2	2:E:104:ASP:CB	2.80	0.64
2:H:101:ASN:O	2:H:104:ASP:OD1	2.15	0.64
3:F:51:TYR:HE1	3:F:55:THR:HG21	1.63	0.64
2:H:154:GLU:N	2:H:155:PRO:CD	2.61	0.63
3:L:92:LEU:O	3:L:99:TRP:CD1	2.52	0.63
1:A:332:ASP:O	1:A:335:SER:OG	2.14	0.63
1:A:539:TYR:CD2	1:B:373:LEU:HD12	2.33	0.63
1:B:161:TYR:CD2	1:B:174:VAL:HG22	2.32	0.63
3:L:93:PHE:CE2	3:L:99:TRP:CD2	2.87	0.63
2:E:102:TRP:HE3	2:E:103:ASP:CA	2.11	0.63
2:K:74:ASN:C	2:K:76:LYS:H	2.07	0.63
3:O:49:LEU:HB3	3:O:50:ILE:HD12	1.79	0.62
1:B:161:TYR:CE1	1:B:203:GLU:CB	2.82	0.62
3:F:85:GLU:OE1	3:F:110:LEU:HD21	2.00	0.62
2:P:74:ASN:C	2:P:76:LYS:H	2.08	0.62
3:J:109:VAL:C	3:J:111:GLY:N	2.55	0.62
1:D:35:ASN:ND2	1:D:282:ALA:HB2	2.15	0.62
2:P:102:TRP:C	2:P:104:ASP:H	2.07	0.62
1:B:161:TYR:CE1	1:B:174:VAL:HG13	2.35	0.62
1:D:36:ASP:OD1	1:D:242:GLY:O	2.17	0.62
1:D:213:ASN:HD22	1:D:297:ARG:HA	1.63	0.62
2:E:52:TRP:HH2	2:E:104:ASP:HB3	1.59	0.62
2:E:53:TYR:CZ	2:E:54:ASP:HB3	2.34	0.62
2:P:35:HIS:ND1	2:P:106:PHE:HZ	1.97	0.62
2:K:57:ASN:HB3	2:K:59:TYR:HE2	1.64	0.61
2:K:102:TRP:C	2:K:104:ASP:H	2.08	0.61
1:D:162:LYS:HE3	1:D:164:LEU:HG	1.82	0.61
1:A:539:TYR:HD2	1:B:373:LEU:HD12	1.65	0.61
3:F:140:ILE:CD1	3:F:199:VAL:HG21	2.30	0.61
2:K:154:GLU:N	2:K:155:PRO:CD	2.63	0.61
2:H:35:HIS:ND1	2:H:106:PHE:HZ	1.98	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:51:CYS:O	1:C:52:VAL:C	2.44	0.61
1:A:213:ASN:HB2	1:A:298:GLY:HA3	1.82	0.61
1:D:201:GLN:HB3	1:D:202:PRO:HD3	1.82	0.61
1:A:122:ASN:HB3	1:A:126:GLY:HA3	1.83	0.61
1:B:440:HIS:O	1:B:465:PRO:HB3	2.01	0.61
2:H:87:SER:O	2:H:117:VAL:HG11	2.01	0.61
1:C:87:TYR:CD2	1:C:119:GLU:HA	2.31	0.61
1:B:278:VAL:HG11	1:B:290:LEU:HD22	1.83	0.60
3:J:33:TYR:CE1	3:J:94:MET:HG3	2.36	0.60
1:A:79:LEU:HD21	1:A:164:LEU:HD12	1.83	0.60
2:E:192:SER:O	2:E:195:LEU:HG	2.01	0.60
2:K:52:TRP:CZ3	2:K:104:ASP:OD1	2.55	0.60
2:E:153:PRO:HD2	2:E:206:HIS:CE1	2.36	0.60
2:P:59:TYR:CG	3:O:96:SER:O	2.54	0.60
1:A:201:GLN:HB3	1:A:202:PRO:HD3	1.84	0.60
1:D:27:TRP:CE2	1:D:77:ASN:ND2	2.70	0.60
1:D:278:VAL:HG11	1:D:290:LEU:HD22	1.84	0.60
2:K:52:TRP:HZ3	2:K:104:ASP:OD1	1.85	0.60
3:F:12:VAL:O	3:F:109:VAL:HG13	2.01	0.60
2:P:7:SER:HB3	2:P:21:SER:HB2	1.83	0.60
2:K:122:THR:HA	2:K:152:PHE:HB3	1.84	0.60
2:K:154:GLU:HG2	2:K:182:TYR:CE2	2.36	0.60
2:K:154:GLU:HA	2:K:182:TYR:CE2	2.36	0.60
1:B:444:GLN:CD	1:B:445:SER:H	2.10	0.60
1:A:355:PHE:CE2	1:C:356:ARG:HD3	2.36	0.59
1:C:201:GLN:HB3	1:C:202:PRO:HD3	1.83	0.59
1:C:332:ASP:O	1:C:335:SER:OG	2.16	0.59
2:H:106:PHE:CD1	3:L:99:TRP:HZ3	2.20	0.59
1:B:201:GLN:HB3	1:B:202:PRO:HD3	1.83	0.59
1:B:437:HIS:CE1	1:B:440:HIS:NE2	2.68	0.59
2:H:152:PHE:N	2:H:153:PRO:CD	2.65	0.59
2:E:199:THR:HG23	2:E:216:LYS:HE3	1.84	0.59
2:K:35:HIS:ND1	2:K:106:PHE:HZ	2.00	0.59
2:K:137:THR:OG1	2:K:142:ALA:HA	2.03	0.59
2:H:137:THR:OG1	2:H:142:ALA:HA	2.02	0.59
3:L:92:LEU:O	3:L:99:TRP:HD1	1.86	0.59
2:H:151:TYR:HD2	2:H:153:PRO:HG2	1.68	0.59
3:J:112:GLN:HG2	3:J:144:TYR:CD2	2.38	0.59
3:L:93:PHE:HE2	3:L:99:TRP:CE3	2.18	0.58
1:C:58:MET:CG	1:C:285:LYS:HG3	2.33	0.58
1:C:27:TRP:CE2	1:C:77:ASN:ND2	2.71	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:J:109:VAL:O	3:J:111:GLY:N	2.37	0.58
1:B:140:LEU:HA	1:B:159:LEU:O	2.03	0.58
1:D:378:GLU:HB2	1:D:381:TRP:CD1	2.38	0.58
3:L:14:PRO:HD3	3:L:110:LEU:CB	2.31	0.58
2:E:153:PRO:HG2	2:E:208:PRO:HB2	1.84	0.58
1:B:213:ASN:HB2	1:B:298:GLY:HA3	1.86	0.58
1:D:220:HIS:CD2	1:D:243:HIS:HB2	2.39	0.58
2:E:190:VAL:HG11	2:E:200:TYR:CE1	2.37	0.58
3:L:93:PHE:CD2	3:L:99:TRP:CD2	2.92	0.58
1:B:445:SER:C	1:B:447:GLY:H	2.07	0.58
3:L:53:THR:HG21	3:L:68:ILE:HG13	1.86	0.58
1:A:385:SER:OG	1:A:477:THR:CG2	2.52	0.58
1:A:385:SER:OG	1:A:477:THR:HG21	2.03	0.58
1:B:478:LYS:O	1:C:401:ARG:CZ	2.52	0.58
1:C:137:PHE:CE2	1:C:139:ILE:HD11	2.39	0.58
1:A:477:THR:C	1:A:479:CYS:H	2.12	0.57
1:C:52:VAL:HG13	1:C:56:ARG:HB2	1.86	0.57
1:A:27:TRP:CE2	1:A:77:ASN:ND2	2.72	0.57
1:B:85:ASP:CG	1:B:220:HIS:CE1	2.83	0.57
3:L:91:VAL:HG22	3:L:99:TRP:NE1	2.20	0.57
1:C:386:MET:HE2	1:C:474:VAL:CB	2.30	0.57
1:D:378:GLU:HB2	1:D:381:TRP:HD1	1.69	0.57
2:H:136:SER:OG	3:L:120:THR:HG23	2.04	0.57
1:A:27:TRP:CZ2	1:A:77:ASN:ND2	2.73	0.57
3:L:108:THR:HG21	3:L:145:PRO:HB3	1.86	0.57
1:D:281:TYR:O	1:D:282:ALA:CB	2.53	0.57
2:H:35:HIS:ND1	2:H:106:PHE:CZ	2.72	0.57
3:O:33:TYR:CE1	3:O:94:MET:HG3	2.38	0.57
1:D:85:ASP:CG	1:D:220:HIS:CE1	2.83	0.57
1:D:27:TRP:CZ2	1:D:77:ASN:ND2	2.72	0.56
1:B:161:TYR:CZ	1:B:203:GLU:CB	2.88	0.56
1:C:206:LYS:HE2	2:K:32:TYR:CE1	2.40	0.56
1:D:32:LEU:HB3	1:D:287:LEU:HD11	1.87	0.56
2:P:106:PHE:CE1	3:O:99:TRP:CH2	2.94	0.56
1:D:448:GLU:HB3	1:D:506:ASP:HB3	1.87	0.56
2:H:154:GLU:HA	2:H:182:TYR:CE2	2.40	0.56
1:B:480:ARG:HG2	1:B:481:VAL:H	1.71	0.56
1:B:44:THR:HB	1:B:49:SER:O	2.06	0.56
1:C:85:ASP:CG	1:C:220:HIS:CE1	2.83	0.56
1:A:85:ASP:CG	1:A:220:HIS:CE1	2.84	0.56
1:C:27:TRP:CZ2	1:C:77:ASN:ND2	2.74	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:160:PRO:HG3	2:K:101:ASN:HB3	1.88	0.56
1:C:453:GLY:O	1:C:455:ILE:HD12	2.05	0.56
1:C:44:THR:HB	1:C:48:SER:HA	1.88	0.55
1:C:162:LYS:HE3	1:C:164:LEU:HG	1.88	0.55
2:H:190:VAL:HG11	2:H:200:TYR:CE1	2.40	0.55
1:A:482:PRO:HG3	1:B:344:VAL:HG11	1.88	0.55
1:B:400:GLU:C	1:B:402:ASN:N	2.64	0.55
1:A:373:LEU:HD12	1:B:539:TYR:HD2	1.71	0.55
1:A:161:TYR:OH	2:E:100:PRO:HB2	2.06	0.55
1:A:476:CYS:CB	1:A:479:CYS:SG	2.94	0.55
1:C:58:MET:SD	1:C:309:LEU:HD11	2.46	0.55
1:C:476:CYS:SG	1:C:477:THR:N	2.80	0.55
2:E:51:PHE:CE1	2:E:55:GLY:HA2	2.41	0.55
1:D:48[B]:SER:HB3	1:D:283:PHE:CZ	2.42	0.55
2:H:152:PHE:O	2:H:152:PHE:HD1	1.90	0.55
2:K:153:PRO:HD3	2:K:206:HIS:CE1	2.42	0.55
2:K:199:THR:HG23	2:K:216:LYS:HE3	1.88	0.55
1:C:59:GLY:HA3	1:C:63:ARG:HD2	1.88	0.55
2:H:95:TYR:CD1	2:H:112:GLY:HA3	2.42	0.55
1:A:444:GLN:O	1:A:446:THR:HG22	2.07	0.55
1:C:539:TYR:N	1:C:540:PRO:CD	2.70	0.54
1:D:44:THR:HB	1:D:49:SER:O	2.07	0.54
2:P:106:PHE:HB2	3:O:38:TYR:HH	1.71	0.54
3:J:52:SER:OG	3:J:55:THR:HB	2.06	0.54
1:C:87:TYR:HE1	1:C:98:GLY:HA2	1.72	0.54
2:H:102:TRP:C	2:H:104:ASP:H	2.14	0.54
2:H:153:PRO:HD3	2:H:206:HIS:CE1	2.41	0.54
1:C:278:VAL:HG11	1:C:290:LEU:HD22	1.88	0.54
1:A:130:PRO:O	1:A:134:GLU:HB2	2.06	0.54
1:A:402:ASN:CG	1:A:403:ASN:H	2.14	0.54
1:A:476:CYS:HB3	1:A:479:CYS:SG	2.48	0.54
1:D:214:LYS:HG2	1:D:294:PHE:CD1	2.42	0.54
2:K:153:PRO:HG3	2:K:206:HIS:NE2	2.22	0.54
1:C:504:GLY:O	1:C:508:PHE:O	2.26	0.54
2:E:194:SER:HB2	2:E:198:GLN:HB3	1.89	0.54
1:A:201:GLN:NE2	1:A:233:VAL:HA	2.22	0.54
1:A:373:LEU:HD12	1:B:539:TYR:CD2	2.42	0.54
1:A:539:TYR:N	1:A:540:PRO:CD	2.70	0.54
1:D:132:LEU:HD13	1:D:157:LEU:HB2	1.90	0.54
2:H:51:PHE:CE2	2:H:58:LYS:HG2	2.42	0.54
3:F:143:PHE:CE2	3:F:146:GLY:HA2	2.43	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:K:152:PHE:N	2:K:153:PRO:CD	2.70	0.54
2:H:153:PRO:HG3	2:H:206:HIS:NE2	2.22	0.54
1:B:321:LYS:HD2	1:B:324:ILE:HG21	1.89	0.54
2:P:28:THR:O	2:P:31:SER:OG	2.25	0.54
1:A:140:LEU:HA	1:A:159:LEU:O	2.07	0.54
2:P:106:PHE:CD1	3:O:99:TRP:CH2	2.95	0.54
2:K:190:VAL:HG11	2:K:200:TYR:CE1	2.42	0.54
1:D:220:HIS:CE1	1:D:243:HIS:HA	2.43	0.53
2:P:103:ASP:OD1	2:P:105:ALA:HB3	2.08	0.53
1:C:292:ILE:CG2	1:C:300:VAL:HG21	2.27	0.53
1:D:253:PRO:HG2	1:D:259:PRO:HD3	1.90	0.53
2:E:125:PRO:HD3	2:E:206:HIS:HD1	1.72	0.53
1:A:32:LEU:HB3	1:A:287:LEU:HD11	1.90	0.53
1:C:32:LEU:HB3	1:C:287:LEU:HD11	1.89	0.53
1:B:161:TYR:CZ	1:B:203:GLU:HB2	2.43	0.53
1:B:539:TYR:N	1:B:540:PRO:CD	2.71	0.53
1:C:267:VAL:O	1:C:274:LYS:O	2.26	0.53
1:C:386:MET:CE	1:C:474:VAL:CB	2.76	0.53
1:D:211:ASN:ND2	3:O:58:SER:OG	2.41	0.53
3:F:12:VAL:HG12	3:F:109:VAL:HG22	1.90	0.53
3:F:85:GLU:OE1	3:F:110:LEU:CD2	2.57	0.53
3:F:112:GLN:N	3:F:113:PRO:CD	2.72	0.53
1:C:385:SER:OG	1:C:477:THR:CG2	2.56	0.53
1:A:539:TYR:HD2	1:B:373:LEU:CD1	2.22	0.53
1:C:45:SER:CB	1:C:51:CYS:HB3	2.35	0.53
1:B:174:VAL:CG1	1:B:200:LEU:HD22	2.39	0.53
2:H:153:PRO:HG3	2:H:206:HIS:CG	2.43	0.53
3:J:113:PRO:O	3:J:115:ALA:N	2.42	0.53
1:A:220:HIS:CE1	1:A:243:HIS:HA	2.44	0.52
2:H:154:GLU:N	2:H:155:PRO:HD2	2.24	0.52
3:L:33:TYR:HB3	3:L:92:LEU:HB3	1.91	0.52
2:E:35:HIS:ND1	2:E:106:PHE:HZ	2.06	0.52
3:L:110:LEU:HD22	3:L:111:GLY:N	2.24	0.52
2:E:83:MET:HB3	2:E:86:LEU:HD21	1.91	0.52
1:B:27:TRP:CZ2	1:B:77:ASN:ND2	2.78	0.52
1:B:120:PHE:CE1	1:B:127:LEU:HD22	2.44	0.52
1:B:220:HIS:CE1	1:B:243:HIS:HA	2.44	0.52
1:C:39:SER:O	1:C:39:SER:OG	2.24	0.52
1:D:118:HIS:CD2	1:D:395:ARG:NH1	2.77	0.52
1:D:539:TYR:N	1:D:540:PRO:CD	2.73	0.52
2:P:106:PHE:CE1	3:O:99:TRP:CZ3	2.97	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:402:ASN:CG	1:A:403:ASN:N	2.67	0.52
1:D:132:LEU:HD13	1:D:157:LEU:CB	2.40	0.52
2:H:12:VAL:O	2:H:117:VAL:HA	2.10	0.52
2:H:83:MET:HB3	2:H:86:LEU:HD21	1.91	0.52
2:E:102:TRP:C	2:E:104:ASP:H	2.17	0.52
2:K:47:TRP:HZ2	2:K:50:VAL:HG12	1.75	0.52
2:H:57:ASN:HB3	2:H:59:TYR:CZ	2.44	0.52
2:E:154:GLU:HG2	2:E:182:TYR:CD2	2.45	0.52
1:A:210:LEU:HD21	2:E:107:ASP:OD2	2.10	0.52
2:H:122:THR:HA	2:H:152:PHE:HB3	1.90	0.52
1:A:478:LYS:HD3	1:A:494:LYS:HZ1	1.75	0.52
1:B:174:VAL:HG11	1:B:200:LEU:HD22	1.92	0.52
1:B:220:HIS:HE2	1:B:243:HIS:HD1	1.58	0.52
1:B:437:HIS:ND1	1:B:440:HIS:CD2	2.78	0.52
1:C:220:HIS:CE1	1:C:243:HIS:HA	2.45	0.52
1:A:206:LYS:HZ2	2:E:101:ASN:H	1.57	0.52
1:B:367:ALA:HA	1:B:531:TYR:CD2	2.45	0.52
1:A:475:LEU:CD2	1:A:482:PRO:HB2	2.41	0.51
1:B:32:LEU:HB3	1:B:287:LEU:HD11	1.92	0.51
1:D:220:HIS:HE2	1:D:243:HIS:HD1	1.57	0.51
1:B:27:TRP:CE2	1:B:77:ASN:ND2	2.78	0.51
2:K:102:TRP:HZ3	3:J:52:SER:HB2	1.76	0.51
2:K:154:GLU:N	2:K:155:PRO:HD2	2.25	0.51
1:B:335:SER:HA	1:B:408:TRP:HB3	1.93	0.51
2:E:47:TRP:HZ2	2:E:50:VAL:HG12	1.75	0.51
3:F:56:ARG:HD3	3:F:64:PHE:O	2.09	0.51
2:P:102:TRP:CE3	2:P:103:ASP:HB3	2.46	0.51
1:B:161:TYR:CE2	1:B:174:VAL:CG2	2.84	0.51
1:B:206:LYS:NZ	2:H:101:ASN:OD1	2.35	0.51
1:D:35:ASN:O	1:D:282:ALA:HB1	2.11	0.51
3:J:52:SER:OG	3:J:55:THR:CB	2.59	0.51
1:C:475:LEU:HD23	1:C:483:SER:O	2.10	0.51
1:A:360:MET:O	1:A:363:LEU:HB3	2.11	0.51
1:B:178:SER:HB2	1:B:221:SER:HA	1.92	0.51
1:B:317:ASP:OD2	1:B:320:ILE:HG12	2.11	0.51
1:C:175:GLY:HA2	1:C:218:LEU:O	2.10	0.51
1:D:152:SER:CA	2:P:57:ASN:HD21	2.24	0.51
1:D:332:ASP:O	1:D:335:SER:OG	2.17	0.51
2:H:184:LEU:HA	3:L:181:TYR:OH	2.11	0.51
2:K:151:TYR:HD2	2:K:153:PRO:HG2	1.75	0.51
1:B:386:MET:CE	1:B:474:VAL:HB	2.41	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:96:TYR:CE2	1:D:327:TRP:CE3	2.99	0.51
1:C:87:TYR:HE1	1:C:98:GLY:CA	2.23	0.51
2:E:12:VAL:O	2:E:117:VAL:HA	2.11	0.51
2:E:102:TRP:CE3	2:E:103:ASP:HB3	2.46	0.51
2:K:12:VAL:O	2:K:117:VAL:HA	2.11	0.51
1:B:117:ASN:OD1	1:B:220:HIS:CD2	2.63	0.50
1:C:484:TYR:CE2	1:D:543:GLU:HA	2.45	0.50
1:D:118:HIS:NE2	1:D:395:ARG:CZ	2.74	0.50
2:P:59:TYR:CE1	3:O:96:SER:O	2.64	0.50
1:A:335:SER:HA	1:A:408:TRP:HB3	1.92	0.50
1:C:58:MET:HE3	1:C:285:LYS:HE3	1.92	0.50
1:C:91:ILE:HG22	1:C:408:TRP:HH2	1.72	0.50
1:D:67:LYS:HB2	1:D:314:ILE:HD13	1.93	0.50
1:D:145:LYS:HE3	1:D:194:GLU:CD	2.36	0.50
2:H:47:TRP:HZ2	2:H:50:VAL:HG12	1.77	0.50
2:K:57:ASN:HD22	2:K:59:TYR:HE2	1.57	0.50
1:B:360:MET:O	1:B:363:LEU:HB3	2.11	0.50
1:C:508:PHE:HA	1:C:510:MET:HE1	1.93	0.50
3:F:24:LEU:HD13	3:F:26:SER:H	1.76	0.50
2:P:6:GLU:OE2	2:P:6:GLU:N	2.44	0.50
3:O:51:TYR:O	3:O:52:SER:OG	2.29	0.50
1:B:480:ARG:HG3	1:B:481:VAL:HG23	1.93	0.50
1:C:96:TYR:CE2	1:C:327:TRP:CE3	2.99	0.50
1:C:385:SER:OG	1:C:477:THR:HG22	2.12	0.50
1:D:175:GLY:HA2	1:D:218:LEU:O	2.12	0.50
3:L:168:PRO:HA	3:L:178:ALA:HB2	1.94	0.50
1:A:162:LYS:HD2	2:E:102:TRP:HZ2	1.72	0.50
1:A:402:ASN:OD1	1:A:403:ASN:N	2.37	0.50
1:D:464:LYS:HD2	1:D:465:PRO:O	2.11	0.50
2:K:153:PRO:HG3	2:K:206:HIS:CG	2.43	0.50
3:F:38:TYR:CD1	3:F:48:THR:HA	2.47	0.50
2:P:47:TRP:HZ2	2:P:50:VAL:HG12	1.76	0.50
1:D:368:MET:HG2	1:D:421:PHE:CD2	2.47	0.50
1:A:175:GLY:HA2	1:A:218:LEU:O	2.11	0.50
1:B:219:GLY:O	1:B:241:GLY:HA2	2.12	0.50
1:C:147:LYS:O	1:C:190:ASN:ND2	2.41	0.50
1:C:360:MET:O	1:C:363:LEU:HB3	2.11	0.50
1:D:206:LYS:HB3	2:P:32:TYR:OH	2.12	0.50
1:D:398:ILE:HD11	1:D:414:VAL:HG21	1.94	0.50
2:E:95:TYR:CD1	2:E:112:GLY:HA3	2.47	0.50
2:E:53:TYR:CE1	2:E:54:ASP:CG	2.90	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:220:HIS:HE2	1:C:243:HIS:HD1	1.59	0.49
3:F:50:ILE:HD12	3:F:66:GLY:HA3	1.94	0.49
2:K:5:VAL:HG12	2:K:22:CYS:HA	1.94	0.49
1:A:220:HIS:HE2	1:A:243:HIS:HD1	1.60	0.49
1:A:381:TRP:CG	1:A:423:LEU:HD13	2.47	0.49
1:D:147:LYS:O	1:D:190:ASN:ND2	2.42	0.49
1:D:317:ASP:OD2	1:D:320:ILE:HG12	2.13	0.49
2:H:38:ARG:HA	2:H:93:VAL:O	2.11	0.49
1:B:96:TYR:CE2	1:B:327:TRP:CE3	3.00	0.49
1:B:175:GLY:HA2	1:B:218:LEU:O	2.11	0.49
2:H:6:GLU:OE2	2:H:6:GLU:N	2.45	0.49
1:A:348:GLY:H	1:A:397:PRO:HB3	1.77	0.49
1:B:445:SER:C	1:B:447:GLY:N	2.66	0.49
1:A:445:SER:O	1:A:446:THR:C	2.56	0.49
1:C:271:ASP:HB2	1:C:273:ARG:HD3	1.93	0.49
1:D:426:LEU:HD22	1:D:430:THR:HB	1.94	0.49
3:L:109:VAL:C	3:L:110:LEU:O	2.54	0.49
1:A:117:ASN:OD1	1:A:220:HIS:CD2	2.65	0.49
1:A:317:ASP:OD2	1:A:320:ILE:HG12	2.13	0.49
1:C:178:SER:HB2	1:C:221:SER:HA	1.94	0.49
1:D:140:LEU:HA	1:D:159:LEU:O	2.13	0.49
1:D:144:ILE:O	1:D:158:TYR:HE2	1.95	0.49
2:K:27:PHE:CD1	2:K:32:TYR:HD2	2.29	0.49
1:A:79:LEU:CD2	1:A:164:LEU:HD12	2.43	0.49
1:A:220:HIS:CD2	1:A:243:HIS:HB2	2.48	0.49
1:B:231:GLN:HA	1:B:269:SER:HA	1.94	0.49
1:C:381:TRP:CG	1:C:423:LEU:HD13	2.47	0.49
1:D:158:TYR:O	1:D:158:TYR:CD1	2.65	0.49
2:H:153:PRO:CG	2:H:206:HIS:NE2	2.76	0.49
2:H:153:PRO:HD3	2:H:206:HIS:NE2	2.28	0.49
2:K:69:THR:HB	2:K:82:GLN:HB3	1.93	0.49
1:B:368:MET:HG2	1:B:421:PHE:CD2	2.47	0.49
1:C:294:PHE:CE2	1:C:295:ASP:O	2.65	0.49
2:E:150:ASP:OD2	3:F:133:LYS:NZ	2.46	0.49
2:K:47:TRP:CZ2	3:J:97:GLY:O	2.66	0.49
1:A:145:LYS:HG3	2:E:53:TYR:OH	2.12	0.49
1:A:373:LEU:CD1	1:B:539:TYR:HD2	2.26	0.49
1:B:158:TYR:O	1:B:158:TYR:CD1	2.66	0.49
1:B:381:TRP:CG	1:B:423:LEU:HD13	2.47	0.49
1:C:117:ASN:OD1	1:C:220:HIS:CD2	2.64	0.49
1:D:67:LYS:HE2	1:D:314:ILE:HG21	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:360:MET:O	1:D:363:LEU:HB3	2.12	0.49
3:J:38:TYR:CD1	3:J:48:THR:HA	2.48	0.49
1:A:476:CYS:HB2	1:A:479:CYS:SG	2.52	0.49
1:B:161:TYR:CE1	1:B:203:GLU:HB2	2.47	0.49
1:B:510:MET:HE3	1:B:514:GLU:OE2	2.13	0.48
1:C:368:MET:HG2	1:C:421:PHE:CD2	2.48	0.48
1:D:152:SER:HA	2:P:57:ASN:ND2	2.28	0.48
1:D:214:LYS:HG2	1:D:294:PHE:CE1	2.48	0.48
2:P:69:THR:HB	2:P:82:GLN:HB3	1.94	0.48
2:P:102:TRP:O	2:P:104:ASP:N	2.46	0.48
2:E:38:ARG:HA	2:E:93:VAL:O	2.14	0.48
2:P:12:VAL:O	2:P:117:VAL:HA	2.12	0.48
3:J:211:ALA:HB1	3:J:212:PRO:HD2	1.95	0.48
2:H:184:LEU:HD12	2:H:184:LEU:C	2.38	0.48
2:K:44:GLY:HA2	3:J:89:TYR:OH	2.13	0.48
3:J:10:PHE:CE2	3:J:18:VAL:CG2	2.96	0.48
3:J:114:LYS:O	3:J:114:LYS:HG2	2.13	0.48
1:B:290:LEU:HD21	1:B:292:ILE:HD11	1.95	0.48
2:E:123:LYS:HB2	2:E:151:TYR:HA	1.95	0.48
2:P:38:ARG:HA	2:P:93:VAL:O	2.13	0.48
1:A:132:LEU:HB3	1:A:157:LEU:HB3	1.95	0.48
2:H:100:PRO:HG2	2:H:105:ALA:HB3	1.94	0.48
2:E:153:PRO:CG	2:E:206:HIS:NE2	2.70	0.48
2:K:153:PRO:CG	2:K:206:HIS:NE2	2.77	0.48
1:A:368:MET:HG2	1:A:421:PHE:CD2	2.48	0.48
1:B:130:PRO:O	1:B:134:GLU:HB2	2.14	0.48
1:B:478:LYS:O	1:C:401:ARG:NH1	2.46	0.48
1:C:426:LEU:HD22	1:C:430:THR:HB	1.95	0.48
2:K:38:ARG:HA	2:K:93:VAL:O	2.14	0.48
1:A:158:TYR:O	1:A:158:TYR:CD1	2.67	0.48
2:E:101:ASN:O	2:E:102:TRP:O	2.31	0.48
2:E:137:THR:OG1	2:E:142:ALA:HA	2.14	0.48
3:F:168:PRO:HA	3:F:178:ALA:HB2	1.96	0.48
2:K:153:PRO:HB2	2:K:155:PRO:HD2	1.96	0.48
1:B:325:ASN:HA	1:B:328:ARG:HB2	1.95	0.48
3:F:51:TYR:HE1	3:F:55:THR:CG2	2.26	0.48
3:J:151:ALA:HB3	3:J:198:GLN:HB2	1.95	0.48
1:B:400:GLU:HG2	1:B:401:ARG:HG3	1.95	0.48
1:B:426:LEU:HD22	1:B:430:THR:HB	1.96	0.48
1:D:377:ASP:O	1:D:379:MET:HG3	2.14	0.48
2:H:106:PHE:CE1	3:L:99:TRP:HZ3	2.32	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:184:LEU:C	2:E:184:LEU:HD12	2.39	0.48
1:D:72:ARG:HA	1:D:78:VAL:HG21	1.95	0.47
1:D:152:SER:HA	2:P:57:ASN:HD21	1.80	0.47
1:D:348:GLY:H	1:D:397:PRO:HB3	1.78	0.47
3:F:11:SER:OG	3:F:202:GLU:OE2	2.23	0.47
2:P:103:ASP:CB	3:O:52:SER:HA	2.40	0.47
2:K:105:ALA:HB1	3:J:36:ASN:HD22	0.68	0.47
2:K:106:PHE:HB2	3:J:38:TYR:OH	2.14	0.47
1:C:220:HIS:CD2	1:C:243:HIS:HB2	2.49	0.47
3:L:56:ARG:HD3	3:L:64:PHE:O	2.14	0.47
3:L:156:SER:HB2	3:F:157:SER:OG	2.14	0.47
3:O:14:PRO:HG3	3:O:109:VAL:HG12	1.95	0.47
2:K:57:ASN:HB3	2:K:59:TYR:CE2	2.47	0.47
2:K:184:LEU:C	2:K:184:LEU:HD12	2.39	0.47
3:J:90:CYS:SG	3:J:91:VAL:N	2.87	0.47
1:A:79:LEU:CD2	1:A:164:LEU:CD1	2.92	0.47
1:A:341:LYS:HB3	1:A:341:LYS:HE2	1.62	0.47
1:B:120:PHE:HE1	1:B:127:LEU:CD2	2.27	0.47
1:C:504:GLY:C	1:C:508:PHE:O	2.57	0.47
1:D:402:ASN:C	1:D:404:GLY:H	2.23	0.47
2:H:153:PRO:HB2	2:H:155:PRO:HD2	1.96	0.47
3:F:12:VAL:CG1	3:F:109:VAL:HG22	2.45	0.47
3:F:110:LEU:N	3:F:144:TYR:OH	2.48	0.47
2:K:102:TRP:CE3	2:K:103:ASP:HB3	2.49	0.47
1:A:96:TYR:CE2	1:A:327:TRP:CE3	3.02	0.47
1:B:161:TYR:CZ	1:B:203:GLU:HB3	2.49	0.47
1:C:290:LEU:HD21	1:C:292:ILE:HD11	1.97	0.47
1:D:35:ASN:ND2	1:D:282:ALA:CB	2.77	0.47
1:D:219:GLY:O	1:D:241:GLY:HA2	2.15	0.47
3:L:112:GLN:OE1	3:L:144:TYR:CE2	2.67	0.47
2:K:102:TRP:CZ3	3:J:52:SER:HB2	2.49	0.47
1:A:144:ILE:O	1:A:158:TYR:HE2	1.97	0.47
1:B:91:ILE:HG12	1:B:331:LEU:HD22	1.97	0.47
2:E:207:LYS:N	2:E:208:PRO:CD	2.78	0.47
1:A:244:SER:O	1:A:279:GLN:OE1	2.32	0.47
1:A:515:LEU:HD12	1:A:515:LEU:C	2.40	0.47
1:B:341:LYS:HE3	1:B:403:ASN:O	2.15	0.47
1:D:539:TYR:N	1:D:539:TYR:CD1	2.82	0.47
3:F:53:THR:HG21	3:F:68:ILE:HG13	1.97	0.47
3:O:52:SER:HG	3:O:55:THR:HB	1.80	0.47
1:B:515:LEU:C	1:B:515:LEU:HD12	2.39	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:J:168:PRO:HA	3:J:178:ALA:HB2	1.97	0.47
1:D:38:HIS:NE2	1:D:395:ARG:NH1	2.62	0.47
1:D:290:LEU:HD21	1:D:292:ILE:HD11	1.97	0.47
1:D:378:GLU:CB	1:D:381:TRP:HB2	2.45	0.47
1:B:441:ARG:O	1:B:444:GLN:NE2	2.47	0.46
1:C:154:ILE:O	1:C:157:LEU:CG	2.53	0.46
3:L:90:CYS:SG	3:L:91:VAL:N	2.88	0.46
2:K:102:TRP:O	2:K:104:ASP:N	2.48	0.46
2:K:207:LYS:N	2:K:208:PRO:CD	2.78	0.46
1:A:486:PRO:O	1:A:488:LYS:HE3	2.16	0.46
1:A:536[A]:LYS:HD2	1:B:186:ASN:OD1	2.15	0.46
1:B:206:LYS:HB2	2:H:102:TRP:CH2	2.50	0.46
1:D:515:LEU:C	1:D:515:LEU:HD12	2.40	0.46
1:B:426:LEU:HD23	1:B:426:LEU:HA	1.72	0.46
1:C:120:PHE:HE1	1:C:127:LEU:CD2	2.28	0.46
1:D:36:ASP:HA	1:D:85:ASP:HB2	1.97	0.46
1:D:161:TYR:CD1	1:D:161:TYR:C	2.92	0.46
1:A:426:LEU:HD22	1:A:430:THR:HB	1.96	0.46
1:B:310:LEU:HA	1:B:314:ILE:HD12	1.97	0.46
1:B:220:HIS:CD2	1:B:243:HIS:HB2	2.51	0.46
1:B:480:ARG:HG2	1:B:481:VAL:N	2.31	0.46
1:D:360:MET:HE1	1:D:398:ILE:HD12	1.98	0.46
2:E:103:ASP:HB2	3:F:34:TYR:HB3	1.97	0.46
1:C:539:TYR:N	1:C:539:TYR:CD1	2.84	0.46
1:D:178:SER:HB2	1:D:221:SER:HA	1.98	0.46
1:A:219:GLY:O	1:A:241:GLY:HA2	2.16	0.46
1:A:292:ILE:HG22	1:A:294:PHE:CE1	2.51	0.46
1:B:197:ILE:HD11	1:B:228:LEU:HD23	1.97	0.46
1:D:398:ILE:HD11	1:D:414:VAL:CG2	2.45	0.46
1:B:120:PHE:HE1	1:B:127:LEU:HD22	1.81	0.46
2:H:73:ASP:OD1	2:H:75:SER:HB2	2.16	0.46
3:F:51:TYR:HD1	3:F:51:TYR:C	2.24	0.46
3:O:52:SER:OG	3:O:55:THR:CB	2.61	0.46
2:K:74:ASN:C	2:K:76:LYS:N	2.74	0.46
2:K:102:TRP:C	2:K:104:ASP:N	2.74	0.46
1:B:278:VAL:O	1:B:279:GLN:HB2	2.15	0.46
1:B:437:HIS:ND1	1:B:440:HIS:NE2	2.64	0.46
2:H:100:PRO:CG	2:H:105:ALA:HB3	2.46	0.46
2:H:207:LYS:N	2:H:208:PRO:CD	2.78	0.46
2:E:153:PRO:CG	2:E:208:PRO:HB2	2.45	0.46
1:A:163:VAL:N	2:E:102:TRP:HE1	2.07	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:281:TYR:O	1:B:282:ALA:O	2.34	0.46
2:H:91:THR:HA	2:H:115:VAL:O	2.16	0.46
3:F:154:ALA:HB2	3:F:195:TYR:CD1	2.50	0.46
2:K:194:SER:HB2	2:K:198:GLN:HB3	1.97	0.46
1:A:508:PHE:HA	1:A:510:MET:HE1	1.97	0.45
1:B:161:TYR:CZ	1:B:174:VAL:HG13	2.50	0.45
1:B:348:GLY:H	1:B:397:PRO:HB3	1.81	0.45
1:D:138:PRO:HG3	1:D:162:LYS:HE2	1.98	0.45
3:L:14:PRO:CD	3:L:110:LEU:HB2	2.39	0.45
1:C:219:GLY:O	1:C:241:GLY:HA2	2.16	0.45
1:D:117:ASN:OD1	1:D:220:HIS:CD2	2.67	0.45
1:D:206:LYS:HD3	2:P:32:TYR:CE2	2.51	0.45
2:H:184:LEU:C	3:L:181:TYR:OH	2.60	0.45
2:E:121:SER:HB2	2:E:123:LYS:HE2	1.99	0.45
2:P:8:GLY:HA2	2:P:113:THR:OG1	2.16	0.45
1:C:161:TYR:CD1	1:C:161:TYR:C	2.95	0.45
2:H:34:MET:HB3	2:H:79:LEU:HD22	1.98	0.45
2:K:35:HIS:HB2	2:K:106:PHE:HE2	1.82	0.45
3:J:41:THR:O	3:J:42:PRO:C	2.59	0.45
1:A:201:GLN:HE22	1:A:233:VAL:HA	1.81	0.45
1:A:475:LEU:HD21	1:A:482:PRO:HB2	1.98	0.45
1:A:548:PHE:HB2	1:C:351:GLN:HE22	1.81	0.45
1:B:321:LYS:C	1:B:324:ILE:HG22	2.37	0.45
1:C:91:ILE:CG2	1:C:408:TRP:HH2	2.23	0.45
1:D:213:ASN:ND2	1:D:297:ARG:HA	2.29	0.45
1:C:437:HIS:CE1	1:C:507:GLY:HA3	2.52	0.45
2:H:206:HIS:CG	2:H:206:HIS:O	2.70	0.45
2:E:106:PHE:CE2	3:F:99:TRP:CE3	3.04	0.45
2:E:141:THR:HA	2:E:191:PRO:HA	1.99	0.45
2:P:102:TRP:C	2:P:104:ASP:N	2.74	0.45
1:B:437:HIS:HA	1:B:440:HIS:CD2	2.52	0.45
2:E:52:TRP:HH2	2:E:104:ASP:HA	1.82	0.45
2:E:91:THR:HA	2:E:115:VAL:O	2.17	0.45
2:K:114:MET:HE2	2:K:154:GLU:CD	2.41	0.45
1:A:397:PRO:O	1:A:398:ILE:HD13	2.17	0.45
1:A:426:LEU:HA	1:A:426:LEU:HD23	1.74	0.45
1:A:539:TYR:N	1:A:539:TYR:CD1	2.83	0.45
1:C:45:SER:N	1:C:51:CYS:SG	2.90	0.45
1:D:38:HIS:NE2	1:D:118:HIS:CD2	2.82	0.45
1:D:206:LYS:NZ	2:P:27:PHE:HB2	2.32	0.45
1:D:282:ALA:O	1:D:284:GLY:N	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:508:PHE:HA	1:D:510:MET:HE1	1.99	0.45
2:E:47:TRP:CD1	3:F:99:TRP:HE3	2.35	0.45
2:H:65:LYS:C	2:H:67:ARG:H	2.24	0.44
2:E:106:PHE:HB2	3:F:38:TYR:CZ	2.50	0.44
1:D:281:TYR:CE1	1:D:351:GLN:NE2	2.85	0.44
2:E:154:GLU:CG	2:E:182:TYR:CE2	3.01	0.44
2:K:206:HIS:CG	2:K:206:HIS:O	2.69	0.44
1:A:349:SER:HB3	1:D:479:CYS:CB	2.47	0.44
1:B:164:LEU:HD23	1:B:164:LEU:HA	1.81	0.44
2:H:100:PRO:HD3	2:H:107:ASP:CG	2.42	0.44
2:H:154:GLU:HA	2:H:182:TYR:CD2	2.52	0.44
3:F:112:GLN:H	3:F:113:PRO:HD2	1.82	0.44
2:P:106:PHE:O	3:O:38:TYR:OH	2.20	0.44
3:O:41:THR:O	3:O:42:PRO:C	2.59	0.44
3:F:144:TYR:HA	3:F:145:PRO:C	2.42	0.44
2:K:65:LYS:C	2:K:67:ARG:H	2.26	0.44
1:A:87:TYR:O	1:A:93:PHE:CD1	2.70	0.44
2:H:185:SER:N	3:L:181:TYR:OH	2.50	0.44
3:F:151:ALA:HB3	3:F:198:GLN:HB2	1.99	0.44
2:K:43:LYS:HE2	2:K:43:LYS:HA	1.99	0.44
1:A:290:LEU:HD21	1:A:292:ILE:HD11	1.98	0.44
1:C:426:LEU:HD23	1:C:426:LEU:HA	1.72	0.44
1:C:476:CYS:C	1:C:478:LYS:H	2.25	0.44
1:C:515:LEU:HD12	1:C:515:LEU:C	2.42	0.44
3:F:145:PRO:HD2	3:F:201:HIS:NE2	2.33	0.44
2:P:59:TYR:CD2	3:O:96:SER:O	2.71	0.44
1:A:244:SER:C	1:A:246:THR:H	2.25	0.44
1:B:539:TYR:N	1:B:539:TYR:CD1	2.84	0.44
1:C:159:LEU:HD22	2:K:102:TRP:CD1	2.53	0.44
1:D:95:VAL:HG12	1:D:96:TYR:CD1	2.52	0.44
1:D:162:LYS:HD3	2:P:102:TRP:NE1	2.32	0.44
1:A:154:ILE:O	1:A:155:SER:C	2.61	0.44
1:C:59:GLY:CA	1:C:63:ARG:HD2	2.48	0.44
1:D:281:TYR:O	1:D:282:ALA:HB2	2.17	0.44
2:H:103:ASP:O	2:H:104:ASP:C	2.61	0.44
3:L:36:ASN:ND2	3:L:38:TYR:OH	2.51	0.44
3:L:49:LEU:HD23	3:L:50:ILE:HG13	2.00	0.44
2:E:118:SER:HB2	2:E:120:ALA:HB2	2.00	0.44
2:K:153:PRO:HD3	2:K:206:HIS:NE2	2.32	0.44
1:A:548:PHE:HB2	1:C:351:GLN:NE2	2.33	0.44
2:E:51:PHE:CE2	2:E:58:LYS:HG2	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:41:THR:O	3:F:42:PRO:C	2.60	0.44
3:F:51:TYR:C	3:F:51:TYR:CD1	2.95	0.44
3:J:179:SER:HB2	3:J:181:TYR:CE1	2.53	0.44
1:A:160:PRO:HG3	2:E:101:ASN:HB3	1.99	0.43
1:B:210:LEU:CD1	2:H:102:TRP:HE1	2.31	0.43
1:B:273:ARG:HB2	1:B:273:ARG:NH1	2.33	0.43
2:H:55:GLY:O	2:H:56:SER:C	2.60	0.43
2:H:132:PRO:HB3	2:H:144:LEU:CB	2.32	0.43
2:H:141:THR:HA	2:H:191:PRO:HA	2.00	0.43
2:E:55:GLY:O	2:E:56:SER:C	2.61	0.43
2:P:91:THR:HA	2:P:115:VAL:O	2.17	0.43
3:O:94:MET:HE2	3:O:94:MET:H	1.83	0.43
1:D:152:SER:N	2:P:57:ASN:HD21	2.15	0.43
3:F:110:LEU:CA	3:F:144:TYR:OH	2.65	0.43
1:B:296:GLU:O	1:B:296:GLU:HG2	2.18	0.43
1:D:378:GLU:CG	1:D:379:MET:N	2.80	0.43
2:E:35:HIS:ND1	2:E:106:PHE:CZ	2.86	0.43
2:E:73:ASP:OD1	2:E:75:SER:HB2	2.18	0.43
2:E:152:PHE:HB3	2:E:153:PRO:HD3	2.01	0.43
3:O:56:ARG:NH1	3:O:62:ASP:HA	2.32	0.43
2:K:7:SER:HB3	2:K:113:THR:HG21	2.00	0.43
1:B:39:SER:O	1:B:39:SER:OG	2.28	0.43
2:E:135:LYS:O	3:F:208:LYS:NZ	2.44	0.43
2:E:152:PHE:O	2:E:152:PHE:CD2	2.71	0.43
2:P:35:HIS:HB2	2:P:106:PHE:CE2	2.54	0.43
2:K:2:VAL:HG23	2:K:27:PHE:CD2	2.53	0.43
2:K:11:VAL:HA	2:K:116:THR:O	2.19	0.43
2:K:14:PRO:HD3	2:K:118:SER:O	2.18	0.43
1:B:64:LEU:HD23	1:B:108:LEU:HD21	2.00	0.43
1:C:293:GLU:O	1:C:300:VAL:HG23	2.18	0.43
3:F:90:CYS:SG	3:F:91:VAL:N	2.91	0.43
2:K:104:ASP:O	3:J:93:PHE:CD2	2.72	0.43
3:J:208:LYS:HD3	3:J:208:LYS:HA	1.93	0.43
3:L:41:THR:O	3:L:42:PRO:C	2.60	0.43
2:K:141:THR:HA	2:K:191:PRO:HA	2.00	0.43
1:A:548:PHE:CB	1:C:351:GLN:NE2	2.82	0.43
1:B:161:TYR:OH	1:B:204:VAL:HG23	2.18	0.43
1:D:111:ASP:HB3	1:D:164:LEU:HD11	2.01	0.43
1:D:144:ILE:O	1:D:158:TYR:CE2	2.71	0.43
1:D:162:LYS:HA	2:P:102:TRP:CD1	2.53	0.43
2:P:11:VAL:HA	2:P:116:THR:O	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:349:SER:HB3	1:D:479:CYS:HB2	2.01	0.43
1:C:210:LEU:HD23	1:C:210:LEU:HA	1.89	0.43
1:C:243:HIS:CE1	1:C:377:ASP:O	2.71	0.43
1:D:280:ALA:O	1:D:281:TYR:C	2.62	0.43
2:P:51:PHE:CE1	2:P:55:GLY:HA2	2.54	0.43
2:K:91:THR:HA	2:K:115:VAL:O	2.18	0.43
1:C:46:GLU:HG3	1:C:256:LYS:HE3	2.01	0.43
1:C:63:ARG:NH1	1:C:63:ARG:HB2	2.34	0.43
1:C:264:PRO:HG3	1:C:303:SER:HB2	1.99	0.43
1:D:435:PHE:O	1:D:438:SER:HB3	2.19	0.43
2:P:35:HIS:CD2	2:P:106:PHE:CZ	3.07	0.43
3:J:10:PHE:CD2	3:J:18:VAL:HG21	2.54	0.43
1:C:278:VAL:O	1:C:279:GLN:HB2	2.19	0.43
2:H:58:LYS:O	2:H:59:TYR:CD1	2.72	0.43
2:E:154:GLU:HG3	2:E:182:TYR:CE2	2.54	0.43
3:O:90:CYS:SG	3:O:91:VAL:N	2.91	0.43
1:A:425:GLN:HG2	1:A:492:VAL:CG1	2.49	0.42
1:D:90:THR:OG1	1:D:92:TRP:HB2	2.19	0.42
1:D:179:LYS:HA	1:D:193:PHE:HB3	2.01	0.42
3:L:91:VAL:HG22	3:L:99:TRP:CE2	2.53	0.42
2:E:65:LYS:C	2:E:67:ARG:H	2.26	0.42
1:C:485:ASP:OD1	1:C:485:ASP:N	2.52	0.42
2:E:174:ALA:HA	2:E:184:LEU:HB3	2.00	0.42
3:F:116:ALA:HB1	3:F:117:PRO:HD2	2.01	0.42
1:A:144:ILE:O	1:A:158:TYR:CE2	2.72	0.42
1:B:64:LEU:CD2	1:B:108:LEU:HD21	2.49	0.42
1:B:437:HIS:CD2	1:B:437:HIS:O	2.72	0.42
2:E:201:ILE:HG12	2:E:216:LYS:HA	2.02	0.42
3:F:24:LEU:CD1	3:F:26:SER:O	2.67	0.42
2:K:35:HIS:CD2	2:K:106:PHE:CZ	3.07	0.42
1:A:548:PHE:CG	1:C:351:GLN:NE2	2.87	0.42
1:C:91:ILE:HG21	1:C:408:TRP:CZ2	2.54	0.42
3:J:63:ARG:HB2	3:J:78:THR:O	2.19	0.42
3:J:211:ALA:HB1	3:J:212:PRO:CD	2.49	0.42
1:B:161:TYR:OH	1:B:203:GLU:HB2	2.20	0.42
1:C:348:GLY:H	1:C:397:PRO:HB3	1.83	0.42
2:K:35:HIS:HB2	2:K:106:PHE:CE2	2.54	0.42
2:K:55:GLY:O	2:K:56:SER:C	2.61	0.42
3:J:146:GLY:HA3	3:J:176:TYR:CD2	2.55	0.42
1:B:243:HIS:CE1	1:B:377:ASP:O	2.73	0.42
1:C:201:GLN:OE1	1:C:233:VAL:HA	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:92:TRP:CE2	1:D:331:LEU:HD11	2.53	0.42
3:L:63:ARG:NH1	3:L:81:GLN:HG3	2.34	0.42
2:P:55:GLY:O	2:P:56:SER:C	2.62	0.42
2:P:65:LYS:C	2:P:67:ARG:H	2.27	0.42
2:K:91:THR:HG23	2:K:116:THR:HA	2.01	0.42
3:J:63:ARG:NH1	3:J:81:GLN:HG3	2.34	0.42
3:J:116:ALA:HB1	3:J:117:PRO:HD2	2.01	0.42
1:A:349:SER:CB	1:D:479:CYS:HB2	2.50	0.42
1:C:33:HIS:CD2	1:C:288:GLY:HA3	2.55	0.42
1:B:444:GLN:CD	1:B:445:SER:N	2.77	0.42
1:C:386:MET:HE3	1:C:474:VAL:HB	1.89	0.42
3:L:94:MET:O	3:L:95:GLY:O	2.38	0.42
2:P:34:MET:HB3	2:P:79:LEU:HD22	2.02	0.42
1:A:243:HIS:CE1	1:A:377:ASP:O	2.73	0.42
1:D:100:GLU:HA	1:D:327:TRP:CZ3	2.55	0.42
1:D:152:SER:HB2	2:P:59:TYR:CE1	2.54	0.42
1:D:152:SER:HB2	2:P:59:TYR:OH	2.19	0.42
2:K:45:LEU:HD13	3:J:89:TYR:CE2	2.54	0.42
2:K:201:ILE:HG12	2:K:216:LYS:HA	2.02	0.42
1:A:344:VAL:HG11	1:B:482:PRO:HG3	2.01	0.42
1:B:108:LEU:HD23	1:B:110:TYR:HE1	1.84	0.42
1:C:164:LEU:HD23	1:C:164:LEU:HA	1.88	0.42
1:D:40:ARG:HD3	1:D:90:THR:HG23	2.02	0.42
1:D:278:VAL:O	1:D:279:GLN:HB2	2.20	0.42
1:D:281:TYR:HE1	1:D:351:GLN:NE2	2.17	0.42
3:L:110:LEU:C	3:L:110:LEU:HD13	2.45	0.42
2:E:103:ASP:O	2:E:104:ASP:C	2.63	0.42
2:P:91:THR:HG23	2:P:116:THR:HA	2.02	0.42
2:P:103:ASP:OD2	3:O:52:SER:HB3	2.19	0.42
1:A:435:PHE:O	1:A:438:SER:HB3	2.20	0.41
1:D:48[A]:SER:HB2	1:D:283:PHE:CD1	2.55	0.41
1:D:425:GLN:HG2	1:D:492:VAL:CG1	2.49	0.41
2:E:91:THR:HG23	2:E:116:THR:HA	2.02	0.41
2:P:74:ASN:C	2:P:76:LYS:N	2.75	0.41
2:K:27:PHE:CD1	2:K:32:TYR:CD2	3.07	0.41
3:J:171:GLN:OE1	3:J:177:ALA:HB2	2.20	0.41
1:C:90:THR:OG1	1:C:92:TRP:HB2	2.20	0.41
1:D:426:LEU:HA	1:D:426:LEU:HD23	1.70	0.41
2:K:174:ALA:HA	2:K:184:LEU:HB3	2.01	0.41
3:J:94:MET:HE2	3:J:94:MET:H	1.84	0.41
1:B:179:LYS:HD2	1:B:196:GLU:OE1	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:541:ALA:HA	1:D:482:PRO:HG2	2.01	0.41
2:H:194:SER:HB2	2:H:198:GLN:HB3	2.01	0.41
3:L:38:TYR:HE1	3:L:91:VAL:CG1	2.33	0.41
1:A:374:ARG:HH12	1:B:536:LYS:HD3	1.85	0.41
1:C:482:PRO:HG2	1:D:541:ALA:HA	2.01	0.41
1:D:154:ILE:O	1:D:155:SER:C	2.63	0.41
2:H:91:THR:HG23	2:H:116:THR:HA	2.00	0.41
3:L:92:LEU:C	3:L:99:TRP:HE1	2.28	0.41
3:L:151:ALA:HB3	3:L:198:GLN:HB2	2.02	0.41
2:E:102:TRP:C	2:E:104:ASP:N	2.78	0.41
1:B:154:ILE:O	1:B:155:SER:C	2.63	0.41
1:C:52:VAL:HG13	1:C:52:VAL:O	2.19	0.41
3:F:52:SER:O	3:F:53:THR:CB	2.68	0.41
2:P:74:ASN:O	2:P:76:LYS:N	2.51	0.41
2:P:103:ASP:O	2:P:104:ASP:C	2.63	0.41
1:B:282:ALA:O	1:B:284:GLY:N	2.53	0.41
2:H:106:PHE:CZ	3:L:99:TRP:CE3	3.08	0.41
2:H:151:TYR:OH	2:H:184:LEU:HD23	2.19	0.41
3:F:110:LEU:CD1	3:F:144:TYR:CE2	2.84	0.41
1:A:123:GLY:HA2	1:A:186:ASN:O	2.19	0.41
1:C:113:MET:HB3	1:C:139:ILE:HG12	2.01	0.41
2:E:11:VAL:HA	2:E:116:THR:O	2.21	0.41
2:K:74:ASN:O	2:K:76:LYS:N	2.51	0.41
1:A:152:SER:O	3:F:96:SER:HA	2.20	0.41
1:C:158:TYR:O	1:C:159:LEU:O	2.38	0.41
2:H:106:PHE:HB2	3:L:38:TYR:OH	2.21	0.41
3:F:110:LEU:CG	3:F:144:TYR:HE2	2.34	0.41
2:P:35:HIS:HB2	2:P:106:PHE:HE2	1.86	0.41
1:B:164:LEU:HD12	1:B:173:ILE:HD11	2.02	0.41
1:B:425:GLN:HG2	1:B:492:VAL:CG1	2.51	0.41
1:C:79:LEU:HD23	1:C:79:LEU:HA	1.96	0.41
2:H:11:VAL:HA	2:H:116:THR:O	2.21	0.41
2:E:34:MET:HB3	2:E:79:LEU:HD22	2.03	0.41
3:F:114:LYS:O	3:F:114:LYS:HG3	2.21	0.41
3:O:52:SER:O	3:O:53:THR:CB	2.68	0.41
1:A:461:LEU:HD11	1:A:548:PHE:CE2	2.56	0.41
1:B:90:THR:OG1	1:B:92:TRP:HB2	2.21	0.41
1:B:400:GLU:O	1:B:402:ASN:N	2.54	0.41
2:E:52:TRP:CH2	2:E:104:ASP:CG	2.99	0.41
3:O:63:ARG:CZ	3:O:81:GLN:HG3	2.50	0.41
1:A:310:LEU:HA	1:A:314:ILE:HD12	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:437:HIS:CD2	1:D:437:HIS:O	2.74	0.40
3:L:116:ALA:HB1	3:L:117:PRO:HD2	2.03	0.40
2:K:103:ASP:O	2:K:104:ASP:C	2.64	0.40
1:A:166:VAL:O	1:A:167:GLY:C	2.64	0.40
1:A:278:VAL:O	1:A:279:GLN:HB2	2.20	0.40
1:B:247:PHE:CE2	1:B:257:GLU:HB3	2.57	0.40
1:B:435:PHE:O	1:B:438:SER:HB3	2.21	0.40
1:C:144:ILE:O	1:C:158:TYR:OH	2.17	0.40
1:D:475:LEU:HD23	1:D:483:SER:O	2.22	0.40
1:A:206:LYS:NZ	2:E:101:ASN:H	2.20	0.40
1:B:123:GLY:HA2	1:B:186:ASN:O	2.21	0.40
1:C:375:HIS:ND1	1:C:375:HIS:N	2.69	0.40
2:E:106:PHE:CE2	3:F:99:TRP:CE2	3.09	0.40
1:B:33:HIS:CD2	1:B:288:GLY:HA3	2.56	0.40
1:B:95:VAL:HG12	1:B:96:TYR:CD1	2.56	0.40
1:D:158:TYR:O	1:D:158:TYR:CG	2.74	0.40
1:D:311:ASN:ND2	6:D:601:NAG:H61	2.37	0.40
2:H:102:TRP:C	2:H:104:ASP:N	2.80	0.40
3:J:136:LEU:HD23	3:J:136:LEU:HA	1.95	0.40
1:C:247:PHE:CE2	1:C:257:GLU:HB3	2.56	0.40
1:D:161:TYR:O	2:P:102:TRP:HD1	2.05	0.40
2:P:100:PRO:CD	2:P:105:ALA:O	2.56	0.40
3:J:127:GLU:O	3:J:130:GLN:HB2	2.20	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:O:58:SER:O	2:K:138:SER:N[8_455]	1.95	0.25
3:O:58:SER:O	2:K:138:SER:CA[8_455]	1.98	0.22

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	522/529 (99%)	451 (86%)	60 (12%)	11 (2%)	5	31
1	B	521/529 (98%)	454 (87%)	58 (11%)	9 (2%)	7	34
1	C	521/529 (98%)	444 (85%)	64 (12%)	13 (2%)	4	28
1	D	524/529 (99%)	456 (87%)	57 (11%)	11 (2%)	5	31
2	E	218/233 (94%)	185 (85%)	27 (12%)	6 (3%)	4	27
2	H	218/233 (94%)	186 (85%)	25 (12%)	7 (3%)	3	25
2	K	218/233 (94%)	187 (86%)	23 (11%)	8 (4%)	2	22
2	P	117/233 (50%)	104 (89%)	10 (8%)	3 (3%)	4	28
3	F	211/216 (98%)	185 (88%)	22 (10%)	4 (2%)	6	33
3	J	210/216 (97%)	184 (88%)	21 (10%)	5 (2%)	4	29
3	L	210/216 (97%)	183 (87%)	23 (11%)	4 (2%)	6	33
3	O	108/216 (50%)	89 (82%)	15 (14%)	4 (4%)	2	22
All	All	3598/3912 (92%)	3108 (86%)	405 (11%)	85 (2%)	4	29

All (85) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	85	ASP
1	A	282	ALA
1	A	442	TYR
1	B	85	ASP
1	B	282	ALA
1	B	446	THR
1	C	52	VAL
1	C	85	ASP
1	C	282	ALA
1	C	283	PHE
1	D	85	ASP
1	D	378	GLU
2	H	102	TRP
3	L	110	LEU
2	E	102	TRP
3	F	112	GLN
2	K	138	SER
3	J	114	LYS
1	A	243	HIS
1	A	279	GLN

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Mol	Chain	Res	Type
1	A	446	THR
1	B	243	HIS
1	B	279	GLN
1	B	283	PHE
1	B	505	GLY
1	C	159	LEU
1	C	243	HIS
1	C	279	GLN
1	C	506	ASP
1	D	279	GLN
1	D	282	ALA
1	D	311	ASN
1	D	337	GLN
1	D	383	HIS
2	H	28	THR
3	L	53	THR
2	E	105	ALA
3	F	53	THR
3	F	156	SER
2	P	28	THR
3	O	53	THR
3	O	85	GLU
2	K	6	GLU
2	K	7	SER
3	J	85	GLU
3	J	95	GLY
3	J	110	LEU
1	A	155	SER
1	A	283	PHE
1	A	402	ASN
1	B	155	SER
1	C	299	ASN
1	C	311	ASN
1	D	283	PHE
1	D	401	ARG
2	H	150	ASP
3	L	111	GLY
2	E	28	THR
2	E	150	ASP
3	O	54	ASN
2	K	28	THR
2	K	103	ASP

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Mol	Chain	Res	Type
2	K	150	ASP
3	J	52	SER
1	B	480	ARG
2	H	30	SER
2	P	75	SER
2	P	103	ASP
2	K	75	SER
1	A	281	TYR
1	A	478	LYS
2	H	138	SER
2	E	152	PHE
1	C	275	VAL
1	C	281	TYR
1	D	297	ARG
3	O	95	GLY
1	D	242	GLY
3	L	95	GLY
3	F	95	GLY
2	H	152	PHE
2	K	152	PHE
2	H	66	GLY
2	E	132	PRO
1	C	160	PRO

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	451/456 (99%)	425 (94%)	26 (6%)	18	45
1	B	450/456 (99%)	430 (96%)	20 (4%)	25	50
1	C	448/456 (98%)	421 (94%)	27 (6%)	17	44
1	D	453/456 (99%)	429 (95%)	24 (5%)	20	46
2	E	183/197 (93%)	166 (91%)	17 (9%)	8	32
2	H	182/197 (92%)	168 (92%)	14 (8%)	12	37

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	K	182/197 (92%)	170 (93%)	12 (7%)	15	42
2	P	97/197 (49%)	94 (97%)	3 (3%)	35	57
3	F	177/182 (97%)	156 (88%)	21 (12%)	5	23
3	J	176/182 (97%)	162 (92%)	14 (8%)	11	36
3	L	176/182 (97%)	154 (88%)	22 (12%)	4	21
3	O	91/182 (50%)	86 (94%)	5 (6%)	19	46
All	All	3066/3340 (92%)	2861 (93%)	205 (7%)	15	42

All (205) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	44	THR
1	A	46	GLU
1	A	75	GLU
1	A	87	TYR
1	A	119	GLU
1	A	125	GLU
1	A	133	LYS
1	A	164	LEU
1	A	252	ASN
1	A	268	THR
1	A	302	SER
1	A	312	SER
1	A	324	ILE
1	A	328	ARG
1	A	336	THR
1	A	341	LYS
1	A	350	SER
1	A	375	HIS
1	A	440	HIS
1	A	475	LEU
1	A	478	LYS
1	A	481	VAL
1	A	488	LYS
1	A	534	LYS
1	A	539	TYR
1	A	549	SER
1	B	39	SER
1	B	43	GLN
1	B	87	TYR

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Mol	Chain	Res	Type
1	B	108	LEU
1	B	119	GLU
1	B	297	ARG
1	B	300	VAL
1	B	302	SER
1	B	323	ASP
1	B	336	THR
1	B	350	SER
1	B	352	SER
1	B	356	ARG
1	B	375	HIS
1	B	377	ASP
1	B	440	HIS
1	B	510	MET
1	B	520	SER
1	B	539	TYR
1	B	549	SER
1	C	30	THR
1	C	39	SER
1	C	43	GLN
1	C	46	GLU
1	C	49	SER
1	C	51	CYS
1	C	87	TYR
1	C	119	GLU
1	C	134	GLU
1	C	140	LEU
1	C	157	LEU
1	C	164	LEU
1	C	169	GLU
1	C	190	ASN
1	C	256	LYS
1	C	312	SER
1	C	324	ILE
1	C	328	ARG
1	C	336	THR
1	C	350	SER
1	C	351	GLN
1	C	365	CYS
1	C	375	HIS
1	C	377	ASP
1	C	503	ASN

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Mol	Chain	Res	Type
1	C	539	TYR
1	C	549	SER
1	D	86	GLN
1	D	87	TYR
1	D	119	GLU
1	D	134	GLU
1	D	164	LEU
1	D	168	ASP
1	D	184	LEU
1	D	296	GLU
1	D	302	SER
1	D	312	SER
1	D	314	ILE
1	D	324	ILE
1	D	328	ARG
1	D	338	GLU
1	D	350	SER
1	D	398	ILE
1	D	440	HIS
1	D	445	SER
1	D	477	THR
1	D	481	VAL
1	D	489	MET
1	D	503	ASN
1	D	539	TYR
1	D	549	SER
2	H	2	VAL
2	H	20	LEU
2	H	102	TRP
2	H	103	ASP
2	H	106	PHE
2	H	119	SER
2	H	122	THR
2	H	134	SER
2	H	135	LYS
2	H	162	SER
2	H	167	SER
2	H	202	CYS
2	H	203	ASN
2	H	206	HIS
3	L	2	THR
3	L	10	PHE

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Mol	Chain	Res	Type
3	L	36	ASN
3	L	49	LEU
3	L	53	THR
3	L	57	SER
3	L	92	LEU
3	L	93	PHE
3	L	96	SER
3	L	108	THR
3	L	110	LEU
3	L	125	SER
3	L	126	SER
3	L	141	SER
3	L	142	ASP
3	L	148	VAL
3	L	149	THR
3	L	157	SER
3	L	159	VAL
3	L	160	LYS
3	L	172	SER
3	L	179	SER
2	E	2	VAL
2	E	6	GLU
2	E	20	LEU
2	E	102	TRP
2	E	104	ASP
2	E	119	SER
2	E	121	SER
2	E	122	THR
2	E	134	SER
2	E	135	LYS
2	E	136	SER
2	E	154	GLU
2	E	162	SER
2	E	167	SER
2	E	185	SER
2	E	202	CYS
2	E	203	ASN
3	F	2	THR
3	F	10	PHE
3	F	51	TYR
3	F	53	THR
3	F	57	SER

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Mol	Chain	Res	Type
3	F	92	LEU
3	F	93	PHE
3	F	96	SER
3	F	108	THR
3	F	109	VAL
3	F	125	SER
3	F	126	SER
3	F	140	ILE
3	F	141[A]	SER
3	F	141[B]	SER
3	F	155	ASP
3	F	157	SER
3	F	159	VAL
3	F	169	SER
3	F	172	SER
3	F	179	SER
2	P	2	VAL
2	P	20	LEU
2	P	76	LYS
3	O	10	PHE
3	O	50	ILE
3	O	57	SER
3	O	93	PHE
3	O	96	SER
2	K	2	VAL
2	K	5	VAL
2	K	20	LEU
2	K	76	LYS
2	K	104	ASP
2	K	119	SER
2	K	121	SER
2	K	135	LYS
2	K	162	SER
2	K	167	SER
2	K	202	CYS
2	K	203	ASN
3	J	1	GLN
3	J	10	PHE
3	J	50	ILE
3	J	57	SER
3	J	93	PHE
3	J	96	SER

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Mol	Chain	Res	Type
3	J	125	SER
3	J	126	SER
3	J	141	SER
3	J	157	SER
3	J	159	VAL
3	J	160	LYS
3	J	172	SER
3	J	179	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (55) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	35	ASN
1	A	38	HIS
1	A	53	ASN
1	A	77	ASN
1	A	88	GLN
1	A	201	GLN
1	A	371	ASN
1	A	372	ASN
1	A	375	HIS
1	A	440	HIS
1	A	456	HIS
1	B	35	ASN
1	B	70	GLN
1	B	190	ASN
1	B	211	ASN
1	B	371	ASN
1	B	437	HIS
1	C	35	ASN
1	C	38	HIS
1	C	279	GLN
1	C	311	ASN
1	C	351	GLN
1	C	370	ASN
1	C	372	ASN
1	C	425	GLN
1	C	440	HIS
1	C	456	HIS
1	D	35	ASN
1	D	53	ASN
1	D	86	GLN

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Mol	Chain	Res	Type
1	D	118	HIS
1	D	153	GLN
1	D	211	ASN
1	D	304	HIS
1	D	351	GLN
1	D	371	ASN
1	D	383	HIS
2	H	77	ASN
2	H	82	GLN
3	L	36	ASN
3	L	112	GLN
3	L	192	HIS
2	E	39	GLN
2	E	57	ASN
2	E	77	ASN
3	F	188	GLN
3	F	192	HIS
2	P	82	GLN
2	P	111	GLN
2	K	57	ASN
2	K	82	GLN
2	K	101	ASN
2	K	111	GLN
3	J	36	ASN
3	J	192	HIS

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 ⓘ

5 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The

Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	G	1	1,4	14,14,15	0.28	0	17,19,21	0.77	0
4	NAG	G	2	4	14,14,15	0.28	0	17,19,21	0.63	0
5	NAG	I	1	1,5	14,14,15	0.29	0	17,19,21	0.90	0
5	NAG	I	2	5	14,14,15	0.28	0	17,19,21	0.58	0
5	BMA	I	3	5	11,11,12	0.26	0	15,15,17	0.72	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	G	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	G	2	4	-	0/6/23/26	0/1/1/1
5	NAG	I	1	1,5	-	0/6/23/26	0/1/1/1
5	NAG	I	2	5	-	2/6/23/26	0/1/1/1
5	BMA	I	3	5	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

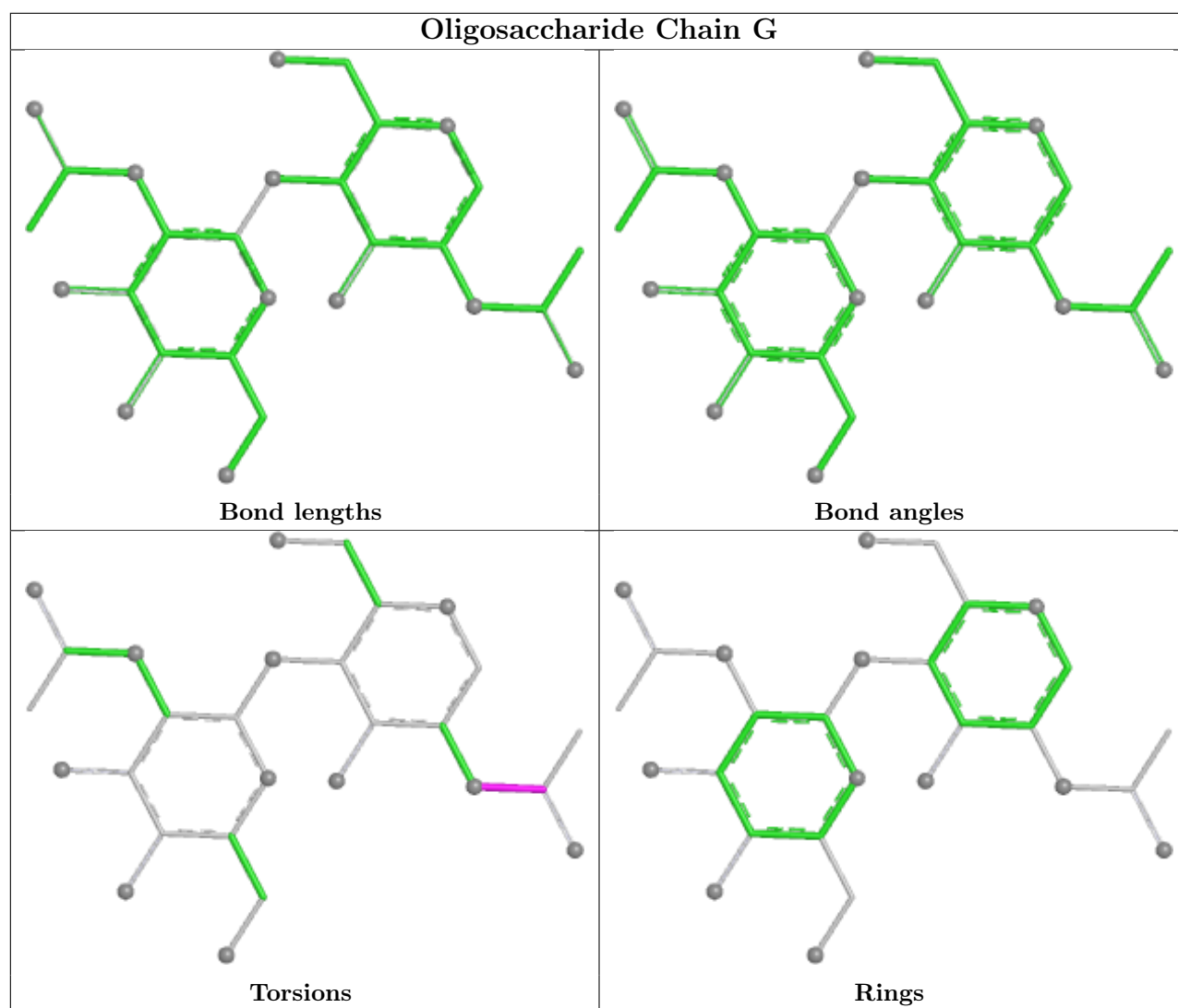
All (4) torsion outliers are listed below:

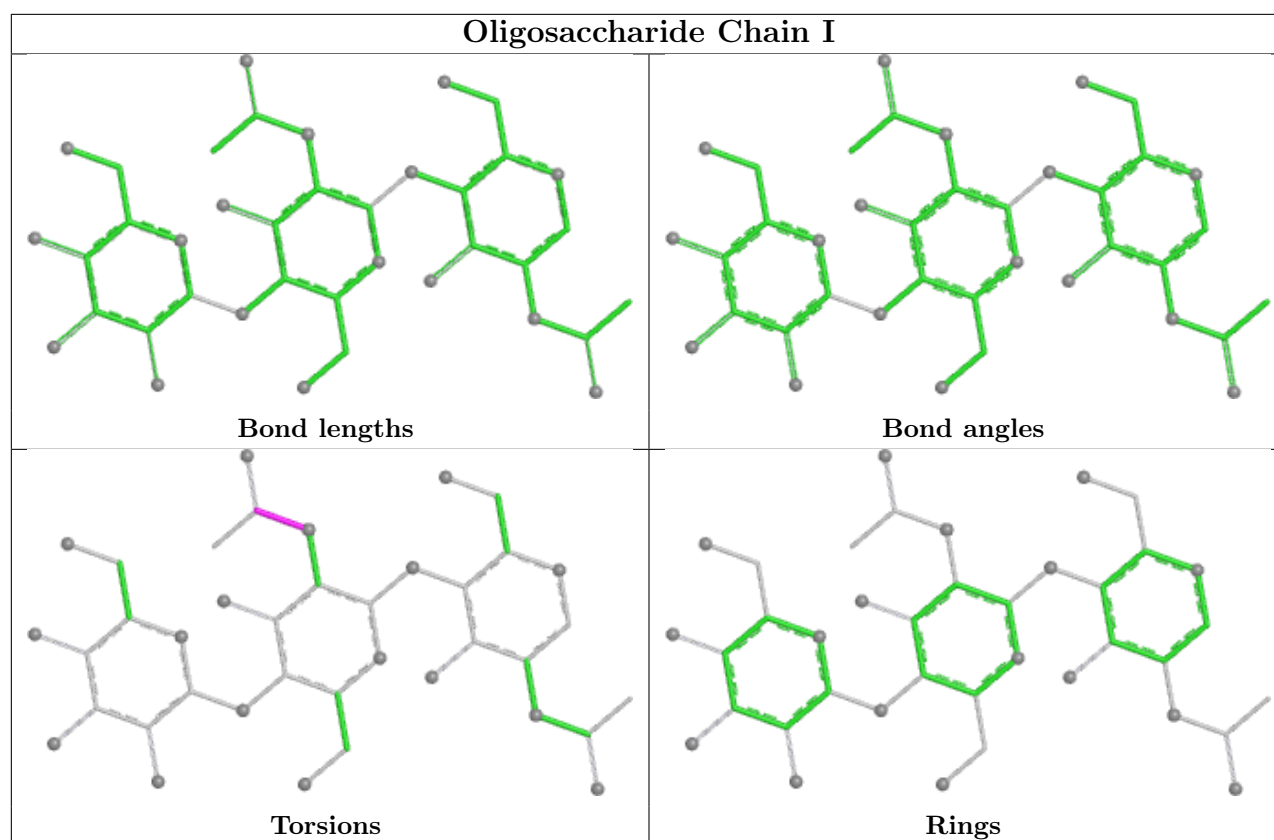
Mol	Chain	Res	Type	Atoms
5	I	2	NAG	O7-C7-N2-C2
5	I	2	NAG	C8-C7-N2-C2
4	G	1	NAG	C8-C7-N2-C2
4	G	1	NAG	O7-C7-N2-C2

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.





5.6 Ligand geometry [i](#)

1 ligand is 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	$\# Z > 2$
6	NAG	D	601	1	14,14,15	0.87	1 (7%)	17,19,21	2.23	5 (29%)

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
6	NAG	D	601	1	-	1/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	D	601	NAG	C1-C2	2.50	1.55	1.52

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	D	601	NAG	C1-O5-C5	5.40	119.42	112.19
6	D	601	NAG	O5-C5-C6	3.57	114.61	107.66
6	D	601	NAG	C4-C3-C2	-3.35	106.11	111.02
6	D	601	NAG	C2-N2-C7	2.97	126.88	122.90
6	D	601	NAG	O7-C7-N2	2.23	125.93	121.98

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	D	601	NAG	C3-C2-N2-C7

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	D	601	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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	523/529 (98%)	-0.44	0 100 100	91, 158, 211, 284	1 (0%)
1	B	523/529 (98%)	-0.39	2 (0%) 88 72	91, 159, 213, 273	0
1	C	523/529 (98%)	-0.36	0 100 100	150, 218, 269, 445	0
1	D	523/529 (98%)	-0.36	4 (0%) 82 60	123, 213, 265, 425	3 (0%)
2	E	219/233 (93%)	-0.37	1 (0%) 87 69	108, 191, 263, 362	1 (0%)
2	H	220/233 (94%)	-0.35	2 (0%) 81 58	102, 160, 233, 333	0
2	K	220/233 (94%)	-0.27	1 (0%) 87 69	181, 241, 311, 489	0
2	P	119/233 (51%)	-0.31	0 100 100	224, 270, 327, 353	0
3	F	212/216 (98%)	-0.32	3 (1%) 73 48	114, 171, 233, 265	1 (0%)
3	J	212/216 (98%)	-0.21	1 (0%) 87 69	174, 238, 296, 321	0
3	L	212/216 (98%)	-0.41	1 (0%) 87 69	96, 160, 212, 251	0
3	O	110/216 (50%)	-0.33	0 100 100	192, 243, 293, 343	0
All	All	3616/3912 (92%)	-0.36	15 (0%) 88 72	91, 194, 275, 489	6 (0%)

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	104	ASP	3.5
3	J	175	LYS	3.4
3	L	133	LYS	3.4
3	F	130	GLN	2.9
2	K	104	ASP	2.8
1	D	392	GLY	2.7
1	D	184	LEU	2.7
3	F	51	TYR	2.5
3	F	106	LYS	2.5
1	B	378	GLU	2.3
1	D	129	GLU	2.2

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Mol	Chain	Res	Type	RSRZ
2	H	102	TRP	2.1
1	B	494	LYS	2.1
1	D	494	LYS	2.0
2	E	101	ASN	2.0

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

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

6.3 Carbohydrates [i](#)

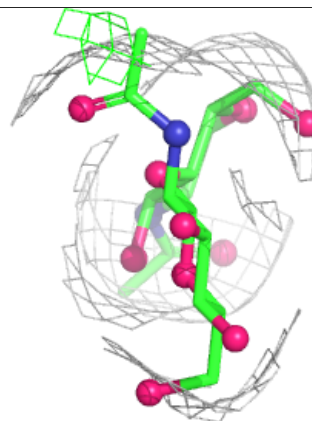
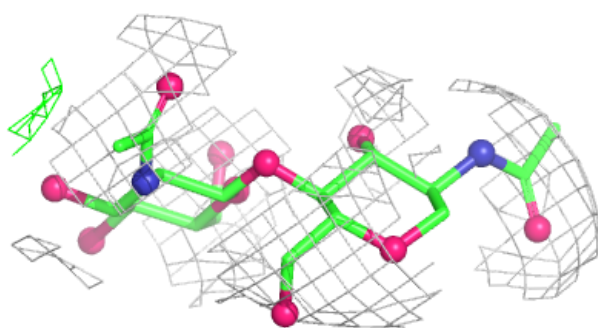
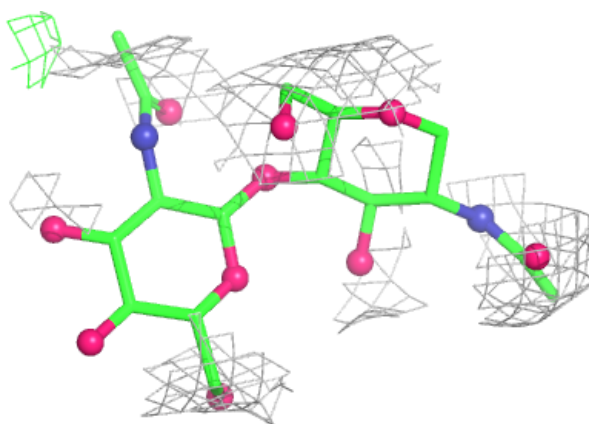
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	NAG	G	1	14/15	-	-	157,195,224,241	0
4	NAG	G	2	14/15	-	-	209,248,270,287	0
5	NAG	I	1	14/15	-	-	172,206,232,244	0
5	NAG	I	2	14/15	-	-	198,224,271,328	0
5	BMA	I	3	11/12	-	-	202,268,299,306	0

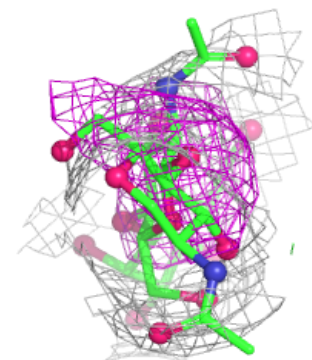
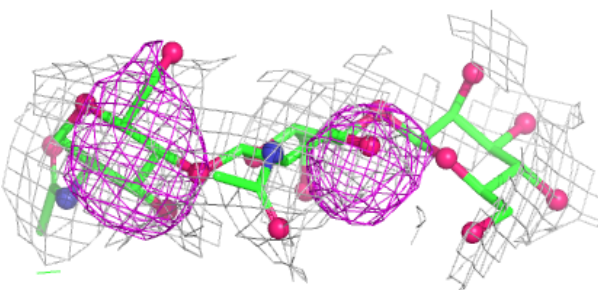
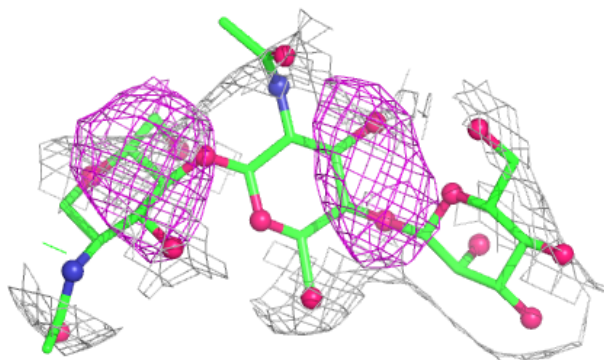
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain G:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around Chain I:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



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($\text{\AA}^2$)	Q<0.9
6	NAG	D	601	14/15	0.77	0.08	155,229,256,274	0

6.5 Other polymers [i](#)

There are no such residues in this entry.