



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 04:33 AM UTC

PDB ID : 3QI9 / pdb_00003qi9
Title : Crystal structure of mouse CD1d-alpha-phosphatidylinositol with mouse Valpha14-Vbeta6 2A3-D NKT TCR
Authors : Clarke, A.J.; Rossjohn, J.
Deposited on : 2011-01-26
Resolution : 2.30 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

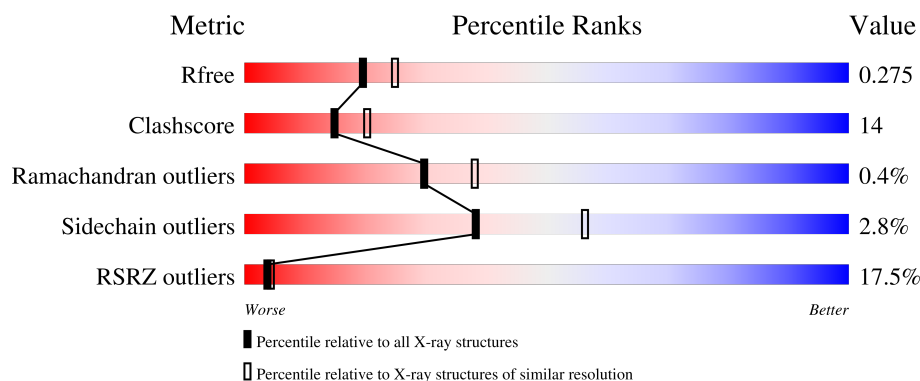
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	302	4% 79% 15% • •
2	B	99	5% 87% 12% •
3	C	207	32% 79% 13% • 5%
4	D	243	24% 83% 15% • •
5	E	2	100%

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 6863 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 Antigen-presenting glycoprotein CD1d1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	290	2333	1486	406	427	14	6	0	0

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	201	HIS	ASP	conflict	UNP P11609
A	280	GLY	-	expression tag	UNP P11609
A	281	SER	-	expression tag	UNP P11609
A	282	LEU	-	expression tag	UNP P11609
A	283	HIS	-	expression tag	UNP P11609
A	284	HIS	-	expression tag	UNP P11609
A	285	ILE	-	expression tag	UNP P11609
A	286	LEU	-	expression tag	UNP P11609
A	287	ASP	-	expression tag	UNP P11609
A	288	ALA	-	expression tag	UNP P11609
A	289	GLN	-	expression tag	UNP P11609
A	290	LYS	-	expression tag	UNP P11609
A	291	MET	-	expression tag	UNP P11609
A	292	VAL	-	expression tag	UNP P11609
A	293	TRP	-	expression tag	UNP P11609
A	294	ASN	-	expression tag	UNP P11609
A	295	HIS	-	expression tag	UNP P11609
A	296	ARG	-	expression tag	UNP P11609
A	297	HIS	-	expression tag	UNP P11609
A	298	HIS	-	expression tag	UNP P11609
A	299	HIS	-	expression tag	UNP P11609
A	300	HIS	-	expression tag	UNP P11609
A	301	HIS	-	expression tag	UNP P11609
A	302	HIS	-	expression tag	UNP P11609

- Molecule 2 is a protein called Beta-2-microglobulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	99	Total	C	N	O	S	1	0	0
			814	520	138	149	7			

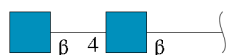
- Molecule 3 is a protein called NKT TCR V alpha 14.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	196	Total	C	N	O	S	3	0	0
			1514	938	260	309	7			

- Molecule 4 is a protein called NKT TCR V beta 6 2A3-D.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	239	Total	C	N	O	S	5	0	0
			1897	1201	326	363	7			

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



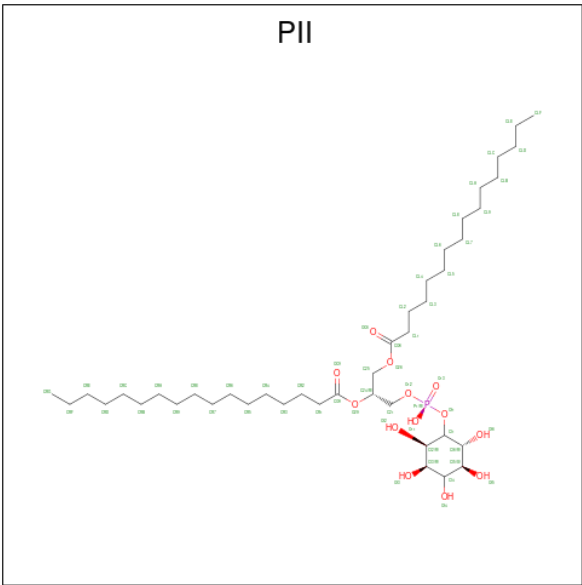
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	2	Total	C	N	O		0	0	0
			28	16	2	10				

- Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



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

- Molecule 7 is 2-[(HYDROXY{(2R,3R,5S,6R)-2,3,4,5,6-PENTAHYDROXYCYCLOHEXYL}OXY}PHOSPHORYL)OXY]-1-[(PALMITOYLOXY)METHYL]ETHYL HEPTADECANOATE (CCD ID: PII) (formula: C₄₂H₈₁O₁₃P).

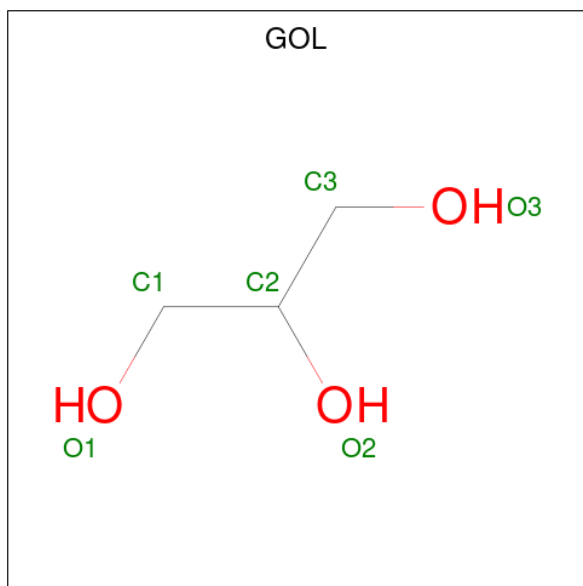


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	1	Total	C	O	P	0	0
			56	42	13	1		

- Molecule 8 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	1	Total	Mg	0	0
			1	1		

- Molecule 9 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	B	1	Total	C	O	0	0
			6	3	3		
9	D	1	Total	C	O	0	0
			6	3	3		

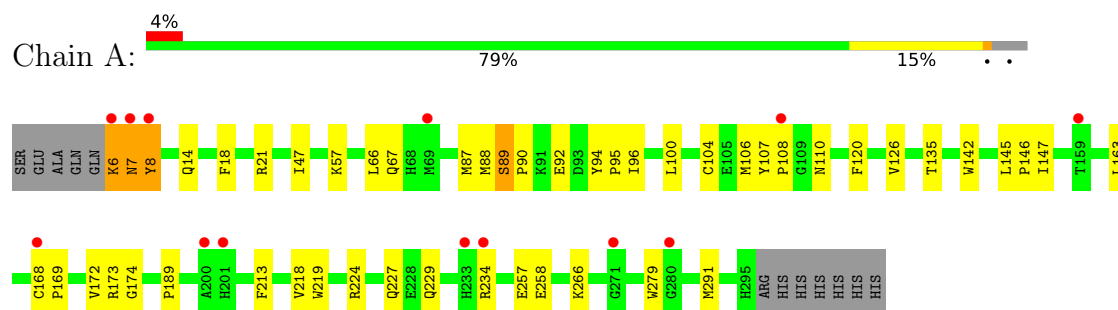
- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	68	Total	O	0	0
			68	68		
10	B	26	Total	O	0	0
			26	26		
10	C	38	Total	O	0	0
			38	38		
10	D	48	Total	O	0	0
			48	48		

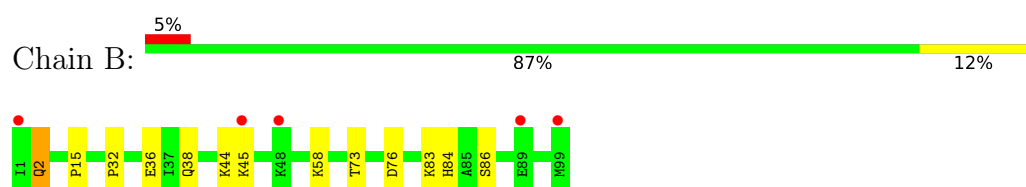
3 Residue-property plots

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.

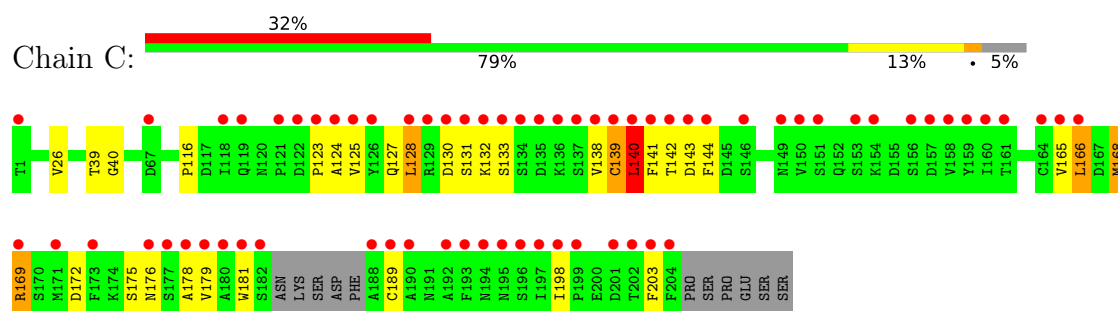
- Molecule 1: Antigen-presenting glycoprotein CD1d1



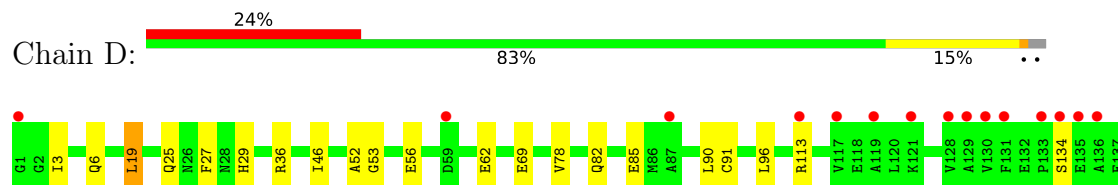
- Molecule 2: Beta-2-microglobulin

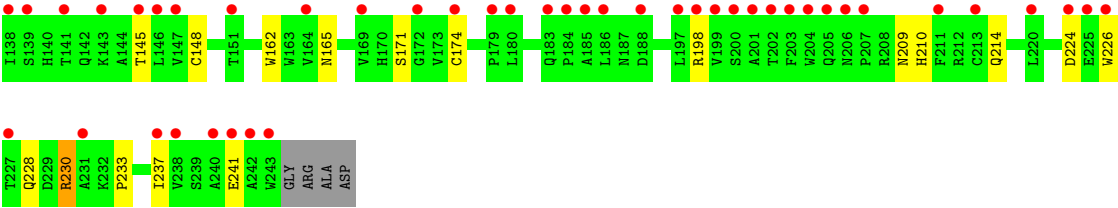


- Molecule 3: NKT TCR V alpha 14



- Molecule 4: NKT TCR V beta 6 2A3-D





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

Chain E: 100%

MAG1
MAG2

4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	95.70Å 95.70Å 289.71Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.31 – 2.30 39.31 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.9 (39.31-2.30) 99.9 (39.31-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.10 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.247 , 0.275 0.249 , 0.275	Depositor DCC
R_{free} test set	3082 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	47.6	Xtriage
Anisotropy	0.038	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 25.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6863	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.44% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PII, MG, GOL, NAG

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.75	0/2403	0.75	2/3266 (0.1%)
2	B	0.67	0/840	0.69	1/1140 (0.1%)
3	C	0.58	0/1539	0.72	0/2089
4	D	0.57	0/1944	0.69	0/2636
All	All	0.65	0/6726	0.72	3/9131 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	7	ASN	N-CA-C	8.52	120.56	111.28
1	A	8	TYR	N-CA-C	5.86	119.02	109.06
2	B	2	GLN	N-CA-C	5.26	117.82	111.40

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	2333	0	2232	45	0
2	B	814	0	788	13	1
3	C	1514	0	1466	99	1
4	D	1897	0	1830	45	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	E	28	0	25	0	0
6	A	28	0	26	0	0
7	A	56	0	80	2	0
8	A	1	0	0	0	0
9	B	6	0	8	0	0
9	D	6	0	8	0	0
10	A	68	0	0	1	0
10	B	26	0	0	0	0
10	C	38	0	0	0	0
10	D	48	0	0	0	0
All	All	6863	0	6463	180	1

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 14.

All (180) 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)
3:C:141:PHE:CZ	3:C:144:PHE:HB3	1.40	1.56
3:C:168:MET:CE	4:D:198:ARG:HG2	1.31	1.54
3:C:40:GLY:CA	4:D:113:ARG:NH2	1.68	1.52
3:C:141:PHE:CE2	3:C:144:PHE:CD1	2.15	1.32
3:C:140:LEU:CD1	3:C:179:VAL:CG2	2.08	1.30
3:C:140:LEU:HD13	3:C:179:VAL:CG2	1.63	1.28
3:C:140:LEU:CD1	3:C:179:VAL:HG22	1.60	1.27
3:C:168:MET:CE	4:D:198:ARG:CG	2.13	1.25
3:C:140:LEU:HB2	3:C:178:ALA:O	1.38	1.22
3:C:168:MET:HE1	4:D:198:ARG:CG	1.73	1.18
3:C:40:GLY:HA2	4:D:113:ARG:NH2	1.33	1.18
3:C:40:GLY:HA3	4:D:113:ARG:NH2	1.52	1.17
3:C:141:PHE:CZ	3:C:144:PHE:CB	2.29	1.14
2:B:38:GLN:NE2	2:B:45:LYS:HE2	1.60	1.14
1:A:89:SER:HB3	1:A:90:PRO:CD	1.72	1.13
3:C:141:PHE:HZ	3:C:144:PHE:CB	1.61	1.12
3:C:140:LEU:HD12	3:C:179:VAL:HG23	1.32	1.12
3:C:168:MET:HE1	4:D:198:ARG:HG2	1.22	1.11
1:A:89:SER:HB3	1:A:90:PRO:HD3	1.14	1.10
1:A:89:SER:CB	1:A:90:PRO:CD	2.29	1.08
3:C:168:MET:HE2	4:D:198:ARG:HG2	1.18	1.08
3:C:123:PRO:HA	3:C:141:PHE:HE1	1.14	1.04
4:D:145:THR:HG22	4:D:198:ARG:HD2	1.39	1.04

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:38:GLN:NE2	2:B:45:LYS:CE	2.21	1.02
3:C:140:LEU:CB	3:C:178:ALA:O	2.08	1.02
3:C:141:PHE:CE2	3:C:144:PHE:CG	2.48	1.00
3:C:140:LEU:CD1	3:C:179:VAL:HG23	1.84	1.00
2:B:38:GLN:HE22	2:B:45:LYS:CE	1.73	0.99
3:C:168:MET:HE2	4:D:198:ARG:CG	1.84	0.99
3:C:139:CYS:O	3:C:140:LEU:HB3	1.65	0.97
3:C:123:PRO:CA	3:C:141:PHE:HE1	1.78	0.96
3:C:125:VAL:HG12	3:C:141:PHE:HB2	1.47	0.96
3:C:140:LEU:HD12	3:C:179:VAL:CG2	1.89	0.95
3:C:166:LEU:HD22	3:C:166:LEU:O	1.65	0.95
2:B:38:GLN:HE22	2:B:45:LYS:NZ	1.65	0.95
3:C:123:PRO:HA	3:C:141:PHE:CE1	2.04	0.93
1:A:6:LYS:HA	1:A:107:TYR:HD1	1.34	0.90
3:C:141:PHE:CE2	3:C:144:PHE:HB3	2.09	0.87
3:C:140:LEU:CB	3:C:179:VAL:HA	2.07	0.84
3:C:40:GLY:CA	4:D:113:ARG:HH22	1.89	0.84
3:C:168:MET:HE1	4:D:198:ARG:HG3	1.58	0.83
3:C:141:PHE:CZ	3:C:144:PHE:CD1	2.67	0.82
3:C:125:VAL:CG1	3:C:141:PHE:HB2	2.08	0.82
3:C:166:LEU:HD13	3:C:166:LEU:H	1.44	0.82
3:C:140:LEU:CG	3:C:178:ALA:O	2.28	0.82
3:C:140:LEU:HB2	3:C:179:VAL:HA	1.62	0.82
1:A:6:LYS:HA	1:A:107:TYR:CD1	2.15	0.81
3:C:141:PHE:CE2	3:C:144:PHE:HD1	1.91	0.81
3:C:40:GLY:HA3	4:D:113:ARG:HH22	1.44	0.80
3:C:140:LEU:HD23	3:C:140:LEU:O	1.82	0.80
1:A:89:SER:CB	1:A:90:PRO:HD3	1.96	0.79
2:B:38:GLN:HE22	2:B:45:LYS:HE2	1.31	0.79
3:C:141:PHE:HE2	3:C:144:PHE:CG	1.98	0.78
3:C:125:VAL:HG12	3:C:141:PHE:CB	2.14	0.78
3:C:168:MET:HE2	4:D:198:ARG:CB	2.13	0.78
4:D:230:ARG:HH21	4:D:230:ARG:HG3	1.47	0.78
2:B:38:GLN:CD	2:B:45:LYS:HE2	2.08	0.78
3:C:140:LEU:HD13	3:C:179:VAL:HG22	0.80	0.77
1:A:168:CYS:O	1:A:172:VAL:HG23	1.86	0.74
1:A:89:SER:CB	1:A:90:PRO:HD2	2.17	0.74
3:C:40:GLY:N	4:D:113:ARG:NH2	2.34	0.74
3:C:141:PHE:CE2	3:C:144:PHE:CB	2.67	0.73
2:B:38:GLN:NE2	2:B:45:LYS:NZ	2.37	0.73
3:C:123:PRO:CA	3:C:141:PHE:CE1	2.67	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:168:MET:HE2	4:D:198:ARG:HB3	1.70	0.72
3:C:166:LEU:HD22	3:C:166:LEU:C	2.17	0.70
3:C:138:VAL:O	3:C:138:VAL:HG13	1.90	0.70
3:C:141:PHE:CZ	3:C:144:PHE:CG	2.79	0.69
1:A:89:SER:HB2	1:A:90:PRO:HD2	1.75	0.67
3:C:169:ARG:HG3	4:D:171:SER:HB3	1.76	0.67
1:A:89:SER:HB2	1:A:90:PRO:CD	2.26	0.66
3:C:140:LEU:HB2	3:C:178:ALA:C	2.20	0.65
1:A:135:THR:HG23	1:A:147:ILE:HD12	1.77	0.65
3:C:141:PHE:CZ	3:C:144:PHE:HD1	2.13	0.64
3:C:141:PHE:HE2	3:C:144:PHE:CD1	2.03	0.64
3:C:168:MET:HE3	4:D:198:ARG:HG2	1.65	0.64
3:C:166:LEU:H	3:C:166:LEU:CD1	2.11	0.63
4:D:6:GLN:NE2	4:D:91:CYS:H	1.96	0.63
4:D:36:ARG:HB2	4:D:46:ILE:HD11	1.81	0.63
1:A:168:CYS:HB3	1:A:169:PRO:HD3	1.78	0.63
3:C:166:LEU:HD13	3:C:175:SER:O	1.99	0.63
4:D:25:GLN:HE22	4:D:29:HIS:H	1.47	0.62
4:D:25:GLN:NE2	4:D:29:HIS:H	1.97	0.62
3:C:166:LEU:CD1	3:C:175:SER:O	2.48	0.61
1:A:219:TRP:HB3	1:A:266:LYS:HB2	1.83	0.60
3:C:141:PHE:CD2	3:C:144:PHE:CD1	2.84	0.60
3:C:140:LEU:HB3	3:C:179:VAL:HA	1.82	0.60
1:A:7:ASN:HB3	1:A:8:TYR:CD2	2.37	0.59
1:A:110:ASN:OD1	1:A:173:ARG:NH1	2.36	0.58
3:C:124:ALA:N	3:C:141:PHE:CD1	2.72	0.58
3:C:139:CYS:CB	3:C:189:CYS:HG	2.12	0.57
3:C:124:ALA:O	3:C:141:PHE:HA	2.05	0.57
3:C:138:VAL:HG23	3:C:181:TRP:HB3	1.84	0.57
3:C:166:LEU:HB3	4:D:174:CYS:HB2	1.87	0.57
1:A:87:MET:HE2	4:D:96:LEU:HD11	1.85	0.56
3:C:139:CYS:CB	3:C:189:CYS:SG	2.92	0.56
1:A:189:PRO:HB3	1:A:213:PHE:HB3	1.88	0.56
1:A:145:LEU:HB3	1:A:146:PRO:HD3	1.88	0.56
3:C:140:LEU:O	3:C:140:LEU:CD2	2.54	0.55
1:A:106:MET:HE3	1:A:173:ARG:HH21	1.71	0.55
4:D:230:ARG:HG3	4:D:230:ARG:NH2	2.21	0.55
2:B:32:PRO:O	2:B:84:HIS:HE1	1.90	0.55
3:C:140:LEU:HD12	3:C:178:ALA:O	2.08	0.54
2:B:38:GLN:HE22	2:B:45:LYS:HZ1	1.48	0.54
3:C:124:ALA:N	3:C:141:PHE:CE1	2.76	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:110:ASN:HA	1:A:173:ARG:NH1	2.23	0.53
1:A:14:GLN:HB3	1:A:100:LEU:HB3	1.90	0.53
3:C:123:PRO:HB3	3:C:141:PHE:CE1	2.43	0.53
1:A:88:MET:HE2	1:A:142:TRP:HE3	1.74	0.52
4:D:6:GLN:HE22	4:D:90:LEU:HA	1.74	0.52
4:D:165:ASN:HD21	4:D:209:ASN:HA	1.75	0.52
2:B:36:GLU:HB3	2:B:83:LYS:HB3	1.91	0.52
1:A:18:PHE:HB2	1:A:96:ILE:HB	1.92	0.52
4:D:25:GLN:HE21	4:D:27:PHE:N	2.08	0.52
3:C:140:LEU:HG	3:C:178:ALA:O	2.07	0.51
3:C:166:LEU:CD1	3:C:166:LEU:N	2.74	0.51
1:A:94:TYR:HB3	1:A:95:PRO:HA	1.92	0.51
1:A:106:MET:HE3	1:A:173:ARG:NH2	2.26	0.51
3:C:140:LEU:CD1	3:C:178:ALA:O	2.59	0.51
1:A:21:ARG:NH1	4:D:56:GLU:OE2	2.44	0.50
4:D:145:THR:CG2	4:D:198:ARG:HD2	2.28	0.50
1:A:104:CYS:SG	1:A:172:VAL:HG21	2.51	0.50
3:C:138:VAL:O	3:C:138:VAL:CG1	2.60	0.49
3:C:172:ASP:OD1	3:C:172:ASP:O	2.30	0.49
4:D:19:LEU:HB2	4:D:78:VAL:HB	1.94	0.49
3:C:141:PHE:HZ	3:C:144:PHE:HB3	0.66	0.48
3:C:140:LEU:HD12	3:C:178:ALA:C	2.38	0.48
4:D:165:ASN:ND2	4:D:210:HIS:H	2.12	0.47
7:A:405:PII:OC6	7:A:405:PII:H212	2.13	0.47
3:C:123:PRO:CB	3:C:141:PHE:CE1	2.97	0.47
4:D:25:GLN:HE21	4:D:27:PHE:H	1.62	0.47
1:A:126:VAL:HG22	1:A:147:ILE:HD11	1.97	0.47
1:A:258:GLU:HB3	1:A:279:TRP:CD1	2.50	0.47
3:C:127:GLN:C	3:C:128:LEU:HD13	2.39	0.47
4:D:224:ASP:OD1	4:D:224:ASP:N	2.48	0.47
1:A:218:VAL:O	1:A:234:ARG:NH2	2.47	0.46
1:A:163:LEU:HD13	7:A:405:PII:HR72	1.98	0.46
1:A:88:MET:HE2	1:A:142:TRP:CE3	2.50	0.46
3:C:140:LEU:HB2	3:C:179:VAL:CA	2.40	0.46
1:A:227:GLN:HE21	1:A:229:GLN:HE22	1.64	0.46
3:C:125:VAL:HG12	3:C:141:PHE:CA	2.45	0.45
3:C:166:LEU:C	3:C:166:LEU:CD2	2.88	0.45
3:C:124:ALA:H	3:C:141:PHE:HD1	1.64	0.45
3:C:165:VAL:HG22	3:C:176:ASN:OD1	2.17	0.45
3:C:116:PRO:HG3	3:C:165:VAL:HG21	1.98	0.45
3:C:125:VAL:HA	3:C:141:PHE:HA	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:230:ARG:HH21	4:D:230:ARG:CG	2.24	0.44
1:A:6:LYS:N	1:A:108:PRO:HD2	2.33	0.44
1:A:47:ILE:H	1:A:67:GLN:HE21	1.66	0.44
3:C:128:LEU:N	3:C:128:LEU:HD22	2.33	0.44
1:A:174:GLY:HA3	10:A:506:HOH:O	2.17	0.44
3:C:124:ALA:O	3:C:141:PHE:HD1	2.00	0.44
4:D:145:THR:HG22	4:D:198:ARG:CD	2.29	0.43
3:C:131:SER:HA	3:C:132:LYS:HA	1.70	0.43
1:A:94:TYR:HA	1:A:95:PRO:C	2.43	0.43
3:C:124:ALA:N	3:C:141:PHE:HD1	2.13	0.43
3:C:127:GLN:O	3:C:128:LEU:HD13	2.18	0.43
3:C:128:LEU:CD1	4:D:134:SER:HB2	2.48	0.43
1:A:224:ARG:NH2	1:A:257:GLU:O	2.51	0.43
4:D:214:GLN:HG3	4:D:237:ILE:HG23	2.01	0.43
1:A:110:ASN:HA	1:A:173:ARG:HH12	1.83	0.43
2:B:84:HIS:HD2	2:B:86:SER:OG	2.01	0.43
1:A:92:GLU:CG	1:A:120:PHE:HZ	2.32	0.43
1:A:106:MET:CG	1:A:172:VAL:HG11	2.49	0.43
1:A:92:GLU:HG2	1:A:120:PHE:CZ	2.54	0.42
4:D:148:CYS:HB2	4:D:162:TRP:CZ2	2.55	0.42
3:C:123:PRO:HB3	3:C:141:PHE:CZ	2.54	0.42
3:C:128:LEU:HD13	4:D:134:SER:HB2	2.01	0.42
4:D:226:TRP:CZ2	4:D:233:PRO:HD3	2.55	0.42
3:C:124:ALA:HB2	3:C:203:PHE:HB3	2.01	0.42
1:A:57:LYS:HD2	1:A:174:GLY:HA2	2.02	0.42
1:A:92:GLU:HG2	1:A:120:PHE:HZ	1.83	0.42
3:C:141:PHE:CE1	3:C:142:THR:O	2.73	0.42
4:D:52:ALA:HA	4:D:53:GLY:HA2	1.76	0.41
3:C:142:THR:OG1	3:C:143:ASP:N	2.53	0.41
1:A:291:MET:HG2	2:B:15:PRO:HG2	2.03	0.41
4:D:230:ARG:NH2	4:D:230:ARG:CG	2.84	0.41
3:C:125:VAL:HG13	3:C:141:PHE:HB2	2.00	0.41
4:D:82:GLN:HB2	4:D:85:GLU:HB2	2.02	0.41
2:B:73:THR:OG1	2:B:76:ASP:HB2	2.21	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:44:LYS:NZ	3:C:133:SER:OG[6_534]	1.89	0.31

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	288/302 (95%)	282 (98%)	5 (2%)	1 (0%)	36	46
2	B	97/99 (98%)	95 (98%)	2 (2%)	0	100	100
3	C	192/207 (93%)	181 (94%)	9 (5%)	2 (1%)	12	15
4	D	237/243 (98%)	229 (97%)	8 (3%)	0	100	100
All	All	814/851 (96%)	787 (97%)	24 (3%)	3 (0%)	30	38

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	89	SER
3	C	140	LEU
3	C	130	ASP

5.3.2 Protein sidechains [i](#)

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	253/264 (96%)	251 (99%)	2 (1%)	73	86
2	B	92/93 (99%)	90 (98%)	2 (2%)	45	65
3	C	175/186 (94%)	166 (95%)	9 (5%)	21	32
4	D	206/208 (99%)	199 (97%)	7 (3%)	32	49
All	All	726/751 (97%)	706 (97%)	20 (3%)	38	56

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	LYS
1	A	66	LEU
2	B	2	GLN
2	B	58	LYS
3	C	26	VAL
3	C	39	THR
3	C	128	LEU
3	C	139	CYS
3	C	140	LEU
3	C	166	LEU
3	C	168	MET
3	C	169	ARG
3	C	198	ILE
4	D	3	ILE
4	D	19	LEU
4	D	62	GLU
4	D	69	GLU
4	D	228	GLN
4	D	230	ARG
4	D	241	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	55	GLN
1	A	60	ASN
1	A	67	GLN
1	A	117	HIS
1	A	154	GLN
1	A	186	GLN
1	A	227	GLN
1	A	295	HIS
2	B	2	GLN
2	B	29	GLN
2	B	38	GLN
2	B	67	HIS
2	B	84	HIS
3	C	14	GLN
3	C	120	ASN
4	D	6	GLN
4	D	25	GLN
4	D	122	ASN
4	D	165	ASN

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Mol	Chain	Res	Type
4	D	170	HIS
4	D	178	GLN
4	D	187	ASN
4	D	209	ASN
4	D	228	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 ⓘ

2 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	$\# Z > 2$
5	NAG	E	1	5,1	14,14,15	0.54	0	17,19,21	1.12	2 (11%)
5	NAG	E	2	5	14,14,15	0.64	0	17,19,21	1.09	1 (5%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	E	1	5,1	-	2/6/23/26	0/1/1/1
5	NAG	E	2	5	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	E	2	NAG	C4-C3-C2	2.74	115.03	111.02
5	E	1	NAG	C4-C3-C2	2.60	114.83	111.02
5	E	1	NAG	O4-C4-C3	-2.58	104.30	110.38

There are no chirality outliers.

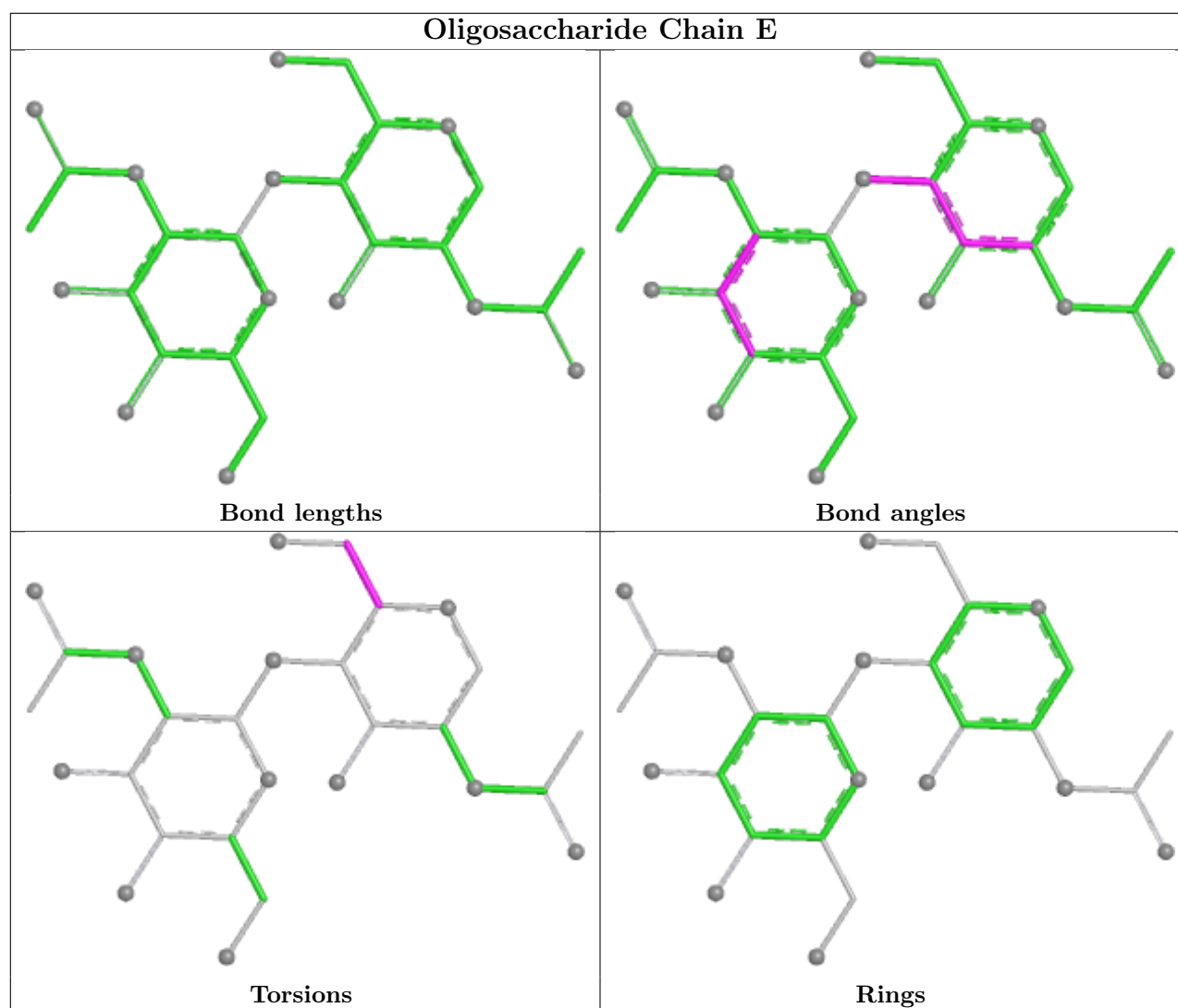
All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	E	1	NAG	C4-C5-C6-O6
5	E	1	NAG	O5-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

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



5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 1 is monoatomic - leaving 5 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
6	NAG	A	402	1	14,14,15	0.45	0	17,19,21	1.14	1 (5%)
9	GOL	B	100	-	5,5,5	0.28	0	5,5,5	0.38	0
9	GOL	D	300	-	5,5,5	0.29	0	5,5,5	0.27	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	NAG	A	401	1	14,14,15	0.43	0	17,19,21	1.70	2 (11%)
7	PII	A	405	-	56,56,56	2.05	3 (5%)	65,68,68	2.51	10 (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
6	NAG	A	402	1	-	4/6/23/26	0/1/1/1
9	GOL	B	100	-	-	2/4/4/4	-
9	GOL	D	300	-	-	0/4/4/4	-
6	NAG	A	401	1	-	4/6/23/26	0/1/1/1
7	PII	A	405	-	-	22/51/75/75	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	A	405	PII	OC9-CO9	12.90	1.61	1.22
7	A	405	PII	C21-C24	-5.95	1.32	1.50
7	A	405	PII	C25-C24	2.80	1.59	1.50

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	A	405	PII	O29-CO9-OC9	-11.89	95.91	123.70
7	A	405	PII	C25-C24-C21	-8.87	91.11	111.78
7	A	405	PII	O29-C24-C21	7.74	136.11	108.34
7	A	405	PII	P1-O12-C21	6.64	159.38	121.35
6	A	401	NAG	C1-O5-C5	4.83	118.66	112.19
7	A	405	PII	O26-C25-C24	4.67	121.86	108.40
7	A	405	PII	O29-CO9-CR1	3.81	119.72	111.48
6	A	401	NAG	O5-C1-C2	-3.46	105.94	111.29
6	A	402	NAG	C1-O5-C5	3.44	116.79	112.19
7	A	405	PII	OC9-CO9-CR1	-3.20	111.25	123.78
7	A	405	PII	OI1-CI1-CI6	2.43	113.87	108.73
7	A	405	PII	O26-CO6-CL1	2.37	119.06	111.83
7	A	405	PII	O29-C24-C25	2.29	116.55	108.34

There are no chirality outliers.

All (32) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	A	405	PII	CI6-CI1-OI1-P1
7	A	405	PII	OC9-CO9-O29-C24
7	A	405	PII	CR1-CO9-O29-C24
9	B	100	GOL	O1-C1-C2-C3
6	A	401	NAG	C8-C7-N2-C2
6	A	401	NAG	O7-C7-N2-C2
6	A	401	NAG	C4-C5-C6-O6
6	A	401	NAG	O5-C5-C6-O6
6	A	402	NAG	O5-C5-C6-O6
6	A	402	NAG	C8-C7-N2-C2
6	A	402	NAG	O7-C7-N2-C2
7	A	405	PII	C25-C24-O29-CO9
7	A	405	PII	CRC-CRD-CRE-CRF
7	A	405	PII	CL8-CL9-CLA-CLB
9	B	100	GOL	O1-C1-C2-O2
7	A	405	PII	CR9-CRA-CRB-CRC
7	A	405	PII	CLA-CLB-CLC-CLD
7	A	405	PII	CR7-CR8-CR9-CRA
7	A	405	PII	CL5-CL6-CL7-CL8
6	A	402	NAG	C4-C5-C6-O6
7	A	405	PII	CR5-CR6-CR7-CR8
7	A	405	PII	CR6-CR7-CR8-CR9
7	A	405	PII	O12-C21-C24-C25
7	A	405	PII	O12-C21-C24-O29
7	A	405	PII	CR4-CR5-CR6-CR7
7	A	405	PII	CR2-CR3-CR4-CR5
7	A	405	PII	CRB-CRC-CRD-CRE
7	A	405	PII	CR1-CR2-CR3-CR4
7	A	405	PII	C21-C24-C25-O26
7	A	405	PII	CL6-CL7-CL8-CL9
7	A	405	PII	CRA-CRB-CRC-CRD
7	A	405	PII	CL2-CL1-CO6-O26

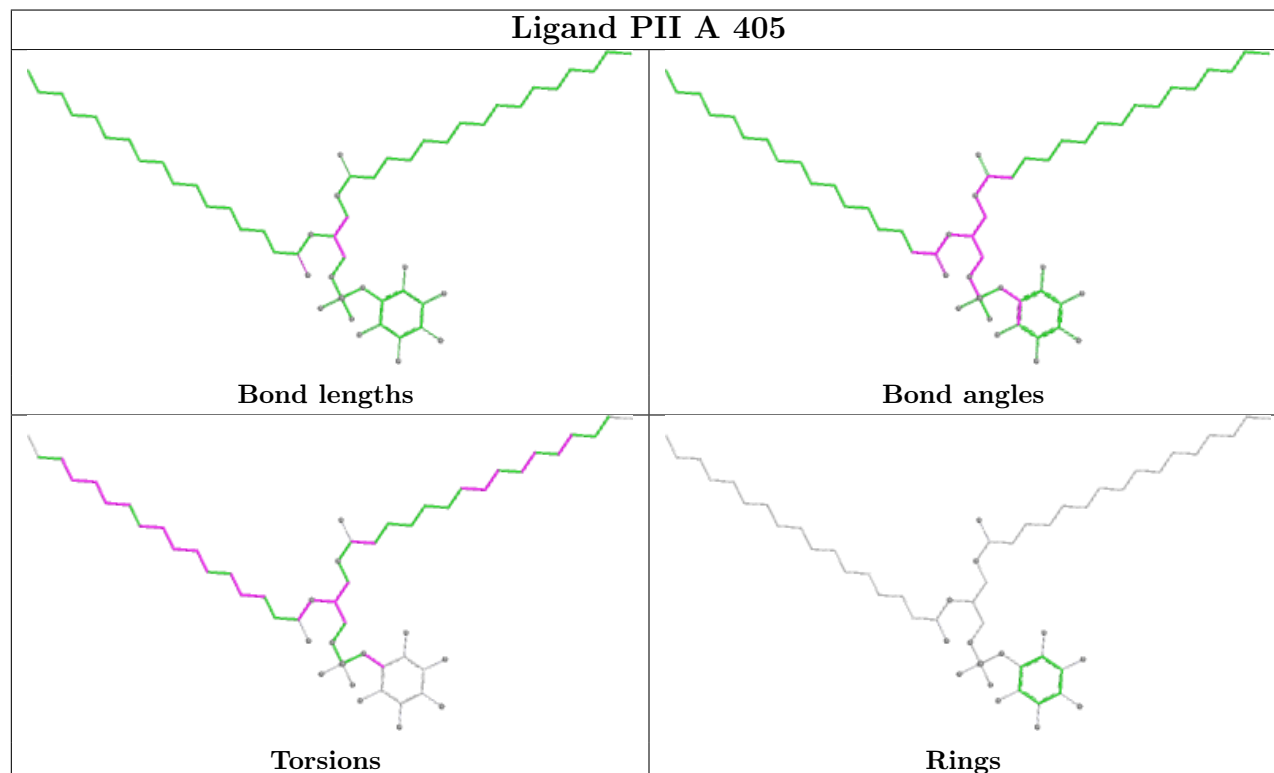
There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	405	PII	2	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.



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	290/302 (96%)	0.56	13 (4%) 38 40	33, 44, 59, 82	5 (1%)
2	B	99/99 (100%)	0.60	5 (5%) 33 35	34, 46, 61, 73	1 (1%)
3	C	196/207 (94%)	1.52	67 (34%) 1 1	29, 49, 113, 124	2 (1%)
4	D	239/243 (98%)	1.26	59 (24%) 2 2	31, 63, 93, 103	3 (1%)
All	All	824/851 (96%)	1.00	144 (17%) 4 4	29, 48, 96, 124	11 (1%)

All (144) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	188	ALA	10.6
3	C	141	PHE	10.5
3	C	140	LEU	9.1
3	C	125	VAL	6.9
3	C	139	CYS	6.1
3	C	128	LEU	5.9
3	C	198	ILE	5.9
3	C	197	ILE	5.7
3	C	134	SER	5.5
3	C	124	ALA	5.3
3	C	138	VAL	5.0
3	C	181	TRP	4.8
4	D	207	PRO	4.5
4	D	211	PHE	4.4
3	C	178	ALA	4.4
3	C	204	PHE	4.4
3	C	190	ALA	4.3
3	C	203	PHE	4.3
3	C	202	THR	4.3
3	C	164	CYS	4.2
4	D	136	ALA	4.2

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Mol	Chain	Res	Type	RSRZ
3	C	132	LYS	4.1
3	C	182	SER	4.0
3	C	126	TYR	4.0
4	D	147	VAL	3.9
3	C	136	LYS	3.9
3	C	131	SER	3.9
4	D	226	TRP	3.9
4	D	130	VAL	3.7
3	C	130	ASP	3.7
4	D	231	ALA	3.7
3	C	199	PRO	3.7
3	C	189	CYS	3.6
4	D	227	THR	3.6
1	A	7	ASN	3.5
4	D	139	SER	3.5
2	B	1	ILE	3.5
4	D	242	ALA	3.5
4	D	119	ALA	3.4
3	C	166	LEU	3.4
1	A	159	THR	3.4
4	D	128	VAL	3.4
3	C	192	ALA	3.4
4	D	113	ARG	3.3
3	C	137	SER	3.3
4	D	186	LEU	3.3
4	D	129	ALA	3.2
3	C	193	PHE	3.2
3	C	157	ASP	3.2
1	A	6	LYS	3.2
4	D	204	TRP	3.2
3	C	142	THR	3.2
3	C	150	VAL	3.1
4	D	145	THR	3.1
3	C	173	PHE	3.1
3	C	195	ASN	3.1
1	A	280	GLY	3.0
3	C	144	PHE	3.0
3	C	133	SER	3.0
1	A	108	PRO	3.0
4	D	240	ALA	3.0
4	D	1	GLY	3.0
3	C	180	ALA	3.0

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Mol	Chain	Res	Type	RSRZ
3	C	123	PRO	3.0
3	C	159	TYR	2.9
2	B	48	LYS	2.9
3	C	119	GLN	2.9
4	D	188	ASP	2.9
3	C	121	PRO	2.9
4	D	237	ILE	2.9
4	D	146	LEU	2.9
4	D	220	LEU	2.8
2	B	45	LYS	2.8
4	D	199	VAL	2.8
4	D	243	TRP	2.8
4	D	202	THR	2.7
4	D	134	SER	2.7
1	A	200	ALA	2.7
4	D	185	ALA	2.7
3	C	161	THR	2.7
3	C	179	VAL	2.6
4	D	201	ALA	2.6
4	D	164	VAL	2.6
3	C	165	VAL	2.6
3	C	118	ILE	2.6
3	C	160	ILE	2.6
4	D	138	ILE	2.6
4	D	205	GLN	2.5
3	C	122	ASP	2.5
4	D	200	SER	2.4
4	D	143	LYS	2.4
1	A	168	CYS	2.4
3	C	154	LYS	2.4
3	C	177	SER	2.4
4	D	241	GLU	2.4
4	D	180	LEU	2.4
4	D	135	GLU	2.3
3	C	67	ASP	2.3
4	D	198	ARG	2.3
3	C	158	VAL	2.3
4	D	169	VAL	2.3
3	C	201	ASP	2.3
4	D	121	LYS	2.3
4	D	133	PRO	2.3
2	B	99	MET	2.3

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Mol	Chain	Res	Type	RSRZ
4	D	203	PHE	2.3
4	D	174	CYS	2.3
4	D	206	ASN	2.3
3	C	1	THR	2.3
3	C	196	SER	2.3
4	D	131	PHE	2.2
3	C	176	ASN	2.2
4	D	59	ASP	2.2
3	C	129	ARG	2.2
3	C	135	ASP	2.2
3	C	143	ASP	2.2
1	A	271	GLY	2.2
4	D	172	GLY	2.2
2	B	89	GLU	2.2
4	D	197	LEU	2.2
3	C	171	MET	2.2
3	C	146	SER	2.2
3	C	151	SER	2.2
3	C	156	SER	2.2
4	D	117	VAL	2.2
4	D	183	GLN	2.2
3	C	149	ASN	2.1
4	D	224	ASP	2.1
1	A	8	TYR	2.1
4	D	225	GLU	2.1
1	A	201	HIS	2.1
3	C	153	SER	2.1
4	D	151	THR	2.1
1	A	69	MET	2.1
1	A	234	ARG	2.1
4	D	141	THR	2.1
4	D	87	ALA	2.1
1	A	233	HIS	2.1
4	D	184	PRO	2.1
4	D	213	CYS	2.0
4	D	238	VAL	2.0
3	C	194	ASN	2.0
3	C	169	ARG	2.0
4	D	179	PRO	2.0

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

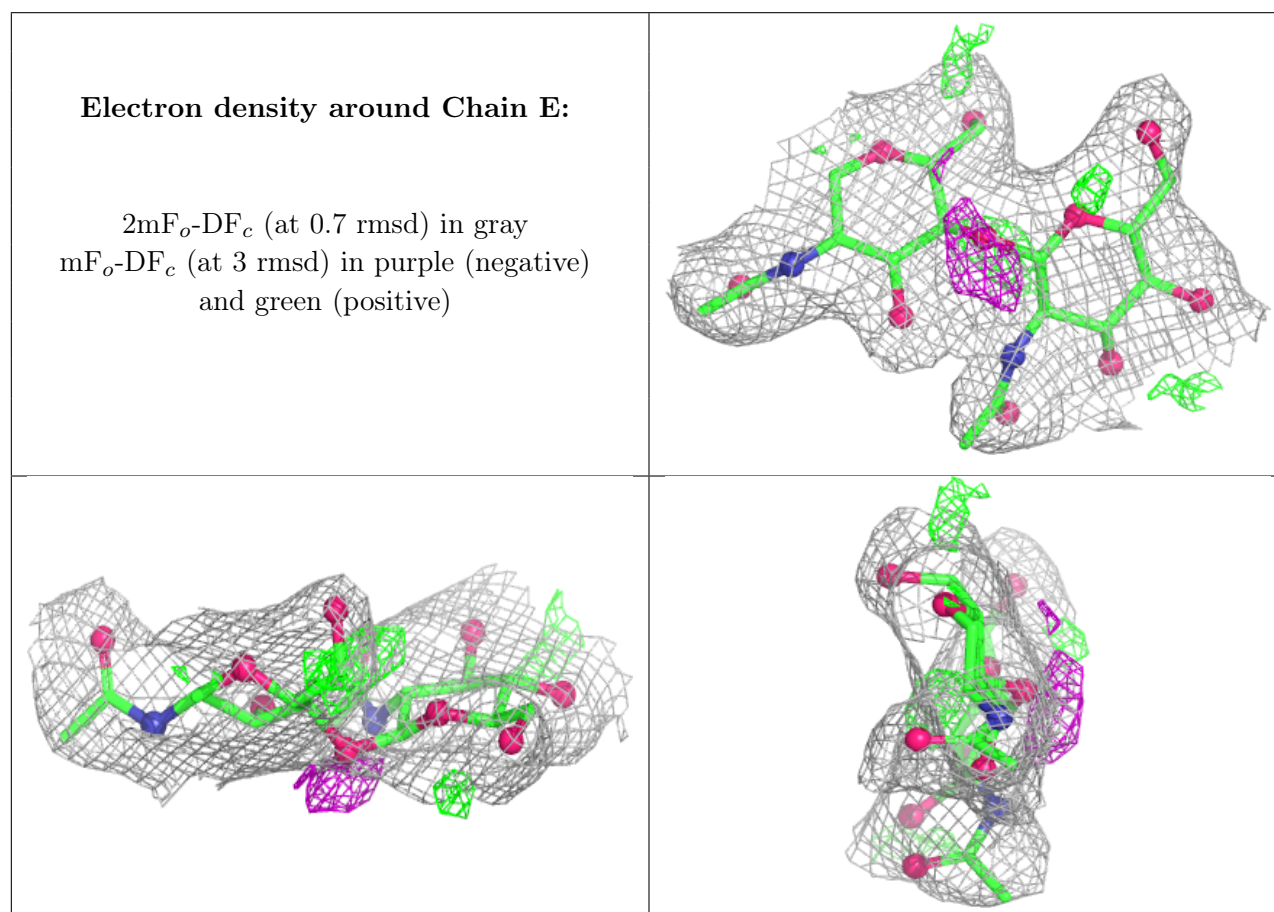
There are no non-standard protein/DNA/RNA residues in this entry.

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
5	NAG	E	2	14/15	0.74	0.17	71,74,79,81	0
5	NAG	E	1	14/15	0.86	0.13	56,58,62,67	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.

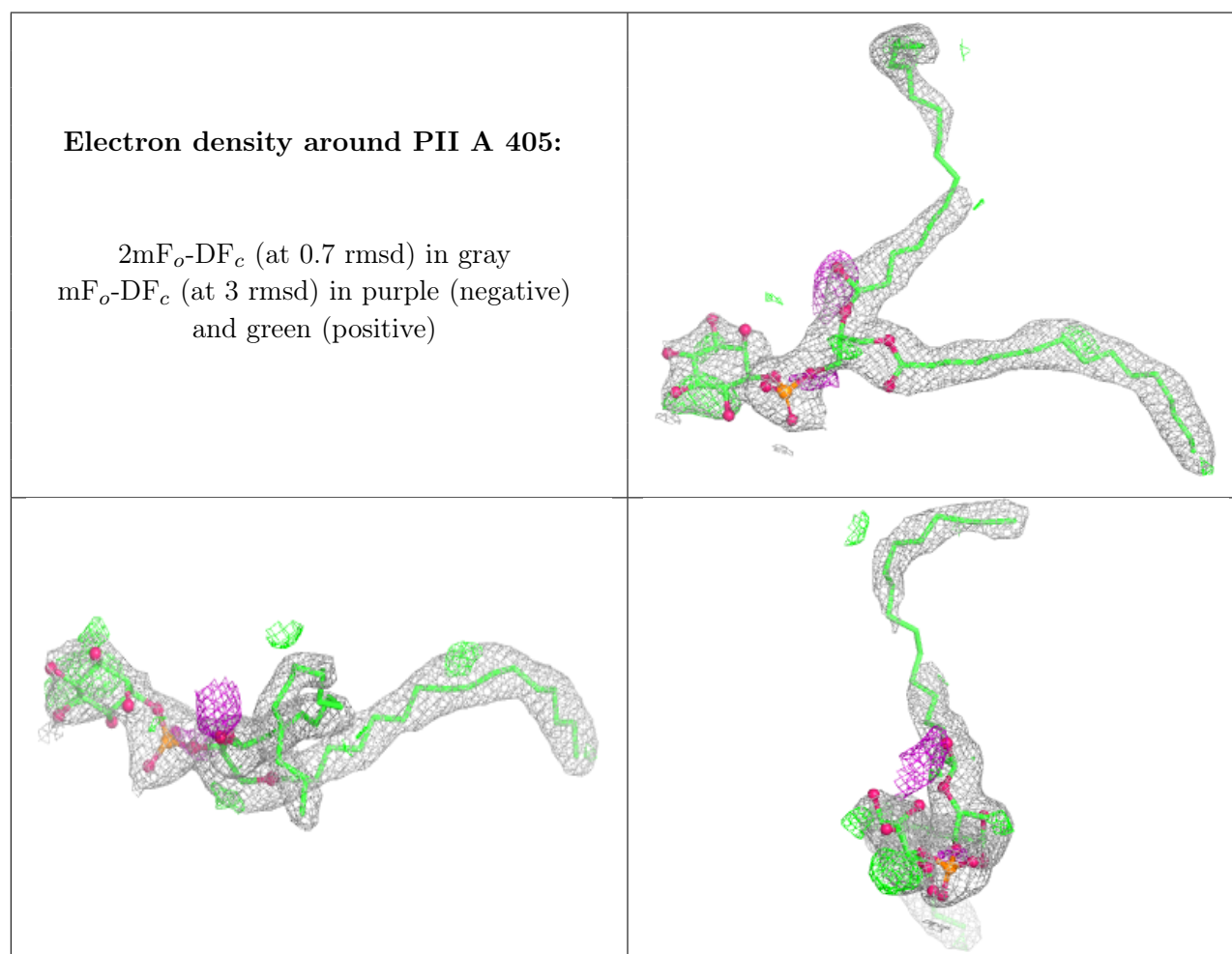


6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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
6	NAG	A	401	14/15	0.56	0.20	52,54,55,55	0
9	GOL	B	100	6/6	0.68	0.27	66,67,69,70	0
9	GOL	D	300	6/6	0.71	0.22	70,76,77,77	0
7	PII	A	405	56/56	0.85	0.22	45,57,64,66	11
6	NAG	A	402	14/15	0.88	0.11	51,53,55,55	0
8	MG	A	406	1/1	0.93	0.33	62,62,62,62	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.



6.5 Other polymers [i](#)

There are no such residues in this entry.