



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:19 PM UTC

PDB ID : 3KG2 / pdb_00003kg2
Title : AMPA subtype ionotropic glutamate receptor in complex with competitive antagonist ZK 200775
Authors : Sobolevsky, A.I.; Rosconi, M.P.; Gouaux, E.
Deposited on : 2009-10-28
Resolution : 3.60 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
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
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

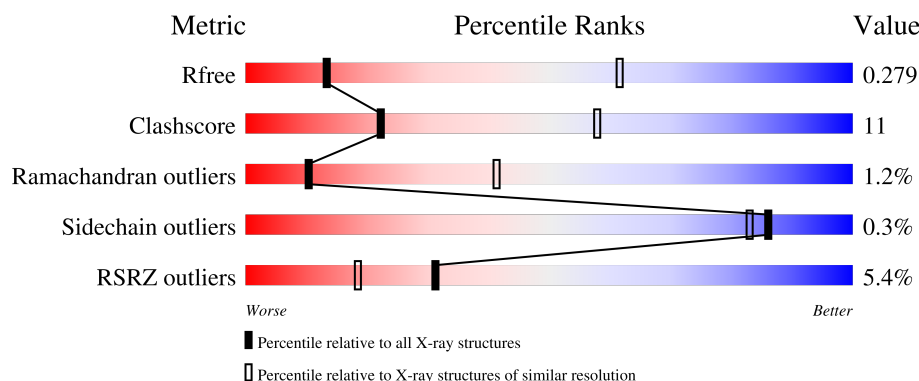
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.60 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.



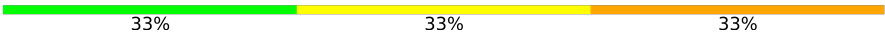
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1747 (3.70-3.50)
Clashscore	190562	1827 (3.70-3.50)
Ramachandran outliers	187476	1773 (3.70-3.50)
Sidechain outliers	187428	1772 (3.70-3.50)
RSRZ outliers	180081	1745 (3.70-3.50)

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	823	6% 75% 19% 5%
1	B	823	6% 75% 19% 5%
1	C	823	4% 75% 19% 5%
1	D	823	5% 75% 19% • 5%
2	E	4	50% 25% 25%

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Mol	Chain	Length	Quality of chain
2	F	4	 50% 25% 25%
3	G	3	 33% 33% 33%
3	H	3	 33% 67%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 22686 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 Glutamate receptor 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	779	Total	C	N	O	S	0	0	0
			5600	3605	891	1077	27			
1	B	779	Total	C	N	O	S	0	0	0
			5600	3605	891	1077	27			
1	C	779	Total	C	N	O	S	0	0	0
			5600	3605	891	1077	27			
1	D	779	Total	C	N	O	S	0	0	0
			5600	3605	891	1077	27			

There are 68 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	241	GLU	ASN	engineered mutation	UNP P19491
A	?	-	VAL	deletion	UNP P19491
A	?	-	THR	deletion	UNP P19491
A	?	-	LEU	deletion	UNP P19491
A	?	-	PRO	deletion	UNP P19491
A	?	-	SER	deletion	UNP P19491
A	?	-	GLY	deletion	UNP P19491
A	385	ASP	ASN	engineered mutation	UNP P19491
A	392	GLN	ASN	engineered mutation	UNP P19491
A	410	ALA	LYS	engineered mutation	UNP P19491
A	413	ALA	GLU	engineered mutation	UNP P19491
A	414	ALA	MET	engineered mutation	UNP P19491
A	416	ALA	GLU	engineered mutation	UNP P19491
A	589	ALA	CYS	engineered mutation	UNP P19491
A	744	THR	ASN	variant	UNP P19491
A	745	PRO	ALA	variant	UNP P19491
A	754	SER	ASN	variant	UNP P19491
B	241	GLU	ASN	engineered mutation	UNP P19491
B	?	-	VAL	deletion	UNP P19491
B	?	-	THR	deletion	UNP P19491
B	?	-	LEU	deletion	UNP P19491

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Chain	Residue	Modelled	Actual	Comment	Reference
B	?	-	PRO	deletion	UNP P19491
B	?	-	SER	deletion	UNP P19491
B	?	-	GLY	deletion	UNP P19491
B	385	ASP	ASN	engineered mutation	UNP P19491
B	392	GLN	ASN	engineered mutation	UNP P19491
B	410	ALA	LYS	engineered mutation	UNP P19491
B	413	ALA	GLU	engineered mutation	UNP P19491
B	414	ALA	MET	engineered mutation	UNP P19491
B	416	ALA	GLU	engineered mutation	UNP P19491
B	589	ALA	CYS	engineered mutation	UNP P19491
B	744	THR	ASN	variant	UNP P19491
B	745	PRO	ALA	variant	UNP P19491
B	754	SER	ASN	variant	UNP P19491
C	241	GLU	ASN	engineered mutation	UNP P19491
C	?	-	VAL	deletion	UNP P19491
C	?	-	THR	deletion	UNP P19491
C	?	-	LEU	deletion	UNP P19491
C	?	-	PRO	deletion	UNP P19491
C	?	-	SER	deletion	UNP P19491
C	?	-	GLY	deletion	UNP P19491
C	385	ASP	ASN	engineered mutation	UNP P19491
C	392	GLN	ASN	engineered mutation	UNP P19491
C	410	ALA	LYS	engineered mutation	UNP P19491
C	413	ALA	GLU	engineered mutation	UNP P19491
C	414	ALA	MET	engineered mutation	UNP P19491
C	416	ALA	GLU	engineered mutation	UNP P19491
C	589	ALA	CYS	engineered mutation	UNP P19491
C	744	THR	ASN	variant	UNP P19491
C	745	PRO	ALA	variant	UNP P19491
C	754	SER	ASN	variant	UNP P19491
D	241	GLU	ASN	engineered mutation	UNP P19491
D	?	-	VAL	deletion	UNP P19491
D	?	-	THR	deletion	UNP P19491
D	?	-	LEU	deletion	UNP P19491
D	?	-	PRO	deletion	UNP P19491
D	?	-	SER	deletion	UNP P19491
D	?	-	GLY	deletion	UNP P19491
D	385	ASP	ASN	engineered mutation	UNP P19491
D	392	GLN	ASN	engineered mutation	UNP P19491
D	410	ALA	LYS	engineered mutation	UNP P19491
D	413	ALA	GLU	engineered mutation	UNP P19491
D	414	ALA	MET	engineered mutation	UNP P19491

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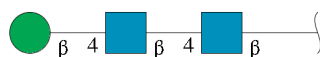
Chain	Residue	Modelled	Actual	Comment	Reference
D	416	ALA	GLU	engineered mutation	UNP P19491
D	589	ALA	CYS	engineered mutation	UNP P19491
D	744	THR	ASN	variant	UNP P19491
D	745	PRO	ALA	variant	UNP P19491
D	754	SER	ASN	variant	UNP P19491

- Molecule 2 is an oligosaccharide called beta-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



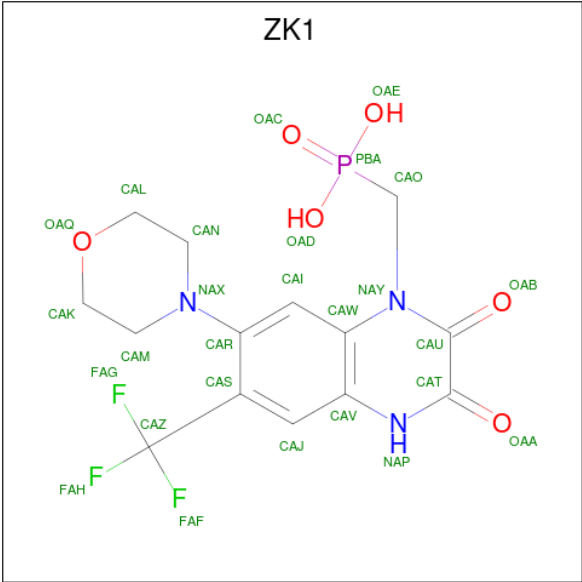
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	4	Total	C	N	O	0	0	0
			50	28	2	20			
2	F	4	Total	C	N	O	0	0	0
			50	28	2	20			

- Molecule 3 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
3	G	3	Total	C	N	O	0	0	0
			39	22	2	15			
3	H	3	Total	C	N	O	0	0	0
			39	22	2	15			

- Molecule 4 is {[7-morpholin-4-yl-2,3-dioxo-6-(trifluoromethyl)-3,4-dihydroquinoxalin-1(2H)-yl]methyl}phosphonic acid (CCD ID: ZK1) (formula: C₁₄H₁₅F₃N₃O₆P).

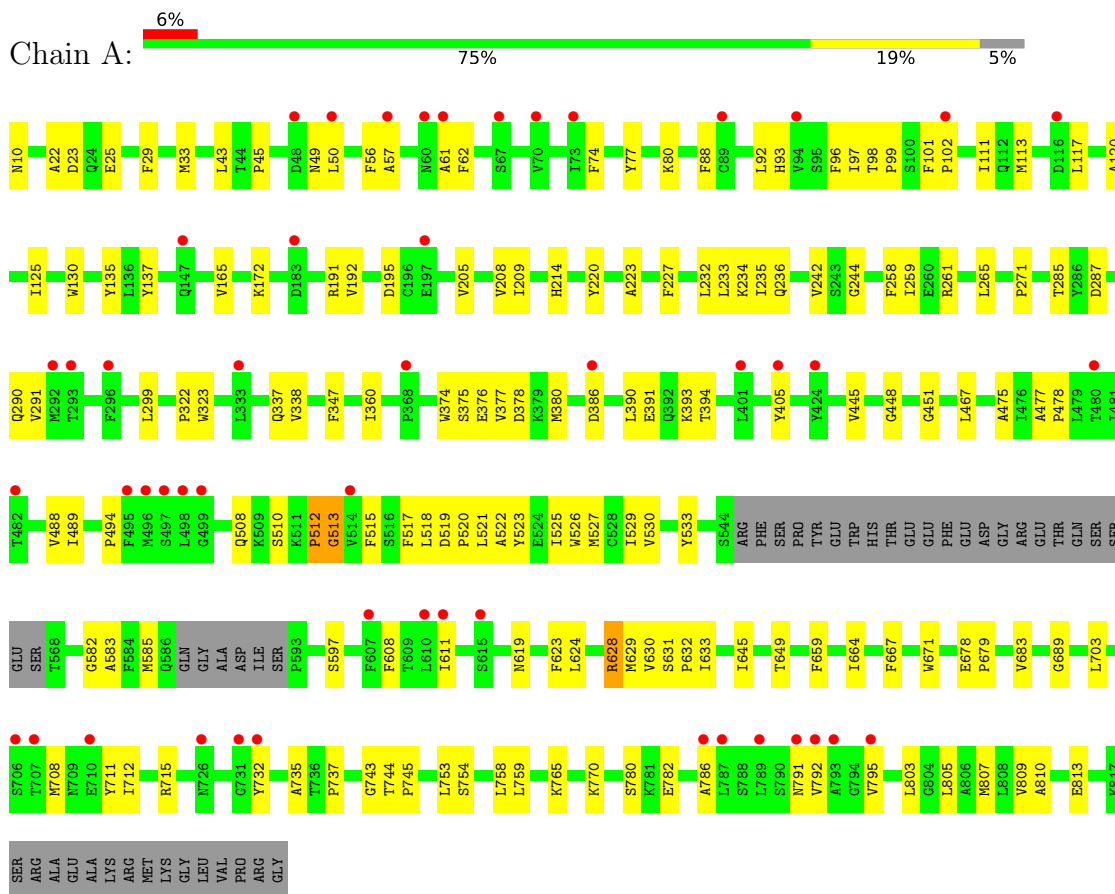


Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
4	A	1	Total	C	F	N	O	P	0	0
			27	14	3	3	6	1		
4	B	1	Total	C	F	N	O	P	0	0
			27	14	3	3	6	1		
4	C	1	Total	C	F	N	O	P	0	0
			27	14	3	3	6	1		
4	D	1	Total	C	F	N	O	P	0	0
			27	14	3	3	6	1		

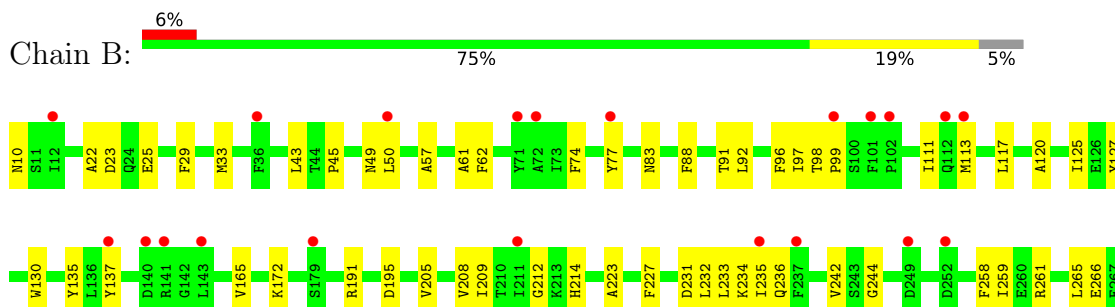
3 Residue-property plots [i](#)

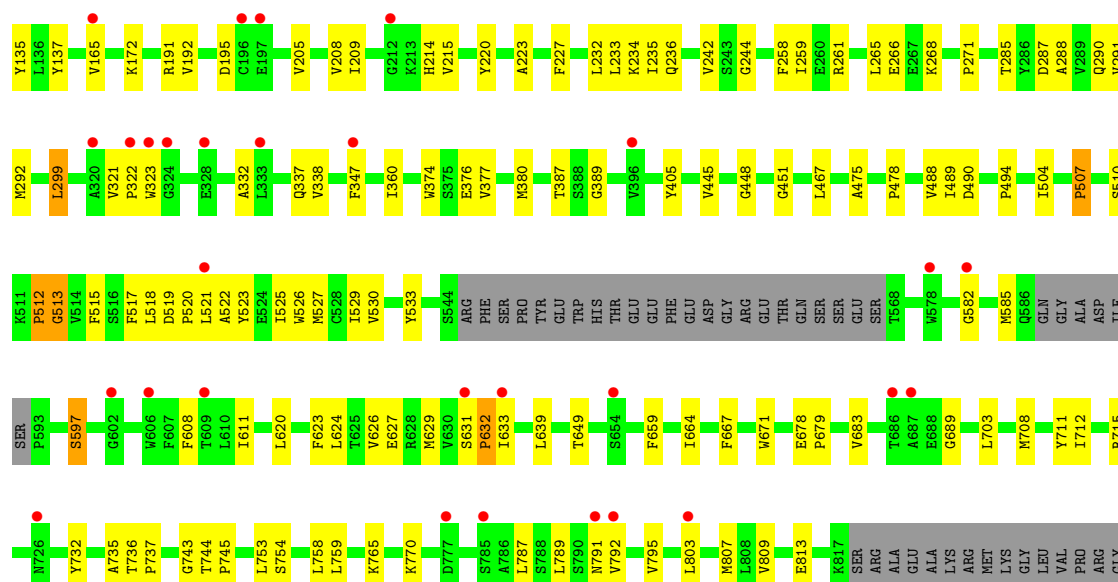
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Glutamate receptor 2



• Molecule 1: Glutamate receptor 2





- Molecule 2: beta-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain E: 50% 25% 25%



- Molecule 2: beta-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F: 50% 25% 25%



- Molecule 3: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain G: 33% 33% 33%



- Molecule 3: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain H: 33% 67%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	90.72Å 109.85Å 161.13Å 85.32° 84.75° 78.92°	Depositor
Resolution (Å)	49.89 – 3.60 49.89 – 3.60	Depositor EDS
% Data completeness (in resolution range)	98.8 (49.89-3.60) 98.9 (49.89-3.60)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.15 (at 3.57Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.286 , 0.296 0.269 , 0.279	Depositor DCC
R_{free} test set	3484 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	136.5	Xtriage
Anisotropy	0.162	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 99.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	22686	wwPDB-VP
Average B, all atoms (Å ²)	179.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.26% 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, ZK1, 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	0.26	0/5721	0.68	5/7837 (0.1%)
1	B	0.25	0/5721	0.67	2/7837 (0.0%)
1	C	0.26	0/5721	0.70	5/7837 (0.1%)
1	D	0.25	0/5721	0.67	2/7837 (0.0%)
All	All	0.25	0/22884	0.68	14/31348 (0.0%)

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	628	ARG	NE-CZ-NH1	-8.12	113.38	121.50
1	C	628	ARG	NE-CZ-NH2	7.82	126.24	119.20
1	C	787	LEU	N-CA-C	6.06	118.37	110.43
1	A	391	GLU	N-CA-C	-5.77	106.58	113.97
1	A	628	ARG	NE-CZ-NH2	-5.76	114.01	119.20
1	B	585	MET	N-CA-C	5.60	117.92	110.53
1	C	585	MET	N-CA-C	5.58	117.90	110.53
1	A	585	MET	N-CA-C	5.50	117.79	110.53
1	D	585	MET	N-CA-C	5.50	117.79	110.53
1	A	582	GLY	N-CA-C	-5.36	108.77	114.67
1	C	582	GLY	N-CA-C	-5.35	108.79	114.67
1	D	582	GLY	N-CA-C	-5.34	108.80	114.67
1	B	582	GLY	N-CA-C	-5.33	108.81	114.67
1	A	628	ARG	CD-NE-CZ	5.13	131.59	124.40

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5600	0	5083	124	0
1	B	5600	0	5084	125	0
1	C	5600	0	5083	120	0
1	D	5600	0	5083	129	0
2	E	50	0	43	3	0
2	F	50	0	43	1	0
3	G	39	0	34	1	0
3	H	39	0	34	2	0
4	A	27	0	13	0	0
4	B	27	0	13	1	0
4	C	27	0	13	1	0
4	D	27	0	13	1	0
All	All	22686	0	20539	456	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 11.

All (456) 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:626:VAL:HB	1:C:628:ARG:HH11	1.24	0.99
1:D:649:THR:HG22	1:D:703:LEU:HB2	1.68	0.76
1:A:525:ILE:HG12	1:B:789:LEU:HD13	1.68	0.76
1:B:649:THR:HG22	1:B:703:LEU:HB2	1.68	0.76
1:A:649:THR:HG22	1:A:703:LEU:HB2	1.68	0.75
1:A:623:PHE:HE1	1:B:785:SER:HA	1.52	0.74
1:C:649:THR:HG22	1:C:703:LEU:HB2	1.68	0.74
1:A:623:PHE:CE1	1:B:785:SER:HA	2.23	0.74
1:C:611:ILE:HG21	1:D:795:VAL:HG21	1.70	0.72
1:A:628:ARG:NH1	1:D:627:GLU:H	1.91	0.68
1:C:521:LEU:HD22	1:C:526:TRP:CD2	2.32	0.64
1:B:521:LEU:HD23	1:B:525:ILE:HB	1.80	0.64
1:A:521:LEU:HD23	1:A:525:ILE:HB	1.80	0.64
1:A:810:ALA:HB2	1:D:597:SER:OG	1.99	0.63
1:D:521:LEU:HD23	1:D:525:ILE:HB	1.80	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:521:LEU:HD22	1:B:526:TRP:CD2	2.35	0.62
1:C:525:ILE:HG12	1:D:789:LEU:HD13	1.81	0.62
1:C:50:LEU:HD23	1:C:57:ALA:HB1	1.82	0.62
1:A:521:LEU:HD22	1:A:526:TRP:CD2	2.35	0.62
1:D:510:SER:O	1:D:512:PRO:HD3	2.00	0.62
1:A:619:ASN:OD1	1:B:787:LEU:HB2	2.00	0.62
1:D:521:LEU:HD22	1:D:526:TRP:CD2	2.35	0.62
1:A:628:ARG:NH2	1:D:623:PHE:HD1	1.98	0.61
1:D:522:ALA:HB3	1:D:525:ILE:HG13	1.83	0.61
1:A:50:LEU:HD23	1:A:57:ALA:HB1	1.82	0.61
1:A:475:ALA:HB3	1:A:735:ALA:HB3	1.82	0.61
1:C:521:LEU:HD23	1:C:525:ILE:HB	1.83	0.61
1:D:50:LEU:HD23	1:D:57:ALA:HB1	1.82	0.61
1:D:489:ILE:HD12	1:D:735:ALA:HB1	1.83	0.61
1:A:522:ALA:HB3	1:A:525:ILE:HG13	1.82	0.60
1:B:50:LEU:HD23	1:B:57:ALA:HB1	1.83	0.60
1:B:522:ALA:HB3	1:B:525:ILE:HG13	1.83	0.60
1:C:522:ALA:HB3	1:C:525:ILE:HG13	1.83	0.60
1:D:517:PHE:HB2	1:D:791:ASN:OD1	2.02	0.60
1:A:795:VAL:HG21	1:D:611:ILE:HG21	1.84	0.59
1:A:235:ILE:HD13	1:A:242:VAL:HG21	1.85	0.59
1:A:628:ARG:HH12	1:D:627:GLU:H	1.50	0.59
1:C:33:MET:HE1	1:C:43:LEU:HB2	1.86	0.58
1:D:235:ILE:HD13	1:D:242:VAL:HG21	1.86	0.58
1:B:10:ASN:HD22	1:B:10:ASN:N	2.01	0.58
1:C:475:ALA:HB3	1:C:735:ALA:HB3	1.84	0.58
1:D:10:ASN:N	1:D:10:ASN:HD22	2.02	0.58
1:B:597:SER:OG	1:C:810:ALA:HB2	2.04	0.58
1:A:33:MET:HE1	1:A:43:LEU:HB2	1.86	0.58
1:D:33:MET:HE1	1:D:43:LEU:HB2	1.86	0.58
1:C:235:ILE:HD13	1:C:242:VAL:HG21	1.86	0.58
1:B:235:ILE:HD13	1:B:242:VAL:HG21	1.85	0.58
1:D:629:MET:HE2	1:D:629:MET:HA	1.84	0.58
1:A:93:HIS:ND1	1:A:322:PRO:HB3	2.19	0.57
1:D:633:ILE:HD13	1:D:639:LEU:HD21	1.86	0.57
1:A:10:ASN:N	1:A:10:ASN:HD22	2.02	0.57
1:A:99:PRO:HA	1:A:113:MET:HB2	1.87	0.57
1:B:629:MET:HE2	1:B:629:MET:HA	1.84	0.57
1:D:113:MET:HE1	1:D:347:PHE:HZ	1.70	0.57
1:B:33:MET:HE1	1:B:43:LEU:HB2	1.86	0.57
1:B:489:ILE:HD12	1:B:735:ALA:HB1	1.87	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:77:TYR:CE2	1:C:98:THR:HG21	2.40	0.56
1:C:521:LEU:HD22	1:C:526:TRP:CG	2.40	0.56
1:D:99:PRO:HA	1:D:113:MET:HB2	1.87	0.56
1:C:10:ASN:N	1:C:10:ASN:HD22	2.02	0.56
1:C:517:PHE:HB2	1:C:791:ASN:OD1	2.05	0.56
1:B:227:PHE:CD1	1:B:244:GLY:HA3	2.41	0.56
1:B:517:PHE:HB2	1:B:791:ASN:OD1	2.06	0.56
1:C:209:ILE:HA	1:C:214:HIS:CD2	2.41	0.56
1:D:77:TYR:CE2	1:D:98:THR:HG21	2.40	0.56
1:D:227:PHE:CD1	1:D:244:GLY:HA3	2.41	0.56
1:D:521:LEU:HD22	1:D:526:TRP:CG	2.41	0.56
1:A:227:PHE:CD1	1:A:244:GLY:HA3	2.41	0.56
1:A:376:GLU:HG3	1:A:377:VAL:HG13	1.88	0.56
1:C:520:PRO:HA	1:C:623:PHE:HE2	1.71	0.56
1:A:77:TYR:CE2	1:A:98:THR:HG21	2.40	0.56
1:B:521:LEU:HD22	1:B:526:TRP:CG	2.41	0.56
1:B:626:VAL:CB	1:C:628:ARG:HH11	2.08	0.56
1:C:227:PHE:CD1	1:C:244:GLY:HA3	2.41	0.56
1:A:209:ILE:HA	1:A:214:HIS:CD2	2.41	0.55
1:B:99:PRO:HA	1:B:113:MET:HB2	1.87	0.55
1:B:209:ILE:HA	1:B:214:HIS:CD2	2.41	0.55
1:B:299:LEU:HD11	1:B:332:ALA:HB2	1.88	0.55
1:B:626:VAL:HB	1:C:628:ARG:NH1	2.08	0.55
1:A:521:LEU:HD22	1:A:526:TRP:CG	2.41	0.55
1:C:91:THR:HG21	1:D:56:PHE:CE2	2.41	0.55
1:C:376:GLU:HG3	1:C:377:VAL:HG13	1.88	0.55
1:D:209:ILE:HA	1:D:214:HIS:CD2	2.41	0.55
1:C:99:PRO:HA	1:C:113:MET:HB2	1.87	0.55
1:C:489:ILE:HD12	1:C:735:ALA:HB1	1.88	0.55
1:B:77:TYR:CE2	1:B:98:THR:HG21	2.41	0.55
1:D:376:GLU:HG3	1:D:377:VAL:HG13	1.88	0.55
1:B:765:LYS:O	1:B:770:LYS:HB2	2.07	0.55
1:B:376:GLU:HG3	1:B:377:VAL:HG13	1.88	0.54
1:D:765:LYS:O	1:D:770:LYS:HB2	2.07	0.54
1:A:753:LEU:HD22	1:A:758:LEU:HD13	1.89	0.54
1:C:753:LEU:HD22	1:C:758:LEU:HD13	1.89	0.54
1:A:62:PHE:HE2	1:A:92:LEU:HD12	1.73	0.54
1:A:765:LYS:O	1:A:770:LYS:HB2	2.07	0.54
1:D:753:LEU:HD22	1:D:758:LEU:HD13	1.89	0.54
1:D:475:ALA:HB3	1:D:735:ALA:HB3	1.89	0.54
1:C:62:PHE:HE2	1:C:92:LEU:HD12	1.73	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:754:SER:HB3	1:A:759:LEU:HD12	1.90	0.54
1:B:753:LEU:HD22	1:B:758:LEU:HD13	1.89	0.53
1:B:754:SER:HB3	1:B:759:LEU:HD12	1.90	0.53
1:D:754:SER:HB3	1:D:759:LEU:HD12	1.90	0.53
1:C:754:SER:HB3	1:C:759:LEU:HD12	1.90	0.53
1:C:765:LYS:O	1:C:770:LYS:HB2	2.08	0.53
1:A:337:GLN:HE22	2:E:1:NAG:H2	1.74	0.53
1:B:633:ILE:HD13	1:B:639:LEU:HD21	1.91	0.53
1:A:113:MET:HE1	1:A:347:PHE:HZ	1.73	0.53
1:D:62:PHE:HE2	1:D:92:LEU:HD12	1.73	0.52
1:B:205:VAL:O	1:B:209:ILE:HG13	2.10	0.52
1:B:523:TYR:O	1:B:527:MET:HG2	2.10	0.52
1:C:205:VAL:O	1:C:209:ILE:HG13	2.09	0.52
1:C:523:TYR:O	1:C:527:MET:HG2	2.10	0.52
1:A:205:VAL:O	1:A:209:ILE:HG13	2.10	0.52
1:C:113:MET:HE2	1:C:113:MET:HA	1.92	0.52
1:C:360:ILE:HG12	1:C:380:MET:HE2	1.91	0.52
1:D:205:VAL:O	1:D:209:ILE:HG13	2.09	0.52
1:A:113:MET:HA	1:A:113:MET:HE2	1.92	0.52
1:D:405:TYR:CD1	1:D:478:PRO:HG3	2.45	0.52
1:D:523:TYR:O	1:D:527:MET:HG2	2.10	0.52
1:D:405:TYR:CG	1:D:478:PRO:HG3	2.44	0.52
1:D:504:ILE:HG21	1:D:633:ILE:HD12	1.90	0.52
1:A:360:ILE:HG12	1:A:380:MET:HE2	1.91	0.52
1:A:611:ILE:HG21	1:B:795:VAL:HG21	1.92	0.52
1:B:62:PHE:HE2	1:B:92:LEU:HD12	1.72	0.52
1:B:360:ILE:HG12	1:B:380:MET:HE2	1.91	0.52
1:B:113:MET:HE2	1:B:113:MET:HA	1.91	0.52
1:C:405:TYR:CG	1:C:478:PRO:HG3	2.45	0.52
1:D:360:ILE:HG12	1:D:380:MET:HE2	1.91	0.52
1:D:337:GLN:HE22	3:H:1:NAG:H2	1.74	0.51
1:B:405:TYR:CG	1:B:478:PRO:HG3	2.45	0.51
1:D:113:MET:HE1	1:D:347:PHE:CZ	2.45	0.51
1:A:337:GLN:NE2	2:E:1:NAG:H2	2.25	0.51
1:A:405:TYR:CD1	1:A:478:PRO:HG3	2.45	0.51
1:A:405:TYR:CG	1:A:478:PRO:HG3	2.45	0.51
1:A:628:ARG:HH11	1:D:626:VAL:HB	1.74	0.51
1:A:633:ILE:HD11	1:A:645:ILE:HD12	1.92	0.51
1:C:83:ASN:ND2	1:D:80:LYS:HA	2.26	0.51
1:A:523:TYR:O	1:A:527:MET:HG2	2.09	0.51
1:B:494:PRO:HA	1:B:732:TYR:O	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:780:SER:C	1:A:782:GLU:H	2.19	0.51
1:C:74:PHE:CZ	1:C:285:THR:HG23	2.46	0.51
1:C:405:TYR:CD1	1:C:478:PRO:HG3	2.45	0.51
1:D:290:GLN:HG2	1:D:338:VAL:HG21	1.93	0.51
1:B:515:PHE:HD1	1:B:518:LEU:HD12	1.76	0.51
1:D:337:GLN:NE2	3:H:1:NAG:H2	2.26	0.51
1:C:494:PRO:HA	1:C:732:TYR:O	2.11	0.51
1:B:74:PHE:CZ	1:B:285:THR:HG23	2.46	0.50
1:D:74:PHE:CZ	1:D:285:THR:HG23	2.46	0.50
1:A:74:PHE:CZ	1:A:285:THR:HG23	2.47	0.50
1:D:113:MET:HA	1:D:113:MET:HE2	1.92	0.50
1:B:611:ILE:HG21	1:C:795:VAL:HG21	1.93	0.50
1:C:515:PHE:HD1	1:C:518:LEU:HD12	1.76	0.50
1:B:405:TYR:CD1	1:B:478:PRO:HG3	2.46	0.50
1:A:608:PHE:HZ	1:B:796:PHE:CZ	2.30	0.50
1:C:43:LEU:O	1:C:45:PRO:HD3	2.12	0.50
1:D:43:LEU:O	1:D:45:PRO:HD3	2.12	0.49
1:B:475:ALA:HB3	1:B:735:ALA:HB3	1.93	0.49
1:A:43:LEU:O	1:A:45:PRO:HD3	2.12	0.49
1:A:489:ILE:HD12	1:A:735:ALA:HB1	1.94	0.49
1:A:494:PRO:HA	1:A:732:TYR:O	2.12	0.49
1:A:515:PHE:HD1	1:A:518:LEU:HD12	1.76	0.49
1:A:56:PHE:CE2	1:B:91:THR:HG21	2.47	0.49
1:A:445:VAL:HG13	1:A:448:GLY:HA2	1.95	0.49
1:C:659:PHE:HB3	1:C:671:TRP:HB2	1.95	0.49
1:C:127:TYR:CE2	1:C:382:LEU:HD21	2.47	0.49
1:B:526:TRP:O	1:B:529:ILE:HG22	2.12	0.49
1:B:536:VAL:HG22	1:C:803:LEU:HD21	1.94	0.49
1:D:49:ASN:C	1:D:50:LEU:HD12	2.38	0.49
1:A:49:ASN:C	1:A:50:LEU:HD12	2.38	0.49
1:A:80:LYS:HA	1:B:83:ASN:ND2	2.27	0.49
1:A:517:PHE:HB2	1:A:791:ASN:OD1	2.12	0.49
1:B:43:LEU:O	1:B:45:PRO:HD3	2.11	0.49
1:A:526:TRP:O	1:A:529:ILE:HG22	2.13	0.49
1:D:515:PHE:HD1	1:D:518:LEU:HD12	1.76	0.49
1:D:526:TRP:O	1:D:529:ILE:HG22	2.13	0.49
1:A:50:LEU:HD22	1:A:61:ALA:HB2	1.95	0.48
1:B:50:LEU:HD22	1:B:61:ALA:HB2	1.95	0.48
1:B:659:PHE:HB3	1:B:671:TRP:HB2	1.95	0.48
1:C:526:TRP:O	1:C:529:ILE:HG22	2.13	0.48
1:C:49:ASN:C	1:C:50:LEU:HD12	2.38	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:494:PRO:HA	1:D:732:TYR:O	2.14	0.48
1:A:659:PHE:HB3	1:A:671:TRP:HB2	1.95	0.48
1:C:619:ASN:HA	1:D:624:LEU:HD13	1.95	0.48
1:A:628:ARG:CZ	1:D:623:PHE:HD1	2.27	0.48
1:B:539:VAL:HG13	1:C:807:MET:SD	2.53	0.48
1:D:93:HIS:ND1	1:D:322:PRO:HB3	2.28	0.48
1:D:664:ILE:HB	1:D:667:PHE:HD2	1.79	0.48
1:C:258:PHE:HD2	1:C:259:ILE:HD12	1.79	0.48
1:C:445:VAL:HG13	1:C:448:GLY:HA2	1.95	0.48
1:C:521:LEU:HD22	1:C:526:TRP:CE2	2.49	0.48
1:A:664:ILE:HB	1:A:667:PHE:HD2	1.78	0.48
1:B:445:VAL:HG13	1:B:448:GLY:HA2	1.95	0.48
1:D:50:LEU:HD22	1:D:61:ALA:HB2	1.95	0.48
1:D:659:PHE:HB3	1:D:671:TRP:HB2	1.95	0.48
1:D:195:ASP:HA	1:D:223:ALA:HB3	1.95	0.47
1:D:258:PHE:HD2	1:D:259:ILE:HD12	1.79	0.47
1:C:664:ILE:HB	1:C:667:PHE:HD2	1.78	0.47
1:A:97:ILE:HG13	1:A:111:ILE:HB	1.97	0.47
1:B:664:ILE:HB	1:B:667:PHE:HD2	1.79	0.47
1:D:711:TYR:O	1:D:715:ARG:HG2	2.14	0.47
1:B:49:ASN:C	1:B:50:LEU:HD12	2.38	0.47
1:B:195:ASP:HA	1:B:223:ALA:HB3	1.96	0.47
1:C:29:PHE:O	1:C:33:MET:HG2	2.15	0.47
1:C:50:LEU:HD22	1:C:61:ALA:HB2	1.95	0.47
1:A:711:TYR:O	1:A:715:ARG:HG2	2.14	0.47
1:B:97:ILE:HG13	1:B:111:ILE:HB	1.96	0.47
1:D:97:ILE:HG13	1:D:111:ILE:HB	1.97	0.47
1:D:530:VAL:O	1:D:533:TYR:HB3	2.15	0.47
1:B:208:VAL:HG12	1:B:214:HIS:HB3	1.97	0.47
1:A:258:PHE:HD2	1:A:259:ILE:HD12	1.79	0.47
1:B:258:PHE:HD2	1:B:259:ILE:HD12	1.79	0.47
1:B:530:VAL:O	1:B:533:TYR:HB3	2.15	0.47
1:C:195:ASP:HA	1:C:223:ALA:HB3	1.96	0.47
1:C:530:VAL:O	1:C:533:TYR:HB3	2.15	0.47
1:A:512:PRO:HB2	1:A:513:GLY:H	1.51	0.47
1:A:525:ILE:HG12	1:B:789:LEU:CD1	2.40	0.47
1:D:445:VAL:HG13	1:D:448:GLY:HA2	1.95	0.47
1:B:711:TYR:O	1:B:715:ARG:HG2	2.14	0.47
1:C:208:VAL:HG12	1:C:214:HIS:HB3	1.97	0.47
1:C:467:LEU:HD22	1:C:737:PRO:HD3	1.96	0.47
1:A:195:ASP:HA	1:A:223:ALA:HB3	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:29:PHE:O	1:B:33:MET:HG2	2.14	0.47
1:C:117:LEU:HD12	1:C:120:ALA:HB3	1.97	0.47
1:C:299:LEU:HD11	1:C:332:ALA:HB2	1.97	0.46
1:D:22:ALA:HB1	1:D:25:GLU:HB2	1.98	0.46
1:C:113:MET:HE1	1:C:347:PHE:HZ	1.79	0.46
1:C:711:TYR:O	1:C:715:ARG:HG2	2.14	0.46
1:A:530:VAL:O	1:A:533:TYR:HB3	2.15	0.46
1:D:29:PHE:O	1:D:33:MET:HG2	2.15	0.46
1:A:62:PHE:CE2	1:A:92:LEU:HD12	2.51	0.46
1:D:135:TYR:CE1	1:D:137:TYR:HB3	2.51	0.46
1:D:208:VAL:HG12	1:D:214:HIS:HB3	1.98	0.46
1:D:23:ASP:HB3	1:D:271:PRO:HB2	1.98	0.46
1:D:490:ASP:HB2	1:D:736:THR:HG23	1.98	0.46
1:C:135:TYR:CE1	1:C:137:TYR:HB3	2.50	0.46
1:A:22:ALA:HB1	1:A:25:GLU:HB2	1.98	0.46
1:A:98:THR:HA	1:A:99:PRO:HD3	1.81	0.46
1:A:521:LEU:CD2	1:A:525:ILE:HB	2.45	0.46
1:A:135:TYR:CE1	1:A:137:TYR:HB3	2.50	0.46
1:A:208:VAL:HG12	1:A:214:HIS:HB3	1.97	0.46
1:D:683:VAL:HG11	1:D:689:GLY:HA2	1.98	0.46
1:A:29:PHE:O	1:A:33:MET:HG2	2.15	0.46
1:A:628:ARG:HH12	1:D:627:GLU:N	2.14	0.46
1:B:135:TYR:CE1	1:B:137:TYR:HB3	2.50	0.46
1:C:97:ILE:HG13	1:C:111:ILE:HB	1.97	0.46
1:D:117:LEU:HD12	1:D:120:ALA:HB3	1.98	0.46
1:C:22:ALA:HB1	1:C:25:GLU:HB2	1.97	0.45
1:C:23:ASP:HB3	1:C:271:PRO:HB2	1.98	0.45
1:C:683:VAL:HG11	1:C:689:GLY:HA2	1.98	0.45
1:D:521:LEU:HD22	1:D:526:TRP:CE2	2.52	0.45
1:A:23:ASP:HB3	1:A:271:PRO:HB2	1.98	0.45
1:A:117:LEU:HD12	1:A:120:ALA:HB3	1.98	0.45
1:A:629:MET:HE2	1:C:629:MET:HE2	1.98	0.45
1:B:113:MET:HE1	1:B:347:PHE:HZ	1.81	0.45
1:A:519:ASP:HB2	1:A:520:PRO:HD3	1.98	0.45
1:A:678:GLU:HA	1:A:679:PRO:C	2.41	0.45
1:B:536:VAL:HG21	1:B:605:TRP:CE3	2.51	0.45
1:C:619:ASN:CA	1:D:624:LEU:HD13	2.46	0.45
1:B:22:ALA:HB1	1:B:25:GLU:HB2	1.97	0.45
1:B:209:ILE:CD1	1:B:234:LYS:HB2	2.47	0.45
1:D:62:PHE:CE2	1:D:92:LEU:HD12	2.51	0.45
1:D:678:GLU:HA	1:D:679:PRO:C	2.41	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:125:ILE:HG23	1:A:130:TRP:HB2	1.99	0.45
1:A:209:ILE:CD1	1:A:234:LYS:HB2	2.46	0.45
1:B:323:TRP:CD1	1:B:323:TRP:H	2.35	0.45
1:C:62:PHE:CE2	1:C:92:LEU:HD12	2.51	0.45
1:A:628:ARG:HB2	1:D:626:VAL:HG21	1.98	0.45
1:B:62:PHE:CE2	1:B:88:PHE:HB3	2.52	0.45
1:B:165:VAL:HG22	1:B:165:VAL:O	2.17	0.45
1:D:519:ASP:HB2	1:D:520:PRO:HD3	1.98	0.45
1:B:23:ASP:HB3	1:B:271:PRO:HB2	1.98	0.45
1:B:521:LEU:HD22	1:B:526:TRP:CE2	2.52	0.45
1:D:261:ARG:O	1:D:265:LEU:HG	2.17	0.45
1:C:209:ILE:CD1	1:C:234:LYS:HB2	2.46	0.45
1:C:323:TRP:H	1:C:323:TRP:CD1	2.35	0.45
1:D:62:PHE:CE2	1:D:88:PHE:HB3	2.52	0.45
1:B:490:ASP:HB2	1:B:736:THR:HG23	1.98	0.45
1:B:521:LEU:CD2	1:B:525:ILE:HB	2.46	0.45
1:B:678:GLU:HA	1:B:679:PRO:C	2.41	0.45
1:B:683:VAL:HG11	1:B:689:GLY:HA2	1.99	0.45
1:B:62:PHE:CE2	1:B:92:LEU:HD12	2.51	0.44
1:B:601:VAL:HG23	1:C:806:ALA:CB	2.46	0.44
1:D:521:LEU:CD2	1:D:525:ILE:HB	2.46	0.44
1:B:117:LEU:HD12	1:B:120:ALA:HB3	1.98	0.44
1:C:98:THR:HA	1:C:99:PRO:HD3	1.80	0.44
1:C:519:ASP:HB2	1:C:520:PRO:HD3	1.98	0.44
1:A:515:PHE:CD1	1:A:518:LEU:HD12	2.52	0.44
1:A:521:LEU:HD22	1:A:526:TRP:CE2	2.52	0.44
1:B:130:TRP:CH2	1:B:191:ARG:HB3	2.53	0.44
1:C:261:ARG:O	1:C:265:LEU:HG	2.17	0.44
1:C:678:GLU:HA	1:C:679:PRO:C	2.41	0.44
1:A:62:PHE:CE2	1:A:88:PHE:HB3	2.52	0.44
1:A:96:PHE:CE2	1:A:98:THR:HB	2.53	0.44
1:A:113:MET:HE1	1:A:347:PHE:CZ	2.51	0.44
1:A:130:TRP:CE2	1:A:191:ARG:HD3	2.53	0.44
1:A:165:VAL:HG22	1:A:165:VAL:O	2.18	0.44
1:A:683:VAL:HG11	1:A:689:GLY:HA2	1.99	0.44
1:B:98:THR:HA	1:B:99:PRO:HD3	1.80	0.44
1:C:96:PHE:CE2	1:C:98:THR:HB	2.53	0.44
1:D:96:PHE:CE2	1:D:98:THR:HB	2.53	0.44
1:D:504:ILE:CG2	1:D:633:ILE:HD12	2.48	0.44
1:C:512:PRO:HB2	1:C:513:GLY:H	1.49	0.44
1:D:209:ILE:CD1	1:D:234:LYS:HB2	2.47	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:125:ILE:HG23	1:D:130:TRP:HB2	1.99	0.44
1:D:287:ASP:O	1:D:291:VAL:HG23	2.18	0.44
2:F:1:NAG:H61	2:F:2:NAG:H83	1.99	0.44
1:C:521:LEU:HB3	1:C:526:TRP:CE2	2.53	0.44
1:A:287:ASP:O	1:A:291:VAL:HG23	2.18	0.44
1:B:96:PHE:CE2	1:B:98:THR:HB	2.53	0.44
1:B:519:ASP:HB2	1:B:520:PRO:HD3	1.98	0.44
4:C:833:ZK1:HAI	4:C:833:ZK1:HAOA	1.82	0.44
1:D:165:VAL:HG22	1:D:165:VAL:O	2.18	0.44
1:A:261:ARG:O	1:A:265:LEU:HG	2.17	0.44
1:C:130:TRP:CE2	1:C:191:ARG:HD3	2.53	0.44
1:B:261:ARG:O	1:B:265:LEU:HG	2.17	0.43
1:C:62:PHE:CE2	1:C:88:PHE:HB3	2.52	0.43
1:C:165:VAL:O	1:C:165:VAL:HG22	2.18	0.43
1:C:787:LEU:HA	1:C:787:LEU:HD22	1.74	0.43
1:D:515:PHE:CD1	1:D:518:LEU:HD12	2.52	0.43
1:B:287:ASP:O	1:B:291:VAL:HG23	2.18	0.43
1:B:525:ILE:HD13	1:C:792:VAL:HG21	1.99	0.43
1:B:623:PHE:CZ	1:C:786:ALA:HB2	2.54	0.43
1:C:125:ILE:HG23	1:C:130:TRP:HB2	1.99	0.43
1:C:521:LEU:CD2	1:C:525:ILE:HB	2.47	0.43
1:B:125:ILE:HG23	1:B:130:TRP:HB2	1.99	0.43
1:B:270:TYR:HA	1:B:271:PRO:HD2	1.87	0.43
1:A:323:TRP:H	1:A:323:TRP:CD1	2.35	0.43
1:B:515:PHE:CD1	1:B:518:LEU:HD12	2.52	0.43
1:C:97:ILE:N	1:C:97:ILE:HD12	2.34	0.43
1:B:774:GLY:O	1:B:778:SER:N	2.47	0.43
1:C:488:VAL:HG23	1:C:489:ILE:HG23	2.00	0.43
1:C:515:PHE:CD1	1:C:518:LEU:HD12	2.52	0.43
1:D:620:LEU:HA	1:D:623:PHE:HD2	1.82	0.43
1:B:488:VAL:HG23	1:B:489:ILE:HG23	2.00	0.43
1:C:93:HIS:ND1	1:C:322:PRO:HB3	2.33	0.43
1:C:287:ASP:O	1:C:291:VAL:HG23	2.18	0.43
1:A:209:ILE:HD11	1:A:235:ILE:HG23	2.01	0.43
1:B:321:VAL:HA	1:B:322:PRO:HD3	1.86	0.43
1:B:360:ILE:HD11	1:B:374:TRP:HB2	2.01	0.43
1:C:360:ILE:HD11	1:C:374:TRP:HB2	2.01	0.43
1:A:130:TRP:CH2	1:A:191:ARG:HB3	2.53	0.43
1:C:321:VAL:HA	1:C:322:PRO:HD3	1.86	0.43
1:D:323:TRP:H	1:D:323:TRP:CD1	2.36	0.43
1:A:97:ILE:N	1:A:97:ILE:HD12	2.34	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:628:ARG:NH1	1:D:623:PHE:HA	2.33	0.43
1:C:525:ILE:HD13	1:D:792:VAL:HG21	2.00	0.43
1:D:97:ILE:HD12	1:D:97:ILE:N	2.34	0.43
1:D:130:TRP:CE2	1:D:191:ARG:HD3	2.53	0.43
1:A:508:GLN:C	1:A:510:SER:H	2.26	0.42
1:B:130:TRP:CE2	1:B:191:ARG:HD3	2.54	0.42
1:B:708:MET:O	1:B:712:ILE:HG12	2.19	0.42
1:D:321:VAL:HA	1:D:322:PRO:HD3	1.86	0.42
1:A:786:ALA:HB1	1:D:519:ASP:O	2.19	0.42
1:B:97:ILE:HD12	1:B:97:ILE:N	2.34	0.42
1:C:130:TRP:CH2	1:C:191:ARG:HB3	2.53	0.42
1:C:508:GLN:C	1:C:510:SER:H	2.27	0.42
1:D:209:ILE:HD11	1:D:235:ILE:HG23	2.01	0.42
1:D:360:ILE:HD11	1:D:374:TRP:HB2	2.01	0.42
1:B:233:LEU:HD23	1:B:236:GLN:OE1	2.19	0.42
1:B:521:LEU:HB3	1:B:526:TRP:CE2	2.55	0.42
1:C:233:LEU:HD23	1:C:236:GLN:OE1	2.20	0.42
1:D:233:LEU:HD23	1:D:236:GLN:OE1	2.20	0.42
1:A:290:GLN:HG2	1:A:338:VAL:HG21	2.00	0.42
1:A:795:VAL:HG12	1:D:608:PHE:CD1	2.55	0.42
1:B:512:PRO:HB2	1:B:513:GLY:H	1.66	0.42
1:B:623:PHE:HZ	1:C:786:ALA:HB2	1.85	0.42
1:C:209:ILE:HD11	1:C:235:ILE:HG23	2.01	0.42
1:C:809:VAL:O	1:C:813:GLU:HG3	2.19	0.42
1:A:101:PHE:HA	1:A:102:PRO:HD3	1.94	0.42
1:A:488:VAL:HG23	1:A:489:ILE:HG23	2.00	0.42
1:C:805:LEU:O	1:C:809:VAL:HG23	2.20	0.42
1:D:507:PRO:HB2	1:D:629:MET:SD	2.60	0.42
1:D:130:TRP:CH2	1:D:191:ARG:HB3	2.54	0.42
1:D:488:VAL:HG23	1:D:489:ILE:HG23	2.00	0.42
1:D:708:MET:O	1:D:712:ILE:HG12	2.20	0.42
1:A:360:ILE:HD11	1:A:374:TRP:HB2	2.01	0.42
1:A:795:VAL:HG12	1:D:608:PHE:HD1	1.85	0.42
1:C:533:TYR:CZ	1:C:583:ALA:HB1	2.54	0.42
1:D:521:LEU:HB3	1:D:526:TRP:CE2	2.54	0.42
1:A:533:TYR:CZ	1:A:583:ALA:HB1	2.55	0.42
1:B:787:LEU:HD13	1:B:788:SER:O	2.19	0.42
1:D:299:LEU:HD11	1:D:332:ALA:HB2	2.01	0.42
1:D:513:GLY:C	1:D:515:PHE:H	2.28	0.42
1:A:809:VAL:O	1:A:813:GLU:HG3	2.19	0.42
1:B:809:VAL:O	1:B:813:GLU:HG3	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:233:LEU:HD23	1:A:236:GLN:OE1	2.19	0.41
1:A:521:LEU:HB3	1:A:526:TRP:CE2	2.54	0.41
1:B:232:LEU:O	1:B:236:GLN:HB2	2.20	0.41
1:D:512:PRO:HB2	1:D:513:GLY:H	1.51	0.41
1:B:633:ILE:HD13	1:B:639:LEU:CD2	2.49	0.41
1:C:232:LEU:O	1:C:236:GLN:HB2	2.20	0.41
1:A:803:LEU:O	1:A:807:MET:HG2	2.21	0.41
1:B:513:GLY:C	1:B:515:PHE:H	2.28	0.41
1:D:803:LEU:O	1:D:807:MET:HG2	2.21	0.41
1:A:232:LEU:O	1:A:236:GLN:HB2	2.20	0.41
1:C:803:LEU:O	1:C:807:MET:HG2	2.21	0.41
1:D:631:SER:HA	1:D:632:PRO:HD3	1.86	0.41
1:D:809:VAL:O	1:D:813:GLU:HG3	2.19	0.41
1:B:803:LEU:O	1:B:807:MET:HG2	2.20	0.41
1:C:288:ALA:O	1:C:292:MET:HG3	2.21	0.41
1:C:624:LEU:O	1:C:628:ARG:HG3	2.21	0.41
4:D:833:ZK1:HAI	4:D:833:ZK1:HAOA	1.82	0.41
1:A:708:MET:O	1:A:712:ILE:HG12	2.20	0.41
1:C:619:ASN:OD1	1:D:787:LEU:HB2	2.20	0.41
1:D:380:MET:HE3	1:D:380:MET:HB2	1.95	0.41
1:A:467:LEU:HD22	1:A:737:PRO:HD3	2.03	0.41
1:B:209:ILE:HD11	1:B:235:ILE:HG23	2.01	0.41
1:B:288:ALA:O	1:B:292:MET:HG3	2.21	0.41
1:B:533:TYR:HA	1:B:605:TRP:CH2	2.56	0.41
1:C:18:PHE:HA	1:C:19:PRO:HD3	1.96	0.41
1:C:519:ASP:N	1:C:520:PRO:CD	2.84	0.41
1:C:520:PRO:HA	1:C:623:PHE:CE2	2.53	0.41
1:C:708:MET:O	1:C:712:ILE:HG12	2.20	0.41
1:D:519:ASP:N	1:D:520:PRO:CD	2.84	0.41
2:E:1:NAG:H61	2:E:2:NAG:H83	2.02	0.41
1:A:624:LEU:O	1:A:628:ARG:HG3	2.20	0.41
1:A:744:THR:HB	1:A:745:PRO:HD3	2.02	0.41
1:A:792:VAL:HG21	1:D:525:ILE:HD13	2.02	0.41
1:B:113:MET:HE1	1:B:347:PHE:CZ	2.55	0.41
1:C:101:PHE:HA	1:C:102:PRO:HD3	1.94	0.41
1:D:744:THR:HB	1:D:745:PRO:HD3	2.02	0.41
1:A:519:ASP:N	1:A:520:PRO:CD	2.84	0.41
1:B:127:TYR:CE2	1:B:382:LEU:HD21	2.56	0.41
1:B:608:PHE:HZ	1:C:796:PHE:CZ	2.38	0.41
1:B:805:LEU:O	1:B:809:VAL:HG23	2.20	0.41
1:C:375:SER:HB3	1:C:378:ASP:HB2	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:192:VAL:HB	1:D:220:TYR:CD1	2.56	0.41
1:A:375:SER:HB3	1:A:378:ASP:HB2	2.03	0.40
1:B:299:LEU:CD1	1:B:332:ALA:HB2	2.51	0.40
1:B:608:PHE:HD1	1:C:795:VAL:HG12	1.85	0.40
1:B:758:LEU:HD23	1:B:758:LEU:O	2.21	0.40
4:B:833:ZK1:HAOA	4:B:833:ZK1:HAI	1.81	0.40
1:D:288:ALA:O	1:D:292:MET:HG3	2.20	0.40
1:D:467:LEU:HD22	1:D:737:PRO:HD3	2.03	0.40
3:G:2:NAG:H83	3:G:2:NAG:H2	1.92	0.40
1:B:744:THR:HB	1:B:745:PRO:HD3	2.02	0.40
1:C:113:MET:HE1	1:C:347:PHE:CZ	2.55	0.40
1:C:513:GLY:C	1:C:515:PHE:H	2.29	0.40
1:D:232:LEU:O	1:D:236:GLN:HB2	2.21	0.40
1:D:235:ILE:H	1:D:235:ILE:HG13	1.73	0.40
1:A:192:VAL:HB	1:A:220:TYR:CD1	2.56	0.40
1:A:513:GLY:C	1:A:515:PHE:H	2.29	0.40
1:A:628:ARG:NH2	1:D:623:PHE:CD1	2.84	0.40
1:B:266:GLU:HG2	1:B:268:LYS:H	1.86	0.40
1:B:477:ALA:HB1	1:B:478:PRO:CD	2.51	0.40
1:C:124:LEU:HA	1:C:380:MET:HE1	2.02	0.40
1:A:477:ALA:HB1	1:A:478:PRO:CD	2.52	0.40
1:A:628:ARG:CZ	1:D:623:PHE:HA	2.52	0.40
1:A:631:SER:HA	1:A:632:PRO:HD3	1.92	0.40
1:B:231:ASP:HB3	1:B:234:LYS:HE2	2.04	0.40
1:B:519:ASP:N	1:B:520:PRO:CD	2.84	0.40
1:D:520:PRO:HA	1:D:623:PHE:HE2	1.86	0.40
1:A:805:LEU:O	1:A:809:VAL:HG23	2.21	0.40
1:B:212:GLY:HA3	1:D:215:VAL:HG12	2.02	0.40
1:B:380:MET:HE3	1:B:380:MET:HB2	1.95	0.40
1:D:266:GLU:HG2	1:D:268:LYS:H	1.86	0.40
1:D:387:THR:C	1:D:389:GLY:H	2.29	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	773/823 (94%)	709 (92%)	54 (7%)	10 (1%)	9	39
1	B	773/823 (94%)	712 (92%)	52 (7%)	9 (1%)	10	41
1	C	773/823 (94%)	714 (92%)	49 (6%)	10 (1%)	9	39
1	D	773/823 (94%)	710 (92%)	55 (7%)	8 (1%)	12	45
All	All	3092/3292 (94%)	2845 (92%)	210 (7%)	37 (1%)	10	41

All (37) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	172	LYS
1	B	172	LYS
1	B	512	PRO
1	C	172	LYS
1	D	172	LYS
1	A	386	ASP
1	A	390	LEU
1	A	393	LYS
1	A	512	PRO
1	A	513	GLY
1	B	513	GLY
1	C	512	PRO
1	C	785	SER
1	D	512	PRO
1	C	390	LEU
1	C	513	GLY
1	A	597	SER
1	B	384	GLU
1	B	597	SER
1	C	389	GLY
1	C	597	SER
1	D	597	SER
1	B	507	PRO
1	B	632	PRO
1	D	507	PRO
1	D	513	GLY
1	D	632	PRO
1	A	630	VAL
1	A	743	GLY
1	B	743	GLY

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Mol	Chain	Res	Type
1	C	630	VAL
1	C	743	GLY
1	D	743	GLY
1	A	451	GLY
1	B	451	GLY
1	C	451	GLY
1	D	451	GLY

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	534/697 (77%)	532 (100%)	2 (0%)	84	81
1	B	534/697 (77%)	533 (100%)	1 (0%)	87	85
1	C	534/697 (77%)	531 (99%)	3 (1%)	78	79
1	D	534/697 (77%)	533 (100%)	1 (0%)	87	85
All	All	2136/2788 (77%)	2129 (100%)	7 (0%)	86	83

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	299	LEU
1	A	394	THR
1	B	299	LEU
1	C	299	LEU
1	C	628	ARG
1	C	787	LEU
1	D	299	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	337	GLN
1	B	619	ASN

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Mol	Chain	Res	Type
1	D	337	GLN
1	D	344	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

14 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	NAG	E	1	2,1	14,14,15	0.48	0	17,19,21	0.71	0
2	NAG	E	2	2	14,14,15	0.50	0	17,19,21	0.76	1 (5%)
2	BMA	E	3	2	11,11,12	0.62	0	15,15,17	0.44	0
2	BMA	E	4	2	11,11,12	0.62	0	15,15,17	0.43	0
2	NAG	F	1	2,1	14,14,15	0.50	0	17,19,21	0.69	0
2	NAG	F	2	2	14,14,15	0.47	0	17,19,21	0.75	1 (5%)
2	BMA	F	3	2	11,11,12	0.62	0	15,15,17	0.50	0
2	BMA	F	4	2	11,11,12	0.65	0	15,15,17	0.54	0
3	NAG	G	1	1,3	14,14,15	0.48	0	17,19,21	0.92	1 (5%)
3	NAG	G	2	3	14,14,15	0.47	0	17,19,21	0.77	1 (5%)
3	BMA	G	3	3	11,11,12	0.63	0	15,15,17	0.47	0
3	NAG	H	1	1,3	14,14,15	0.50	0	17,19,21	0.67	0
3	NAG	H	2	3	14,14,15	0.48	0	17,19,21	0.74	1 (5%)
3	BMA	H	3	3	11,11,12	0.62	0	15,15,17	0.48	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
2	NAG	E	1	2,1	-	2/6/23/26	0/1/1/1
2	NAG	E	2	2	-	2/6/23/26	0/1/1/1
2	BMA	E	3	2	-	0/2/19/22	0/1/1/1
2	BMA	E	4	2	-	1/2/19/22	0/1/1/1
2	NAG	F	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	F	2	2	-	2/6/23/26	0/1/1/1
2	BMA	F	3	2	-	0/2/19/22	0/1/1/1
2	BMA	F	4	2	-	1/2/19/22	0/1/1/1
3	NAG	G	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	G	2	3	-	2/6/23/26	0/1/1/1
3	BMA	G	3	3	-	0/2/19/22	0/1/1/1
3	NAG	H	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	H	2	3	-	2/6/23/26	0/1/1/1
3	BMA	H	3	3	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	G	2	NAG	C1-O5-C5	2.29	115.26	112.19
3	G	1	NAG	C1-O5-C5	2.18	115.11	112.19
2	F	2	NAG	C1-O5-C5	2.14	115.05	112.19
3	H	2	NAG	C1-O5-C5	2.05	114.93	112.19
2	E	2	NAG	C1-O5-C5	2.04	114.93	112.19

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	E	1	NAG	C8-C7-N2-C2
2	E	1	NAG	O7-C7-N2-C2
2	E	2	NAG	C8-C7-N2-C2
2	E	2	NAG	O7-C7-N2-C2
2	F	2	NAG	C8-C7-N2-C2
2	F	2	NAG	O7-C7-N2-C2
3	G	2	NAG	C8-C7-N2-C2

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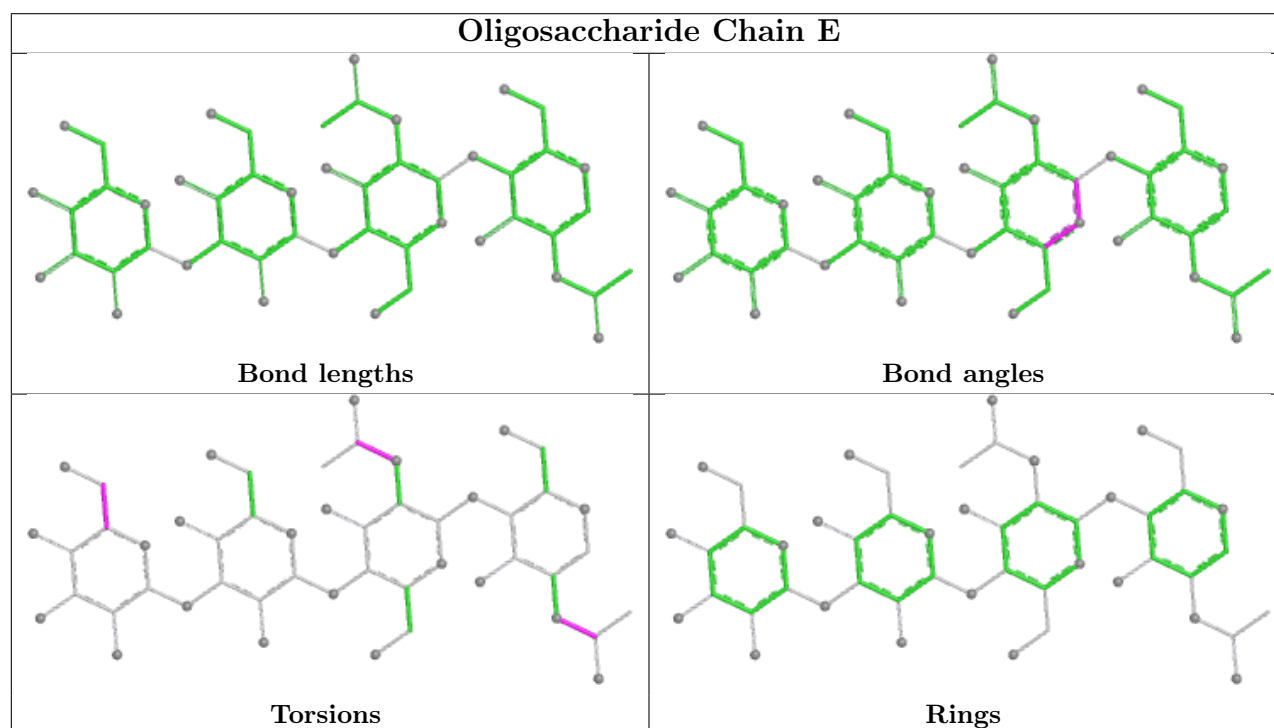
Mol	Chain	Res	Type	Atoms
3	G	2	NAG	O7-C7-N2-C2
3	H	2	NAG	C8-C7-N2-C2
3	H	2	NAG	O7-C7-N2-C2
3	H	1	NAG	C8-C7-N2-C2
3	H	1	NAG	O7-C7-N2-C2
3	G	1	NAG	C8-C7-N2-C2
3	G	1	NAG	O7-C7-N2-C2
2	F	4	BMA	O5-C5-C6-O6
2	E	4	BMA	O5-C5-C6-O6

There are no ring outliers.

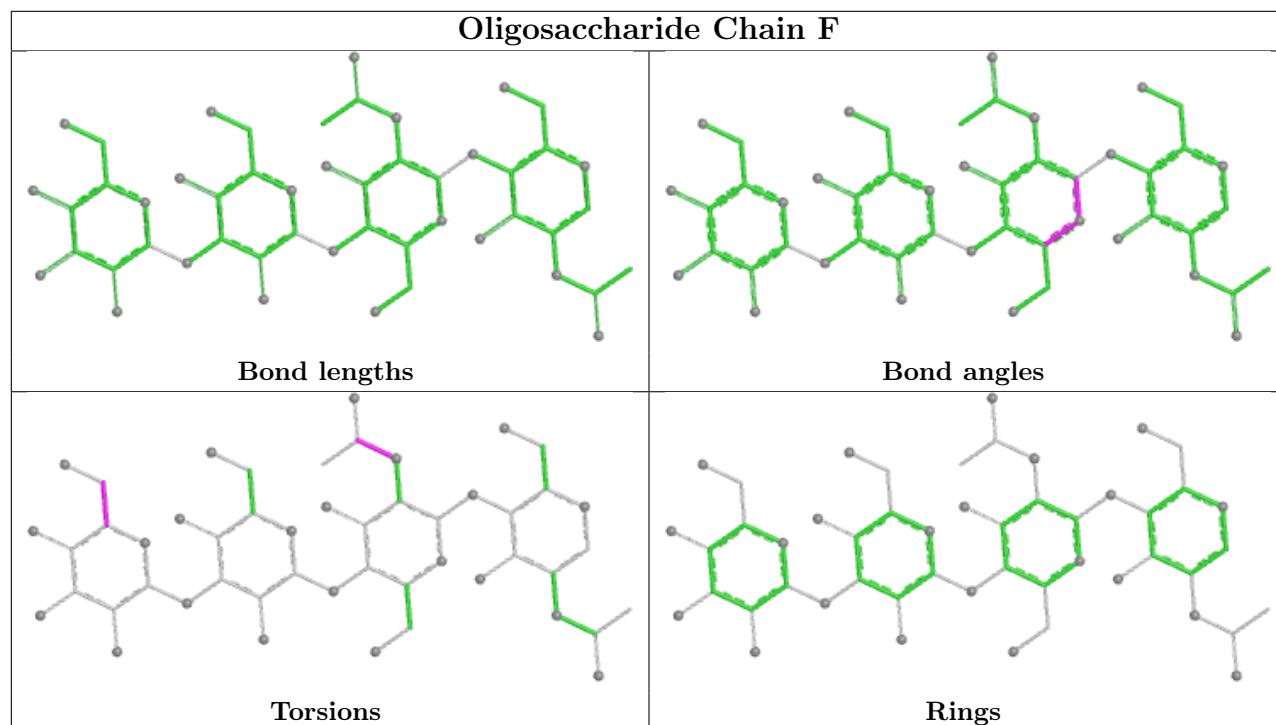
6 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	H	1	NAG	2	0
2	F	1	NAG	1	0
2	E	2	NAG	1	0
2	E	1	NAG	3	0
3	G	2	NAG	1	0
2	F	2	NAG	1	0

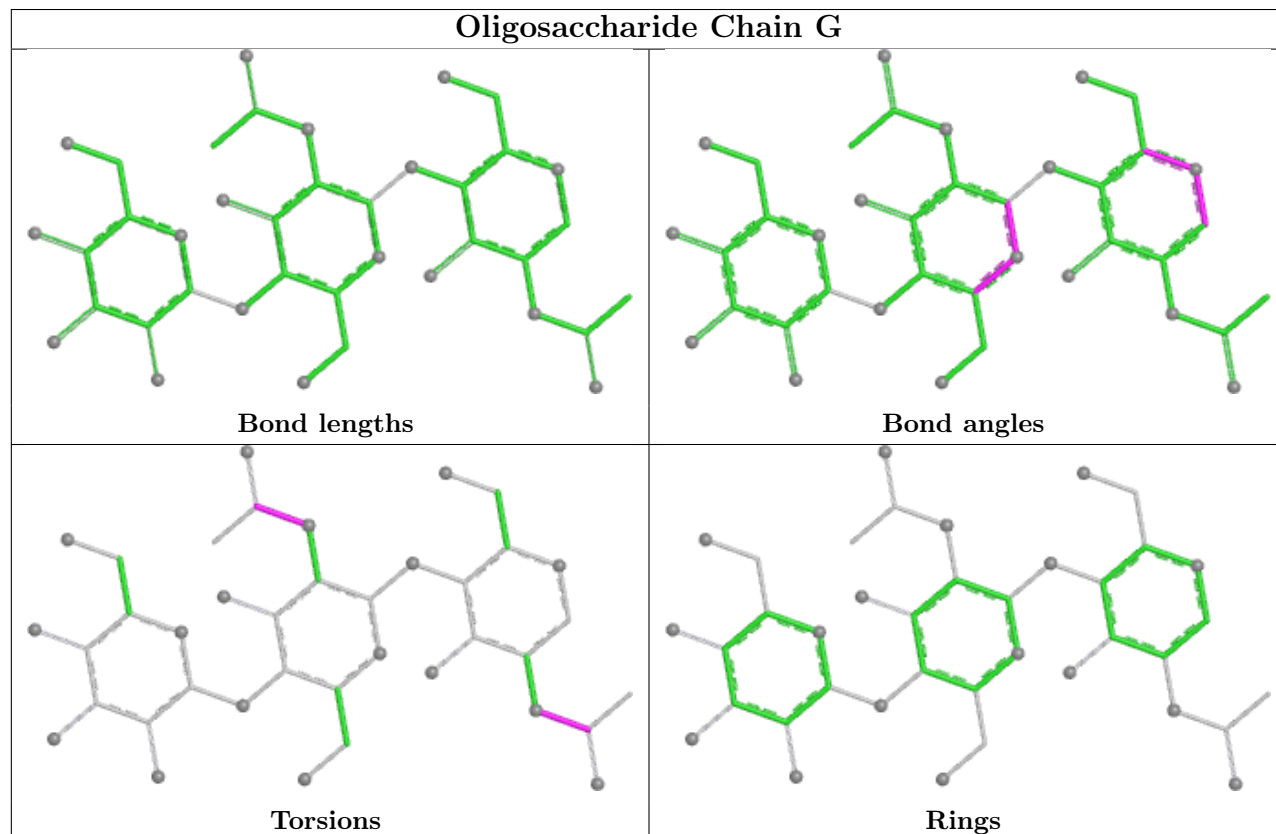
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

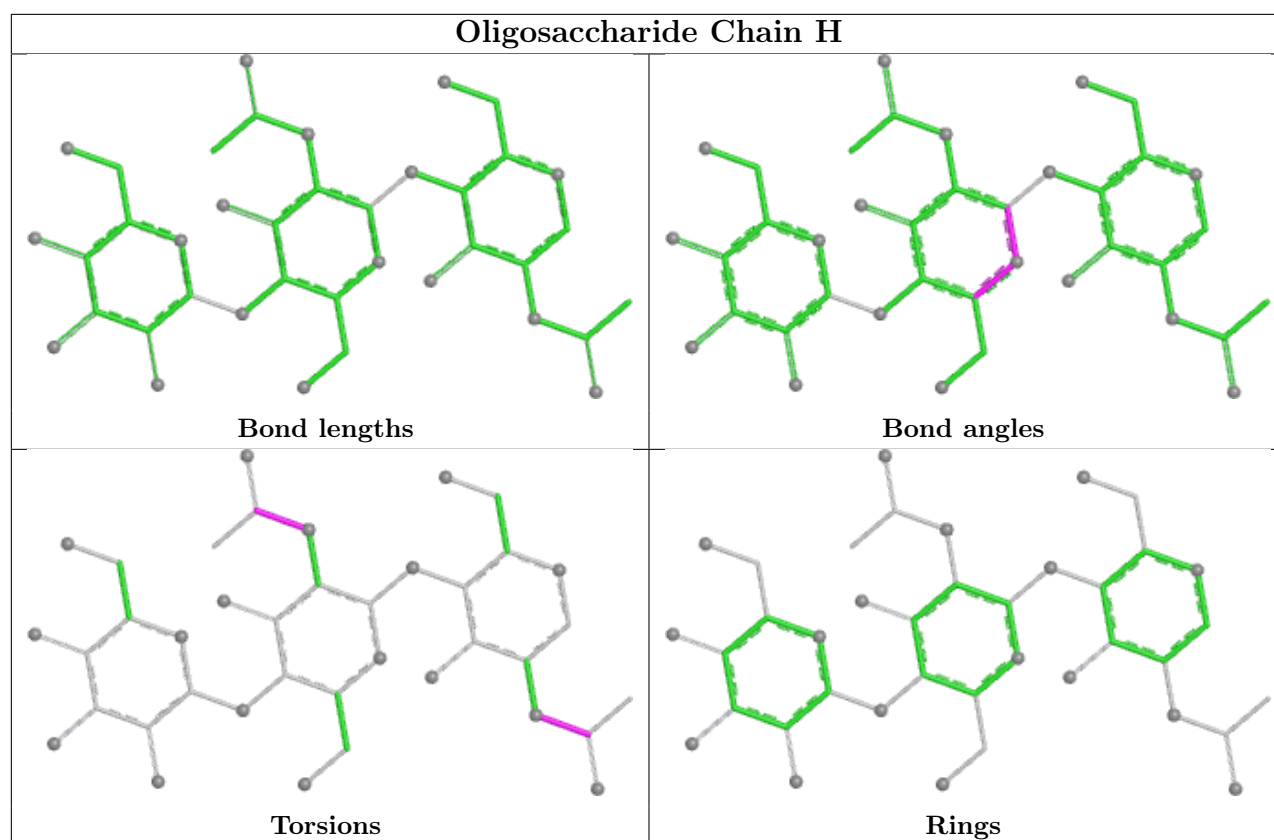


Oligosaccharide Chain F



Oligosaccharide Chain G





5.6 Ligand geometry [i](#)

4 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	ZK1	A	833	-	29,29,29	3.42	11 (37%)	45,45,45	1.58	7 (15%)
4	ZK1	C	833	-	29,29,29	3.41	11 (37%)	45,45,45	1.57	7 (15%)
4	ZK1	D	833	-	29,29,29	3.42	11 (37%)	45,45,45	1.57	7 (15%)
4	ZK1	B	833	-	29,29,29	3.42	10 (34%)	45,45,45	1.56	7 (15%)

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	ZK1	A	833	-	-	5/13/23/23	0/3/3/3
4	ZK1	C	833	-	-	5/13/23/23	0/3/3/3
4	ZK1	D	833	-	-	5/13/23/23	0/3/3/3
4	ZK1	B	833	-	-	5/13/23/23	0/3/3/3

All (43) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	833	ZK1	OAA-CAT	9.37	1.41	1.23
4	D	833	ZK1	OAA-CAT	9.34	1.41	1.23
4	B	833	ZK1	OAA-CAT	9.32	1.41	1.23
4	C	833	ZK1	OAA-CAT	9.25	1.41	1.23
4	B	833	ZK1	OAB-CAU	8.74	1.40	1.23
4	C	833	ZK1	OAB-CAU	8.74	1.40	1.23
4	D	833	ZK1	OAB-CAU	8.72	1.40	1.23
4	A	833	ZK1	OAB-CAU	8.70	1.40	1.23
4	A	833	ZK1	CAT-NAP	5.95	1.44	1.35
4	D	833	ZK1	CAT-NAP	5.90	1.44	1.35
4	C	833	ZK1	CAT-NAP	5.90	1.44	1.35
4	B	833	ZK1	CAT-NAP	5.88	1.44	1.35
4	B	833	ZK1	CAU-CAT	-5.55	1.45	1.53
4	A	833	ZK1	CAU-CAT	-5.50	1.45	1.53
4	C	833	ZK1	CAU-CAT	-5.49	1.45	1.53
4	D	833	ZK1	CAU-CAT	-5.47	1.45	1.53
4	D	833	ZK1	PBA-OAE	4.78	1.65	1.55
4	A	833	ZK1	PBA-OAE	4.76	1.65	1.55
4	B	833	ZK1	PBA-OAE	4.76	1.65	1.55
4	C	833	ZK1	PBA-OAE	4.73	1.65	1.55
4	A	833	ZK1	PBA-OAD	4.65	1.65	1.55
4	D	833	ZK1	PBA-OAD	4.64	1.65	1.55
4	C	833	ZK1	PBA-OAD	4.62	1.65	1.55
4	B	833	ZK1	PBA-OAD	4.61	1.65	1.55
4	B	833	ZK1	CAJ-CAS	3.59	1.44	1.39
4	A	833	ZK1	CAJ-CAS	3.56	1.44	1.39
4	C	833	ZK1	CAJ-CAS	3.56	1.44	1.39
4	D	833	ZK1	CAJ-CAS	3.56	1.44	1.39
4	B	833	ZK1	CAW-NAY	3.34	1.46	1.41
4	D	833	ZK1	CAW-NAY	3.34	1.46	1.41
4	A	833	ZK1	CAW-NAY	3.33	1.46	1.41
4	C	833	ZK1	CAW-NAY	3.28	1.46	1.41
4	D	833	ZK1	PBA-CAO	2.40	1.86	1.81

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	833	ZK1	PBA-CAO	2.37	1.86	1.81
4	A	833	ZK1	PBA-CAO	2.36	1.86	1.81
4	B	833	ZK1	PBA-CAO	2.36	1.86	1.81
4	A	833	ZK1	CAR-NAX	2.13	1.46	1.41
4	D	833	ZK1	CAR-NAX	2.12	1.46	1.41
4	C	833	ZK1	CAR-NAX	2.12	1.46	1.41
4	B	833	ZK1	CAR-NAX	2.10	1.45	1.41
4	A	833	ZK1	CAI-CAW	2.04	1.42	1.39
4	C	833	ZK1	CAI-CAW	2.01	1.42	1.39
4	D	833	ZK1	CAI-CAW	2.01	1.42	1.39

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	833	ZK1	CAN-NAX-CAM	4.43	121.54	111.57
4	D	833	ZK1	CAN-NAX-CAM	4.41	121.48	111.57
4	A	833	ZK1	CAN-NAX-CAM	4.40	121.47	111.57
4	B	833	ZK1	CAN-NAX-CAM	4.40	121.47	111.57
4	A	833	ZK1	CAO-NAY-CAU	3.78	120.39	116.55
4	B	833	ZK1	CAO-NAY-CAU	3.75	120.36	116.55
4	D	833	ZK1	CAO-NAY-CAU	3.75	120.36	116.55
4	C	833	ZK1	CAO-NAY-CAU	3.67	120.28	116.55
4	A	833	ZK1	CAV-NAP-CAT	-3.44	120.00	124.82
4	D	833	ZK1	CAV-NAP-CAT	-3.40	120.05	124.82
4	B	833	ZK1	CAV-NAP-CAT	-3.40	120.06	124.82
4	C	833	ZK1	CAV-NAP-CAT	-3.38	120.08	124.82
4	B	833	ZK1	CAU-CAT-NAP	2.54	120.06	117.46
4	D	833	ZK1	CAU-CAT-NAP	2.54	120.05	117.46
4	A	833	ZK1	CAU-CAT-NAP	2.53	120.05	117.46
4	A	833	ZK1	CAI-CAR-NAX	-2.51	118.94	122.59
4	C	833	ZK1	CAU-CAT-NAP	2.50	120.01	117.46
4	C	833	ZK1	CAI-CAR-NAX	-2.45	119.02	122.59
4	B	833	ZK1	CAI-CAR-NAX	-2.43	119.05	122.59
4	D	833	ZK1	CAI-CAR-NAX	-2.43	119.06	122.59
4	A	833	ZK1	CAT-CAU-NAY	2.34	119.95	117.41
4	D	833	ZK1	CAW-NAY-CAU	-2.32	119.98	122.84
4	A	833	ZK1	CAW-NAY-CAU	-2.31	120.00	122.84
4	D	833	ZK1	CAT-CAU-NAY	2.30	119.91	117.41
4	C	833	ZK1	CAT-CAU-NAY	2.29	119.89	117.41
4	C	833	ZK1	CAW-NAY-CAU	-2.26	120.05	122.84
4	B	833	ZK1	CAT-CAU-NAY	2.24	119.84	117.41
4	B	833	ZK1	CAW-NAY-CAU	-2.23	120.10	122.84

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	833	ZK1	NAY-CAO-PBA-OAE
4	B	833	ZK1	NAY-CAO-PBA-OAE
4	C	833	ZK1	NAY-CAO-PBA-OAE
4	D	833	ZK1	NAY-CAO-PBA-OAE
4	C	833	ZK1	CAI-CAR-NAX-CAN
4	A	833	ZK1	CAI-CAR-NAX-CAN
4	B	833	ZK1	CAI-CAR-NAX-CAN
4	D	833	ZK1	CAI-CAR-NAX-CAN
4	A	833	ZK1	NAY-CAO-PBA-OAC
4	A	833	ZK1	NAY-CAO-PBA-OAD
4	B	833	ZK1	NAY-CAO-PBA-OAC
4	B	833	ZK1	NAY-CAO-PBA-OAD
4	C	833	ZK1	NAY-CAO-PBA-OAC
4	C	833	ZK1	NAY-CAO-PBA-OAD
4	D	833	ZK1	NAY-CAO-PBA-OAC
4	D	833	ZK1	NAY-CAO-PBA-OAD
4	A	833	ZK1	CAS-CAR-NAX-CAN
4	C	833	ZK1	CAS-CAR-NAX-CAN
4	D	833	ZK1	CAS-CAR-NAX-CAN
4	B	833	ZK1	CAS-CAR-NAX-CAN

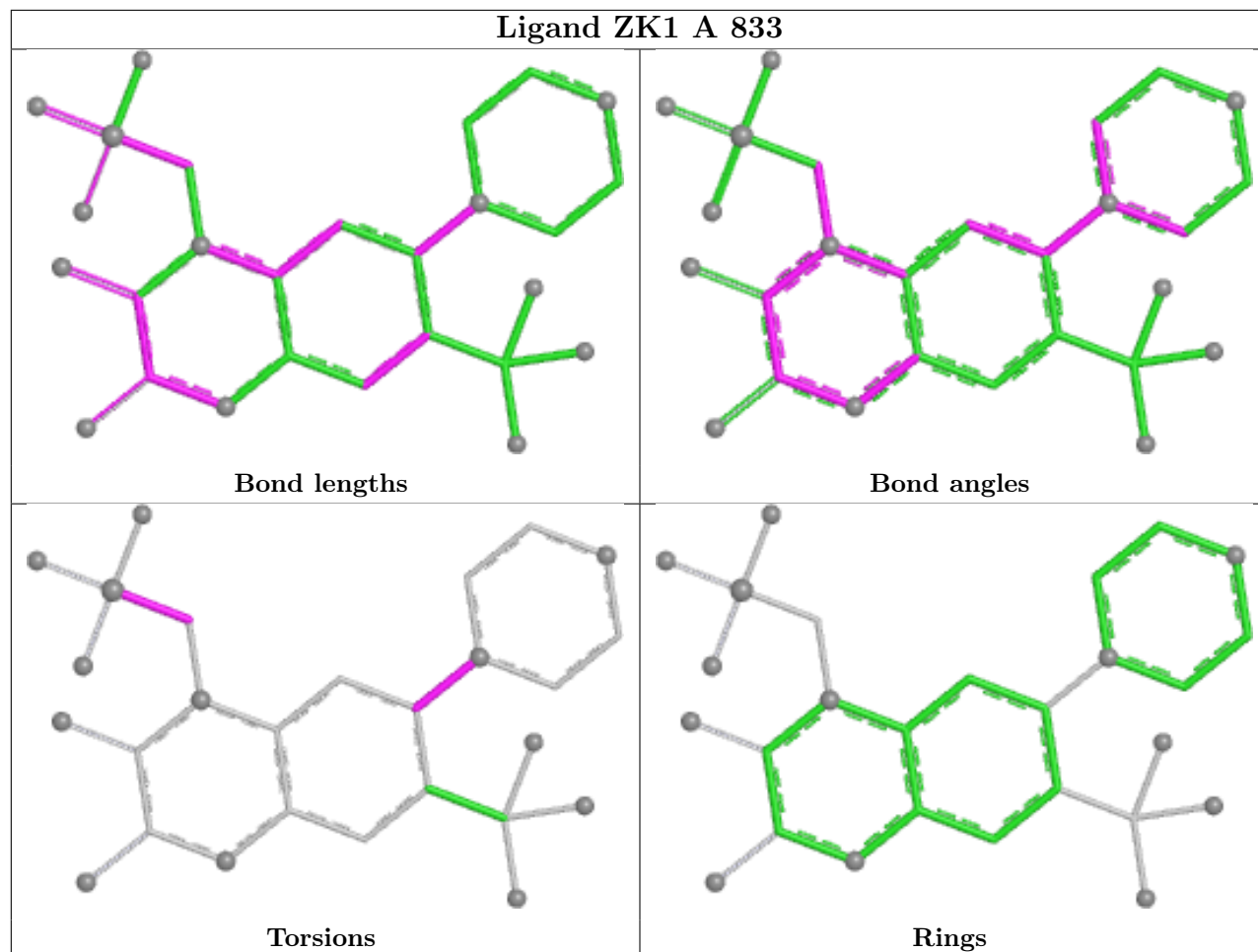
There are no ring outliers.

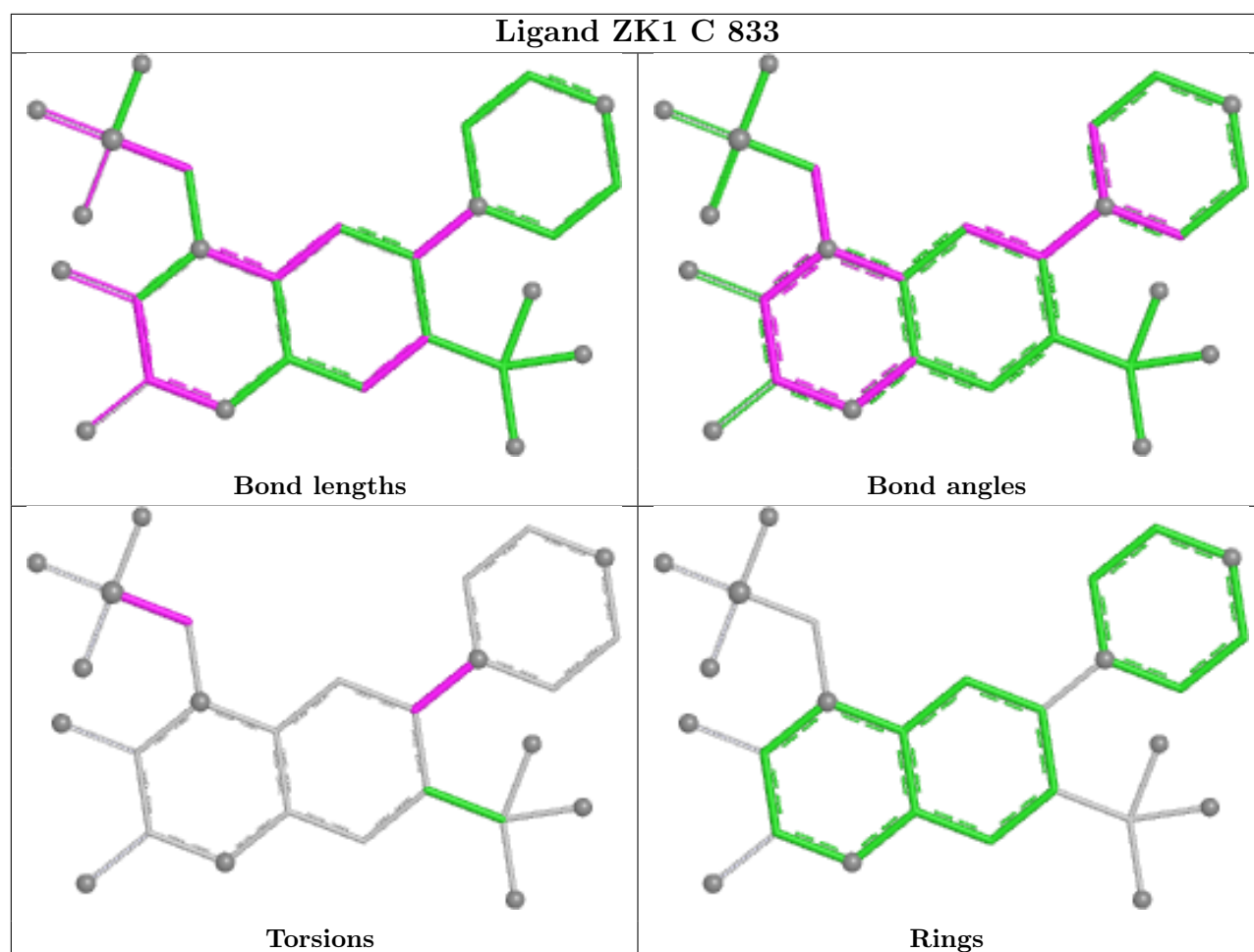
3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	833	ZK1	1	0
4	D	833	ZK1	1	0
4	B	833	ZK1	1	0

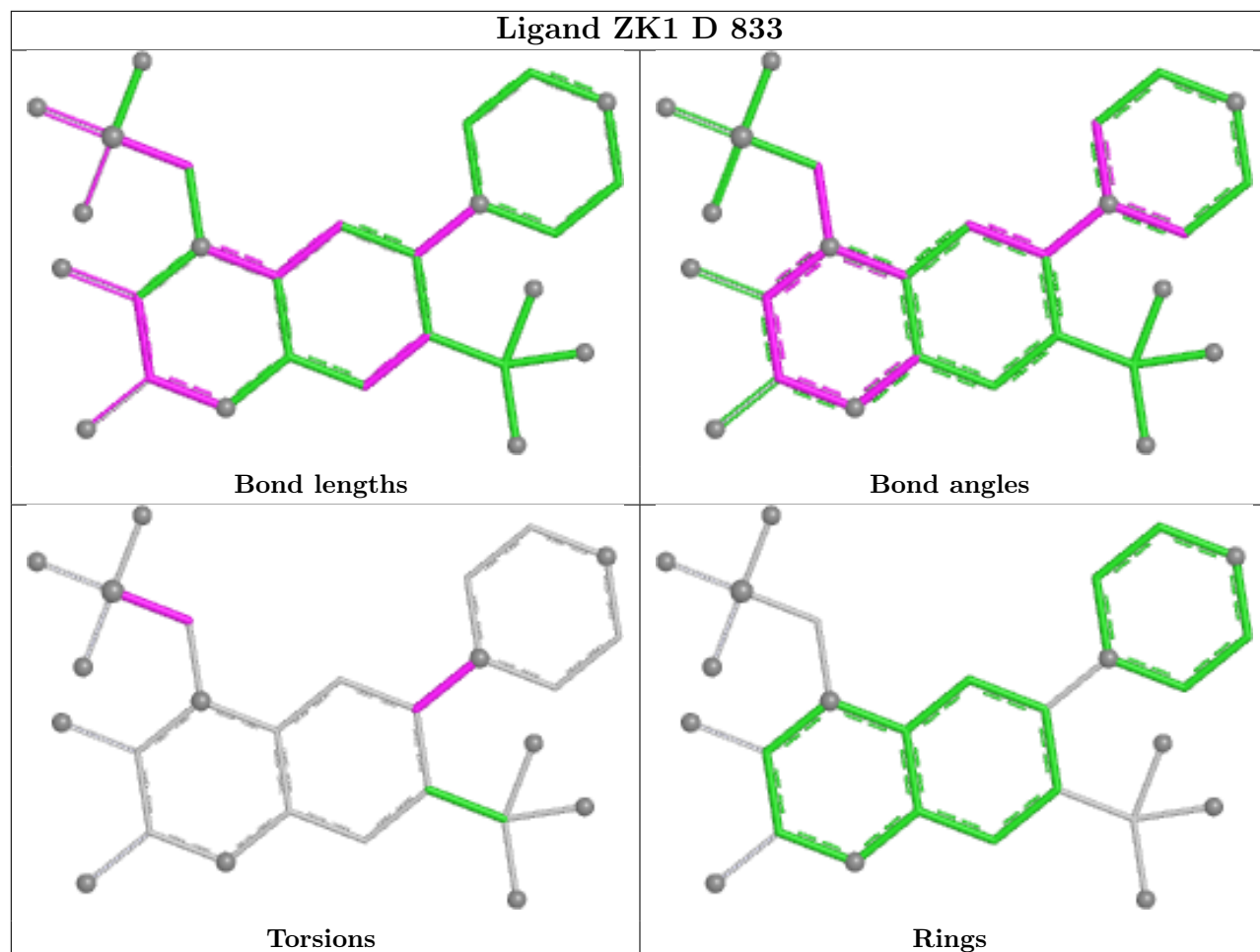
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient

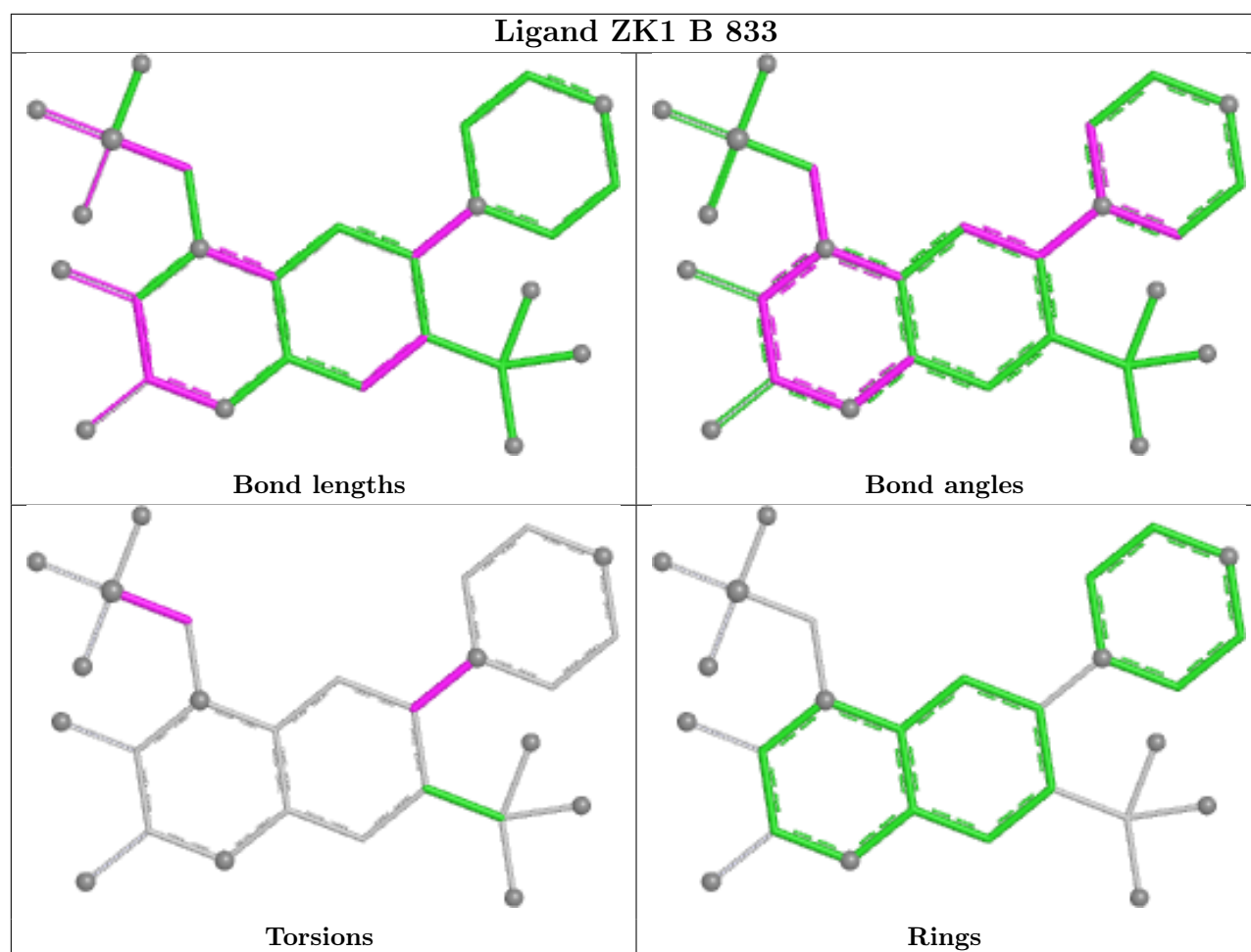
equivalents in the CSD to analyse the geometry.





Ligand ZK1 D 833





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

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

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2			OWAB(Å ²)	Q < 0.9
1	A	779/823 (94%)	0.36	49 (6%)	26	16	108, 154, 266, 372	0
1	B	779/823 (94%)	0.38	47 (6%)	27	17	105, 152, 270, 342	0
1	C	779/823 (94%)	0.22	34 (4%)	39	22	125, 196, 287, 349	0
1	D	779/823 (94%)	0.26	38 (4%)	35	20	115, 177, 287, 356	0
All	All	3116/3292 (94%)	0.30	168 (5%)	31	18	105, 169, 281, 372	0

All (168) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	792	VAL	7.4
1	A	497	SER	7.2
1	D	602	GLY	6.6
1	D	792	VAL	6.0
1	B	725	GLY	6.0
1	A	791	ASN	5.5
1	C	510	SER	5.4
1	D	72	ALA	5.3
1	C	72	ALA	4.9
1	A	710	GLU	4.8
1	A	498	LEU	4.8
1	B	101	PHE	4.8
1	B	77	TYR	4.5
1	B	792	VAL	4.4
1	A	611	ILE	4.4
1	A	496	MET	4.4
1	A	732	TYR	4.3
1	B	665	ALA	4.1
1	A	731	GLY	4.1
1	C	779	GLY	3.9
1	A	610	LEU	3.8

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Mol	Chain	Res	Type	RSRZ
1	D	687	ALA	3.8
1	C	186	LEU	3.7
1	B	235	ILE	3.7
1	B	12	ILE	3.6
1	A	615	SER	3.6
1	B	724	GLY	3.6
1	B	113	MET	3.5
1	A	293	THR	3.5
1	B	726	ASN	3.5
1	B	295	ALA	3.4
1	B	633	ILE	3.4
1	D	803	LEU	3.4
1	B	211	ILE	3.3
1	B	50	LEU	3.3
1	C	630	VAL	3.3
1	A	50	LEU	3.3
1	B	516	SER	3.3
1	B	102	PRO	3.3
1	B	616	TYR	3.3
1	C	803	LEU	3.2
1	D	631	SER	3.2
1	D	686	THR	3.2
1	D	324	GLY	3.2
1	D	785	SER	3.2
1	D	197	GLU	3.1
1	A	197	GLU	3.1
1	C	724	GLY	3.1
1	D	165	VAL	3.1
1	A	296	PHE	3.0
1	B	71	TYR	3.0
1	C	578	TRP	3.0
1	B	666	VAL	3.0
1	D	606	TRP	2.9
1	A	707	THR	2.9
1	C	248	VAL	2.9
1	B	36	PHE	2.9
1	A	726	ASN	2.9
1	B	791	ASN	2.9
1	C	107	HIS	2.9
1	D	322	PRO	2.9
1	B	306	ILE	2.8
1	D	609	THR	2.8

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Mol	Chain	Res	Type	RSRZ
1	C	787	LEU	2.8
1	B	112	GLN	2.8
1	B	513	GLY	2.8
1	A	94	VAL	2.7
1	D	328	GLU	2.7
1	C	187	LYS	2.7
1	B	291	VAL	2.7
1	B	518	LEU	2.7
1	A	147	GLN	2.7
1	B	137	TYR	2.7
1	C	15	GLY	2.7
1	D	212	GLY	2.7
1	C	50	LEU	2.7
1	A	514	VAL	2.6
1	B	140	ASP	2.6
1	B	615	SER	2.6
1	A	183	ASP	2.6
1	A	333	LEU	2.6
1	D	791	ASN	2.6
1	A	89	CYS	2.6
1	C	292	MET	2.6
1	A	607	PHE	2.6
1	B	623	PHE	2.6
1	A	482	THR	2.5
1	D	113	MET	2.5
1	D	347	PHE	2.5
1	A	706	SER	2.5
1	B	252	ASP	2.5
1	D	726	ASN	2.5
1	A	499	GLY	2.5
1	A	789	LEU	2.5
1	B	519	ASP	2.5
1	A	67	SER	2.5
1	D	62	PHE	2.5
1	D	633	ILE	2.5
1	C	527	MET	2.4
1	D	94	VAL	2.4
1	A	787	LEU	2.4
1	A	793	ALA	2.4
1	C	725	GLY	2.4
1	B	456	ASP	2.4
1	C	777	ASP	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	514	VAL	2.4
1	A	786	ALA	2.4
1	B	468	VAL	2.4
1	A	61	ALA	2.4
1	A	73	ILE	2.3
1	D	50	LEU	2.3
1	B	249	ASP	2.3
1	B	658	PHE	2.3
1	D	323	TRP	2.3
1	D	578	TRP	2.3
1	D	333	LEU	2.3
1	C	708	MET	2.3
1	B	542	LEU	2.2
1	C	526	TRP	2.2
1	C	181	PHE	2.2
1	C	295	ALA	2.2
1	A	405	TYR	2.2
1	D	71	TYR	2.2
1	C	291	VAL	2.2
1	A	116	ASP	2.2
1	A	480	THR	2.2
1	B	299	LEU	2.2
1	D	84	THR	2.2
1	A	292	MET	2.2
1	D	396	VAL	2.2
1	D	196	CYS	2.2
1	C	517	PHE	2.2
1	D	654	SER	2.2
1	A	70	VAL	2.2
1	B	99	PRO	2.2
1	C	791	ASN	2.2
1	A	424	TYR	2.2
1	B	141	ARG	2.2
1	C	71	TYR	2.2
1	B	143	LEU	2.2
1	B	179	SER	2.2
1	D	777	ASP	2.1
1	A	495	PHE	2.1
1	C	333	LEU	2.1
1	C	220	TYR	2.1
1	D	582	GLY	2.1
1	B	72	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	102	PRO	2.1
1	C	609	THR	2.1
1	C	594	ARG	2.1
1	D	31	VAL	2.1
1	C	51	GLU	2.1
1	A	57	ALA	2.1
1	B	514	VAL	2.1
1	A	401	LEU	2.1
1	B	284	LEU	2.1
1	A	368	PRO	2.1
1	C	385	ASP	2.1
1	B	784	THR	2.1
1	A	48	ASP	2.0
1	D	92	LEU	2.0
1	A	795	VAL	2.0
1	A	60	ASN	2.0
1	D	521	LEU	2.0
1	A	386	ASP	2.0
1	B	237	PHE	2.0
1	D	320	ALA	2.0
1	C	110	VAL	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.

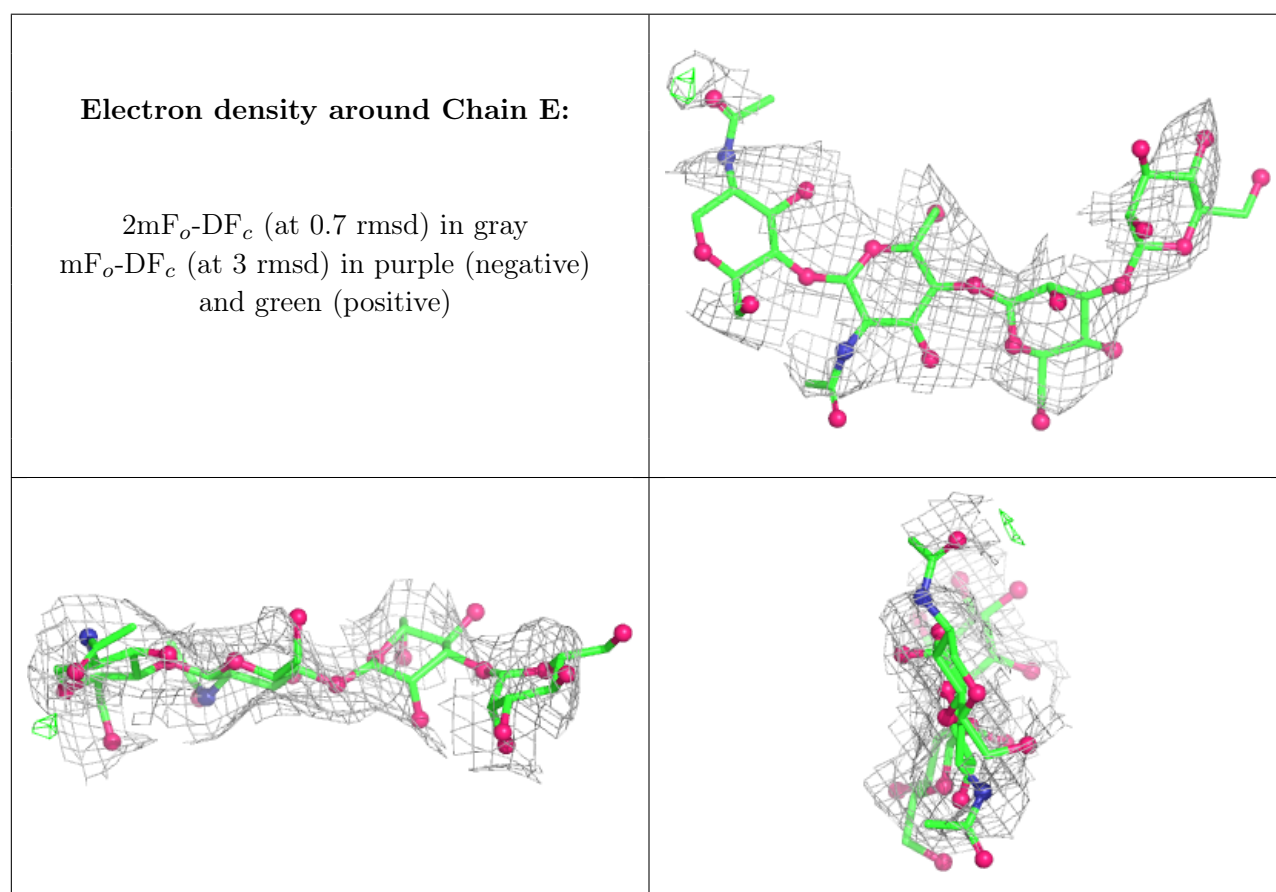
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	NAG	G	1	14/15	-0.11	0.19	415,415,415,415	0
2	NAG	E	1	14/15	0.25	0.14	253,253,253,253	0
3	NAG	G	2	14/15	0.27	0.16	274,274,274,274	0
2	NAG	E	2	14/15	0.34	0.16	277,277,277,277	0
3	BMA	G	3	11/12	0.34	0.11	345,345,345,345	0
2	BMA	E	3	11/12	0.40	0.11	303,303,303,303	0
3	BMA	H	3	11/12	0.48	0.08	265,265,265,265	0

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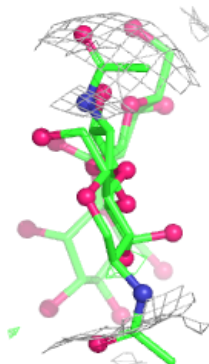
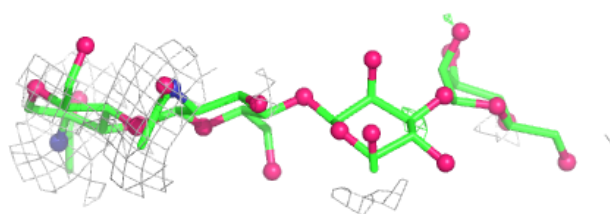
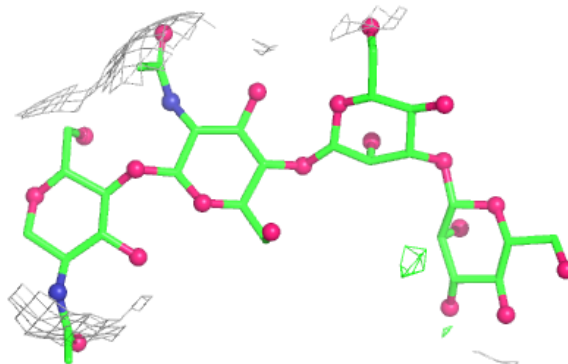
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	BMA	E	4	11/12	0.53	0.07	292,292,292,292	0
2	BMA	F	4	11/12	0.54	0.10	242,242,242,242	0
3	NAG	H	2	14/15	0.58	0.10	283,283,283,283	0
2	NAG	F	2	14/15	0.63	0.09	187,187,187,187	0
2	BMA	F	3	11/12	0.73	0.06	209,209,209,209	0
3	NAG	H	1	14/15	0.81	0.11	267,267,267,267	0
2	NAG	F	1	14/15	0.82	0.10	195,195,195,195	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.

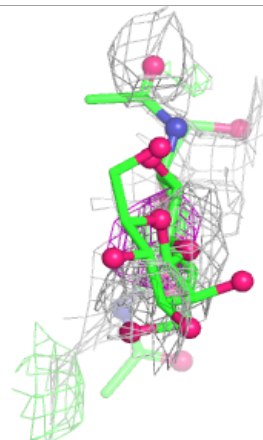
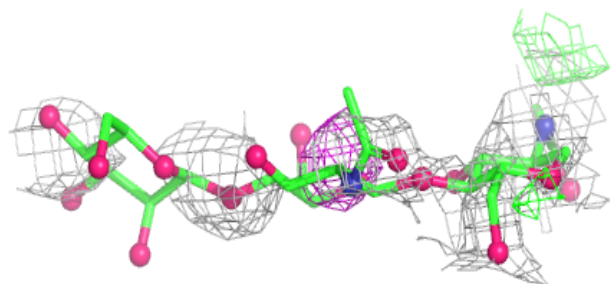
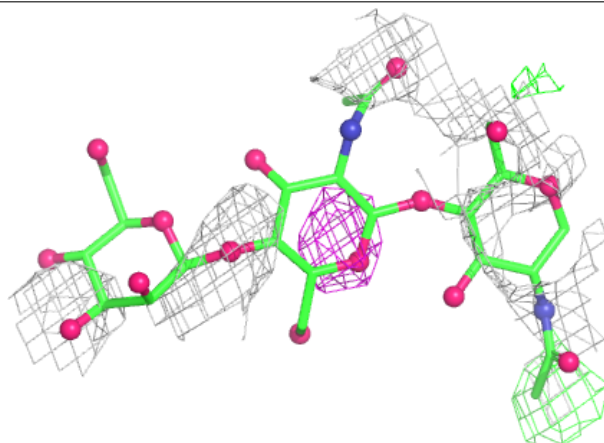


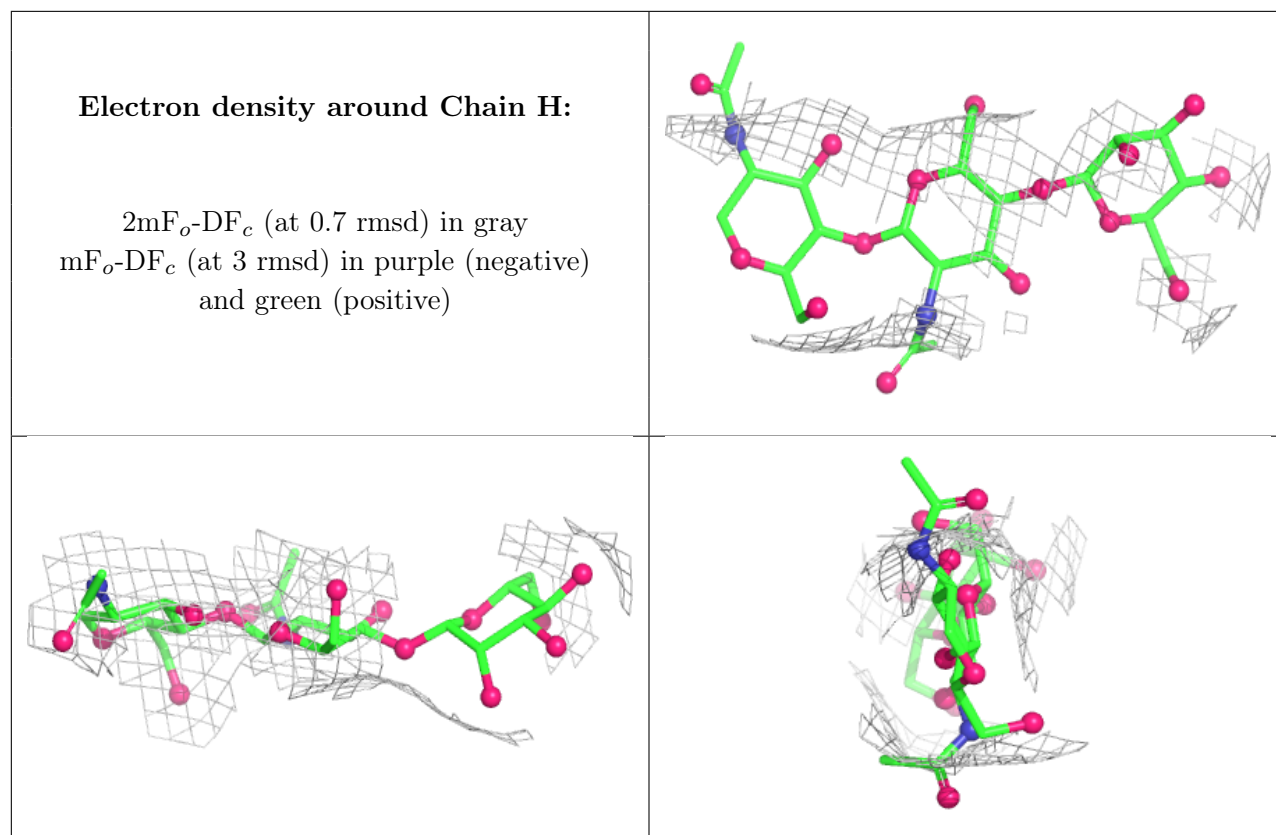
Electron density around Chain F:

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

**Electron density around Chain G:**

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





6.4 Ligands [i](#)

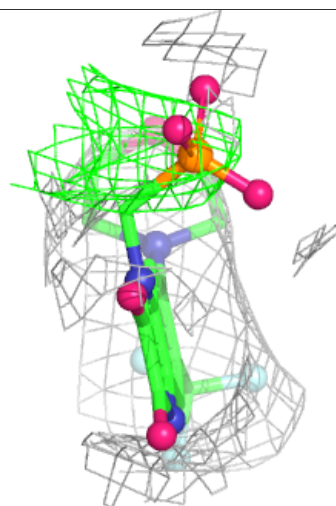
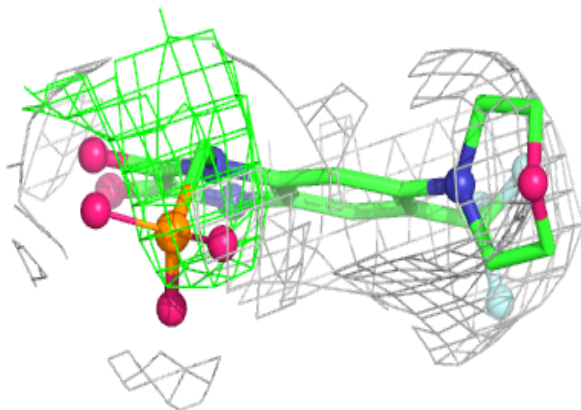
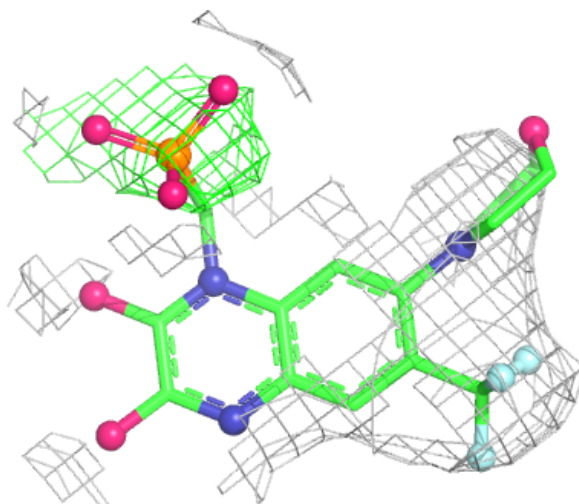
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	ZK1	B	833	27/27	0.87	0.13	160,160,160,160	0
4	ZK1	C	833	27/27	0.88	0.09	162,162,162,162	0
4	ZK1	A	833	27/27	0.92	0.08	108,108,108,108	0
4	ZK1	D	833	27/27	0.92	0.09	162,162,162,162	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

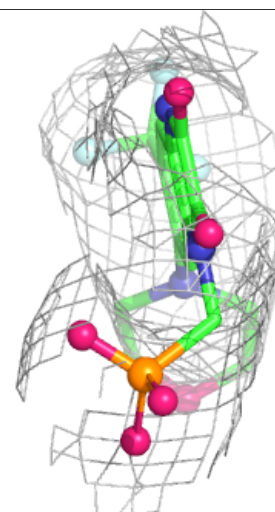
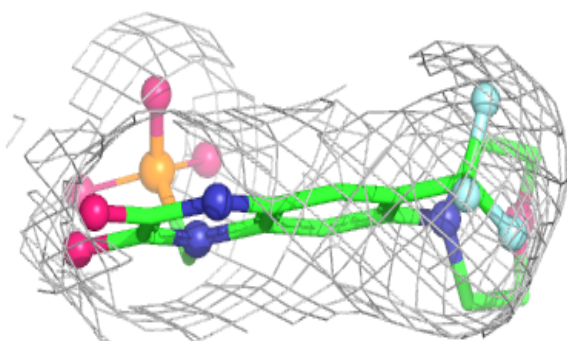
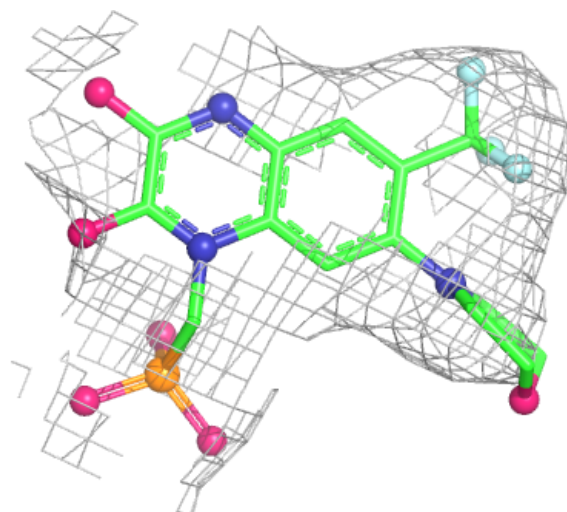
Electron density around ZK1 B 833:

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



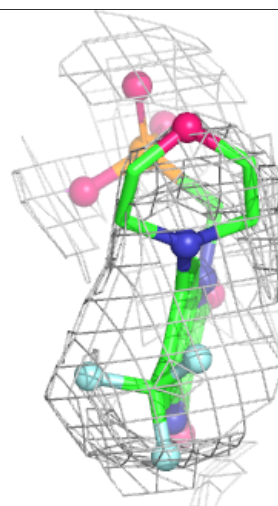
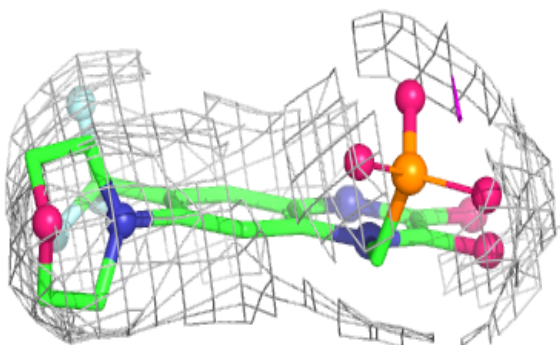
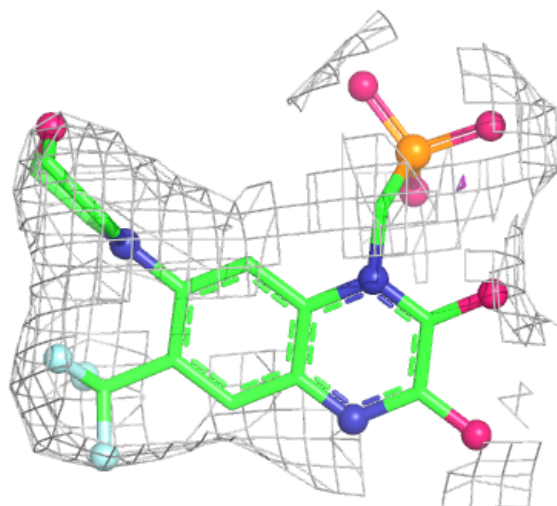
Electron density around ZK1 C 833:

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



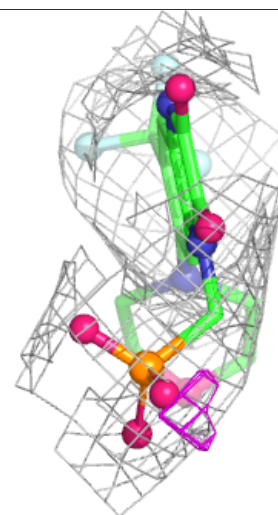
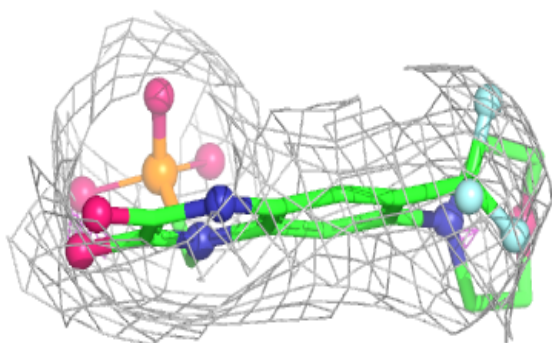
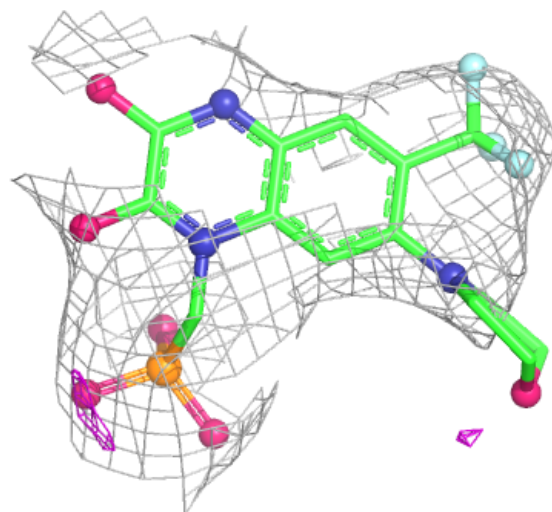
Electron density around ZK1 A 833:

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



Electron density around ZK1 D 833:

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



6.5 Other polymers [i](#)

There are no such residues in this entry.