



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 08:01 PM UTC

PDB ID : 2WRE / pdb_00002wre
Title : structure of H2 japan hemagglutinin with human receptor
Authors : Liu, J.; Stevens, D.J.; Haire, L.F.; Walker, P.A.; Coombs, P.J.; Russell, R.J.;
Gamblin, S.J.; Skehel, J.J.
Deposited on : 2009-09-01
Resolution : 3.00 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

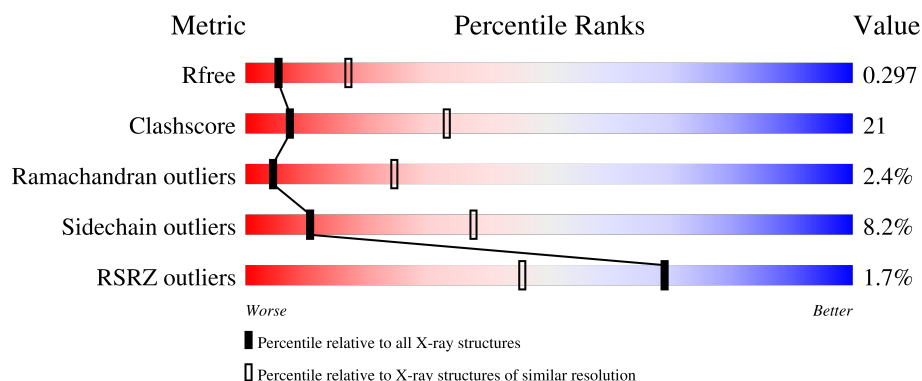
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	505	2% 52% 37% 7%
1	B	505	2% 58% 34%
1	C	505	2% 55% 35% 5%
2	D	2	50% 50%
2	E	2	50% 50%

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Mol	Chain	Length	Quality of chain
2	F	2	 50%50%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 11443 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.

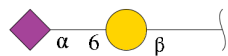
- Molecule 1 is a protein called HEMAGGLUTININ.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	471	Total	C	N	O	S	0	0	0
			3708	2335	633	716	24			
1	B	486	Total	C	N	O	S	0	0	0
			3841	2417	658	742	24			
1	C	481	Total	C	N	O	S	0	0	0
			3798	2392	652	730	24			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	107	HIS	TYR	conflict	UNP Q67085
A	136	GLN	ARG	conflict	UNP Q67085
A	159	ASN	ASP	conflict	UNP Q67085
A	185	ILE	ASN	conflict	UNP Q67085
A	188	LYS	THR	conflict	UNP Q67085
A	307	LYS	ARG	conflict	UNP Q67085
B	107	HIS	TYR	conflict	UNP Q67085
B	136	GLN	ARG	conflict	UNP Q67085
B	159	ASN	ASP	conflict	UNP Q67085
B	185	ILE	ASN	conflict	UNP Q67085
B	188	LYS	THR	conflict	UNP Q67085
B	307	LYS	ARG	conflict	UNP Q67085
C	107	HIS	TYR	conflict	UNP Q67085
C	136	GLN	ARG	conflict	UNP Q67085
C	159	ASN	ASP	conflict	UNP Q67085
C	185	ILE	ASN	conflict	UNP Q67085
C	188	LYS	THR	conflict	UNP Q67085
C	307	LYS	ARG	conflict	UNP Q67085

- Molecule 2 is an oligosaccharide called N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose.

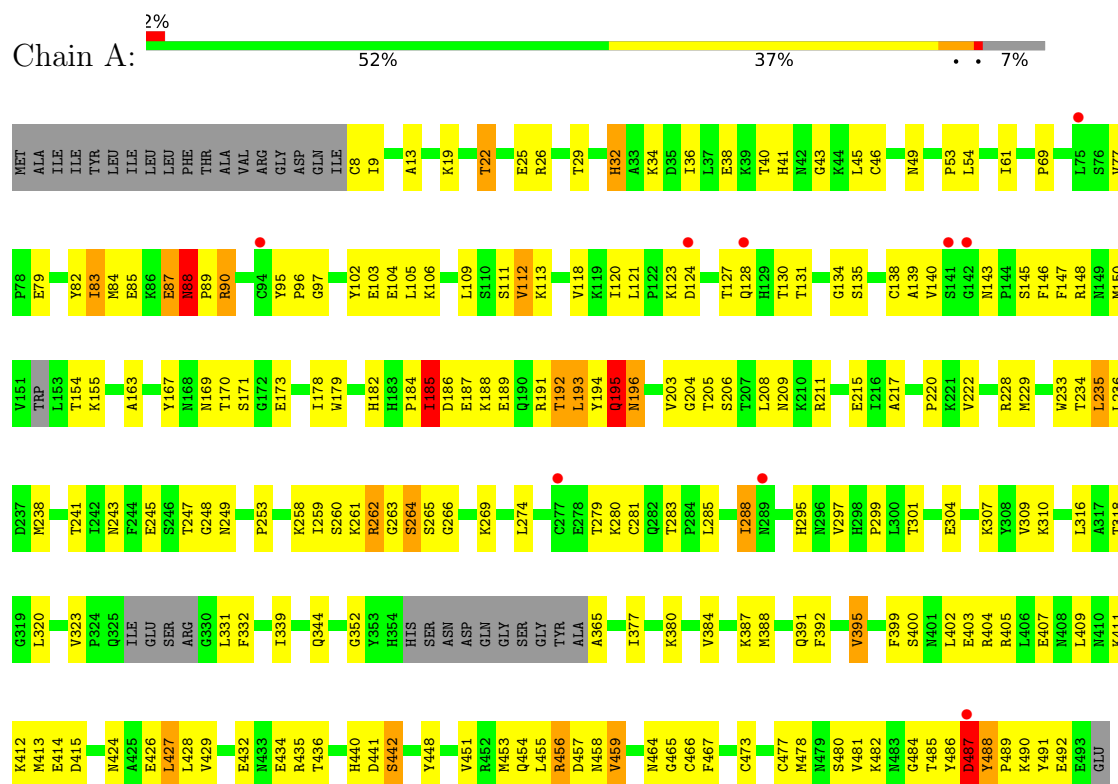


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	D	2	Total	C	N	O	0	0	0
			32	17	1	14			
2	E	2	Total	C	N	O	0	0	0
			32	17	1	14			
2	F	2	Total	C	N	O	0	0	0
			32	17	1	14			

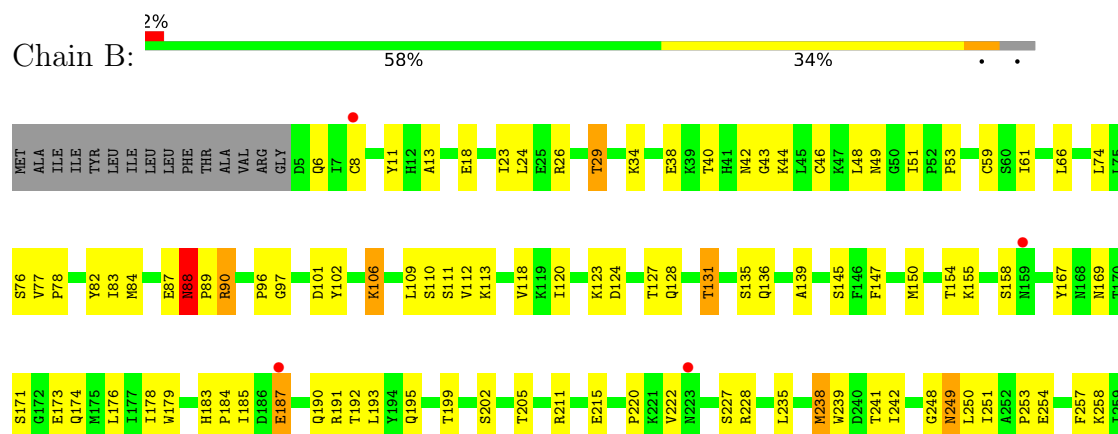
3 Residue-property plots

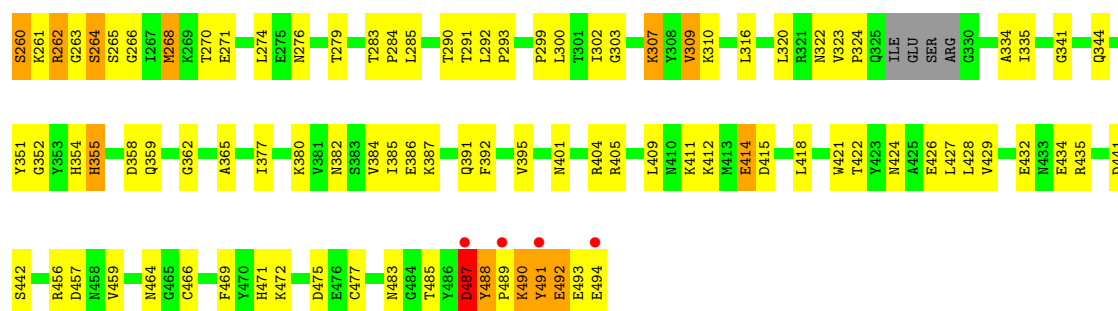
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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

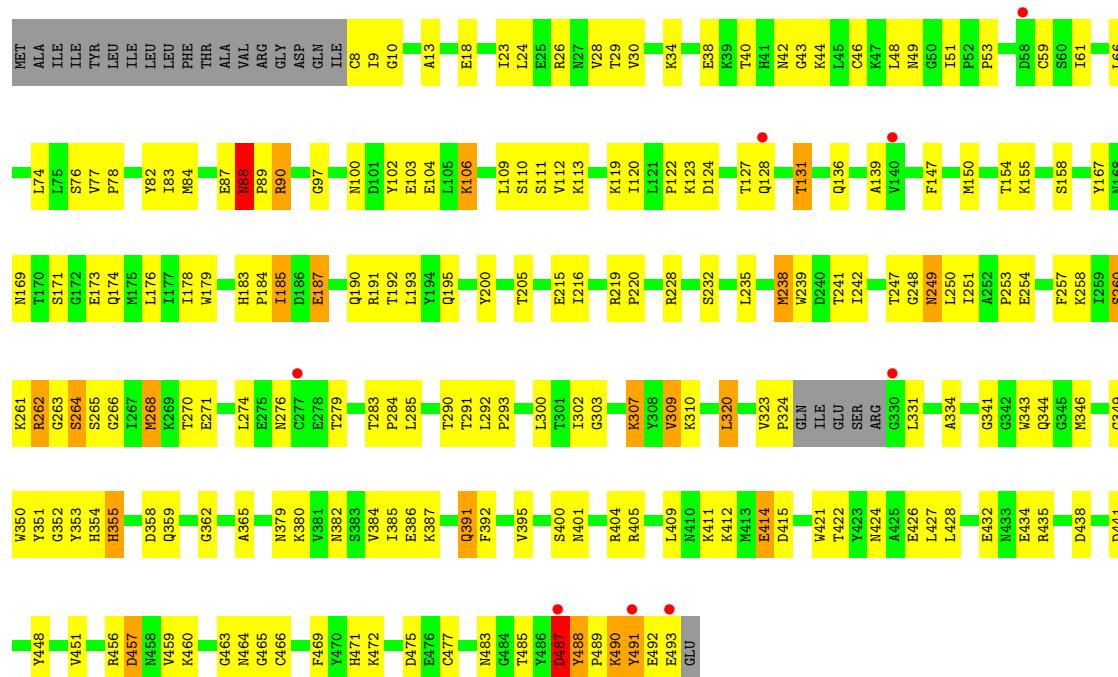


• Molecule 1: HEMAGGLUTININ





• Molecule 1: HEMAGGLUTININ



• Molecule 2: N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose



• Molecule 2: N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose



• Molecule 2: N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	75.97Å 117.46Å 225.37Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.98 – 3.00 19.98 – 3.00	Depositor EDS
% Data completeness (in resolution range)	96.8 (19.98-3.00) 96.3 (19.98-3.00)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.16 (at 3.01Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.239 , 0.295 0.239 , 0.297	Depositor DCC
R_{free} test set	2006 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	68.9	Xtriage
Anisotropy	0.014	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 52.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	11443	wwPDB-VP
Average B, all atoms (Å ²)	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.70% 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: GAL, SIA

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	0.53	0/3788	0.96	6/5121 (0.1%)
1	B	0.55	0/3928	0.92	5/5313 (0.1%)
1	C	0.55	0/3885	0.92	5/5255 (0.1%)
All	All	0.54	0/11601	0.93	16/15689 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1
1	C	0	1
All	All	0	2

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	88	ASN	CA-C-N	9.97	130.25	119.28
1	A	88	ASN	C-N-CA	9.97	130.25	119.28
1	C	292	LEU	CA-C-N	6.46	126.04	119.19
1	C	292	LEU	C-N-CA	6.46	126.04	119.19
1	A	185	ILE	CB-CA-C	-6.19	104.10	111.65
1	B	88	ASN	CA-C-N	6.05	125.60	119.19
1	B	88	ASN	C-N-CA	6.05	125.60	119.19
1	C	88	ASN	CA-C-N	5.96	125.51	119.19
1	C	88	ASN	C-N-CA	5.96	125.51	119.19
1	C	341	GLY	N-CA-C	5.85	119.64	111.10
1	B	341	GLY	N-CA-C	5.76	119.50	111.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	112	VAL	N-CA-C	5.53	115.84	108.27
1	B	292	LEU	CA-C-N	5.49	124.84	118.85
1	B	292	LEU	C-N-CA	5.49	124.84	118.85
1	A	297	VAL	N-CA-C	5.35	115.56	110.53
1	A	384	VAL	CB-CA-C	-5.16	105.28	111.88

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	490	LYS	Peptide
1	C	490	LYS	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	3708	0	3583	177	0
1	B	3841	0	3704	161	0
1	C	3798	0	3669	165	0
2	D	32	0	28	7	0
2	E	32	0	28	0	0
2	F	32	0	28	1	0
All	All	11443	0	11040	464	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 21.

All (464) 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:53:PRO:HG2	1:B:84:MET:HE3	1.40	1.03
1:B:487:ASP:HA	1:B:488:TYR:HB2	1.43	0.99
1:C:487:ASP:HA	1:C:488:TYR:HB2	1.46	0.95
1:C:492:GLU:HA	1:C:493:GLU:C	1.91	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:53:PRO:HG2	1:C:84:MET:HE3	1.45	0.95
1:C:488:TYR:CD2	1:C:488:TYR:O	2.23	0.91
1:A:140:VAL:HG21	1:A:148:ARG:HH22	1.34	0.91
1:B:488:TYR:O	1:B:488:TYR:CD2	2.25	0.89
1:C:9:ILE:HD11	1:C:451:VAL:HG21	1.53	0.89
1:C:261:LYS:HA	1:C:262:ARG:HB2	1.55	0.89
1:A:435:ARG:NH1	1:C:435:ARG:HH11	1.71	0.88
1:B:261:LYS:HA	1:B:262:ARG:HB2	1.56	0.88
1:A:487:ASP:HA	1:A:488:TYR:HB2	1.58	0.86
1:C:488:TYR:N	1:C:489:PRO:HD3	1.89	0.86
1:B:488:TYR:N	1:B:489:PRO:HD3	1.91	0.85
1:A:69:PRO:HB3	1:A:140:VAL:HG22	1.59	0.85
1:B:83:ILE:HD12	1:B:110:SER:HA	1.60	0.83
1:A:488:TYR:N	1:A:489:PRO:HD3	1.92	0.83
1:C:173:GLU:OE1	1:C:260:SER:HB2	1.81	0.81
1:C:488:TYR:H	1:C:489:PRO:HD3	1.42	0.81
1:B:434:GLU:CD	1:C:435:ARG:HH21	1.89	0.80
1:C:83:ILE:HD12	1:C:110:SER:HA	1.61	0.80
1:B:173:GLU:OE1	1:B:260:SER:HB2	1.82	0.80
1:A:261:LYS:HB2	1:A:262:ARG:CB	2.12	0.79
1:B:303:GLY:O	1:B:392:PHE:HA	1.84	0.78
1:B:490:LYS:C	1:B:491:TYR:CD2	2.61	0.78
1:A:435:ARG:HH21	1:C:434:GLU:CD	1.90	0.78
1:B:359:GLN:HE22	1:B:475:ASP:H	1.32	0.77
1:C:483:ASN:HB3	1:C:485:THR:OG1	1.83	0.77
1:B:487:ASP:HA	1:B:488:TYR:CB	2.15	0.77
1:A:61:ILE:HD12	1:A:106:LYS:HD2	1.65	0.77
1:B:261:LYS:HB2	1:B:262:ARG:HB3	1.66	0.77
1:C:261:LYS:HB2	1:C:262:ARG:HB3	1.65	0.76
1:B:8:CYS:HB2	1:B:354:HIS:HB3	1.68	0.76
1:A:448:TYR:CE1	1:A:465:GLY:HA2	2.21	0.75
1:B:488:TYR:H	1:B:489:PRO:HD3	1.47	0.75
1:B:483:ASN:HB3	1:B:485:THR:OG1	1.85	0.75
1:B:459:VAL:HG12	1:B:469:PHE:HA	1.68	0.74
1:C:359:GLN:HE22	1:C:475:ASP:H	1.35	0.74
1:C:155:LYS:NZ	1:C:195:GLN:HG2	2.03	0.73
1:B:307:LYS:HG2	1:B:421:TRP:CE2	2.23	0.73
1:A:53:PRO:HD2	1:A:274:LEU:HD22	1.69	0.73
1:A:179:TRP:HZ3	1:A:234:THR:HG22	1.52	0.73
1:C:459:VAL:HG12	1:C:469:PHE:HA	1.69	0.73
1:A:424:ASN:HD21	1:C:424:ASN:HD21	1.36	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:487:ASP:HA	1:C:488:TYR:CB	2.17	0.72
1:A:189:GLU:OE2	2:D:2:SIA:H92	1.89	0.72
1:C:488:TYR:O	1:C:488:TYR:HD2	1.73	0.72
1:A:217:ALA:HB2	1:B:202:SER:OG	1.90	0.71
1:B:155:LYS:NZ	1:B:195:GLN:HG2	2.05	0.71
1:A:123:LYS:HG2	1:A:131:THR:HG21	1.72	0.71
1:A:189:GLU:OE2	2:D:2:SIA:C9	2.39	0.71
1:C:111:SER:HB2	1:C:265:SER:HB3	1.71	0.71
1:C:488:TYR:H	1:C:489:PRO:CD	2.04	0.71
1:B:111:SER:HB2	1:B:265:SER:HB3	1.72	0.71
1:A:380:LYS:HE2	1:A:432:GLU:OE2	1.89	0.70
1:A:127:THR:O	1:A:128:GLN:HB2	1.91	0.70
1:A:487:ASP:OD1	1:A:487:ASP:C	2.34	0.70
1:B:489:PRO:C	1:B:491:TYR:CZ	2.69	0.70
1:C:489:PRO:C	1:C:491:TYR:CZ	2.70	0.70
1:C:307:LYS:HD2	1:C:307:LYS:N	2.06	0.70
1:B:489:PRO:HA	1:B:491:TYR:OH	1.93	0.69
1:A:283:THR:HG22	1:A:301:THR:HG22	1.73	0.69
1:C:464:ASN:OD1	1:C:466:CYS:HB2	1.93	0.69
1:A:435:ARG:HH12	1:C:435:ARG:HH11	1.41	0.68
1:A:411:LYS:HE2	1:A:415:ASP:OD1	1.94	0.68
1:B:489:PRO:CA	1:B:491:TYR:OH	2.41	0.68
1:A:299:PRO:O	1:A:395:VAL:HG11	1.94	0.68
1:C:490:LYS:C	1:C:491:TYR:CD2	2.72	0.68
1:B:488:TYR:O	1:B:488:TYR:HD2	1.77	0.68
1:C:187:GLU:O	1:C:191:ARG:HG3	1.93	0.68
1:A:173:GLU:OE1	1:A:260:SER:HB2	1.94	0.68
1:C:123:LYS:HG2	1:C:131:THR:HG21	1.76	0.68
1:A:487:ASP:HA	1:A:488:TYR:CB	2.24	0.68
1:C:489:PRO:C	1:C:490:LYS:HG2	2.18	0.67
1:A:488:TYR:N	1:A:489:PRO:CD	2.58	0.67
1:C:261:LYS:HA	1:C:262:ARG:CB	2.24	0.67
1:A:43:GLY:HA2	1:A:285:LEU:O	1.94	0.67
1:B:488:TYR:H	1:B:489:PRO:CD	2.07	0.66
1:A:154:THR:HG21	2:D:2:SIA:H111	1.75	0.66
1:A:146:PHE:HE2	1:A:150:MET:HB2	1.60	0.66
1:B:471:HIS:ND1	1:B:490:LYS:HD2	2.10	0.66
1:B:187:GLU:O	1:B:191:ARG:HG3	1.96	0.66
1:B:261:LYS:HA	1:B:262:ARG:CB	2.25	0.65
1:B:53:PRO:CG	1:B:84:MET:HE3	2.24	0.64
1:C:8:CYS:O	1:C:353:TYR:HA	1.96	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:155:LYS:HZ3	1:C:195:GLN:HG2	1.63	0.64
1:B:123:LYS:HG2	1:B:131:THR:HG21	1.79	0.64
1:A:208:LEU:HG	1:A:209:ASN:N	2.12	0.64
1:B:53:PRO:HD2	1:B:274:LEU:HD22	1.80	0.64
1:C:53:PRO:HD2	1:C:274:LEU:HD22	1.80	0.64
1:B:471:HIS:CE1	1:B:490:LYS:HD2	2.33	0.64
1:A:454:GLN:OE1	1:A:484:GLY:HA2	1.98	0.63
1:C:184:PRO:HG2	1:C:190:GLN:HE21	1.62	0.63
1:A:167:TYR:CE2	1:A:169:ASN:HA	2.33	0.63
1:A:169:ASN:ND2	1:A:238:MET:HA	2.14	0.63
1:C:489:PRO:HD2	1:C:490:LYS:HE2	1.80	0.63
1:A:455:LEU:HD13	1:A:459:VAL:HG11	1.81	0.62
1:A:189:GLU:O	1:A:193:LEU:HB2	1.99	0.62
1:B:307:LYS:HD2	1:B:307:LYS:N	2.14	0.62
1:B:489:PRO:C	1:B:490:LYS:HG2	2.24	0.62
1:B:490:LYS:O	1:B:491:TYR:CD2	2.51	0.62
1:A:488:TYR:H	1:A:489:PRO:CD	2.11	0.62
1:B:487:ASP:C	1:B:487:ASP:OD1	2.42	0.62
1:B:489:PRO:HD2	1:B:490:LYS:HE2	1.81	0.62
1:A:120:ILE:HG22	1:A:167:TYR:CE1	2.34	0.62
1:B:66:LEU:O	1:B:147:PHE:HB3	2.00	0.62
1:C:334:ALA:HB3	1:C:441:ASP:OD2	2.00	0.62
1:A:54:LEU:HB3	1:A:83:ILE:HG12	1.81	0.61
1:A:388:MET:O	1:A:388:MET:HG3	1.99	0.61
1:A:171:SER:HB2	1:A:258:LYS:HD2	1.82	0.61
1:A:352:GLY:HA3	1:A:365:ALA:HA	1.81	0.61
1:C:265:SER:OG	1:C:266:GLY:N	2.34	0.61
1:A:323:VAL:HG12	1:A:323:VAL:O	2.01	0.60
1:B:261:LYS:CA	1:B:262:ARG:CB	2.80	0.60
1:C:176:LEU:HB3	1:C:257:PHE:HB2	1.84	0.60
1:A:169:ASN:HB2	1:A:236:LEU:HD23	1.82	0.60
1:C:261:LYS:CA	1:C:262:ARG:CB	2.79	0.60
1:C:487:ASP:OD1	1:C:487:ASP:C	2.45	0.60
1:C:8:CYS:HB2	1:C:354:HIS:HB3	1.84	0.60
1:C:112:VAL:N	1:C:262:ARG:HD3	2.16	0.60
1:C:489:PRO:CA	1:C:491:TYR:OH	2.49	0.60
1:B:265:SER:OG	1:B:266:GLY:N	2.34	0.59
1:B:155:LYS:HZ3	1:B:195:GLN:HG2	1.65	0.59
1:B:488:TYR:N	1:B:489:PRO:CD	2.64	0.59
1:C:53:PRO:HG3	1:C:82:TYR:CZ	2.38	0.59
1:A:9:ILE:HD11	1:A:451:VAL:HG21	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:323:VAL:N	1:B:324:PRO:HD3	2.18	0.58
1:B:489:PRO:CB	1:B:491:TYR:OH	2.51	0.58
1:C:489:PRO:HA	1:C:491:TYR:OH	2.03	0.58
1:A:205:THR:OG1	1:A:208:LEU:HB3	2.03	0.58
1:A:488:TYR:H	1:A:489:PRO:HD3	1.65	0.58
1:A:194:TYR:O	1:A:196:ASN:N	2.30	0.58
1:B:89:PRO:O	1:B:90:ARG:HB3	2.04	0.58
1:B:173:GLU:OE2	1:B:258:LYS:HE2	2.04	0.58
1:A:435:ARG:NH2	1:C:434:GLU:CD	2.60	0.57
1:C:471:HIS:ND1	1:C:490:LYS:HD2	2.19	0.57
1:B:184:PRO:HG2	1:B:190:GLN:HE21	1.68	0.57
1:C:173:GLU:OE2	1:C:258:LYS:HE2	2.03	0.57
1:C:184:PRO:HG2	1:C:190:GLN:NE2	2.20	0.57
1:A:97:GLY:HA2	1:A:228:ARG:HB3	1.85	0.57
1:A:211:ARG:NH1	1:C:216:ILE:O	2.37	0.57
1:A:95:TYR:CE1	1:A:229:MET:HE3	2.39	0.57
1:B:43:GLY:HA2	1:B:285:LEU:O	2.05	0.57
1:C:13:ALA:O	1:C:344:GLN:HA	2.05	0.57
1:A:102:TYR:CZ	1:A:106:LYS:HD3	2.39	0.57
1:C:127:THR:O	1:C:128:GLN:HB2	2.05	0.57
1:B:405:ARG:HH21	1:C:104:GLU:HG2	1.69	0.57
1:C:155:LYS:HE3	1:C:192:THR:O	2.04	0.57
1:B:171:SER:HB2	1:B:258:LYS:HD2	1.87	0.57
1:A:111:SER:HB2	1:A:265:SER:HB3	1.86	0.56
1:A:265:SER:OG	1:A:266:GLY:N	2.37	0.56
1:A:400:SER:OG	1:A:403:GLU:HG3	2.04	0.56
1:C:23:ILE:HG22	1:C:24:LEU:HD23	1.86	0.56
1:A:89:PRO:O	1:A:90:ARG:HB3	2.06	0.56
1:B:248:GLY:O	1:B:249:ASN:HB2	2.05	0.56
1:B:334:ALA:HB3	1:B:441:ASP:OD2	2.06	0.56
1:C:53:PRO:CG	1:C:84:MET:HE3	2.28	0.56
1:C:261:LYS:CA	1:C:262:ARG:HB2	2.33	0.56
1:A:248:GLY:C	1:A:249:ASN:HD22	2.13	0.56
1:A:186:ASP:HB3	1:A:189:GLU:H	1.69	0.56
1:A:424:ASN:ND2	1:C:424:ASN:HD21	2.04	0.56
1:B:23:ILE:HG22	1:B:24:LEU:HD23	1.88	0.56
1:A:178:ILE:O	1:A:253:PRO:HG3	2.06	0.56
1:B:112:VAL:N	1:B:262:ARG:HD3	2.21	0.55
1:B:490:LYS:C	1:B:491:TYR:CG	2.84	0.55
1:A:261:LYS:CB	1:A:262:ARG:CB	2.85	0.55
1:A:456:ARG:HH12	1:C:460:LYS:HE3	1.70	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:51:ILE:HD12	1:C:262:ARG:HH22	1.71	0.55
1:A:13:ALA:O	1:A:344:GLN:HA	2.05	0.55
1:A:464:ASN:OD1	1:A:466:CYS:HB2	2.06	0.55
1:A:22:THR:HB	1:A:434:GLU:HB3	1.89	0.55
1:A:179:TRP:CE2	1:A:203:VAL:HG21	2.42	0.55
1:A:488:TYR:O	1:A:488:TYR:CD2	2.59	0.55
1:B:493:GLU:H	1:B:494:GLU:HA	1.72	0.55
1:A:46:CYS:HB2	1:A:279:THR:HG22	1.89	0.55
1:B:176:LEU:HB3	1:B:257:PHE:HB2	1.88	0.55
1:C:89:PRO:O	1:C:90:ARG:HB3	2.05	0.55
1:A:307:LYS:HD3	1:A:388:MET:SD	2.46	0.54
1:B:53:PRO:HG3	1:B:82:TYR:CZ	2.42	0.54
1:B:464:ASN:OD1	1:B:466:CYS:HB2	2.06	0.54
1:B:155:LYS:HE3	1:B:192:THR:O	2.07	0.54
1:A:123:LYS:CG	1:A:131:THR:HG21	2.37	0.54
1:B:42:ASN:O	1:B:44:LYS:HG3	2.08	0.54
1:C:171:SER:HB2	1:C:258:LYS:HD2	1.88	0.54
1:B:359:GLN:NE2	1:B:475:ASP:H	2.02	0.54
1:B:489:PRO:CA	1:B:491:TYR:CZ	2.90	0.54
1:B:491:TYR:HB2	1:B:492:GLU:OE2	2.07	0.54
1:A:473:CYS:SG	1:A:478:MET:HE2	2.47	0.54
1:A:387:LYS:HD2	1:C:426:GLU:HB3	1.88	0.54
1:B:268:MET:HE3	1:B:302:ILE:HD12	1.89	0.54
1:B:113:LYS:N	1:B:262:ARG:HD2	2.22	0.54
1:B:184:PRO:HG2	1:B:190:GLN:NE2	2.21	0.54
1:A:377:ILE:HD11	1:A:436:THR:HG23	1.89	0.53
1:B:154:THR:HG21	1:B:193:LEU:HD22	1.90	0.53
1:A:89:PRO:O	1:A:90:ARG:CB	2.55	0.53
1:C:42:ASN:O	1:C:44:LYS:HG3	2.09	0.53
1:C:183:HIS:CE1	1:C:215:GLU:H	2.26	0.53
1:A:188:LYS:O	1:A:192:THR:HB	2.08	0.53
1:C:489:PRO:CA	1:C:491:TYR:CZ	2.91	0.53
1:C:489:PRO:CB	1:C:491:TYR:OH	2.55	0.53
1:A:53:PRO:HG2	1:A:84:MET:HE3	1.91	0.53
1:C:150:MET:HE1	1:C:179:TRP:HA	1.90	0.53
1:C:89:PRO:O	1:C:90:ARG:CB	2.56	0.53
1:B:89:PRO:O	1:B:90:ARG:CB	2.56	0.53
1:A:261:LYS:CA	1:A:262:ARG:CB	2.87	0.52
1:C:307:LYS:HG2	1:C:421:TRP:CE2	2.44	0.52
1:A:140:VAL:HG21	1:A:148:ARG:NH2	2.14	0.52
1:A:458:ASN:ND2	1:A:490:LYS:CB	2.73	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:478:MET:O	1:A:482:LYS:HG3	2.10	0.52
1:C:66:LEU:O	1:C:147:PHE:HB3	2.09	0.52
1:A:102:TYR:CE2	1:A:106:LYS:HD3	2.43	0.52
1:B:250:LEU:HD12	1:B:251:ILE:N	2.24	0.52
1:B:493:GLU:N	1:B:494:GLU:HA	2.25	0.52
1:B:382:ASN:O	1:B:386:GLU:HG2	2.09	0.52
1:C:276:ASN:CG	1:C:276:ASN:O	2.51	0.52
1:A:79:GLU:HG3	1:A:112:VAL:O	2.10	0.52
1:B:127:THR:O	1:B:128:GLN:HB2	2.09	0.52
1:A:192:THR:HG23	1:A:192:THR:O	2.10	0.51
1:A:455:LEU:C	1:A:456:ARG:HG3	2.35	0.51
1:A:146:PHE:CG	1:A:147:PHE:N	2.78	0.51
1:A:424:ASN:O	1:A:428:LEU:HB2	2.09	0.51
1:A:105:LEU:HB2	1:A:233:TRP:CE2	2.45	0.51
1:A:134:GLY:O	2:D:2:SIA:H4	2.10	0.51
1:A:146:PHE:CD2	1:A:147:PHE:N	2.78	0.51
1:B:51:ILE:HD12	1:B:262:ARG:HH22	1.75	0.51
1:B:380:LYS:HE2	1:B:432:GLU:OE2	2.11	0.51
1:A:424:ASN:HD21	1:B:424:ASN:HD21	1.57	0.51
1:A:435:ARG:NH1	1:C:435:ARG:HD2	2.25	0.51
1:A:485:THR:O	1:A:485:THR:HG22	2.11	0.51
1:A:426:GLU:CG	1:B:387:LYS:HD2	2.39	0.51
1:B:150:MET:HE1	1:B:179:TRP:HA	1.93	0.51
1:C:87:GLU:O	1:C:88:ASN:HB2	2.11	0.51
1:A:53:PRO:HB3	1:A:82:TYR:CE1	2.45	0.51
1:A:459:VAL:HG21	1:A:467:PHE:CD2	2.46	0.51
1:C:250:LEU:HD12	1:C:251:ILE:N	2.25	0.51
1:C:43:GLY:HA2	1:C:285:LEU:O	2.10	0.50
1:C:113:LYS:N	1:C:262:ARG:HD2	2.26	0.50
1:C:293:PRO:HG3	1:C:385:ILE:HA	1.92	0.50
1:C:359:GLN:NE2	1:C:475:ASP:H	2.05	0.50
1:C:471:HIS:CE1	1:C:490:LYS:HD2	2.46	0.50
1:A:130:THR:HB	1:A:154:THR:OG1	2.12	0.50
1:A:196:ASN:N	1:A:196:ASN:ND2	2.59	0.50
1:B:205:THR:HG22	1:B:242:ILE:HG13	1.94	0.50
1:A:104:GLU:HG2	1:C:405:ARG:HH21	1.75	0.50
1:A:120:ILE:HG13	1:A:121:LEU:N	2.26	0.50
1:B:238:MET:HB3	1:B:239:TRP:CD1	2.46	0.50
1:A:434:GLU:CD	1:B:435:ARG:HH21	2.20	0.50
1:C:392:PHE:C	1:C:392:PHE:CD2	2.89	0.50
1:A:288:ILE:HG22	1:A:288:ILE:O	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:276:ASN:CG	1:B:276:ASN:O	2.54	0.50
1:C:380:LYS:HE2	1:C:432:GLU:OE2	2.13	0.49
1:B:392:PHE:C	1:B:392:PHE:CD2	2.89	0.49
1:A:69:PRO:HB3	1:A:140:VAL:CG2	2.36	0.49
1:B:174:GLN:OE1	1:B:235:LEU:HD13	2.11	0.49
1:B:183:HIS:CE1	1:B:215:GLU:H	2.30	0.49
1:B:316:LEU:HD13	1:B:429:VAL:HG22	1.95	0.49
1:A:211:ARG:HB3	1:C:215:GLU:HG2	1.95	0.49
1:B:489:PRO:HB3	1:B:491:TYR:OH	2.12	0.49
1:C:205:THR:HG22	1:C:242:ILE:HG13	1.94	0.49
1:B:471:HIS:CG	1:B:490:LYS:HD2	2.48	0.49
1:A:87:GLU:O	1:A:88:ASN:HB2	2.13	0.48
1:C:248:GLY:C	1:C:249:ASN:HD22	2.21	0.48
1:B:136:GLN:O	1:B:139:ALA:HB2	2.13	0.48
1:C:136:GLN:O	1:C:139:ALA:HB2	2.12	0.48
1:C:103:GLU:OE1	1:C:400:SER:HB3	2.13	0.48
1:A:392:PHE:C	1:A:392:PHE:CD2	2.91	0.48
1:A:481:VAL:HG22	1:A:486:TYR:HB2	1.94	0.48
1:C:284:PRO:HD3	1:C:300:LEU:O	2.13	0.48
1:B:120:ILE:HG22	1:B:167:TYR:CE1	2.48	0.48
1:C:102:TYR:CZ	1:C:106:LYS:HD3	2.47	0.48
1:A:41:HIS:H	1:A:41:HIS:CD2	2.30	0.48
1:A:120:ILE:HG13	1:A:121:LEU:H	1.78	0.48
1:A:316:LEU:HD13	1:A:429:VAL:HG22	1.96	0.48
1:B:299:PRO:O	1:B:395:VAL:HG11	2.14	0.48
1:A:491:TYR:C	1:A:492:GLU:OE2	2.57	0.48
1:C:123:LYS:HB2	1:C:254:GLU:CD	2.38	0.48
1:A:188:LYS:H	1:A:188:LYS:HD2	1.79	0.48
1:A:243:ASN:HD22	1:C:219:ARG:HA	1.79	0.48
1:A:487:ASP:OD1	1:A:487:ASP:O	2.32	0.47
1:B:316:LEU:CD1	1:B:429:VAL:HG22	2.44	0.47
1:C:270:THR:OG1	1:C:271:GLU:N	2.47	0.47
1:A:95:TYR:O	1:A:96:PRO:C	2.57	0.47
1:A:448:TYR:HE1	1:A:465:GLY:HA2	1.75	0.47
1:A:163:ALA:O	1:A:245:GLU:HA	2.13	0.47
1:B:13:ALA:O	1:B:344:GLN:HA	2.15	0.47
1:A:440:HIS:O	1:A:441:ASP:C	2.57	0.47
1:A:261:LYS:HA	1:A:262:ARG:CB	2.45	0.47
1:A:146:PHE:CE2	1:A:150:MET:HB2	2.45	0.47
1:A:235:LEU:HD23	1:A:235:LEU:HA	1.56	0.47
1:A:399:PHE:CD2	1:A:407:GLU:HB2	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:87:GLU:O	1:B:88:ASN:HB2	2.14	0.47
1:B:401:ASN:O	1:B:404:ARG:NH1	2.47	0.47
1:C:87:GLU:O	1:C:88:ASN:CB	2.62	0.47
1:C:174:GLN:OE1	1:C:235:LEU:HD13	2.13	0.47
1:C:268:MET:HE3	1:C:302:ILE:HD12	1.97	0.47
1:A:203:VAL:HG12	1:A:204:GLY:N	2.29	0.47
1:B:489:PRO:HA	1:B:491:TYR:CZ	2.50	0.47
1:B:123:LYS:HB2	1:B:254:GLU:CD	2.39	0.47
1:B:238:MET:HE3	1:B:238:MET:HB2	1.79	0.47
1:C:320:LEU:HD21	1:C:350:TRP:CD1	2.49	0.47
1:A:488:TYR:O	1:A:488:TYR:CG	2.69	0.46
1:C:193:LEU:HD23	1:C:193:LEU:HA	1.68	0.46
1:C:238:MET:HB2	1:C:238:MET:HE3	1.69	0.46
1:C:489:PRO:HB3	1:C:491:TYR:OH	2.15	0.46
1:C:28:VAL:O	1:C:30:VAL:HG13	2.16	0.46
1:A:96:PRO:HB2	1:A:228:ARG:HD3	1.97	0.46
1:B:167:TYR:CE2	1:B:169:ASN:HA	2.50	0.46
1:C:261:LYS:HB2	1:C:262:ARG:CB	2.41	0.46
1:B:485:THR:HG22	1:B:485:THR:O	2.16	0.46
1:A:32:HIS:ND1	1:A:32:HIS:N	2.64	0.46
1:A:53:PRO:HD2	1:A:274:LEU:CD2	2.43	0.46
1:B:358:ASP:OD1	1:B:472:LYS:HE2	2.16	0.46
1:B:426:GLU:HB3	1:C:387:LYS:HD2	1.97	0.46
1:C:238:MET:HB3	1:C:239:TRP:CD1	2.50	0.46
1:C:77:VAL:HB	1:C:78:PRO:HD2	1.98	0.46
1:A:45:LEU:HB3	1:A:84:MET:HE2	1.98	0.45
1:B:352:GLY:HA3	1:B:365:ALA:HA	1.97	0.45
1:A:426:GLU:HB3	1:B:387:LYS:HD2	1.98	0.45
1:C:490:LYS:N	1:C:491:TYR:CE2	2.84	0.45
1:A:139:ALA:HA	1:A:143:ASN:O	2.15	0.45
1:A:186:ASP:OD1	1:A:188:LYS:HD2	2.17	0.45
1:A:331:LEU:HD12	1:B:442:SER:CB	2.45	0.45
1:B:102:TYR:CZ	1:B:106:LYS:HD3	2.52	0.45
1:B:270:THR:OG1	1:B:271:GLU:N	2.49	0.45
1:C:401:ASN:O	1:C:404:ARG:NH1	2.49	0.45
1:C:310:LYS:HG2	1:C:310:LYS:O	2.17	0.45
1:A:135:SER:HA	2:D:2:SIA:O1B	2.17	0.45
1:B:123:LYS:HG2	1:B:131:THR:CG2	2.46	0.45
1:B:310:LYS:O	1:B:310:LYS:HG2	2.17	0.45
1:C:123:LYS:HG2	1:C:131:THR:CG2	2.45	0.45
1:A:95:TYR:CZ	1:A:229:MET:HE3	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:382:ASN:O	1:C:386:GLU:HG2	2.16	0.45
1:C:490:LYS:O	1:C:491:TYR:CD2	2.70	0.45
1:C:38:GLU:HG2	1:C:290:THR:HB	1.99	0.45
1:C:100:ASN:ND2	1:C:232:SER:OG	2.35	0.45
1:C:471:HIS:CG	1:C:490:LYS:HD2	2.52	0.45
1:A:316:LEU:CD1	1:A:429:VAL:HG22	2.47	0.45
1:A:402:LEU:HD13	1:B:101:ASP:OD2	2.17	0.45
1:B:284:PRO:HD3	1:B:300:LEU:O	2.17	0.45
1:C:248:GLY:O	1:C:249:ASN:HB2	2.16	0.45
1:A:38:GLU:O	1:A:295:HIS:HA	2.17	0.45
1:A:103:GLU:HA	1:A:106:LYS:HG3	1.98	0.45
1:C:309:VAL:HG22	1:C:422:THR:HA	1.98	0.45
1:C:411:LYS:O	1:C:415:ASP:HB2	2.17	0.45
1:C:489:PRO:HA	1:C:491:TYR:CZ	2.52	0.45
1:C:200:TYR:CE2	1:C:247:THR:HG23	2.53	0.44
1:C:331:LEU:HB2	1:C:438:ASP:OD1	2.17	0.44
1:C:10:GLY:HA3	1:C:343:TRP:CH2	2.53	0.44
1:A:54:LEU:HD23	1:A:83:ILE:HG12	2.00	0.44
1:B:418:LEU:HD23	1:C:391:GLN:OE1	2.18	0.44
1:C:48:LEU:HD12	1:C:279:THR:HG23	2.00	0.44
1:B:261:LYS:HB2	1:B:262:ARG:CB	2.41	0.44
1:A:69:PRO:CB	1:A:140:VAL:HG22	2.37	0.44
1:C:120:ILE:HG22	1:C:167:TYR:CE1	2.51	0.44
1:C:346:MET:HE2	1:C:346:MET:HB3	1.71	0.44
1:A:155:LYS:NZ	1:A:195:GLN:HG2	2.32	0.44
1:B:135:SER:N	1:B:145:SER:O	2.50	0.44
1:C:349:GLY:HA3	1:C:365:ALA:HB1	2.00	0.44
1:B:48:LEU:HD12	1:B:279:THR:HG23	1.99	0.43
1:B:87:GLU:O	1:B:88:ASN:CB	2.66	0.43
1:C:323:VAL:N	1:C:324:PRO:HD3	2.32	0.43
1:B:261:LYS:HB2	1:B:261:LYS:HE3	1.79	0.43
1:B:354:HIS:O	1:B:362:GLY:O	2.36	0.43
1:C:119:LYS:HD3	1:C:122:PRO:HA	2.00	0.43
1:A:426:GLU:HG3	1:B:387:LYS:HD2	2.00	0.43
1:A:480:SER:O	1:A:485:THR:N	2.49	0.43
1:B:77:VAL:HB	1:B:78:PRO:HD2	2.01	0.43
1:B:97:GLY:HA2	1:B:228:ARG:HB3	1.99	0.43
1:B:412:LYS:NZ	1:C:395:VAL:HG22	2.33	0.43
1:B:24:LEU:O	1:C:379:ASN:ND2	2.36	0.43
1:B:322:ASN:C	1:B:324:PRO:HD3	2.43	0.43
1:B:412:LYS:HE2	1:C:414:GLU:OE1	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:87:GLU:O	1:A:88:ASN:CB	2.66	0.43
1:A:182:HIS:O	1:A:184:PRO:HD3	2.18	0.43
1:C:167:TYR:CE2	1:C:169:ASN:HA	2.53	0.43
1:A:318:THR:O	1:A:377:ILE:HG13	2.19	0.43
1:A:428:LEU:HD12	1:A:428:LEU:HA	1.61	0.43
1:B:456:ARG:HB2	1:B:457:ASP:H	1.48	0.43
1:C:358:ASP:OD1	1:C:472:LYS:HE2	2.18	0.43
1:C:456:ARG:HB2	1:C:457:ASP:H	1.51	0.43
1:A:49:ASN:HD21	1:A:280:LYS:HG2	1.83	0.43
1:A:113:LYS:HB3	1:A:260:SER:O	2.18	0.43
1:A:85:GLU:O	1:A:269:LYS:HA	2.18	0.43
1:B:316:LEU:HD12	1:B:429:VAL:HG13	2.01	0.43
1:C:241:THR:HG22	1:C:242:ILE:N	2.33	0.43
1:B:96:PRO:HB3	1:B:222:VAL:CG1	2.49	0.43
1:C:352:GLY:HA3	1:C:365:ALA:HA	2.00	0.43
1:A:220:PRO:HG3	1:B:241:THR:HB	2.01	0.42
1:B:112:VAL:C	1:B:262:ARG:CD	2.92	0.42
1:A:19:LYS:HE3	1:A:29:THR:OG1	2.19	0.42
1:B:74:LEU:HD23	1:B:74:LEU:HA	1.76	0.42
1:C:485:THR:HG22	1:C:485:THR:O	2.19	0.42
1:B:155:LYS:CE	1:B:195:GLN:HG2	2.48	0.42
1:B:359:GLN:NE2	1:B:475:ASP:N	2.68	0.42
1:A:435:ARG:HH11	1:C:435:ARG:HH11	1.60	0.42
1:B:178:ILE:O	1:B:253:PRO:HG3	2.19	0.42
1:B:309:VAL:HG22	1:B:422:THR:HA	2.02	0.42
1:C:354:HIS:O	1:C:362:GLY:O	2.37	0.42
1:C:205:THR:CG2	1:C:242:ILE:HG13	2.50	0.42
1:B:154:THR:CG2	1:B:193:LEU:HD22	2.50	0.42
1:C:97:GLY:HA2	1:C:228:ARG:HB3	2.02	0.42
1:A:211:ARG:HD3	1:C:215:GLU:HA	2.00	0.42
1:A:310:LYS:O	1:A:310:LYS:HG2	2.19	0.42
1:A:459:VAL:HG21	1:A:467:PHE:HD2	1.85	0.42
1:B:158:SER:O	1:B:195:GLN:HG3	2.20	0.42
1:B:227:SER:C	1:B:228:ARG:HG2	2.45	0.42
1:C:74:LEU:HD23	1:C:74:LEU:HA	1.75	0.42
1:A:145:SER:HB3	1:A:146:PHE:H	1.42	0.42
1:A:215:GLU:HB3	1:B:211:ARG:CB	2.50	0.42
1:A:243:ASN:HD22	1:C:220:PRO:HD3	1.85	0.42
1:B:112:VAL:C	1:B:262:ARG:HD2	2.44	0.42
1:B:384:VAL:HG22	1:B:428:LEU:HD21	2.02	0.42
1:B:491:TYR:HB2	1:B:492:GLU:H	1.64	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:18:GLU:O	1:C:29:THR:HA	2.20	0.42
1:B:11:TYR:CE2	1:B:335:ILE:HA	2.55	0.41
1:B:199:THR:OG1	1:B:249:ASN:ND2	2.53	0.41
1:B:248:GLY:C	1:B:249:ASN:HD22	2.28	0.41
1:C:48:LEU:O	1:C:49:ASN:HB2	2.19	0.41
1:C:193:LEU:HD21	2:F:2:SIA:O10	2.20	0.41
1:C:303:GLY:O	1:C:392:PHE:HA	2.20	0.41
1:C:448:TYR:CE1	1:C:465:GLY:HA2	2.55	0.41
1:A:189:GLU:OE2	2:D:2:SIA:O9	2.36	0.41
1:C:384:VAL:HG22	1:C:428:LEU:HD21	2.02	0.41
1:A:203:VAL:HA	1:A:243:ASN:O	2.20	0.41
1:C:158:SER:O	1:C:195:GLN:HG3	2.20	0.41
2:D:2:SIA:O1A	2:D:2:SIA:H6	2.20	0.41
1:C:185:ILE:H	1:C:185:ILE:HG12	1.67	0.41
1:B:38:GLU:HG2	1:B:290:THR:HB	2.02	0.41
1:B:48:LEU:O	1:B:49:ASN:HB2	2.21	0.41
1:B:307:LYS:CG	1:B:421:TRP:CE2	2.97	0.41
1:B:323:VAL:HG12	1:B:323:VAL:O	2.20	0.41
1:A:96:PRO:HG3	1:A:222:VAL:O	2.21	0.41
1:A:154:THR:HG22	1:A:193:LEU:HD22	2.02	0.41
1:B:220:PRO:HD2	1:C:205:THR:O	2.21	0.41
1:A:404:ARG:O	1:A:405:ARG:C	2.63	0.41
1:A:412:LYS:HE2	1:B:414:GLU:OE2	2.21	0.41
1:A:453:MET:CE	1:C:463:GLY:HA2	2.51	0.41
1:B:248:GLY:O	1:B:249:ASN:CB	2.68	0.41
1:C:155:LYS:CE	1:C:195:GLN:HG2	2.51	0.41
1:A:427:LEU:O	1:A:428:LEU:C	2.63	0.41
1:B:18:GLU:O	1:B:29:THR:HA	2.21	0.41
1:B:120:ILE:HG22	1:B:167:TYR:CZ	2.56	0.41
1:B:491:TYR:O	1:B:494:GLU:OE1	2.37	0.41
1:A:281:CYS:HB2	1:A:304:GLU:O	2.21	0.41
1:A:402:LEU:HD23	1:A:402:LEU:HA	1.79	0.41
1:B:293:PRO:HG3	1:B:385:ILE:HA	2.02	0.41
1:B:411:LYS:O	1:B:415:ASP:HB2	2.21	0.41
1:C:178:ILE:O	1:C:253:PRO:HG3	2.20	0.41
1:A:332:PHE:CE1	1:A:442:SER:HB2	2.56	0.41
1:B:377:ILE:HD13	1:B:377:ILE:HA	1.83	0.41
1:A:155:LYS:HZ3	1:A:195:GLN:HG2	1.85	0.40
1:A:413:MET:HE1	1:C:412:LYS:HG2	2.02	0.40
1:A:426:GLU:CB	1:B:387:LYS:HD2	2.51	0.40
1:C:123:LYS:HB2	1:C:254:GLU:OE1	2.20	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:489:PRO:C	1:C:491:TYR:CE1	2.98	0.40
1:A:485:THR:O	1:A:486:TYR:C	2.64	0.40
1:C:154:THR:HG21	1:C:193:LEU:HD22	2.03	0.40
1:A:105:LEU:HB2	1:A:233:TRP:CZ2	2.55	0.40
1:A:185:ILE:CG1	1:A:189:GLU:OE1	2.70	0.40
1:A:413:MET:HE1	1:C:412:LYS:CG	2.51	0.40
1:C:490:LYS:C	1:C:491:TYR:CG	2.99	0.40
1:A:211:ARG:HB3	1:C:215:GLU:CG	2.51	0.40
1:B:241:THR:HG22	1:B:242:ILE:N	2.37	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	463/505 (92%)	389 (84%)	63 (14%)	11 (2%)	4	24
1	B	482/505 (95%)	433 (90%)	37 (8%)	12 (2%)	4	23
1	C	477/505 (94%)	429 (90%)	37 (8%)	11 (2%)	5	25
All	All	1422/1515 (94%)	1251 (88%)	137 (10%)	34 (2%)	4	24

All (34) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	262	ARG
1	B	262	ARG
1	B	487	ASP
1	B	488	TYR
1	C	262	ARG
1	C	488	TYR
1	C	491	TYR
1	A	487	ASP

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Mol	Chain	Res	Type
1	B	264	SER
1	C	90	ARG
1	C	264	SER
1	C	487	ASP
1	A	87	GLU
1	A	90	ARG
1	A	263	GLY
1	B	90	ARG
1	B	249	ASN
1	B	491	TYR
1	B	492	GLU
1	C	249	ASN
1	A	264	SER
1	A	488	TYR
1	B	76	SER
1	B	88	ASN
1	B	355	HIS
1	C	76	SER
1	C	88	ASN
1	C	355	HIS
1	A	88	ASN
1	A	195	GLN
1	A	288	ILE
1	C	263	GLY
1	B	263	GLY
1	A	395	VAL

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	409/442 (92%)	368 (90%)	41 (10%)	7	29
1	B	423/442 (96%)	391 (92%)	32 (8%)	12	41
1	C	418/442 (95%)	388 (93%)	30 (7%)	13	43
All	All	1250/1326 (94%)	1147 (92%)	103 (8%)	10	37

All (103) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	CYS
1	A	22	THR
1	A	25	GLU
1	A	26	ARG
1	A	32	HIS
1	A	34	LYS
1	A	36	ILE
1	A	40	THR
1	A	77	VAL
1	A	83	ILE
1	A	109	LEU
1	A	118	VAL
1	A	124	ASP
1	A	138	CYS
1	A	170	THR
1	A	185	ILE
1	A	187	GLU
1	A	191	ARG
1	A	192	THR
1	A	193	LEU
1	A	195	GLN
1	A	196	ASN
1	A	206	SER
1	A	235	LEU
1	A	241	THR
1	A	247	THR
1	A	259	ILE
1	A	264	SER
1	A	309	VAL
1	A	320	LEU
1	A	339	ILE
1	A	391	GLN
1	A	409	LEU
1	A	414	GLU
1	A	427	LEU
1	A	442	SER
1	A	456	ARG
1	A	457	ASP
1	A	459	VAL
1	A	477	CYS
1	A	487	ASP
1	B	6	GLN

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Mol	Chain	Res	Type
1	B	26	ARG
1	B	29	THR
1	B	34	LYS
1	B	40	THR
1	B	46	CYS
1	B	59	CYS
1	B	61	ILE
1	B	106	LYS
1	B	109	LEU
1	B	118	VAL
1	B	124	ASP
1	B	131	THR
1	B	185	ILE
1	B	187	GLU
1	B	238	MET
1	B	260	SER
1	B	264	SER
1	B	268	MET
1	B	283	THR
1	B	291	THR
1	B	307	LYS
1	B	309	VAL
1	B	320	LEU
1	B	351	TYR
1	B	355	HIS
1	B	391	GLN
1	B	409	LEU
1	B	414	GLU
1	B	427	LEU
1	B	477	CYS
1	B	487	ASP
1	C	26	ARG
1	C	34	LYS
1	C	40	THR
1	C	46	CYS
1	C	59	CYS
1	C	61	ILE
1	C	106	LYS
1	C	109	LEU
1	C	124	ASP
1	C	131	THR
1	C	185	ILE

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Mol	Chain	Res	Type
1	C	187	GLU
1	C	238	MET
1	C	260	SER
1	C	264	SER
1	C	268	MET
1	C	283	THR
1	C	291	THR
1	C	307	LYS
1	C	309	VAL
1	C	320	LEU
1	C	351	TYR
1	C	355	HIS
1	C	391	GLN
1	C	409	LEU
1	C	414	GLU
1	C	427	LEU
1	C	457	ASP
1	C	477	CYS
1	C	487	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	49	ASN
1	A	100	ASN
1	A	128	GLN
1	A	129	HIS
1	A	190	GLN
1	A	196	ASN
1	A	243	ASN
1	A	249	ASN
1	A	322	ASN
1	A	424	ASN
1	A	440	HIS
1	B	129	HIS
1	B	136	GLN
1	B	190	GLN
1	B	243	ASN
1	B	249	ASN
1	B	359	GLN
1	B	371	GLN
1	B	424	ASN

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Mol	Chain	Res	Type
1	B	454	GLN
1	C	129	HIS
1	C	136	GLN
1	C	190	GLN
1	C	249	ASN
1	C	359	GLN
1	C	371	GLN
1	C	424	ASN
1	C	454	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

6 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	$\# Z > 2$
2	GAL	D	1	2	12,12,12	0.37	0	17,17,17	0.90	0
2	SIA	D	2	2	20,20,21	0.78	0	21,28,31	1.16	2 (9%)
2	GAL	E	1	2	12,12,12	0.67	0	17,17,17	0.48	0
2	SIA	E	2	2	20,20,21	1.18	2 (10%)	21,28,31	1.51	4 (19%)
2	GAL	F	1	2	12,12,12	0.53	0	17,17,17	0.97	1 (5%)
2	SIA	F	2	2	20,20,21	0.92	1 (5%)	21,28,31	1.43	3 (14%)

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	GAL	D	1	2	-	1/2/22/22	0/1/1/1
2	SIA	D	2	2	-	1/18/34/38	0/1/1/1
2	GAL	E	1	2	-	2/2/22/22	0/1/1/1
2	SIA	E	2	2	-	1/18/34/38	0/1/1/1
2	GAL	F	1	2	-	0/2/22/22	0/1/1/1
2	SIA	F	2	2	-	2/18/34/38	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	2	SIA	C4-C5	-2.83	1.50	1.53
2	F	2	SIA	C2-C1	2.24	1.55	1.52
2	E	2	SIA	C2-C1	2.15	1.54	1.52

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	2	SIA	O9-C9-C8	-3.60	103.59	111.16
2	E	2	SIA	C4-C5-N5	-3.37	103.80	110.44
2	E	2	SIA	C8-C7-C6	-2.92	107.57	113.05
2	E	2	SIA	O1B-C1-C2	2.85	120.12	112.71
2	D	2	SIA	C8-C7-C6	-2.44	108.48	113.05
2	F	1	GAL	O5-C5-C4	2.42	114.06	109.70
2	E	2	SIA	O4-C4-C5	-2.39	104.41	109.84
2	D	2	SIA	O1B-C1-C2	2.22	118.48	112.71
2	F	2	SIA	C9-C8-C7	-2.22	107.66	112.17
2	F	2	SIA	O1B-C1-C2	2.16	118.33	112.71

There are no chirality outliers.

All (7) torsion outliers are listed below:

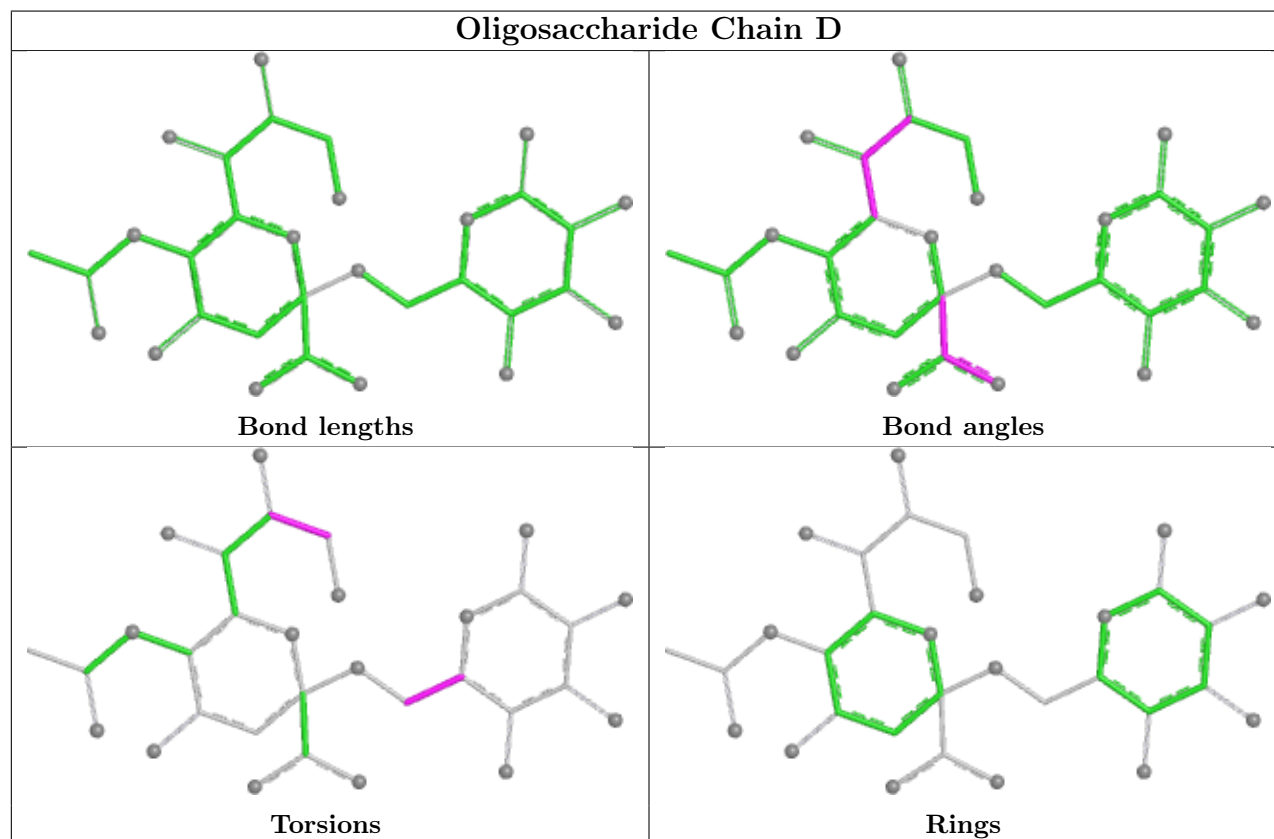
Mol	Chain	Res	Type	Atoms
2	F	2	SIA	C7-C8-C9-O9
2	F	2	SIA	O8-C8-C9-O9
2	E	1	GAL	O5-C5-C6-O6
2	D	2	SIA	O8-C8-C9-O9
2	E	1	GAL	C4-C5-C6-O6
2	E	2	SIA	O8-C8-C9-O9
2	D	1	GAL	O5-C5-C6-O6

There are no ring outliers.

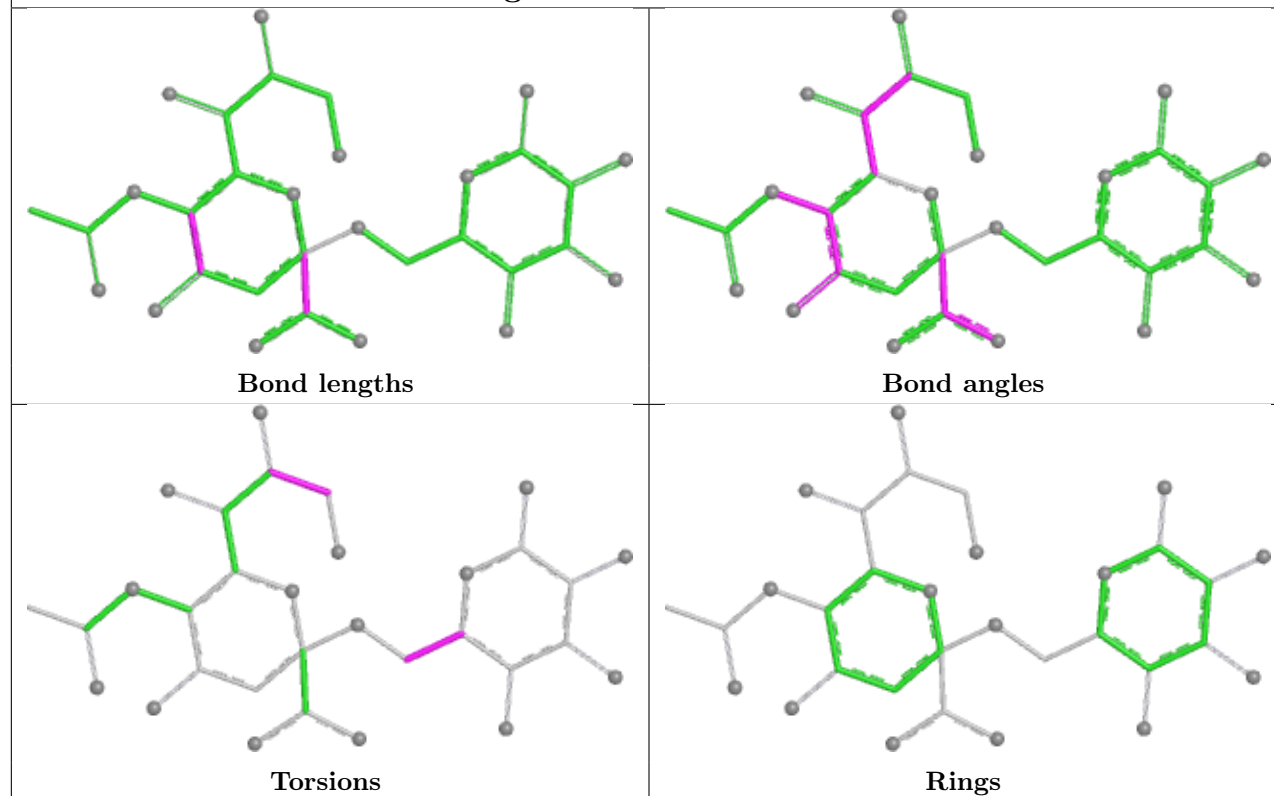
2 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	2	SIA	7	0
2	F	2	SIA	1	0

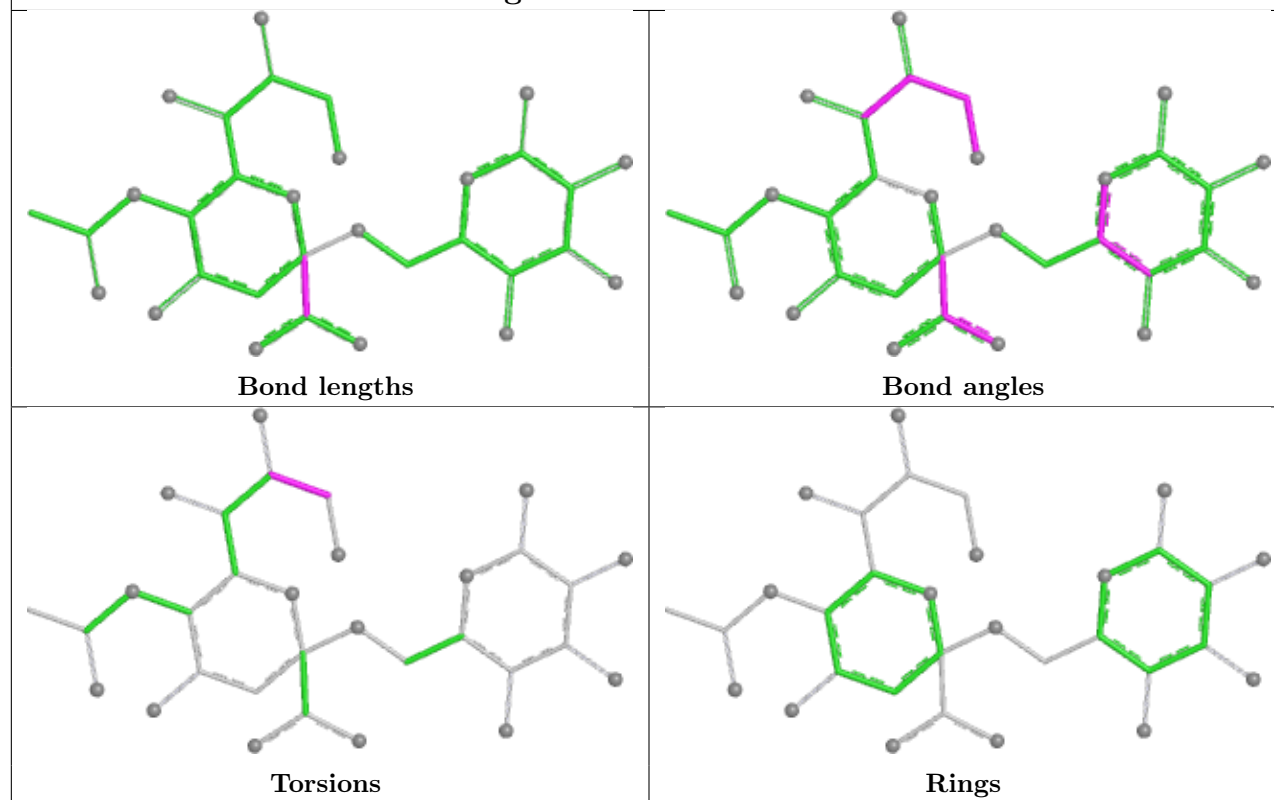
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.



Oligosaccharide Chain E



Oligosaccharide Chain F



5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	471/505 (93%)	-0.21	9 (1%) 66 43	34, 72, 122, 234	0
1	B	486/505 (96%)	-0.10	8 (1%) 70 47	34, 62, 103, 186	0
1	C	481/505 (95%)	-0.11	8 (1%) 69 45	35, 59, 106, 206	0
All	All	1438/1515 (94%)	-0.14	25 (1%) 69 45	34, 64, 112, 234	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	491	TYR	5.3
1	C	128	GLN	4.3
1	C	330	GLY	3.9
1	C	487	ASP	3.6
1	A	141	SER	3.3
1	B	489	PRO	3.0
1	B	487	ASP	2.9
1	C	58	ASP	2.8
1	B	159	ASN	2.8
1	B	187	GLU	2.7
1	C	493	GLU	2.6
1	A	487	ASP	2.6
1	C	277	CYS	2.4
1	A	289	ASN	2.3
1	B	223	ASN	2.3
1	A	142	GLY	2.3
1	A	94	CYS	2.3
1	A	75	LEU	2.3
1	B	491	TYR	2.3
1	B	8	CYS	2.2
1	B	494	GLU	2.2
1	A	124	ASP	2.1
1	A	128	GLN	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	140	VAL	2.1
1	A	277	CYS	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](#)

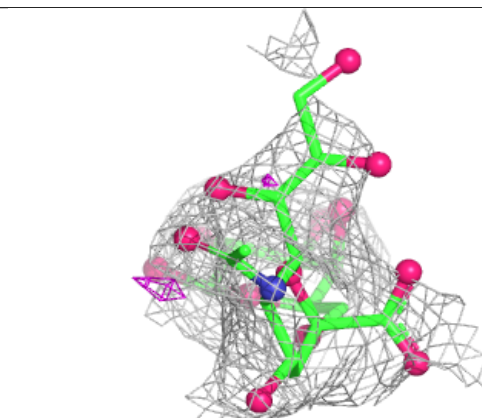
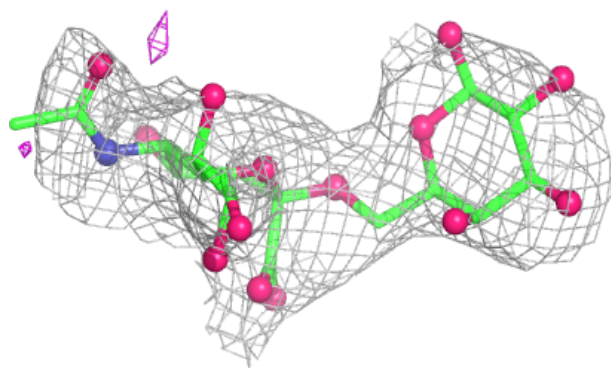
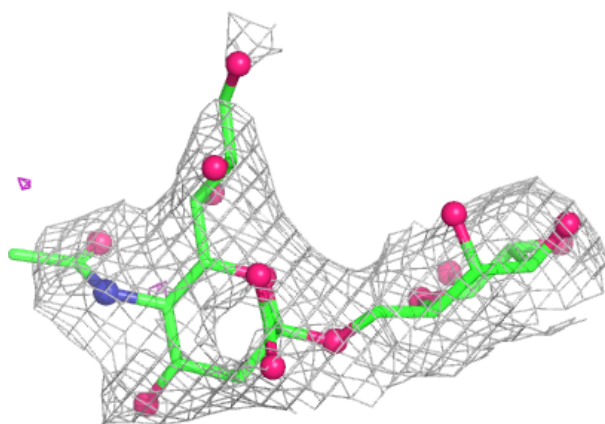
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists 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
2	GAL	E	1	12/12	0.66	0.13	86,86,86,86	0
2	GAL	D	1	12/12	0.76	0.10	110,110,110,110	0
2	GAL	F	1	12/12	0.77	0.08	78,78,78,78	0
2	SIA	D	2	20/21	0.87	0.12	110,110,110,110	0
2	SIA	F	2	20/21	0.88	0.12	78,78,78,78	0
2	SIA	E	2	20/21	0.90	0.10	86,86,86,86	0

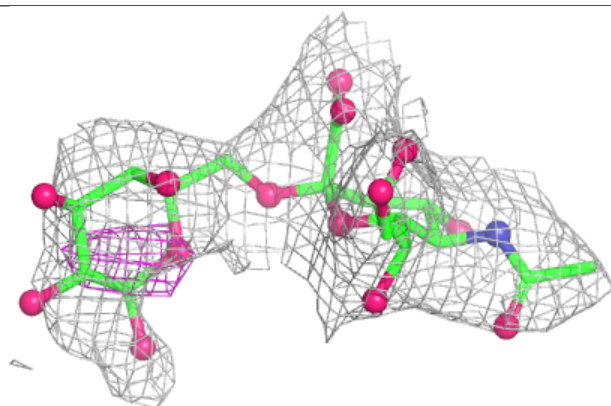
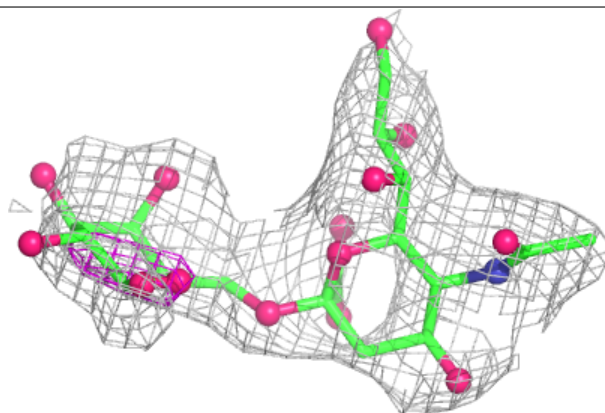
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

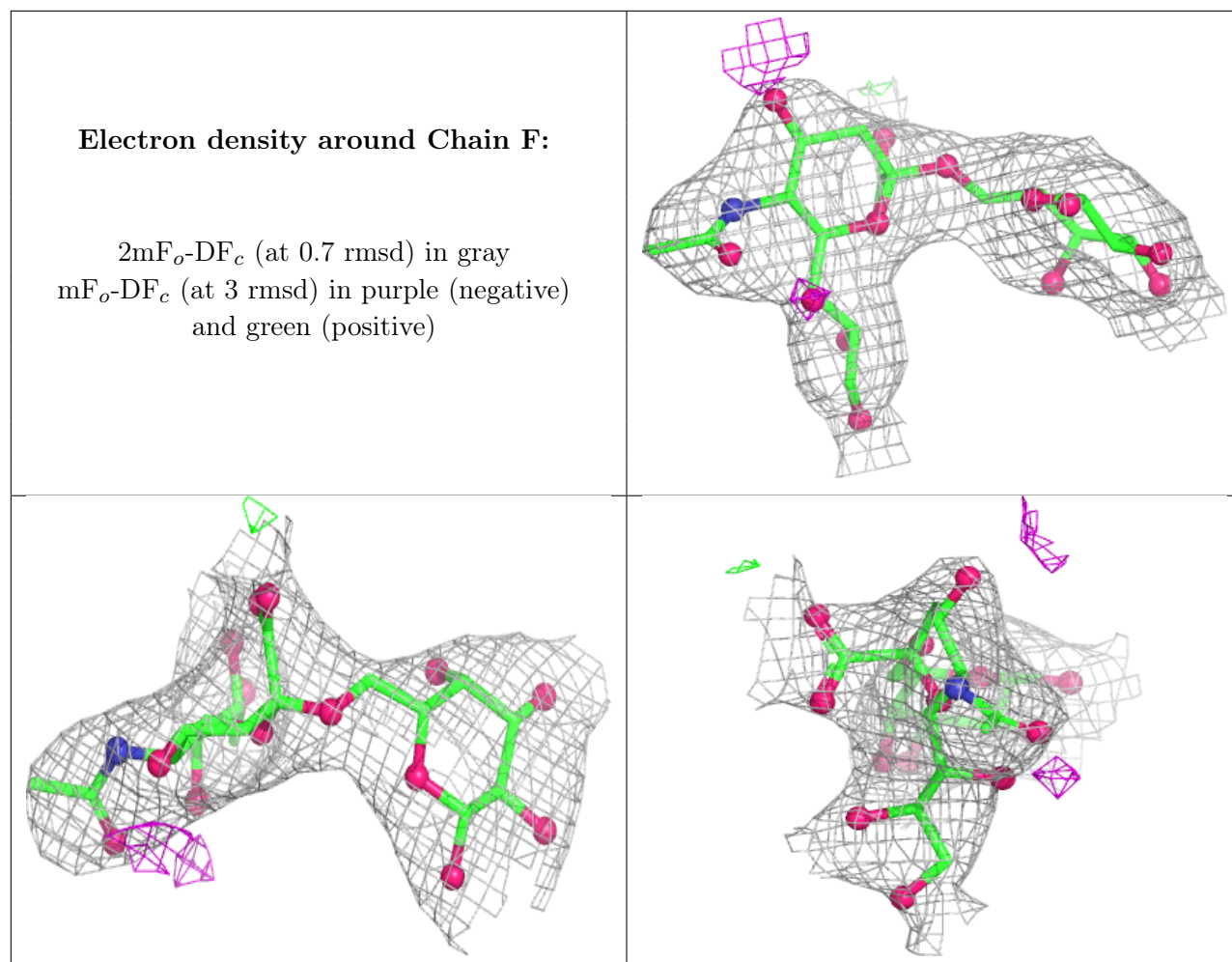
Electron density around Chain D:

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

**Electron density around Chain E:**

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





6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.