



wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 9, 2020 – 06:05 PM BST

PDB ID : 2GRX
Title : Crystal structure of TonB in complex with FhuA, E. coli outer membrane receptor for ferrichrome
Authors : Pawelek, P.D.; Allaire, M.; Coulton, J.W.
Deposited on : 2006-04-25
Resolution : 3.30 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.13.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.13.1

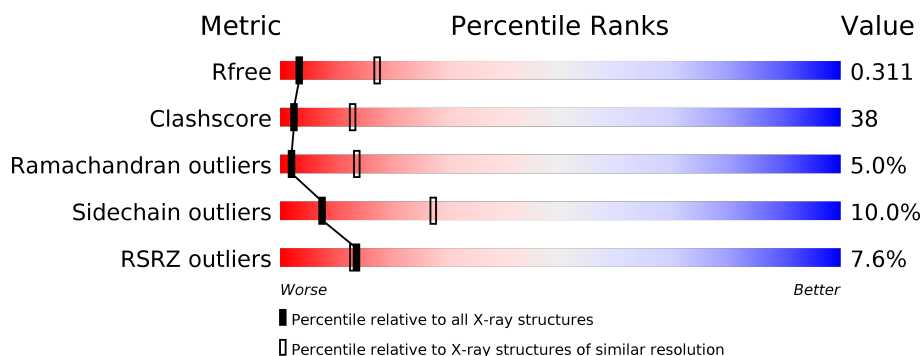
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.30 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}	130704	1149 (3.34-3.26)
Clashscore	141614	1205 (3.34-3.26)
Ramachandran outliers	138981	1183 (3.34-3.26)
Sidechain outliers	138945	1182 (3.34-3.26)
RSRZ outliers	127900	1115 (3.34-3.26)

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 on 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	725	
1	B	725	
2	C	229	
2	D	229	
3	E	5	
3	F	5	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
10	MYR	B	940	-	-	-	X
3	GCN	E	1	-	-	X	-
3	GCN	F	1	-	-	X	-
3	GCN	F	2	-	-	X	-
3	KDO	F	4	-	-	X	-
5	FTT	A	900	-	-	X	-
5	FTT	A	901	-	-	X	X
5	FTT	A	903	-	-	-	X
5	FTT	B	900	-	-	X	-
5	FTT	B	901	-	-	X	X
5	FTT	B	902	-	-	X	-
5	FTT	B	903	-	-	X	X
6	DPO	A	910	-	-	X	-
6	DPO	B	910	-	-	X	-
7	DAO	A	930	-	-	-	X
7	DAO	B	930	-	-	X	-

2 Entry composition [i](#)

There are 10 unique types of molecules in this entry. The entry contains 12533 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 Ferrichrome-iron receptor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	702	Total	C	N	O	S	0	0	0
			5466	3444	926	1082	14			
1	B	691	Total	C	N	O	S	0	0	0
			5394	3398	914	1068	14			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	406	SER	-	SEE REMARK 999	UNP P06971
A	407	SER	-	SEE REMARK 999	UNP P06971
A	408	HIS	-	SEE REMARK 999	UNP P06971
A	409	HIS	-	SEE REMARK 999	UNP P06971
A	410	HIS	-	SEE REMARK 999	UNP P06971
A	411	HIS	-	SEE REMARK 999	UNP P06971
A	412	HIS	-	SEE REMARK 999	UNP P06971
A	413	HIS	-	SEE REMARK 999	UNP P06971
A	414	GLY	-	SEE REMARK 999	UNP P06971
A	415	SER	-	SEE REMARK 999	UNP P06971
A	416	SER	-	SEE REMARK 999	UNP P06971
B	406	SER	-	SEE REMARK 999	UNP P06971
B	407	SER	-	SEE REMARK 999	UNP P06971
B	408	HIS	-	SEE REMARK 999	UNP P06971
B	409	HIS	-	SEE REMARK 999	UNP P06971
B	410	HIS	-	SEE REMARK 999	UNP P06971
B	411	HIS	-	SEE REMARK 999	UNP P06971
B	412	HIS	-	SEE REMARK 999	UNP P06971
B	413	HIS	-	SEE REMARK 999	UNP P06971
B	414	GLY	-	SEE REMARK 999	UNP P06971
B	415	SER	-	SEE REMARK 999	UNP P06971
B	416	SER	-	SEE REMARK 999	UNP P06971

- Molecule 2 is a protein called Protein tonB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	78	Total 624	C 393	N 120	O 109	S 2	0	0	0
2	D	78	Total 624	C 393	N 120	O 109	S 2	0	0	0

There are 44 discrepancies between the modelled and reference sequences:

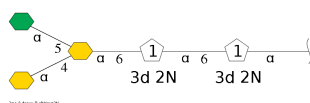
Chain	Residue	Modelled	Actual	Comment	Reference
C	11	GLY	-	SEE REMARK 999	UNP P94739
C	12	SER	-	SEE REMARK 999	UNP P94739
C	13	SER	-	SEE REMARK 999	UNP P94739
C	14	HIS	-	SEE REMARK 999	UNP P94739
C	15	HIS	-	SEE REMARK 999	UNP P94739
C	16	HIS	-	SEE REMARK 999	UNP P94739
C	17	HIS	-	SEE REMARK 999	UNP P94739
C	18	HIS	-	SEE REMARK 999	UNP P94739
C	19	HIS	-	SEE REMARK 999	UNP P94739
C	20	SER	-	SEE REMARK 999	UNP P94739
C	21	SER	-	SEE REMARK 999	UNP P94739
C	22	GLY	-	SEE REMARK 999	UNP P94739
C	23	LEU	-	SEE REMARK 999	UNP P94739
C	24	VAL	-	SEE REMARK 999	UNP P94739
C	25	PRO	-	SEE REMARK 999	UNP P94739
C	26	ARG	-	SEE REMARK 999	UNP P94739
C	27	GLY	-	SEE REMARK 999	UNP P94739
C	28	SER	-	SEE REMARK 999	UNP P94739
C	29	HIS	-	SEE REMARK 999	UNP P94739
C	30	MET	-	SEE REMARK 999	UNP P94739
C	31	SER	-	SEE REMARK 999	UNP P94739
C	32	VAL	-	SEE REMARK 999	UNP P94739
D	11	GLY	-	SEE REMARK 999	UNP P94739
D	12	SER	-	SEE REMARK 999	UNP P94739
D	13	SER	-	SEE REMARK 999	UNP P94739
D	14	HIS	-	SEE REMARK 999	UNP P94739
D	15	HIS	-	SEE REMARK 999	UNP P94739
D	16	HIS	-	SEE REMARK 999	UNP P94739
D	17	HIS	-	SEE REMARK 999	UNP P94739
D	18	HIS	-	SEE REMARK 999	UNP P94739
D	19	HIS	-	SEE REMARK 999	UNP P94739
D	20	SER	-	SEE REMARK 999	UNP P94739
D	21	SER	-	SEE REMARK 999	UNP P94739
D	22	GLY	-	SEE REMARK 999	UNP P94739
D	23	LEU	-	SEE REMARK 999	UNP P94739

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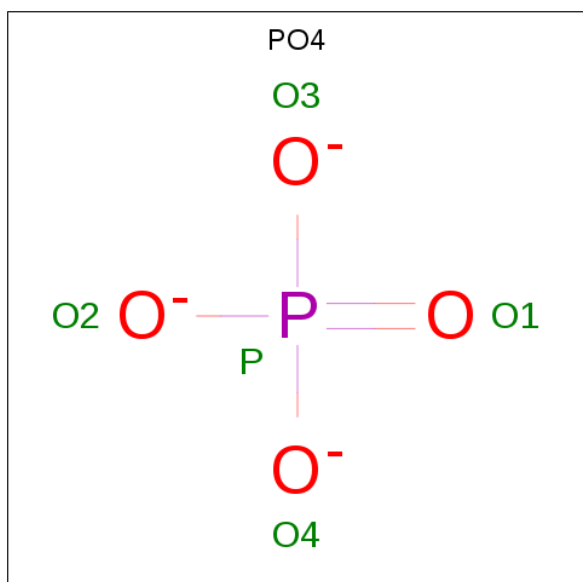
Chain	Residue	Modelled	Actual	Comment	Reference
D	24	VAL	-	SEE REMARK 999	UNP P94739
D	25	PRO	-	SEE REMARK 999	UNP P94739
D	26	ARG	-	SEE REMARK 999	UNP P94739
D	27	GLY	-	SEE REMARK 999	UNP P94739
D	28	SER	-	SEE REMARK 999	UNP P94739
D	29	HIS	-	SEE REMARK 999	UNP P94739
D	30	MET	-	SEE REMARK 999	UNP P94739
D	31	SER	-	SEE REMARK 999	UNP P94739
D	32	VAL	-	SEE REMARK 999	UNP P94739

- Molecule 3 is an oligosaccharide called 3-deoxy-alpha-D-manno-oct-2-ulopyranosonic acid-(2-4)-[L-glycero-alpha-D-manno-heptopyranose-(1-5)]3-deoxy-alpha-D-manno-oct-2-ulopyranosonic acid-(2-6)-2-amino-2,3-dideoxy-alpha-D-glucopyranose-(1-6)-2-amino-2,3-dideoxy-alpha-D-glucopyranose.



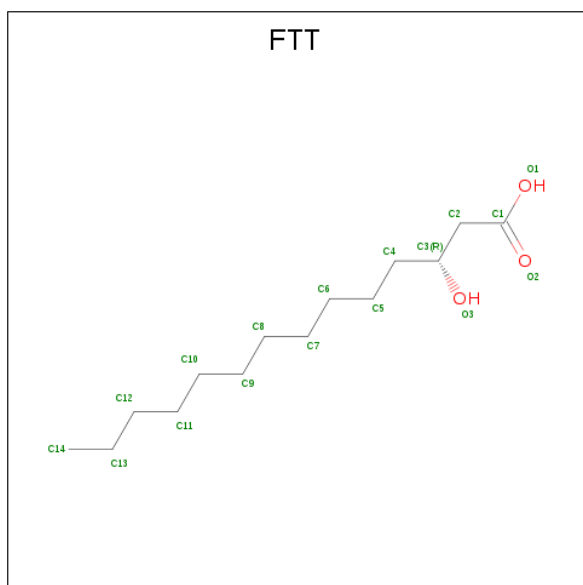
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	E	5	Total	C	N	O	0	0	0
			63	35	2	26			
3	F	5	Total	C	N	O	0	0	0
			63	35	2	26			

- Molecule 4 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



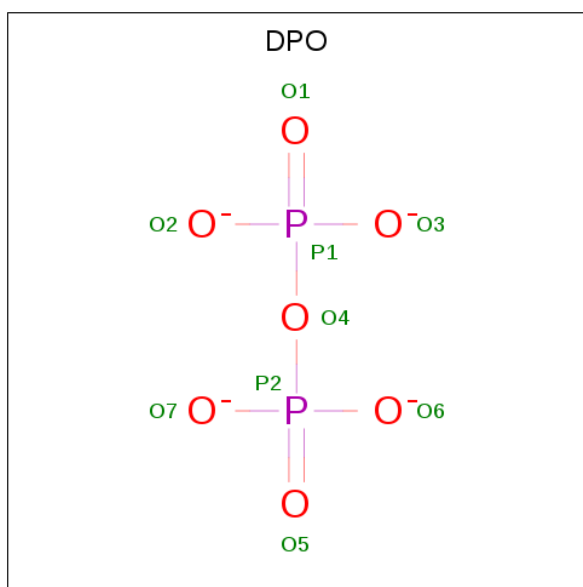
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total O P 4 3 1	0	0
4	B	1	Total O P 4 3 1	0	0

- Molecule 5 is 3-HYDROXY-TETRADECANOIC ACID (three-letter code: FTT) (formula: $C_{14}H_{28}O_3$).



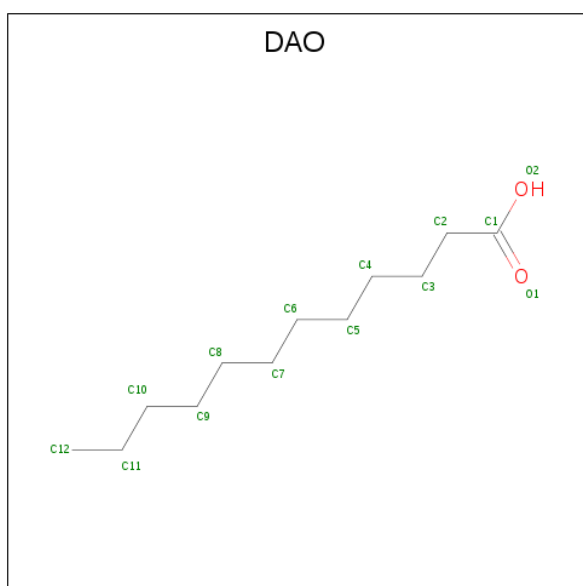
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C O 16 14 2	0	0
5	A	1	Total C O 16 14 2	0	0
5	A	1	Total C O 16 14 2	0	0
5	A	1	Total C O 16 14 2	0	0
5	B	1	Total C O 16 14 2	0	0
5	B	1	Total C O 16 14 2	0	0
5	B	1	Total C O 16 14 2	0	0
5	B	1	Total C O 16 14 2	0	0

- Molecule 6 is DIPHOSPHATE (three-letter code: DPO) (formula: O_7P_2).



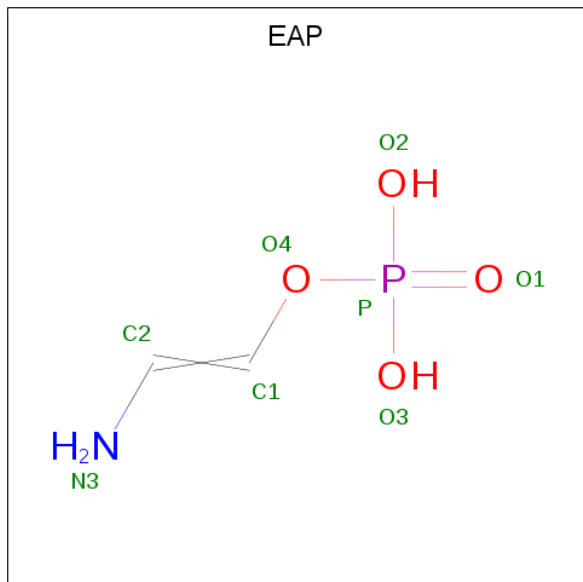
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	O	P	0	0
			8	6	2		
6	B	1	Total	O	P	0	0
			8	6	2		

- Molecule 7 is LAURIC ACID (three-letter code: DAO) (formula: C₁₂H₂₄O₂).



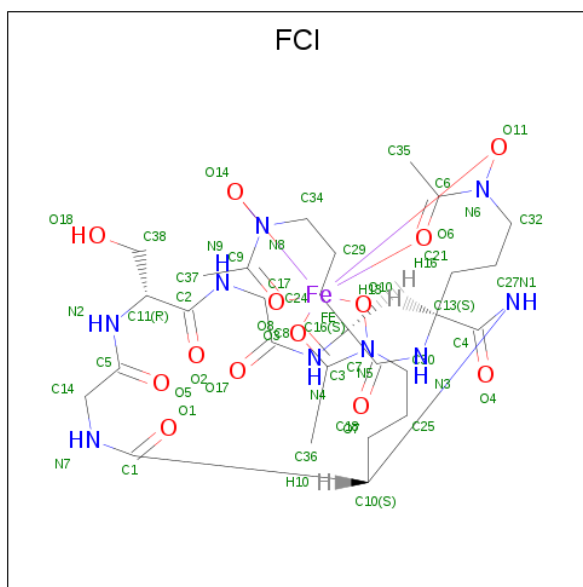
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			13	12	1		
7	B	1	Total	C	O	0	0
			13	12	1		

- Molecule 8 is 2-AMINO-VINYLP-HOSPHATE (three-letter code: EAP) (formula: $C_2H_6NO_4P$).



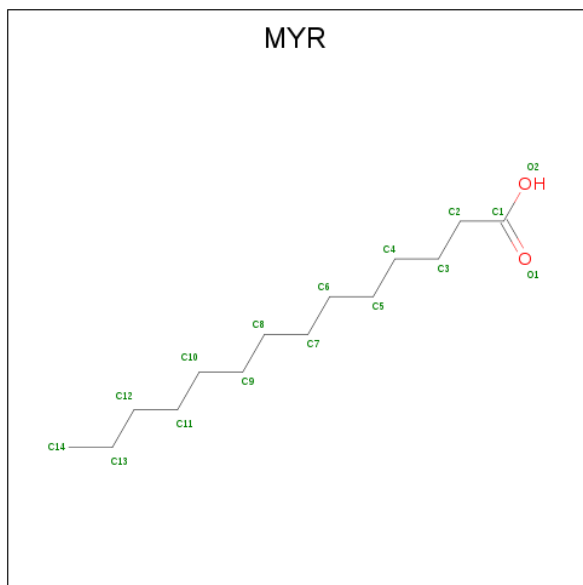
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
8	A	1	Total	C	N	O	P	0	0
			7	2	1	3	1		
8	B	1	Total	C	N	O	P	0	0
			7	2	1	3	1		

- Molecule 9 is FERRICROCIN-IRON (three-letter code: FCI) (formula: $C_{28}H_{44}FeN_9O_{13}$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
9	A	1	Total	C	Fe	N	O	
			46	24	1	9	12	
9	B	1	Total	C	Fe	N	O	
			46	24	1	9	12	

- Molecule 10 is MYRISTIC ACID (three-letter code: MYR) (formula: $C_{14}H_{28}O_2$).

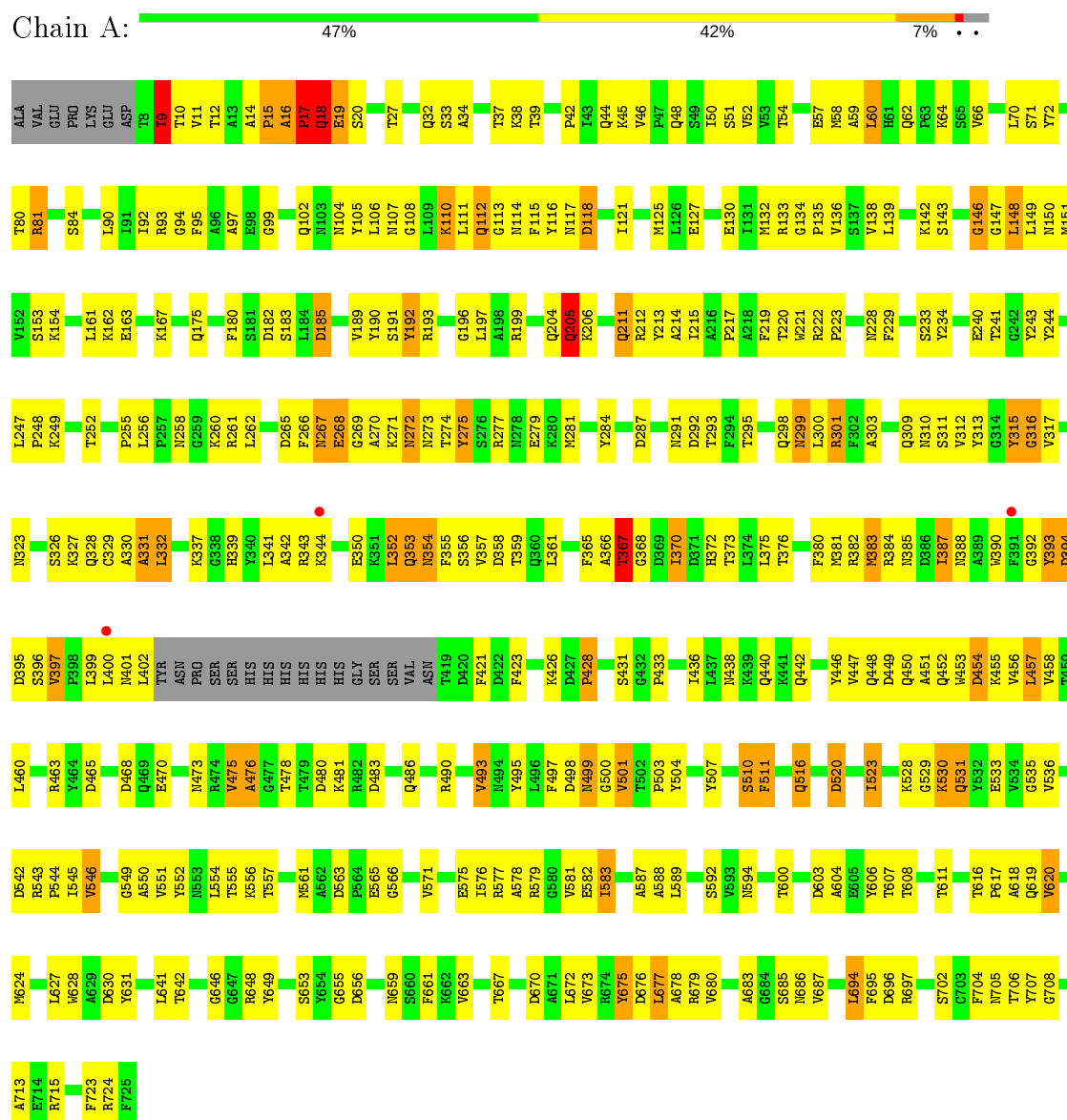


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	B	1	Total	C	O		
			15	14	1	0	0

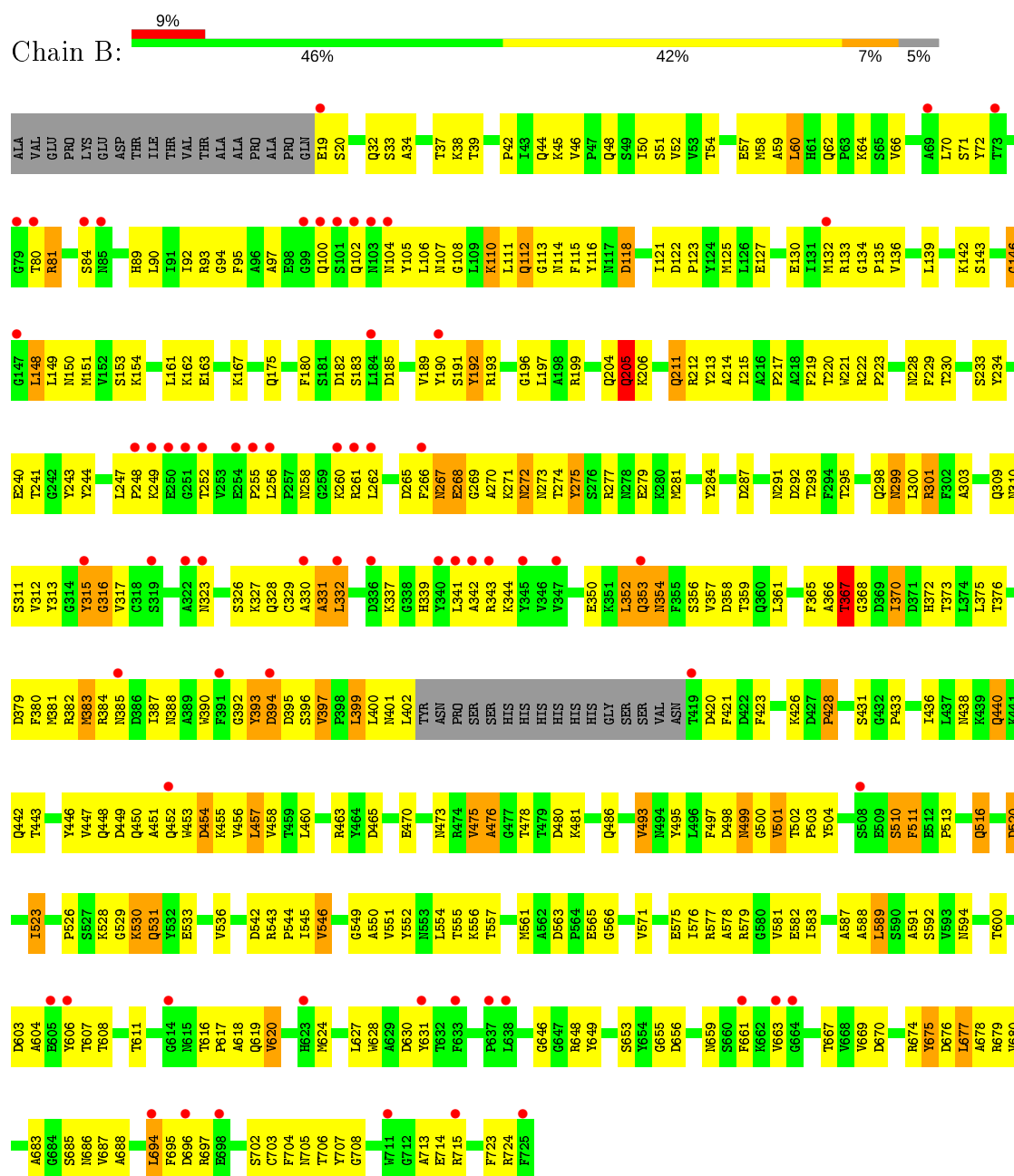
3 Residue-property plots

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.

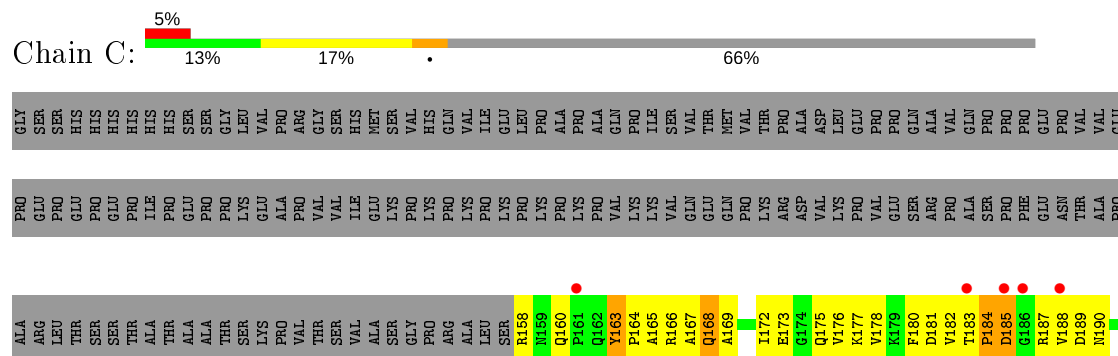
- Molecule 1: Ferrichrome-iron receptor



- Molecule 1: Ferrichrome-iron receptor



- Molecule 2: Protein tonB





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	234.32Å 91.84Å 138.51Å 90.00° 118.86° 90.00°	Depositor
Resolution (Å)	50.00 – 3.30 45.94 – 3.30	Depositor EDS
% Data completeness (in resolution range)	99.2 (50.00-3.30) 99.2 (45.94-3.30)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.33 (at 3.32Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.284 , 0.329 0.260 , 0.311	Depositor DCC
R_{free} test set	2325 reflections (6.01%)	wwPDB-VP
Wilson B-factor (Å ²)	93.6	Xtriage
Anisotropy	0.301	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.25 , 68.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	12533	wwPDB-VP
Average B, all atoms (Å ²)	98.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.62% 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: KDO, FTT, PO4, MYR, DAO, GCN, GMH, FCI, DPO, EAP

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	0.50	0/5599	0.69	1/7609 (0.0%)
1	B	0.40	0/5525	0.66	0/7504
2	C	0.35	0/636	0.56	0/857
2	D	0.31	0/636	0.54	0/857
All	All	0.44	0/12396	0.66	1/16827 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	17	PRO	N-CA-C	5.33	125.96	112.10

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	5466	0	5193	358	0
1	B	5394	0	5120	344	0
2	C	624	0	640	53	0
2	D	624	0	640	48	0
3	E	63	0	50	18	0
3	F	63	0	51	24	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	4	0	0	0	0
4	B	4	0	0	0	0
5	A	64	0	107	30	0
5	B	64	0	107	100	0
6	A	8	0	0	8	0
6	B	8	0	0	8	0
7	A	13	0	23	5	0
7	B	13	0	23	7	0
8	A	7	0	4	2	0
8	B	7	0	4	3	0
9	A	46	0	31	2	0
9	B	46	0	31	1	0
10	B	15	0	27	4	0
All	All	12533	0	12051	922	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 38.

The worst 5 of 922 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)
5:B:902:FTT:H92	5:B:903:FTT:C8	1.28	1.55
3:F:4:KDO:O1B	3:F:4:KDO:C1	1.70	1.39
5:B:902:FTT:C9	5:B:903:FTT:H82	1.52	1.39
5:B:902:FTT:C13	5:B:903:FTT:H131	1.55	1.34
5:B:901:FTT:C14	5:B:903:FTT:H112	1.58	1.33

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	698/725 (96%)	590 (84%)	77 (11%)	31 (4%)	2	16
1	B	687/725 (95%)	582 (85%)	79 (12%)	26 (4%)	3	19
2	C	76/229 (33%)	55 (72%)	11 (14%)	10 (13%)	0	1
2	D	76/229 (33%)	53 (70%)	13 (17%)	10 (13%)	0	1
All	All	1537/1908 (81%)	1280 (83%)	180 (12%)	77 (5%)	2	14

5 of 77 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	16	ALA
1	A	18	GLN
1	A	316	GLY
1	A	331	ALA
1	A	367	THR

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	579/601 (96%)	518 (90%)	61 (10%)	7	25
1	B	572/601 (95%)	516 (90%)	56 (10%)	8	29
2	C	66/200 (33%)	60 (91%)	6 (9%)	9	31
2	D	66/200 (33%)	61 (92%)	5 (8%)	13	39
All	All	1283/1602 (80%)	1155 (90%)	128 (10%)	7	27

5 of 128 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	620	VAL
1	B	192	TYR
2	C	168	GLN
1	A	679	ARG
1	B	80	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 73 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	692	ASN
1	B	205	GLN
2	C	208	ASN
1	B	107	ASN
1	B	267	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 ⓘ

10 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	$\# Z > 2$
3	GCN	E	1	3,5	10,10,11	0.81	0	13,13,15	0.76	0
3	GCN	E	2	3,5,4	10,10,11	1.03	1 (10%)	13,13,15	2.64	2 (15%)
3	KDO	E	3	3	12,15,16	0.73	0	16,21,24	0.86	0
3	KDO	E	4	3	12,15,16	0.47	0	16,21,24	0.89	0
3	GMH	E	5	8,3	13,13,14	0.80	0	17,18,20	0.67	0
3	GCN	F	1	3,5	10,10,11	0.81	0	13,13,15	0.76	0
3	GCN	F	2	3,5,4	10,10,11	1.04	1 (10%)	13,13,15	2.63	2 (15%)
3	KDO	F	3	3	12,15,16	0.74	0	16,21,24	0.85	0
3	KDO	F	4	3	12,15,16	0.46	0	16,21,24	0.89	0
3	GMH	F	5	8,3	13,13,14	0.80	0	17,18,20	0.68	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
3	GCN	E	1	3,5	-	2/2/15/18	0/1/1/1
3	GCN	E	2	3,5,4	-	0/2/15/18	0/1/1/1
3	KDO	E	3	3	-	0/6/26/30	0/1/1/1
3	KDO	E	4	3	-	6/6/26/30	0/1/1/1
3	GMH	E	5	8,3	-	2/6/23/26	0/1/1/1
3	GCN	F	1	3,5	-	2/2/15/18	0/1/1/1
3	GCN	F	2	3,5,4	-	0/2/15/18	0/1/1/1
3	KDO	F	3	3	-	0/6/26/30	0/1/1/1
3	KDO	F	4	3	-	6/6/26/30	0/1/1/1
3	GMH	F	5	8,3	-	2/6/23/26	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	F	2	GCN	C1-C2	2.72	1.55	1.52
3	E	2	GCN	C1-C2	2.68	1.55	1.52

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	2	GCN	C3-C2-C1	8.75	116.92	109.82
3	F	2	GCN	C3-C2-C1	8.71	116.90	109.82
3	E	2	GCN	C1-O5-C5	2.55	115.65	112.19
3	F	2	GCN	C1-O5-C5	2.52	115.61	112.19

There are no chirality outliers.

5 of 20 torsion outliers are listed below:

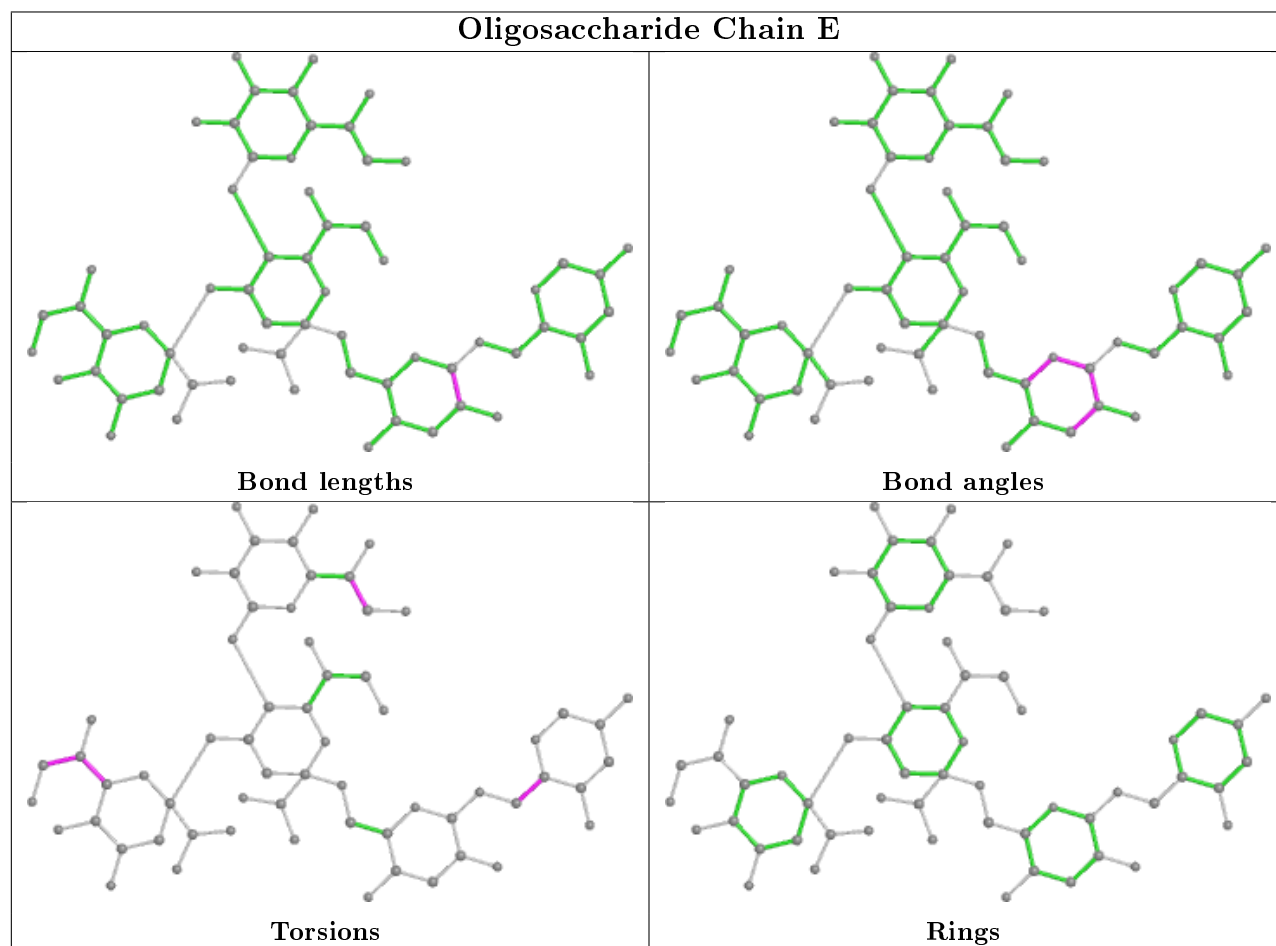
Mol	Chain	Res	Type	Atoms
3	E	4	KDO	C5-C6-C7-O7
3	E	4	KDO	C5-C6-C7-C8
3	E	4	KDO	O6-C6-C7-O7
3	E	4	KDO	O6-C6-C7-C8
3	F	4	KDO	C5-C6-C7-O7

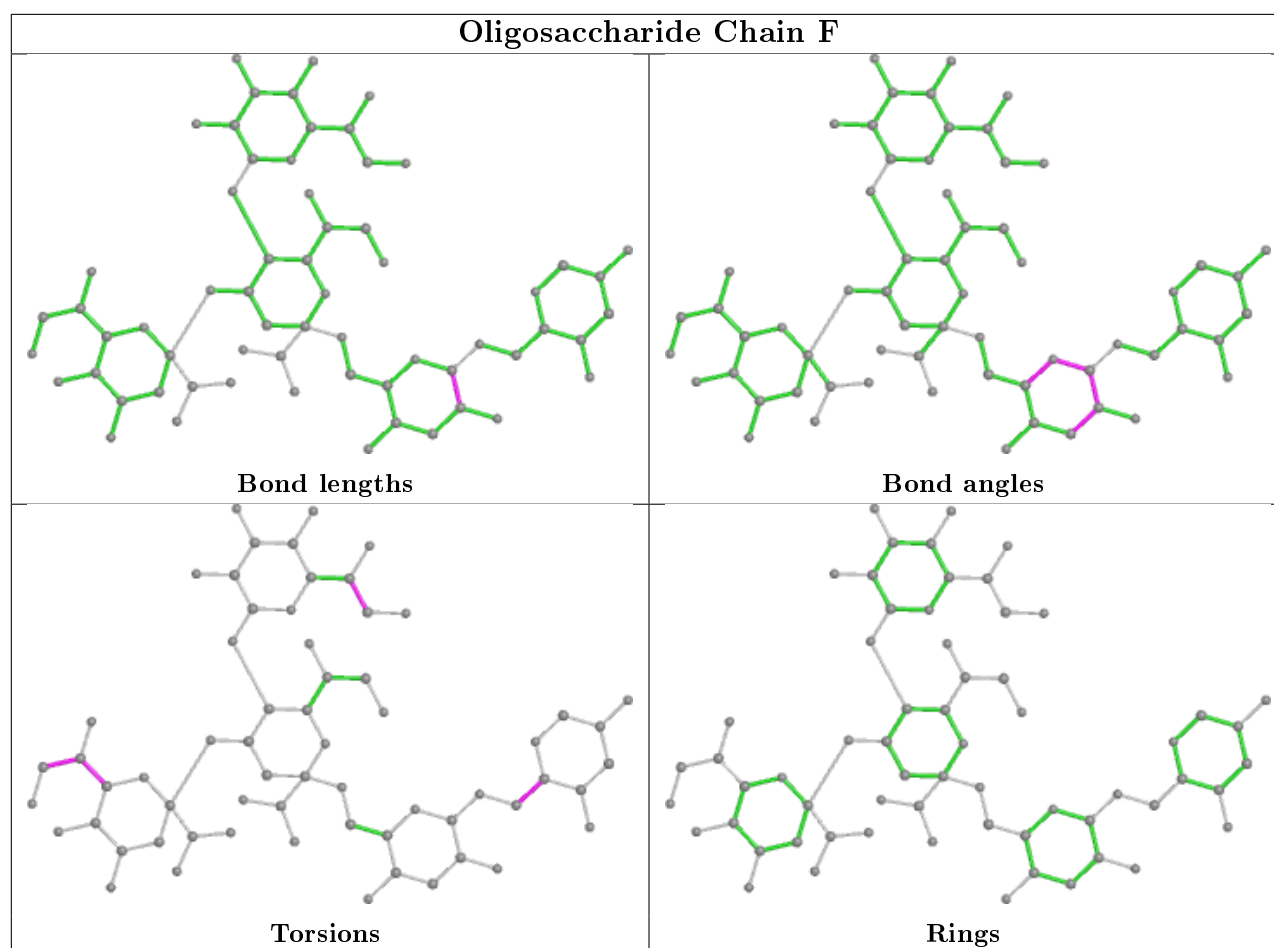
There are no ring outliers.

10 monomers are involved in 42 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	E	4	KDO	3	0
3	F	2	GCN	7	0
3	E	2	GCN	5	0
3	E	3	KDO	3	0
3	E	1	GCN	7	0
3	F	4	KDO	7	0
3	F	3	KDO	3	0
3	E	5	GMH	4	0
3	F	5	GMH	4	0
3	F	1	GCN	7	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.





5.6 Ligand geometry [i](#)

19 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
9	FCI	A	1050	-	45,51,56	1.43	4 (8%)	51,78,87	1.15	4 (7%)
10	MYR	B	940	5	14,14,15	0.99	1 (7%)	13,13,15	0.81	1 (7%)
8	EAP	B	980	3	3,6,7	2.86	1 (33%)	0,6,9	0.00	-
5	FTT	B	901	-	15,15,16	0.33	0	15,15,17	0.80	0
5	FTT	B	900	3	15,15,16	4.13	1 (6%)	15,15,17	2.26	3 (20%)
4	PO4	B	950	3	0,3,4	0.00	-	0,3,6	0.00	-

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	FTT	A	900	3	15,15,16	0.43	0	15,15,17	0.70	1 (6%)
4	PO4	A	950	3	0,3,4	0.00	-	0,3,6	0.00	-
6	DPO	A	910	-	4,7,8	2.44	1 (25%)	5,10,13	0.96	0
7	DAO	A	930	5	12,12,13	1.05	1 (8%)	11,11,13	0.80	1 (9%)
9	FCI	B	1050	-	45,51,56	1.44	4 (8%)	51,78,87	1.15	4 (7%)
5	FTT	B	903	10	15,15,16	0.51	0	15,15,17	0.51	0
5	FTT	A	901	-	15,15,16	0.33	0	15,15,17	0.80	0
8	EAP	A	980	3	3,6,7	2.87	1 (33%)	0,6,9	0.00	-
6	DPO	B	910	-	4,7,8	2.44	1 (25%)	5,10,13	0.96	0
5	FTT	A	902	3,7	15,15,16	0.33	0	15,15,17	0.80	0
5	FTT	B	902	3,7	15,15,16	0.33	0	15,15,17	0.80	0
7	DAO	B	930	5	12,12,13	1.05	1 (8%)	11,11,13	0.80	1 (9%)
5	FTT	A	903	-	15,15,16	0.51	0	15,15,17	0.51	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
9	FCI	A	1050	-	-	11/57/102/116	0/3/6/6
10	MYR	B	940	5	-	2/12/12/13	-
8	EAP	B	980	3	-	0/0/4/5	-
5	FTT	B	901	-	-	5/14/14/15	-
5	FTT	B	900	3	-	2/14/14/15	-
5	FTT	A	900	3	-	0/14/14/15	-
6	DPO	A	910	-	-	0/2/5/6	-
7	DAO	A	930	5	-	1/10/10/11	-
9	FCI	B	1050	-	-	11/57/102/116	0/3/6/6
5	FTT	B	903	10	-	2/14/14/15	-
5	FTT	A	901	-	-	2/14/14/15	-
8	EAP	A	980	3	-	0/0/4/5	-
6	DPO	B	910	-	-	0/2/5/6	-
5	FTT	A	902	3,7	-	3/14/14/15	-
5	FTT	B	902	3,7	-	3/14/14/15	-
7	DAO	B	930	5	-	1/10/10/11	-
5	FTT	A	903	-	-	2/14/14/15	-

The worst 5 of 16 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	900	FTT	C9-C8	15.93	2.41	1.51
9	A	1050	FCI	O10-FE	5.97	2.12	1.99
9	B	1050	FCI	O10-FE	5.95	2.12	1.99
9	B	1050	FCI	O3-FE	5.15	2.19	2.04
9	A	1050	FCI	O3-FE	5.10	2.19	2.04

The worst 5 of 15 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	900	FTT	C9-C8-C7	7.96	154.83	114.42
9	A	1050	FCI	C1-C10-N1	3.46	120.58	111.16
9	B	1050	FCI	C1-C10-N1	3.46	120.57	111.16
9	A	1050	FCI	C34-N8-C9	2.20	130.56	125.65
9	B	1050	FCI	C34-N8-C9	2.20	130.56	125.65

There are no chirality outliers.

5 of 45 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	901	FTT	C1-C2-C3-O3
5	A	901	FTT	C2-C3-C4-C5
9	A	1050	FCI	C25-C30-N5-C3
9	B	1050	FCI	C25-C30-N5-C3
5	A	901	FTT	O3-C3-C4-C5

There are no ring outliers.

17 monomers are involved in 138 short contacts:

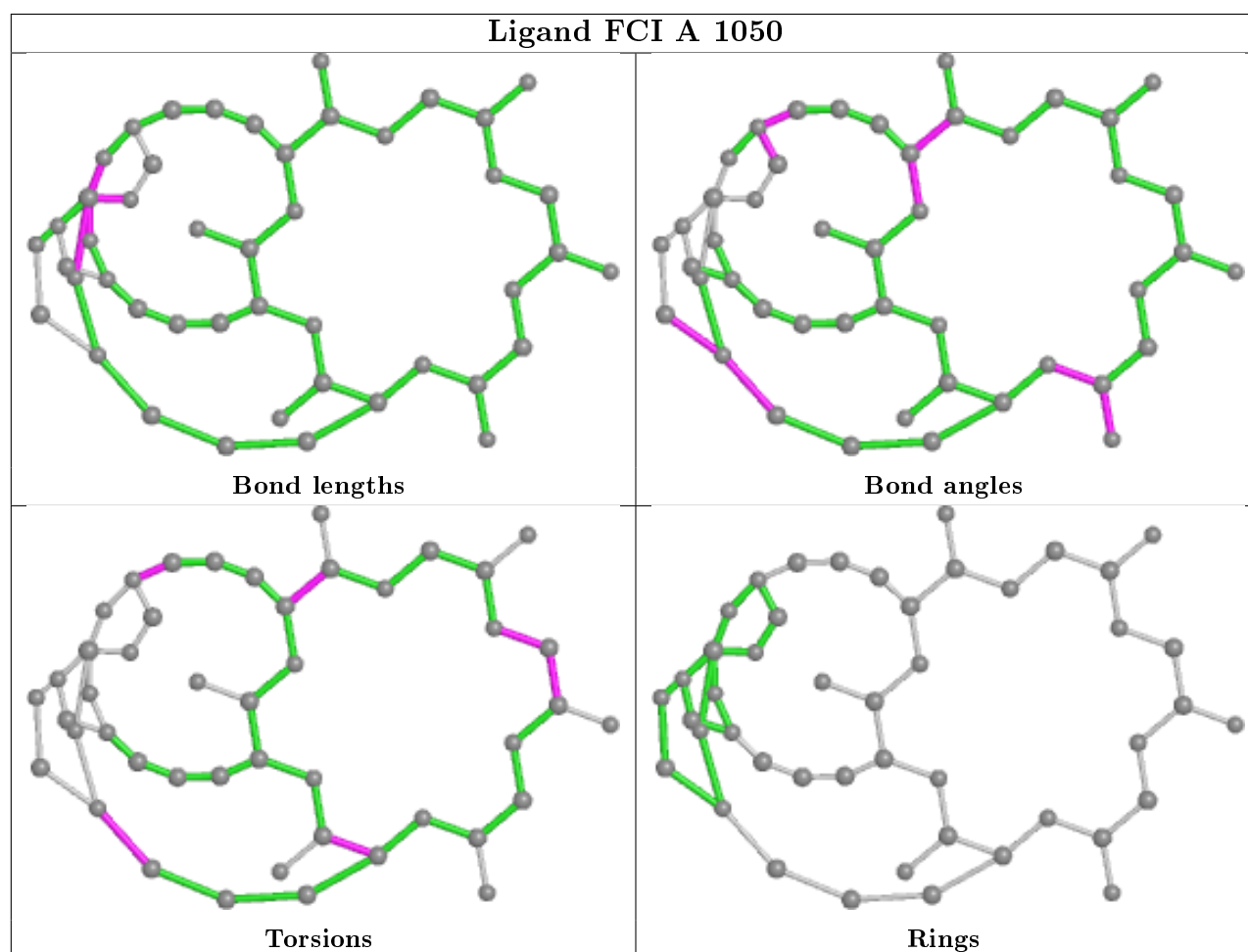
Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	A	1050	FCI	2	0
10	B	940	MYR	4	0
8	B	980	EAP	3	0
5	B	901	FTT	39	0
5	B	900	FTT	17	0
5	A	900	FTT	12	0
6	A	910	DPO	8	0
7	A	930	DAO	5	0
9	B	1050	FCI	1	0
5	B	903	FTT	64	0
5	A	901	FTT	9	0
8	A	980	EAP	2	0
6	B	910	DPO	8	0

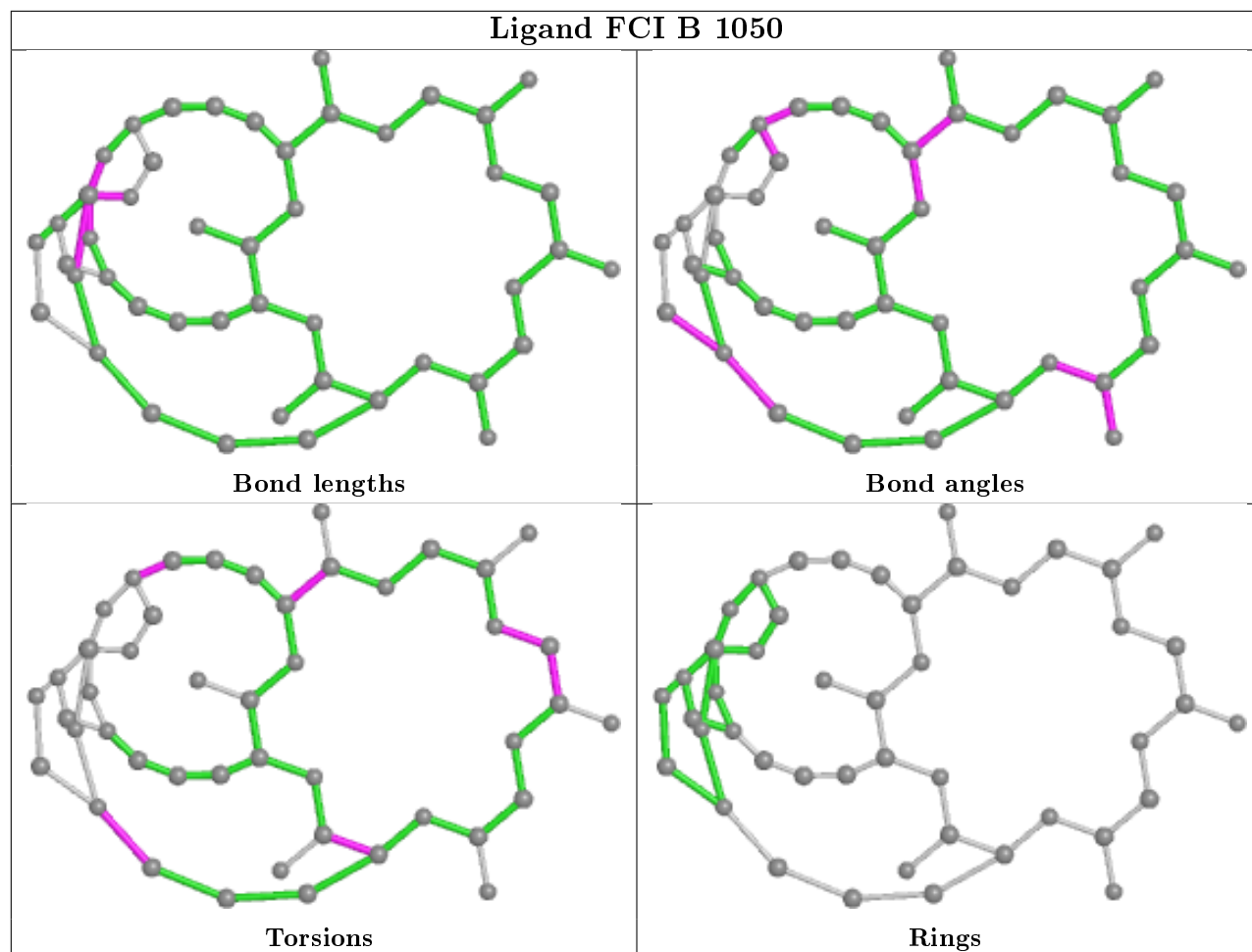
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	902	FTT	6	0
5	B	902	FTT	39	0
7	B	930	DAO	7	0
5	A	903	FTT	3	0

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 [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	702/725 (96%)	-0.16	3 (0%) 92 93	24, 61, 110, 177	0
1	B	691/725 (95%)	0.45	66 (9%) 8 8	44, 115, 186, 210	0
2	C	78/229 (34%)	0.76	11 (14%) 2 2	43, 112, 179, 209	0
2	D	78/229 (34%)	2.38	38 (48%) 0 0	102, 187, 210, 210	0
All	All	1549/1908 (81%)	0.29	118 (7%) 13 13	24, 88, 186, 210	0

The worst 5 of 118 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	178	VAL	7.9
1	B	255	PRO	7.8
2	D	225	VAL	7.6
2	D	159	ASN	7.3
1	B	250	GLU	6.8

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.

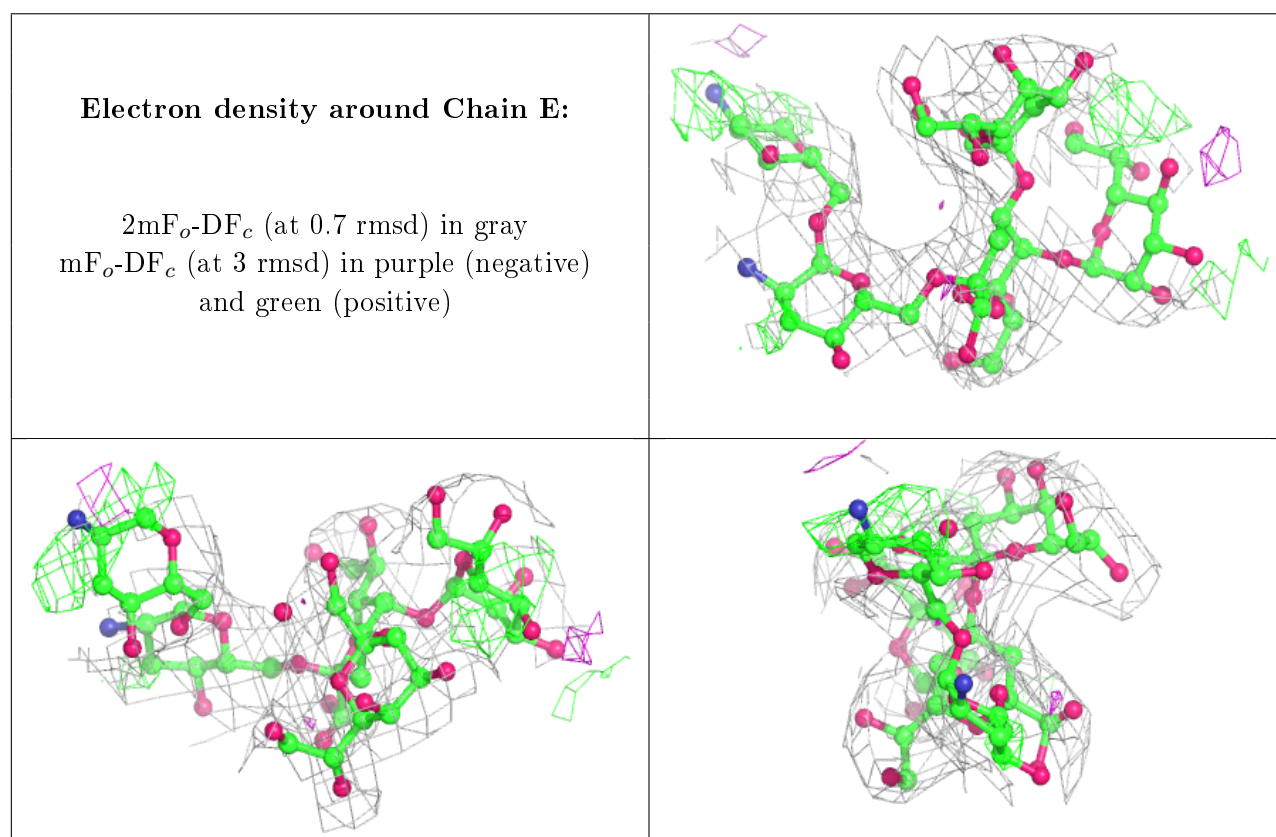
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	KDO	F	4	15/16	0.77	0.24	73,80,98,98	0
3	GMH	F	5	13/14	0.87	0.19	62,67,79,81	0

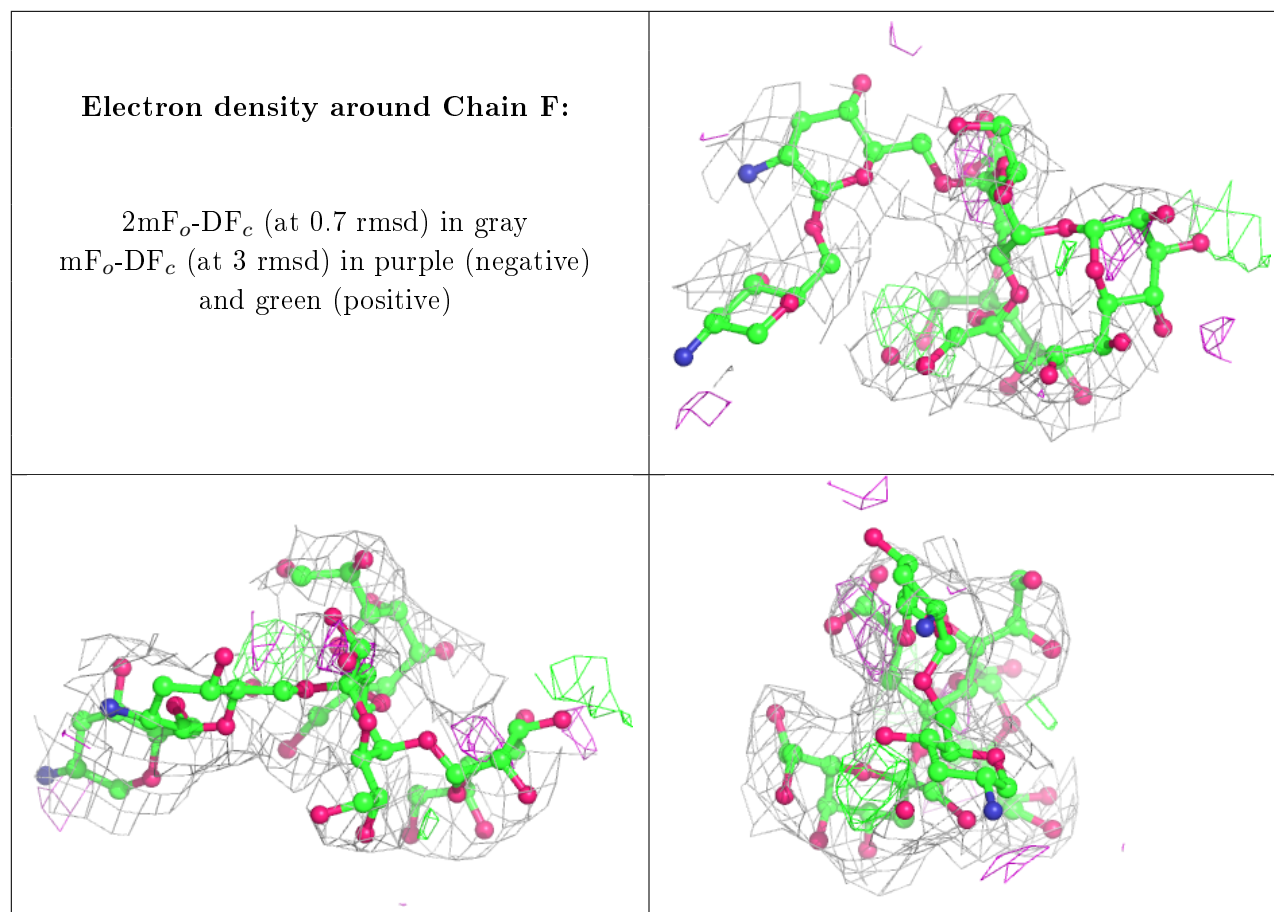
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	KDO	E	4	15/16	0.88	0.20	73,80,98,98	0
3	GMH	E	5	13/14	0.88	0.14	62,67,79,81	0
3	GCN	E	1	10/11	0.88	0.23	62,74,79,87	0
3	GCN	F	2	10/11	0.89	0.08	49,60,64,74	0
3	GCN	E	2	10/11	0.91	0.14	49,60,64,74	0
3	KDO	E	3	15/16	0.92	0.17	65,72,98,98	0
3	KDO	F	3	15/16	0.93	0.21	65,72,98,98	0
3	GCN	F	1	10/11	0.94	0.11	62,74,79,87	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.

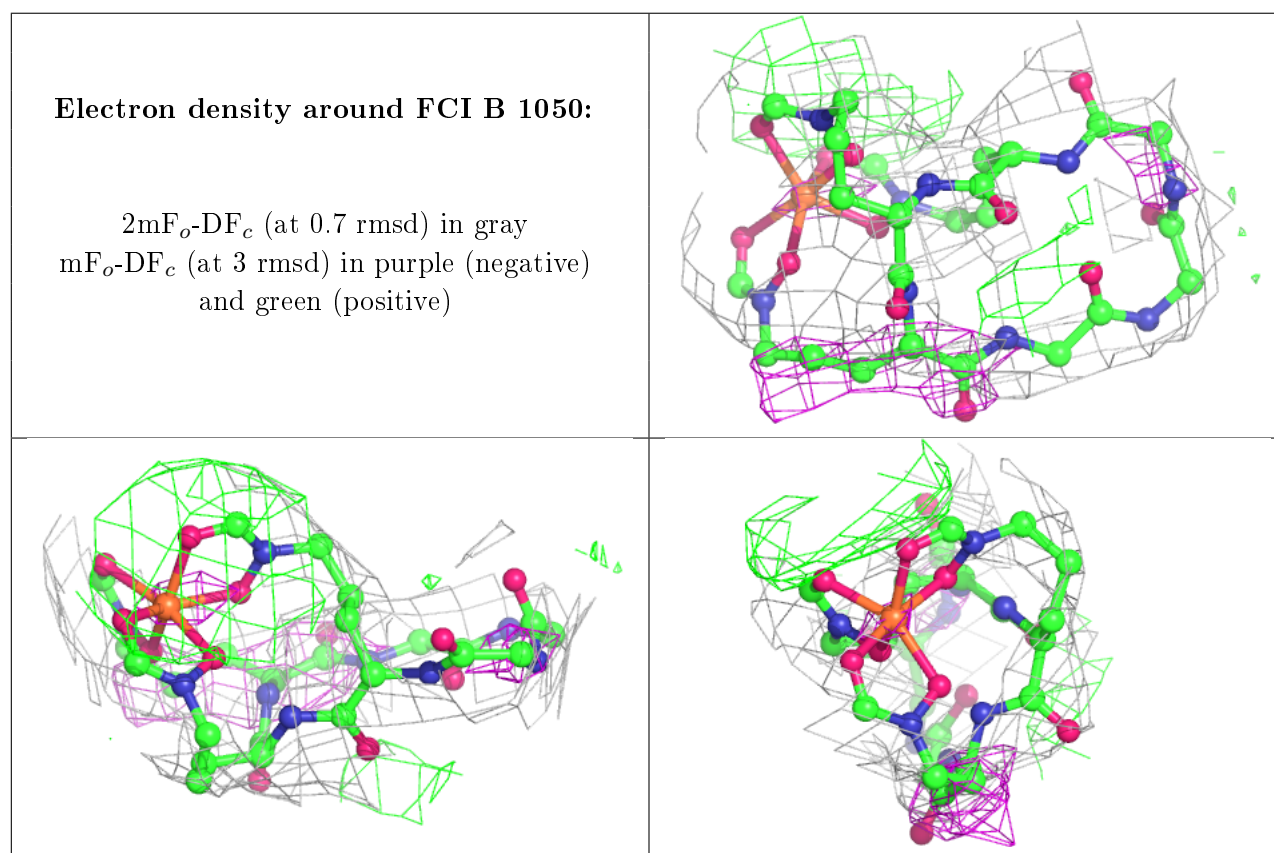
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	FTT	B	903	16/17	0.42	0.88	75,88,92,97	0
5	FTT	B	901	16/17	0.63	0.73	63,70,78,79	0
6	DPO	B	910	8/9	0.69	0.22	62,80,93,95	0
6	DPO	A	910	8/9	0.70	0.22	62,80,93,95	0
10	MYR	B	940	15/16	0.73	1.04	69,86,94,95	0
5	FTT	A	903	16/17	0.74	0.47	75,88,92,97	0
7	DAO	A	930	13/14	0.78	0.60	69,75,82,86	0
5	FTT	A	901	16/17	0.79	0.43	63,70,78,79	0
5	FTT	A	900	16/17	0.80	0.46	76,87,96,102	0
9	FCI	B	1050	46/51	0.81	0.25	44,57,70,81	0
7	DAO	B	930	13/14	0.82	0.40	69,75,82,86	0
5	FTT	B	900	16/17	0.83	0.41	84,95,98,102	0

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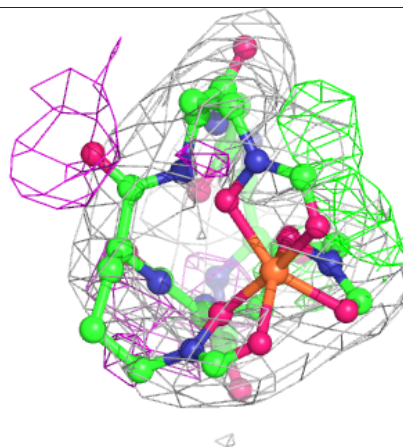
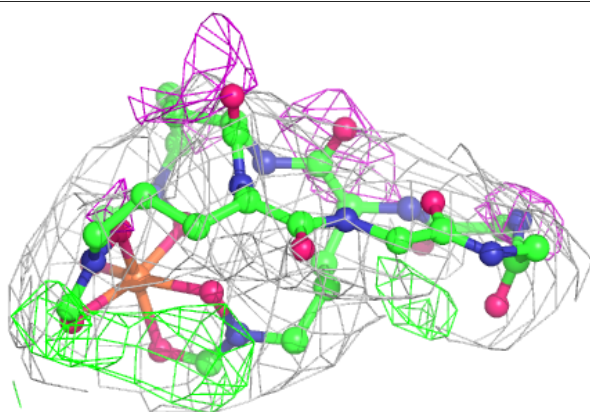
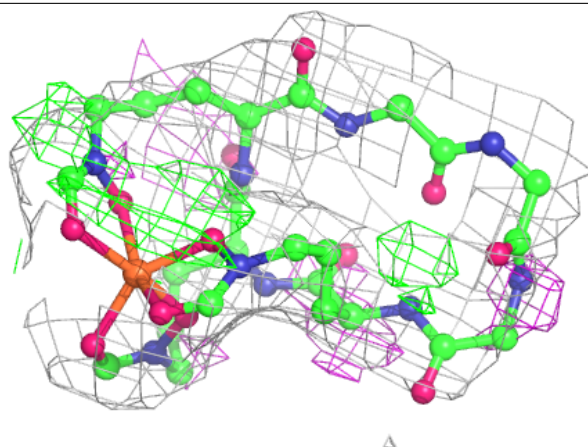
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	FTT	B	902	16/17	0.83	0.57	63,70,78,79	0
9	FCI	A	1050	46/51	0.87	0.28	44,57,70,81	0
5	FTT	A	902	16/17	0.88	0.47	63,70,78,79	0
8	EAP	A	980	7/8	0.89	0.14	81,87,93,95	0
8	EAP	B	980	7/8	0.89	0.16	81,87,93,95	0
4	PO4	A	950	4/5	0.94	0.10	71,74,76,81	0
4	PO4	B	950	4/5	0.95	0.13	71,74,76,81	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 FCI A 1050:

$2mF_o-DF_c$ (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.