



wwPDB X-ray Structure Validation Summary Report ⓘ

May 13, 2020 – 11:09 am BST

PDB ID : 3V83
Title : The 2.1 angstrom crystal structure of diferric human transferrin
Authors : Noinaj, N.; Steere, A.; Mason, A.B.; Buchanan, S.K.
Deposited on : 2011-12-22
Resolution : 2.10 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 : 2.11
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) : 2.11

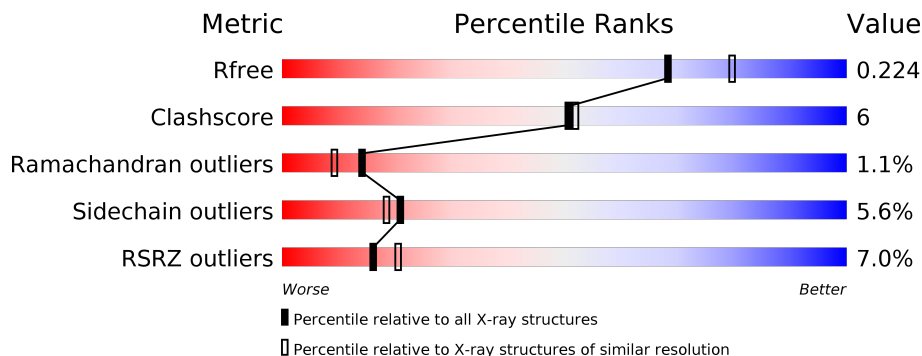
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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	698	
1	B	698	
1	C	698	
1	D	698	
1	E	698	
1	F	698	

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
2	BCT	C	702	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 33928 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.

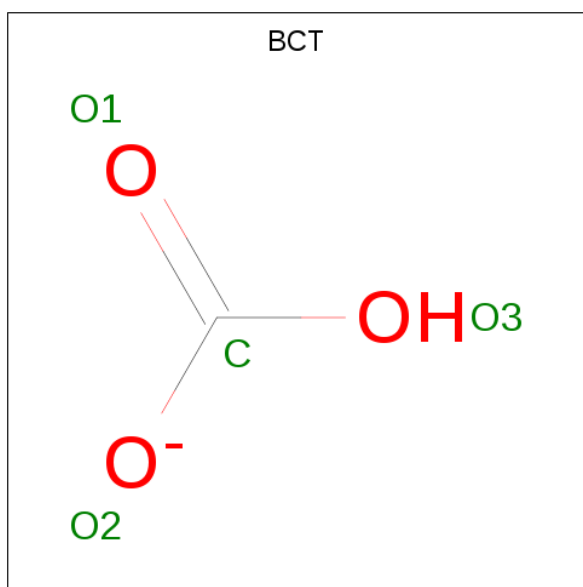
- Molecule 1 is a protein called Serotransferrin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	674	5160	3247	892	974	47	0	0	0
1	B	677	5179	3260	898	974	47	0	0	0
1	C	674	5124	3221	888	968	47	0	0	0
1	D	675	5193	3263	897	986	47	0	0	0
1	E	674	5202	3267	898	990	47	0	0	0
1	F	676	5191	3263	900	981	47	0	0	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	429	VAL	ILE	VARIANT	UNP P02787
B	429	VAL	ILE	VARIANT	UNP P02787
C	429	VAL	ILE	VARIANT	UNP P02787
D	429	VAL	ILE	VARIANT	UNP P02787
E	429	VAL	ILE	VARIANT	UNP P02787
F	429	VAL	ILE	VARIANT	UNP P02787

- Molecule 2 is BICARBONATE ION (three-letter code: BCT) (formula: CHO_3).

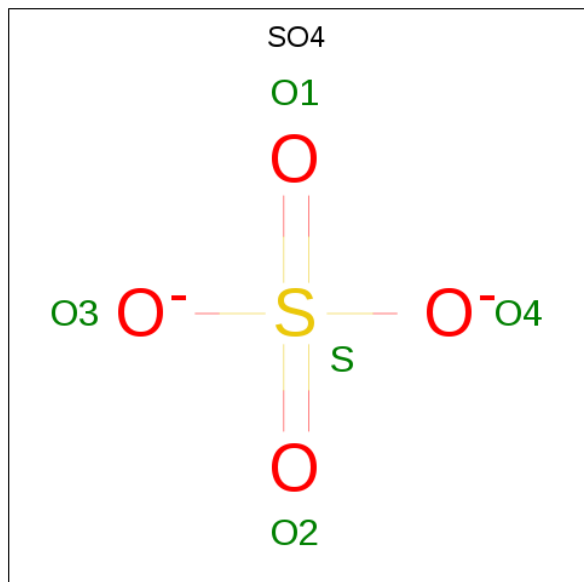


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 4 1 3	0	0
2	A	1	Total C O 4 1 3	0	0
2	B	1	Total C O 4 1 3	0	0
2	B	1	Total C O 4 1 3	0	0
2	C	1	Total C O 4 1 3	0	0
2	C	1	Total C O 4 1 3	0	0
2	D	1	Total C O 4 1 3	0	0
2	D	1	Total C O 4 1 3	0	0
2	E	1	Total C O 4 1 3	0	0
2	E	1	Total C O 4 1 3	0	0
2	F	1	Total C O 4 1 3	0	0
2	F	1	Total C O 4 1 3	0	0

- Molecule 3 is FE (III) ION (three-letter code: FE) (formula: Fe).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	D	2	Total Fe 2 2	0	0
3	E	2	Total Fe 2 2	0	0
3	B	2	Total Fe 2 2	0	0
3	C	2	Total Fe 2 2	0	0
3	A	2	Total Fe 2 2	0	0
3	F	2	Total Fe 2 2	0	0

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total O S 5 4 1	0	0
4	A	1	Total O S 5 4 1	0	0
4	A	1	Total O S 5 4 1	0	0
4	A	1	Total O S 5 4 1	0	0
4	A	1	Total O S 5 4 1	0	0
4	B	1	Total O S 5 4 1	0	0

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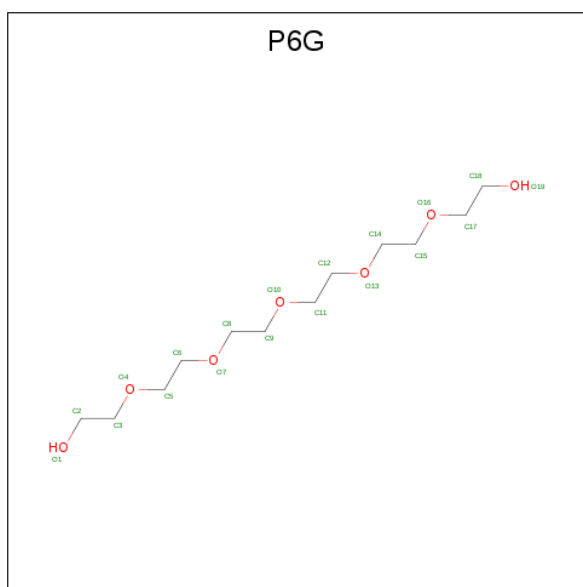
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	D	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	E	1	Total	O	S	0	0
			5	4	1		
4	E	1	Total	O	S	0	0
			5	4	1		
4	E	1	Total	O	S	0	0
			5	4	1		
4	E	1	Total	O	S	0	0
			5	4	1		
4	E	1	Total	O	S	0	0
			5	4	1		
4	E	1	Total	O	S	0	0
			5	4	1		
4	E	1	Total	O	S	0	0
			5	4	1		
4	E	1	Total	O	S	0	0
			5	4	1		
4	F	1	Total	O	S	0	0
			5	4	1		
4	F	1	Total	O	S	0	0
			5	4	1		
4	F	1	Total	O	S	0	0
			5	4	1		
4	F	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is HEXAETHYLENE GLYCOL (three-letter code: P6G) (formula: C₁₂H₂₆O₇).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			19	12	7		
5	A	1	Total	C	O	0	0
			19	12	7		
5	A	1	Total	C	O	0	0
			19	12	7		
5	B	1	Total	C	O	0	0
			19	12	7		
5	B	1	Total	C	O	0	0
			19	12	7		
5	C	1	Total	C	O	0	0
			19	12	7		
5	C	1	Total	C	O	0	0
			19	12	7		
5	C	1	Total	C	O	0	0
			19	12	7		
5	D	1	Total	C	O	0	0
			19	12	7		
5	D	1	Total	C	O	0	0
			19	12	7		
5	E	1	Total	C	O	0	0
			19	12	7		
5	E	1	Total	C	O	0	0
			19	12	7		
5	F	1	Total	C	O	0	0
			19	12	7		
5	F	1	Total	C	O	0	0
			19	12	7		

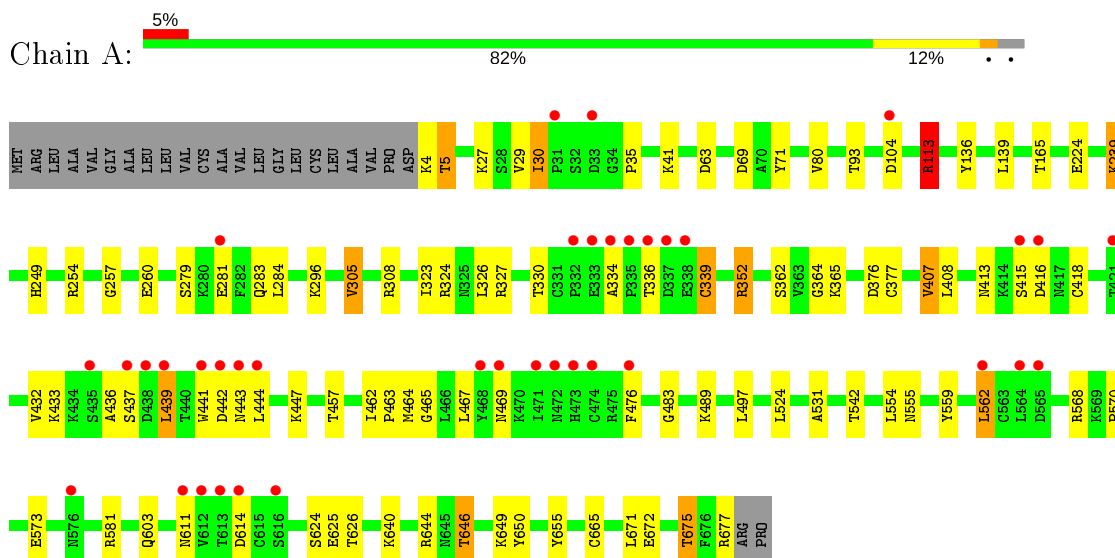
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	352	Total 352	O 352	0	0
6	B	454	Total 454	O 454	0	0
6	C	305	Total 305	O 305	0	0
6	D	398	Total 398	O 398	0	0
6	E	431	Total 431	O 431	0	0
6	F	393	Total 393	O 393	0	0

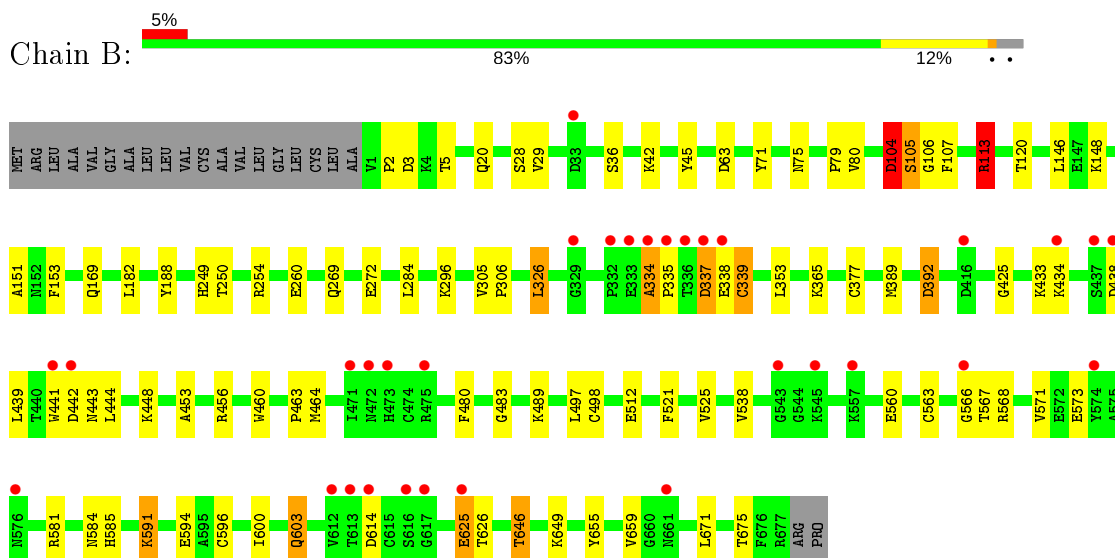
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

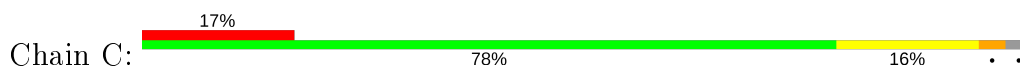
- Molecule 1: Serotransferrin

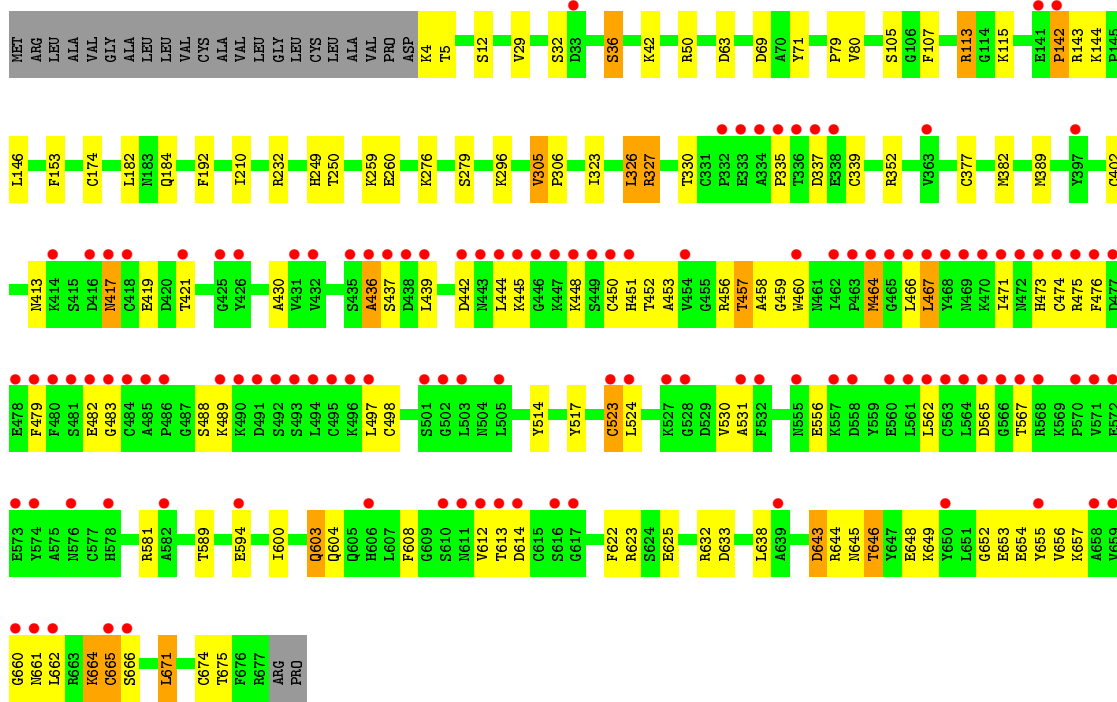


- Molecule 1: Serotransferrin

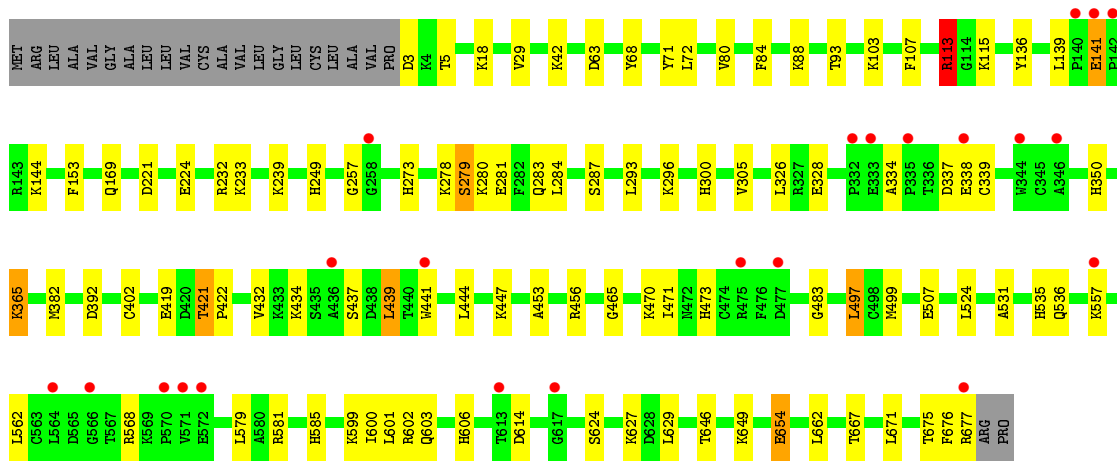
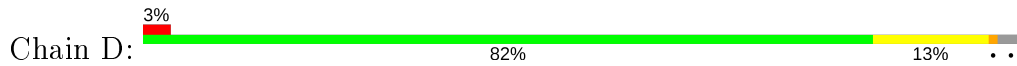


- Molecule 1: Serotransferrin

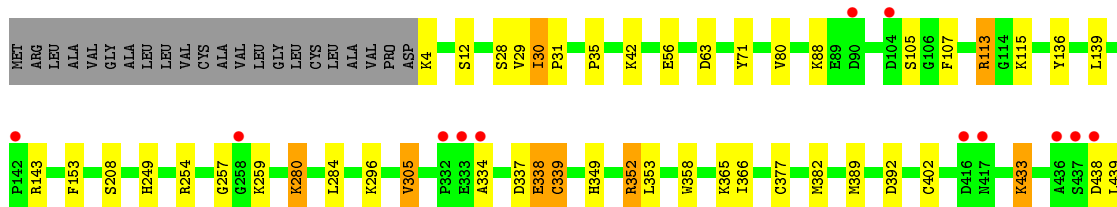
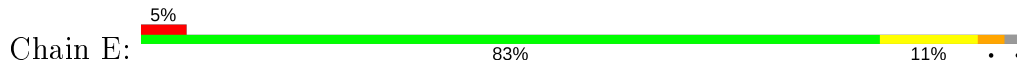


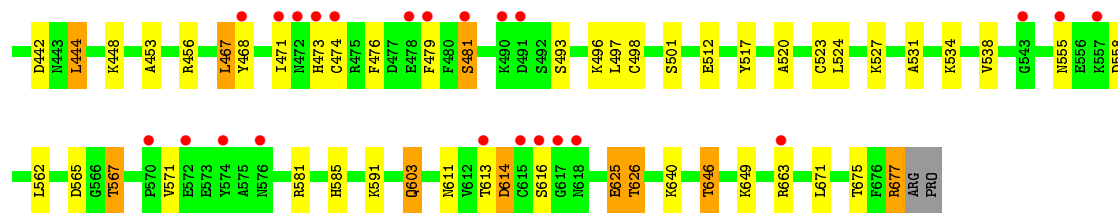


• Molecule 1: Serotransferrin

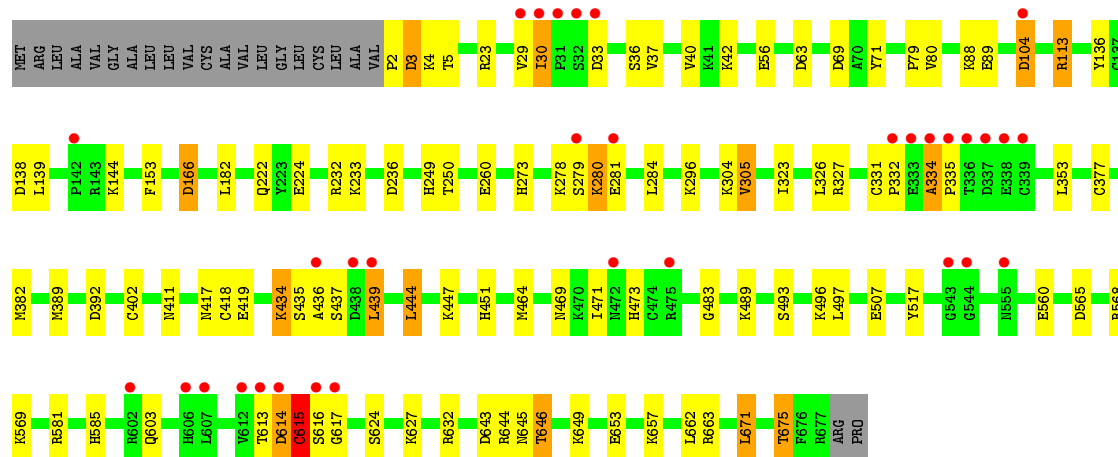
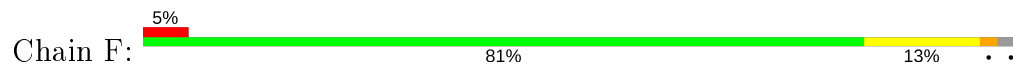


• Molecule 1: Serotransferrin





● Molecule 1: Serotransferrin



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	254.53Å 173.00Å 150.15Å 90.00° 123.26° 90.00°	Depositor
Resolution (Å)	29.95 – 2.10 49.25 – 2.10	Depositor EDS
% Data completeness (in resolution range)	94.4 (29.95-2.10) 94.4 (49.25-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.83 (at 2.10Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, R_{free}	0.180 , 0.230 0.181 , 0.224	Depositor DCC
R_{free} test set	15031 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	27.0	Xtrriage
Anisotropy	0.250	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 60.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.014 for -h-2*k,l	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	33928	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.33% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical 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: BCT, FE, P6G, SO4

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	0.38	0/5278	0.54	1/7144 (0.0%)
1	B	0.42	0/5298	0.57	2/7169 (0.0%)
1	C	0.39	0/5240	0.54	0/7097
1	D	0.41	0/5311	0.57	2/7184 (0.0%)
1	E	0.42	0/5320	0.57	2/7196 (0.0%)
1	F	0.40	0/5309	0.58	2/7183 (0.0%)
All	All	0.40	0/31756	0.56	9/42973 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	D	0	2
1	E	0	2
All	All	0	4

There are no bond length outliers.

The worst 5 of 9 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	113	ARG	NE-CZ-NH1	6.35	123.47	120.30
1	E	113	ARG	NE-CZ-NH1	6.23	123.41	120.30
1	D	113	ARG	NE-CZ-NH1	6.04	123.32	120.30
1	B	113	ARG	NE-CZ-NH2	-5.60	117.50	120.30
1	F	113	ARG	NE-CZ-NH2	-5.23	117.69	120.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	278	LYS	Peptide
1	D	279	SER	Peptide
1	E	613	THR	Peptide
1	E	614	ASP	Peptide

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5160	0	4928	50	0
1	B	5179	0	4949	54	0
1	C	5124	0	4868	84	0
1	D	5193	0	4970	57	0
1	E	5202	0	4992	51	0
1	F	5191	0	4969	62	0
2	A	8	0	1	0	0
2	B	8	0	0	0	0
2	C	8	0	1	3	0
2	D	8	0	1	0	0
2	E	8	0	1	0	0
2	F	8	0	1	0	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
3	C	2	0	0	0	0
3	D	2	0	0	0	0
3	E	2	0	0	0	0
3	F	2	0	0	0	0
4	A	25	0	0	1	0
4	B	45	0	0	1	0
4	C	35	0	0	0	0
4	D	45	0	0	1	0
4	E	45	0	0	0	0
4	F	25	0	0	0	0
5	A	57	0	78	4	0
5	B	38	0	52	6	0
5	C	57	0	78	4	0
5	D	38	0	52	7	0
5	E	38	0	52	4	0
5	F	38	0	52	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	A	352	0	0	6	0
6	B	454	0	0	9	0
6	C	305	0	0	10	0
6	D	398	0	0	10	0
6	E	431	0	0	9	0
6	F	393	0	0	5	0
All	All	33928	0	30045	353	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 6.

The worst 5 of 353 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:C:436:ALA:H	1:C:437:SER:HA	1.23	1.00
1:B:566:GLY:H	1:B:567:THR:HA	1.39	0.85
1:A:113:ARG:HD2	1:E:254:ARG:HD2	1.59	0.82
1:A:646:THR:HG22	1:A:649:LYS:H	1.48	0.79
1:E:646:THR:HG22	1:E:649:LYS:H	1.48	0.78

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	672/698 (96%)	621 (92%)	42 (6%)	9 (1%)	12 7
1	B	675/698 (97%)	630 (93%)	39 (6%)	6 (1%)	17 12
1	C	672/698 (96%)	617 (92%)	46 (7%)	9 (1%)	12 7
1	D	673/698 (96%)	630 (94%)	37 (6%)	6 (1%)	17 12
1	E	672/698 (96%)	632 (94%)	35 (5%)	5 (1%)	22 18

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	F	674/698 (97%)	623 (92%)	43 (6%)	8 (1%)	13	8
All	All	4038/4188 (96%)	3753 (93%)	242 (6%)	43 (1%)	14	9

5 of 43 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	416	ASP
1	A	442	ASP
1	A	614	ASP
1	B	104	ASP
1	B	105	SER

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	547/585 (94%)	514 (94%)	33 (6%)	19	16
1	B	547/585 (94%)	519 (95%)	28 (5%)	24	22
1	C	539/585 (92%)	503 (93%)	36 (7%)	16	13
1	D	555/585 (95%)	535 (96%)	20 (4%)	35	36
1	E	560/585 (96%)	525 (94%)	35 (6%)	18	15
1	F	553/585 (94%)	521 (94%)	32 (6%)	20	17
All	All	3301/3510 (94%)	3117 (94%)	184 (6%)	21	18

5 of 184 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	475	ARG
1	D	141	GLU
1	F	489	LYS
1	C	488	SER
1	C	646	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	469	ASN
1	A	603	GLN
1	F	222	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 82 ligands modelled in this entry, 12 are monoatomic - leaving 70 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	$\# Z > 2$
4	SO4	B	713	-	4,4,4	0.15	0	6,6,6	0.11	0
2	BCT	D	702	3	0,3,3	0.00	-	0,3,3	0.00	-
4	SO4	E	706	-	4,4,4	0.78	0	6,6,6	1.14	1 (16%)
4	SO4	D	705	-	4,4,4	0.13	0	6,6,6	0.08	0
2	BCT	F	701	3	0,3,3	0.00	-	0,3,3	0.00	-
4	SO4	D	712	-	4,4,4	0.12	0	6,6,6	0.12	0
2	BCT	C	701	3	0,3,3	0.00	-	0,3,3	0.00	-
4	SO4	D	706	-	4,4,4	0.15	0	6,6,6	0.11	0
4	SO4	C	708	-	4,4,4	0.13	0	6,6,6	0.15	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	BCT	C	702	3	0,3,3	0.00	-	0,3,3	0.00	-
4	SO4	C	707	-	4,4,4	0.15	0	6,6,6	0.22	0
4	SO4	B	709	-	4,4,4	0.14	0	6,6,6	0.15	0
4	SO4	E	711	-	4,4,4	0.13	0	6,6,6	0.16	0
5	P6G	B	714	-	18,18,18	0.73	0	17,17,17	1.33	2 (11%)
5	P6G	C	713	-	18,18,18	0.67	0	17,17,17	1.57	1 (5%)
4	SO4	C	706	-	4,4,4	0.14	0	6,6,6	0.10	0
2	BCT	A	702	3	0,3,3	0.00	-	0,3,3	0.00	-
2	BCT	E	701	3	0,3,3	0.00	-	0,3,3	0.00	-
4	SO4	B	708	-	4,4,4	0.14	0	6,6,6	0.22	0
4	SO4	E	712	-	4,4,4	0.14	0	6,6,6	0.17	0
2	BCT	A	701	3	0,3,3	0.00	-	0,3,3	0.00	-
4	SO4	C	705	-	4,4,4	0.14	0	6,6,6	0.10	0
5	P6G	B	715	-	18,18,18	0.71	0	17,17,17	1.52	1 (5%)
5	P6G	A	710	-	18,18,18	0.70	0	17,17,17	1.52	0
4	SO4	F	705	-	4,4,4	0.14	0	6,6,6	0.11	0
5	P6G	E	715	-	18,18,18	0.73	0	17,17,17	1.43	0
4	SO4	E	713	-	4,4,4	0.14	0	6,6,6	0.08	0
5	P6G	A	711	-	18,18,18	0.65	0	17,17,17	1.57	2 (11%)
4	SO4	D	710	-	4,4,4	0.12	0	6,6,6	0.19	0
4	SO4	E	707	-	4,4,4	0.14	0	6,6,6	0.16	0
4	SO4	E	710	-	4,4,4	0.14	0	6,6,6	0.11	0
4	SO4	B	707	-	4,4,4	0.16	0	6,6,6	0.12	0
4	SO4	A	706	-	4,4,4	0.16	0	6,6,6	0.22	0
4	SO4	C	709	-	4,4,4	0.16	0	6,6,6	0.12	0
4	SO4	C	711	-	4,4,4	0.13	0	6,6,6	0.06	0
5	P6G	F	710	-	18,18,18	0.71	0	17,17,17	1.55	1 (5%)
2	BCT	B	702	3	0,3,3	0.00	-	0,3,3	0.00	-
2	BCT	D	701	3	0,3,3	0.00	-	0,3,3	0.00	-
5	P6G	D	715	-	18,18,18	0.61	0	17,17,17	1.79	3 (17%)
4	SO4	A	705	-	4,4,4	0.17	0	6,6,6	0.16	0
4	SO4	F	708	-	4,4,4	0.15	0	6,6,6	0.14	0
4	SO4	E	708	-	4,4,4	0.15	0	6,6,6	0.08	0
4	SO4	F	706	-	4,4,4	0.18	0	6,6,6	0.15	0
5	P6G	F	711	-	18,18,18	0.70	0	17,17,17	1.67	2 (11%)
4	SO4	B	706	-	4,4,4	0.18	0	6,6,6	0.13	0
4	SO4	B	712	-	4,4,4	0.15	0	6,6,6	0.06	0
4	SO4	D	711	-	4,4,4	0.15	0	6,6,6	0.12	0
5	P6G	C	714	-	18,18,18	0.66	0	17,17,17	1.57	0
4	SO4	C	710	-	4,4,4	0.15	0	6,6,6	0.08	0
4	SO4	F	707	-	4,4,4	0.14	0	6,6,6	0.15	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	P6G	E	714	-	18,18,18	0.68	0	17,17,17	1.59	2 (11%)
4	SO4	A	707	-	4,4,4	0.11	0	6,6,6	0.15	0
4	SO4	A	709	-	4,4,4	0.15	0	6,6,6	0.08	0
2	BCT	F	702	3	0,3,3	0.00	-	0,3,3	0.00	-
5	P6G	A	712	-	18,18,18	0.70	0	17,17,17	1.50	0
4	SO4	D	713	-	4,4,4	0.16	0	6,6,6	0.08	0
4	SO4	D	707	-	4,4,4	0.12	0	6,6,6	0.15	0
4	SO4	E	705	-	4,4,4	0.15	0	6,6,6	0.07	0
5	P6G	D	714	-	18,18,18	0.69	0	17,17,17	1.59	0
4	SO4	A	708	-	4,4,4	0.18	0	6,6,6	0.12	0
4	SO4	D	708	-	4,4,4	0.12	0	6,6,6	0.14	0
4	SO4	B	705	-	4,4,4	0.15	0	6,6,6	0.32	0
4	SO4	B	711	-	4,4,4	0.18	0	6,6,6	0.13	0
4	SO4	F	709	-	4,4,4	0.15	0	6,6,6	0.13	0
2	BCT	E	702	3	0,3,3	0.00	-	0,3,3	0.00	-
4	SO4	D	709	-	4,4,4	0.18	0	6,6,6	0.14	0
4	SO4	B	710	-	4,4,4	0.15	0	6,6,6	0.19	0
5	P6G	C	712	-	18,18,18	0.72	0	17,17,17	1.46	0
2	BCT	B	701	3	0,3,3	0.00	-	0,3,3	0.00	-
4	SO4	E	709	-	4,4,4	0.14	0	6,6,6	0.16	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
5	P6G	A	712	-	-	11/16/16/16	-
5	P6G	B	714	-	-	4/16/16/16	-
5	P6G	C	713	-	-	7/16/16/16	-
5	P6G	D	715	-	-	9/16/16/16	-
5	P6G	D	714	-	-	13/16/16/16	-
5	P6G	F	710	-	-	8/16/16/16	-
5	P6G	F	711	-	-	13/16/16/16	-
5	P6G	E	714	-	-	9/16/16/16	-
5	P6G	E	715	-	-	12/16/16/16	-
5	P6G	C	714	-	-	11/16/16/16	-
5	P6G	A	711	-	-	10/16/16/16	-
5	P6G	A	710	-	-	10/16/16/16	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	P6G	B	715	-	-	11/16/16/16	-
5	P6G	C	712	-	-	11/16/16/16	-

There are no bond length outliers.

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	E	706	SO4	O4-S-O3	-2.55	98.17	109.06
5	E	714	P6G	C14-O13-C12	2.48	124.04	113.29
5	F	711	P6G	O4-C5-C6	2.35	121.00	110.39
5	F	711	P6G	C8-O7-C6	2.24	123.00	113.29
5	D	715	P6G	O7-C6-C5	2.23	120.46	110.39

There are no chirality outliers.

5 of 139 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	E	715	P6G	C12-C11-O10-C9
5	D	714	P6G	C12-C11-O10-C9
5	C	712	P6G	C14-C15-O16-C17
5	E	714	P6G	C2-C3-O4-C5
5	F	710	P6G	C6-C5-O4-C3

There are no ring outliers.

14 monomers are involved in 37 short contacts:

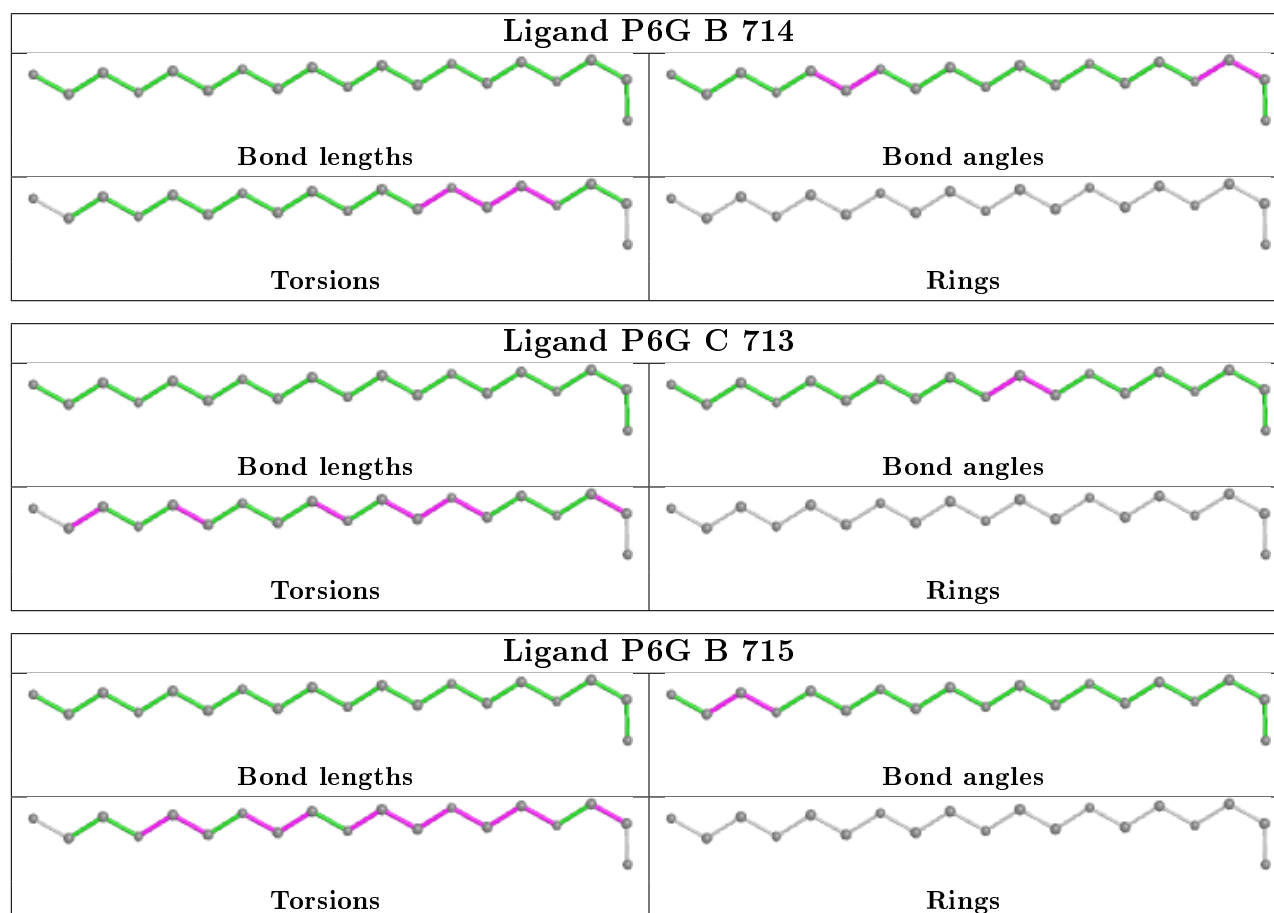
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	702	BCT	3	0
5	C	713	P6G	3	0
5	B	715	P6G	6	0
5	A	710	P6G	1	0
5	E	715	P6G	3	0
4	D	710	SO4	1	0
4	B	707	SO4	1	0
5	D	715	P6G	3	0
4	A	705	SO4	1	0
5	F	711	P6G	6	0
5	E	714	P6G	1	0
5	A	712	P6G	3	0
5	D	714	P6G	4	0

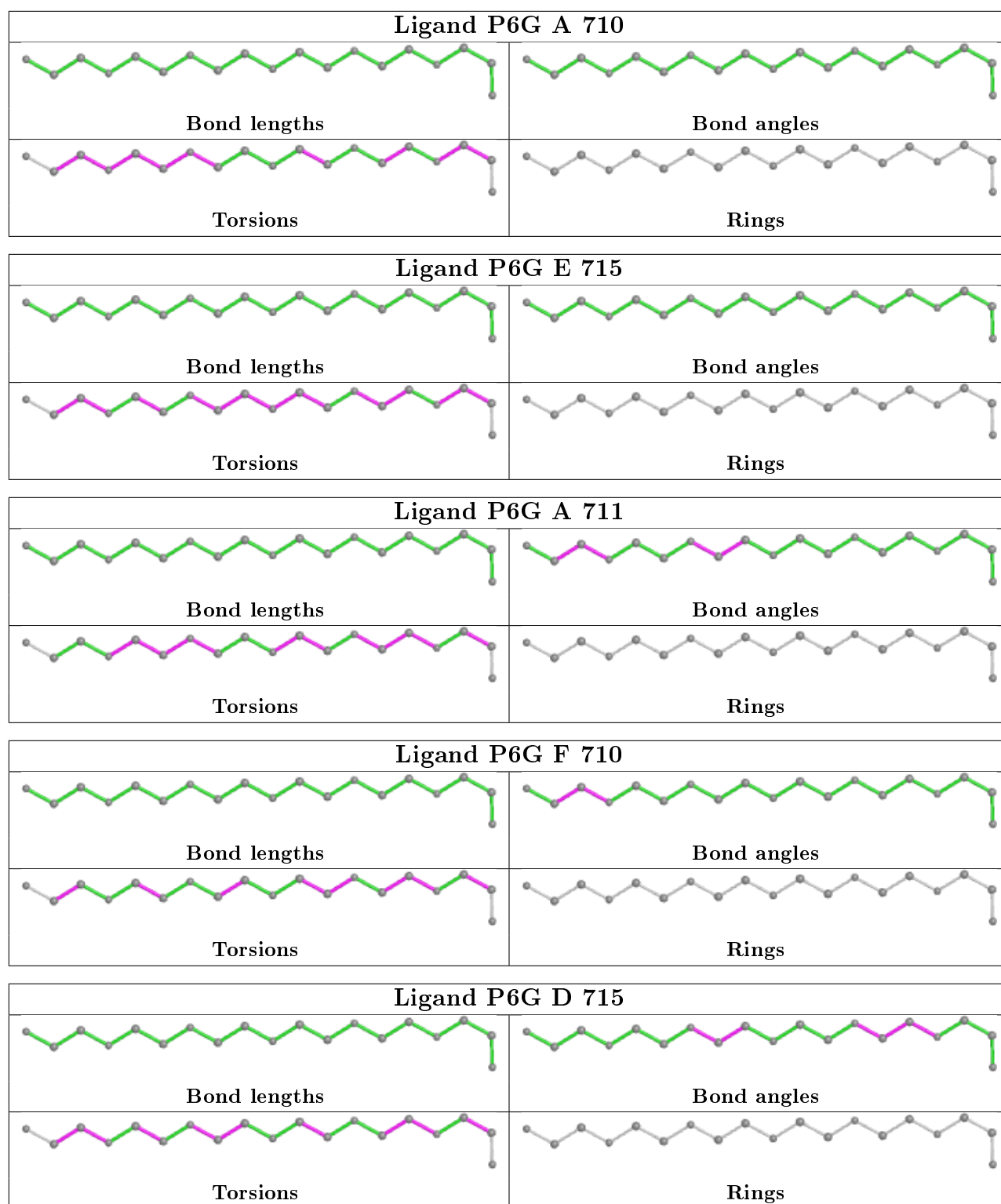
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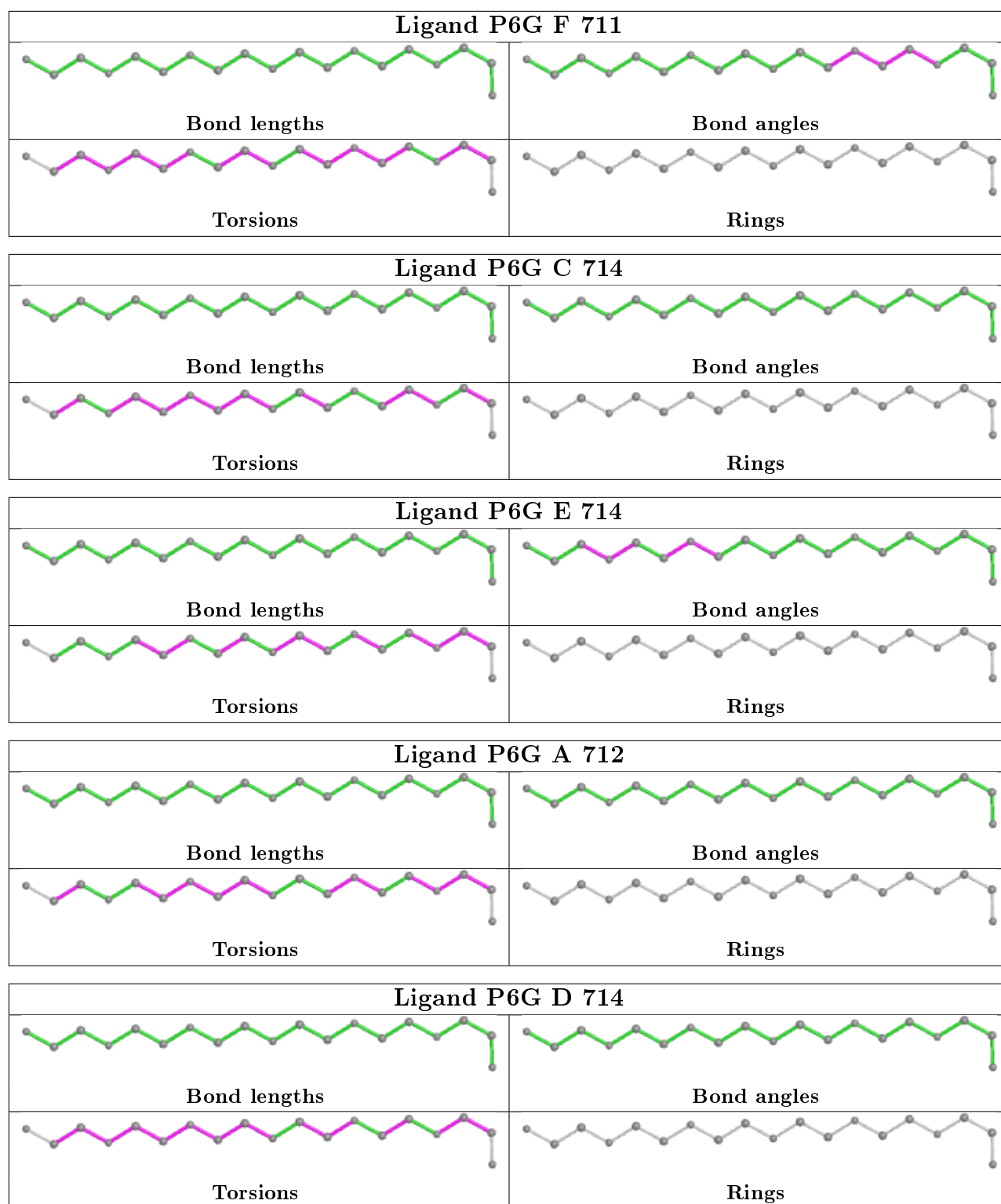
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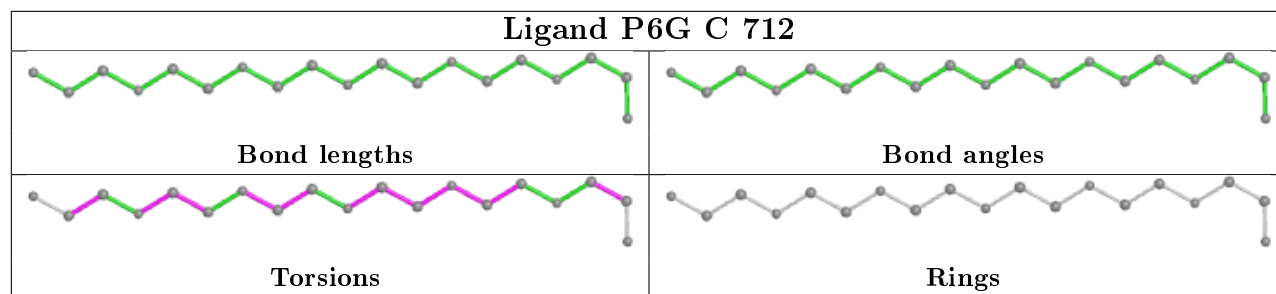
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	C	712	P6G	1	0

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.









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	674/698 (96%)	0.03	38 (5%) 24 29	15, 39, 77, 97	0
1	B	677/698 (96%)	-0.09	32 (4%) 31 37	13, 33, 72, 97	0
1	C	674/698 (96%)	0.81	121 (17%) 1 1	13, 46, 117, 151	0
1	D	675/698 (96%)	-0.04	23 (3%) 45 51	16, 34, 66, 94	0
1	E	674/698 (96%)	0.02	35 (5%) 27 32	13, 35, 73, 93	0
1	F	676/698 (96%)	-0.04	33 (4%) 29 35	13, 36, 70, 98	0
All	All	4050/4188 (96%)	0.12	282 (6%) 16 20	13, 36, 83, 151	0

The worst 5 of 282 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	479	PHE	14.5
1	C	474	CYS	13.6
1	C	478	GLU	13.0
1	C	564	LEU	11.7
1	C	445	LYS	11.2

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](#)

There are no carbohydrates in this entry.

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
5	P6G	D	715	19/19	0.79	0.20	56,74,84,86	0
4	SO4	B	708	5/5	0.81	0.21	86,88,92,96	0
4	SO4	F	706	5/5	0.83	0.16	89,92,95,98	0
5	P6G	A	712	19/19	0.83	0.20	74,81,91,92	0
4	SO4	A	705	5/5	0.84	0.19	93,94,95,97	0
4	SO4	E	706	5/5	0.84	0.42	111,112,113,113	5
4	SO4	C	711	5/5	0.84	0.20	106,109,109,110	0
5	P6G	C	714	19/19	0.85	0.20	56,79,99,100	0
5	P6G	A	711	19/19	0.85	0.16	46,66,84,85	0
4	SO4	F	705	5/5	0.87	0.14	97,97,98,100	0
4	SO4	D	712	5/5	0.88	0.16	95,95,99,100	0
4	SO4	B	713	5/5	0.88	0.20	105,107,108,109	0
4	SO4	E	713	5/5	0.89	0.16	93,95,95,96	0
4	SO4	A	709	5/5	0.89	0.13	97,97,99,99	0
5	P6G	D	714	19/19	0.89	0.18	41,54,79,79	0
4	SO4	A	708	5/5	0.89	0.18	95,98,99,100	0
5	P6G	F	711	19/19	0.90	0.19	28,48,98,100	0
4	SO4	A	707	5/5	0.90	0.18	77,87,88,92	0
4	SO4	D	706	5/5	0.90	0.14	101,103,106,109	0
5	P6G	C	713	19/19	0.90	0.16	43,57,83,83	0
5	P6G	E	714	19/19	0.90	0.21	28,45,83,84	0
4	SO4	C	706	5/5	0.91	0.12	94,95,97,98	0
4	SO4	E	705	5/5	0.91	0.16	107,108,109,110	0
4	SO4	B	707	5/5	0.91	0.12	85,87,90,92	0
4	SO4	F	709	5/5	0.91	0.13	89,89,92,93	0
5	P6G	F	710	19/19	0.92	0.16	37,48,68,71	0
4	SO4	C	709	5/5	0.92	0.13	76,77,80,80	0
5	P6G	A	710	19/19	0.93	0.17	37,48,77,77	0
5	P6G	B	714	19/19	0.93	0.14	22,32,87,87	0
5	P6G	E	715	19/19	0.93	0.17	44,56,71,71	0
4	SO4	E	711	5/5	0.94	0.17	73,77,84,87	0
4	SO4	F	707	5/5	0.94	0.15	85,87,87,90	0
4	SO4	E	708	5/5	0.94	0.12	94,95,96,96	0
4	SO4	B	706	5/5	0.94	0.15	72,79,81,82	0
5	P6G	C	712	19/19	0.94	0.14	28,48,79,81	0
4	SO4	B	712	5/5	0.95	0.11	77,80,83,84	0
4	SO4	A	706	5/5	0.95	0.15	51,66,69,77	0

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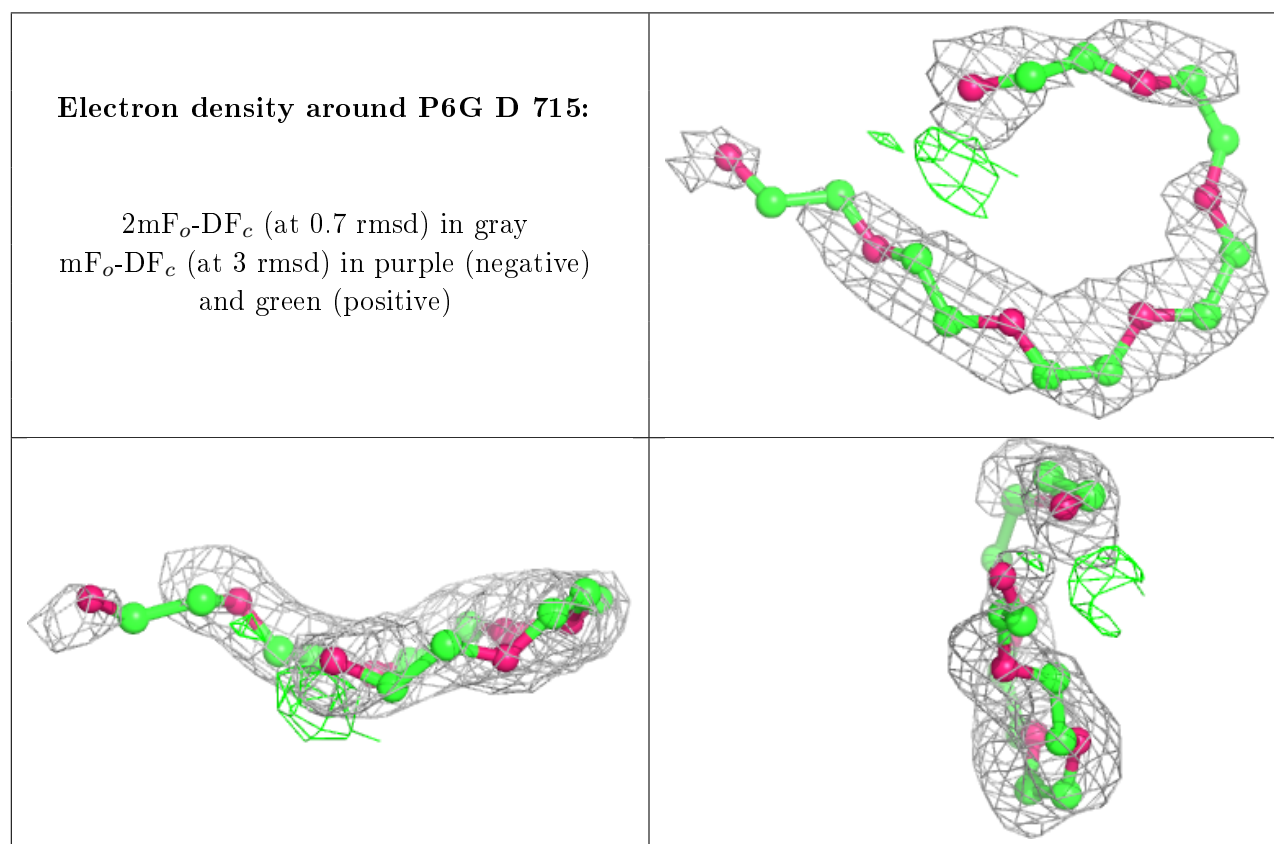
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	BCT	C	702	4/4	0.95	0.13	32,36,38,44	0
4	SO4	E	709	5/5	0.95	0.11	76,77,79,81	0
2	BCT	E	702	4/4	0.95	0.13	26,28,29,30	0
4	SO4	D	709	5/5	0.95	0.11	70,72,75,76	0
4	SO4	F	708	5/5	0.95	0.14	80,85,87,90	0
5	P6G	B	715	19/19	0.95	0.15	32,46,94,95	0
4	SO4	C	710	5/5	0.96	0.15	73,77,79,83	0
3	FE	C	703	1/1	0.96	0.12	38,38,38,38	0
4	SO4	B	711	5/5	0.96	0.19	63,68,74,77	0
4	SO4	D	713	5/5	0.96	0.12	72,73,74,78	0
4	SO4	B	710	5/5	0.97	0.11	44,61,65,72	0
4	SO4	C	707	5/5	0.97	0.10	38,55,63,63	0
4	SO4	B	709	5/5	0.97	0.10	66,67,68,73	0
4	SO4	D	705	5/5	0.97	0.19	68,71,72,78	0
4	SO4	E	710	5/5	0.97	0.12	73,73,75,79	0
4	SO4	D	711	5/5	0.97	0.21	62,67,73,74	0
4	SO4	C	705	5/5	0.97	0.21	77,78,83,83	0
4	SO4	C	708	5/5	0.97	0.10	70,75,76,79	0
4	SO4	D	710	5/5	0.98	0.16	55,55,65,66	0
2	BCT	A	702	4/4	0.98	0.10	28,31,34,38	0
4	SO4	D	708	5/5	0.98	0.12	46,49,58,58	0
4	SO4	B	705	5/5	0.98	0.10	44,45,49,50	0
4	SO4	E	707	5/5	0.98	0.09	41,55,61,63	0
3	FE	E	704	1/1	0.98	0.14	32,32,32,32	0
2	BCT	B	702	4/4	0.98	0.10	27,27,28,31	0
2	BCT	D	701	4/4	0.98	0.11	15,17,21,24	0
4	SO4	E	712	5/5	0.98	0.11	44,48,49,50	0
2	BCT	D	702	4/4	0.98	0.12	23,26,29,32	0
4	SO4	D	707	5/5	0.98	0.09	57,58,59,65	0
3	FE	F	704	1/1	0.99	0.14	31,31,31,31	0
2	BCT	F	702	4/4	0.99	0.13	19,20,21,24	0
2	BCT	A	701	4/4	0.99	0.10	14,14,16,19	0
2	BCT	F	701	4/4	0.99	0.09	8,9,10,11	0
3	FE	D	703	1/1	0.99	0.15	26,26,26,26	0
2	BCT	C	701	4/4	0.99	0.09	14,15,15,17	0
2	BCT	E	701	4/4	0.99	0.10	14,17,17,18	0
3	FE	A	704	1/1	0.99	0.12	34,34,34,34	0
2	BCT	B	701	4/4	0.99	0.11	13,14,15,16	0
3	FE	E	703	1/1	1.00	0.14	19,19,19,19	0
3	FE	C	704	1/1	1.00	0.14	16,16,16,16	0
3	FE	D	704	1/1	1.00	0.15	20,20,20,20	0
3	FE	B	703	1/1	1.00	0.12	29,29,29,29	0

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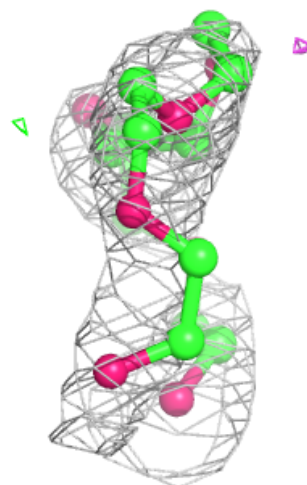
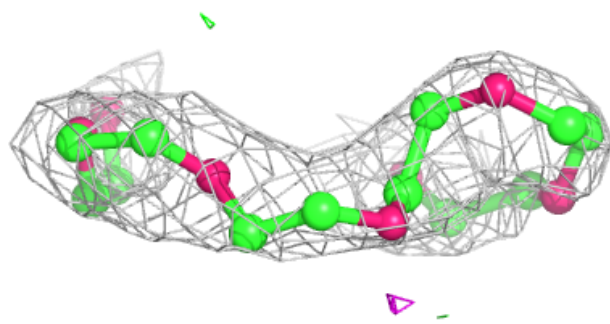
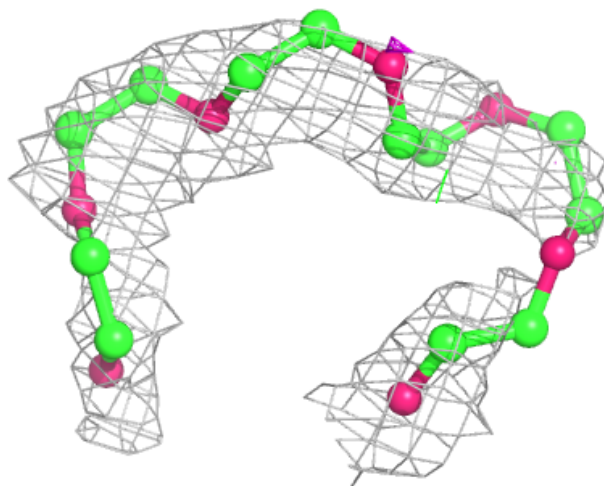
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	FE	F	703	1/1	1.00	0.14	16,16,16,16	0
3	FE	A	703	1/1	1.00	0.15	19,19,19,19	0
3	FE	B	704	1/1	1.00	0.14	15,15,15,15	0

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.



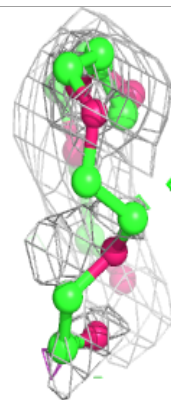
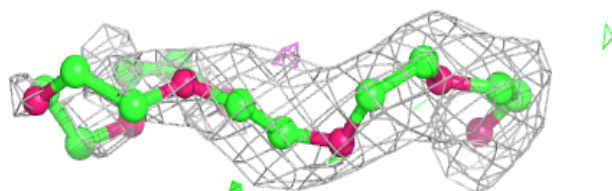
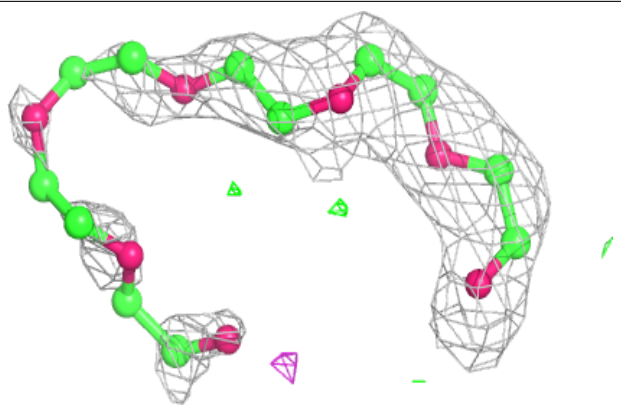
Electron density around P6G A 712:

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



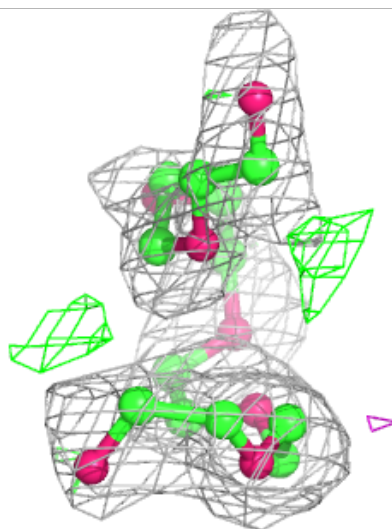
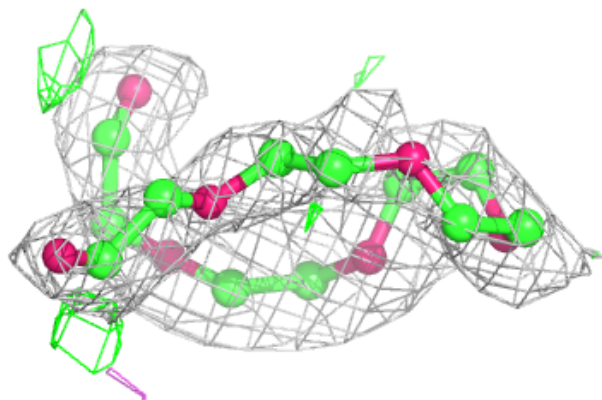
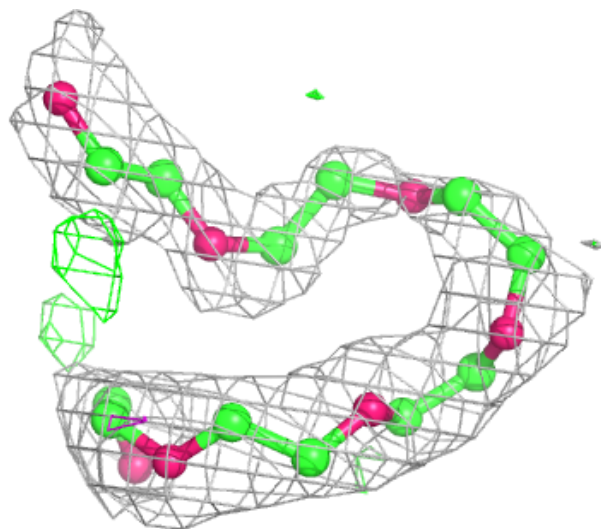
Electron density around P6G C 714:

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



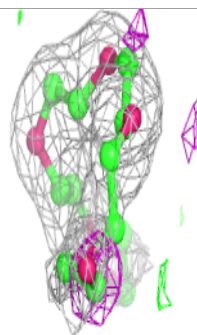
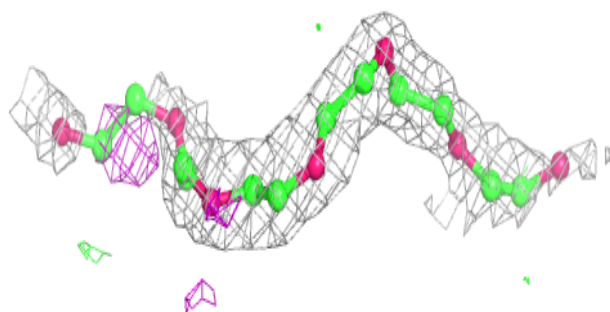
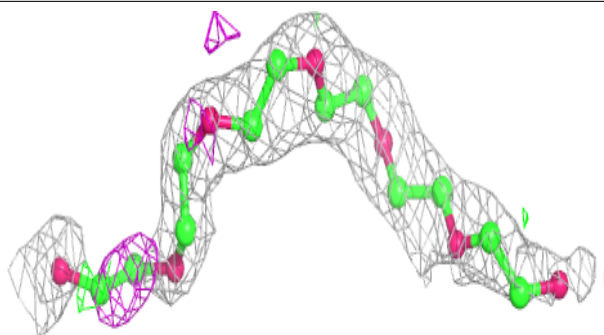
Electron density around P6G A 711:

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

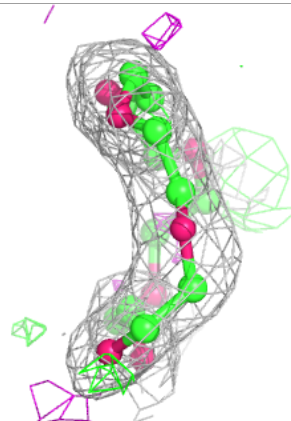
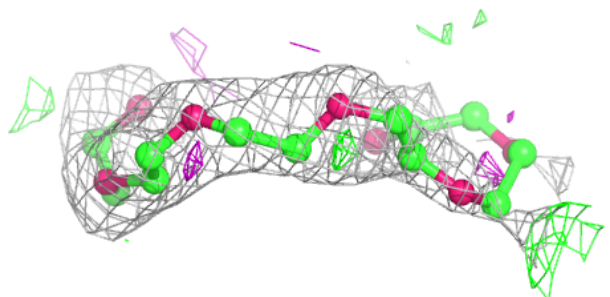
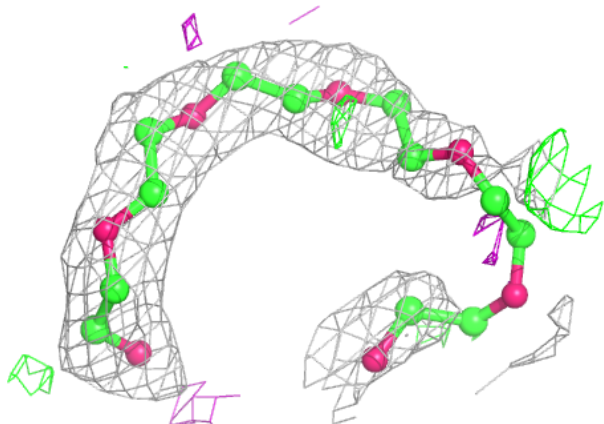


Electron density around P6G D 714:

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

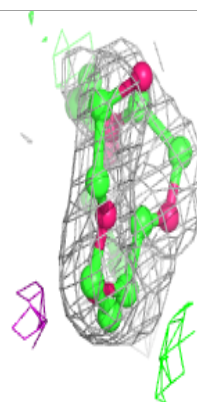
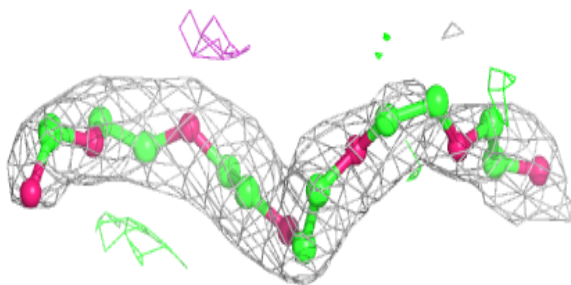
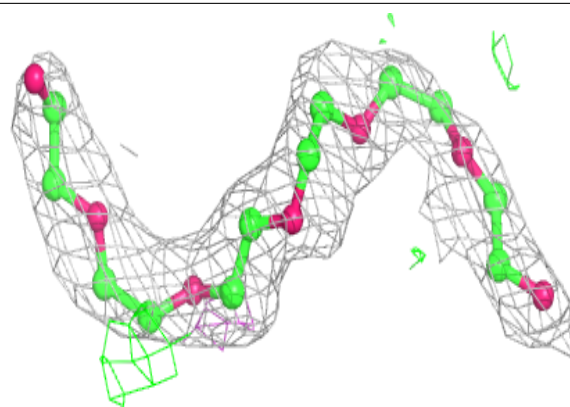
**Electron density around P6G F 711:**

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



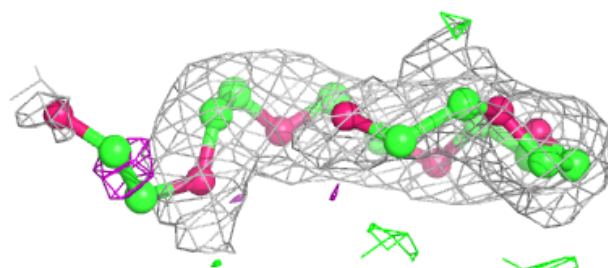
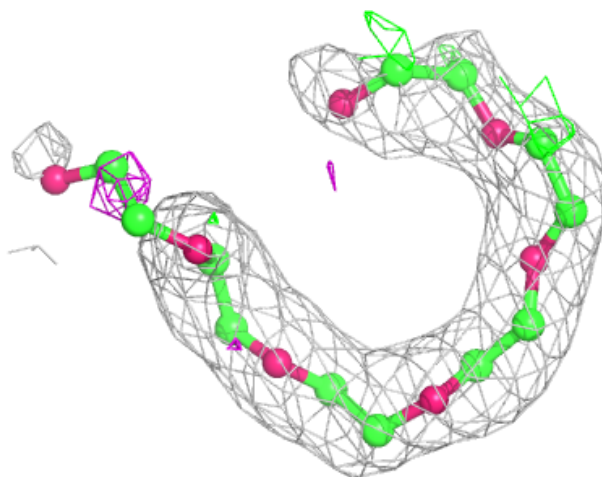
Electron density around P6G C 713:

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



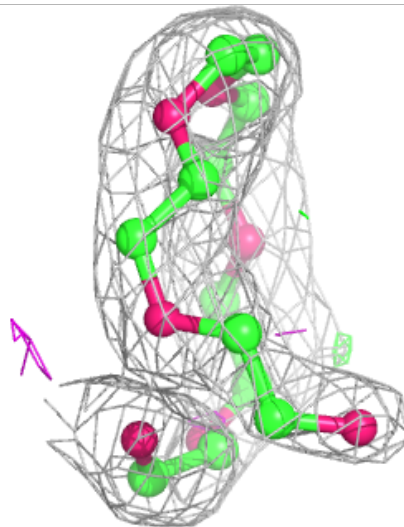
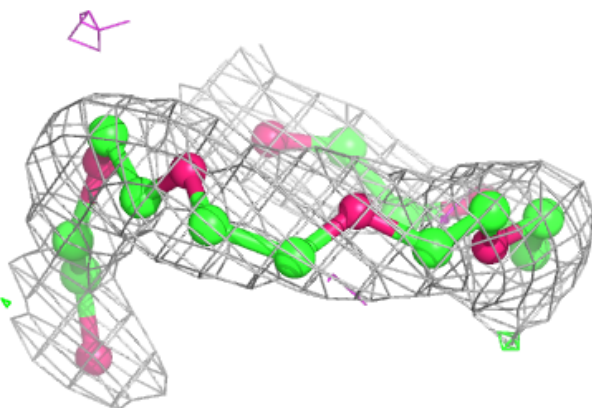
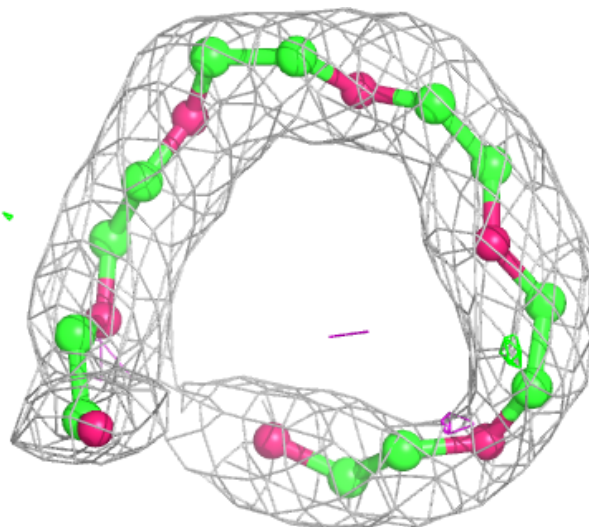
Electron density around P6G E 714:

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



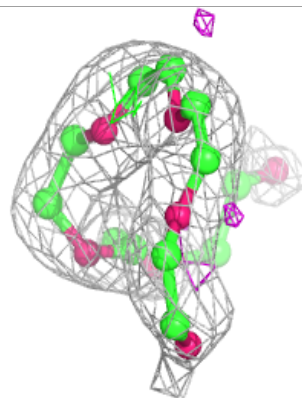
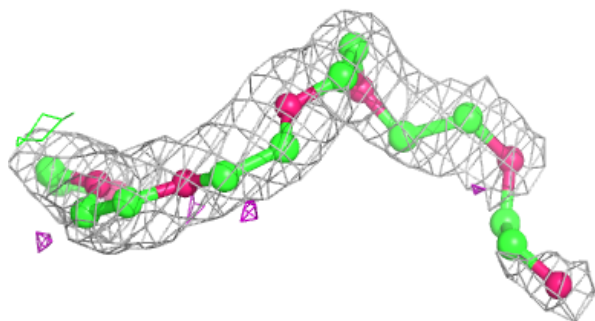
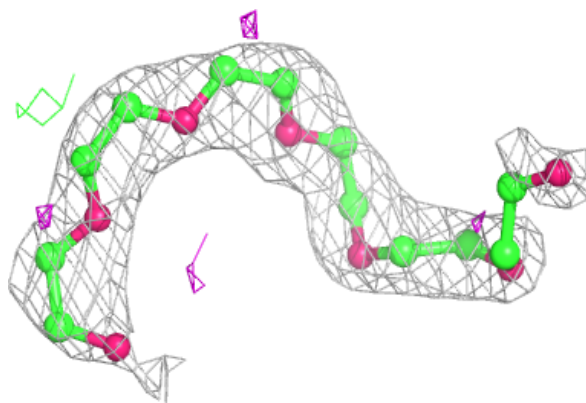
Electron density around P6G F 710:

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



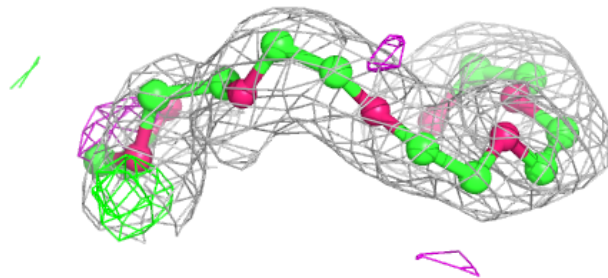
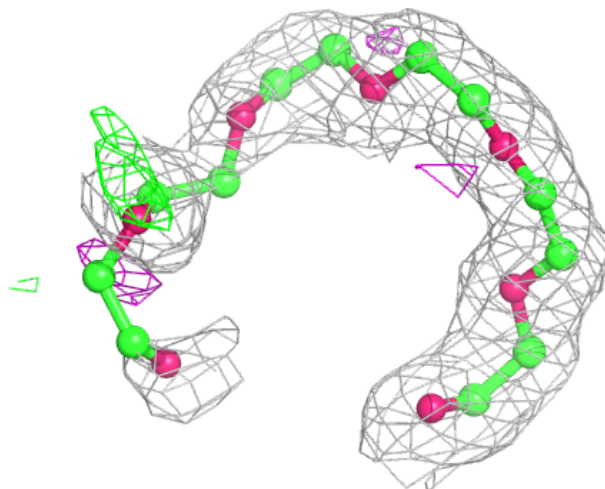
Electron density around P6G A 710:

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



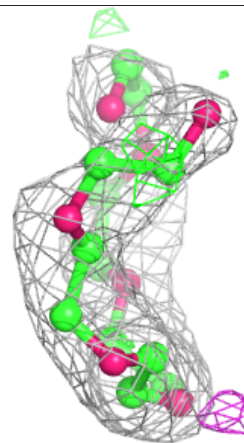
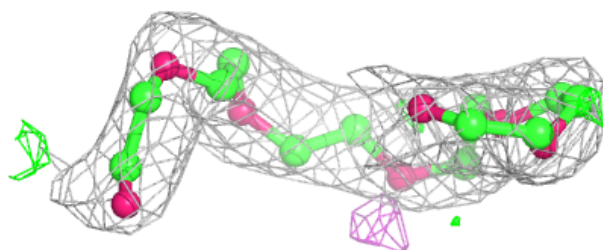
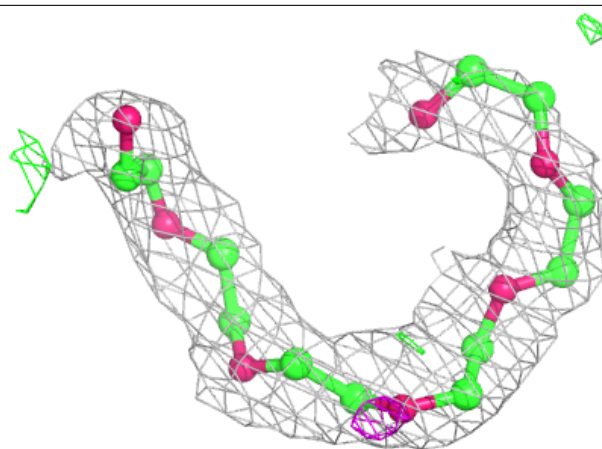
Electron density around P6G B 714:

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



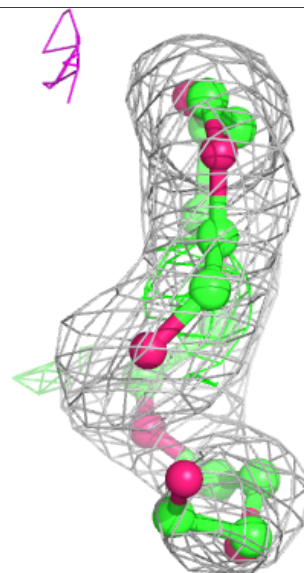
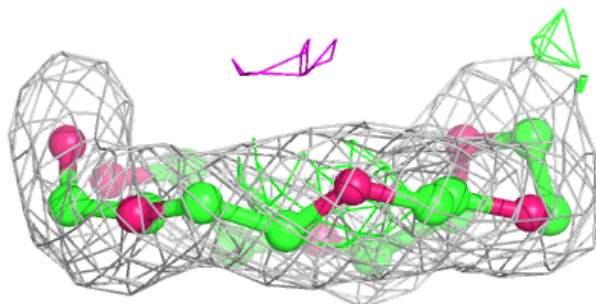
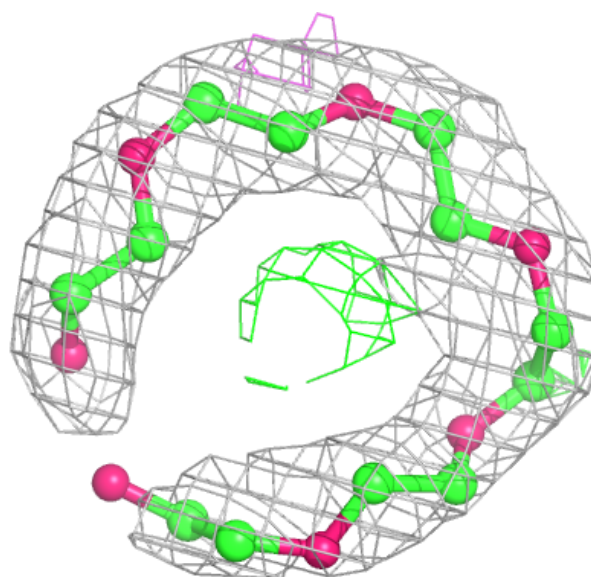
Electron density around P6G E 715:

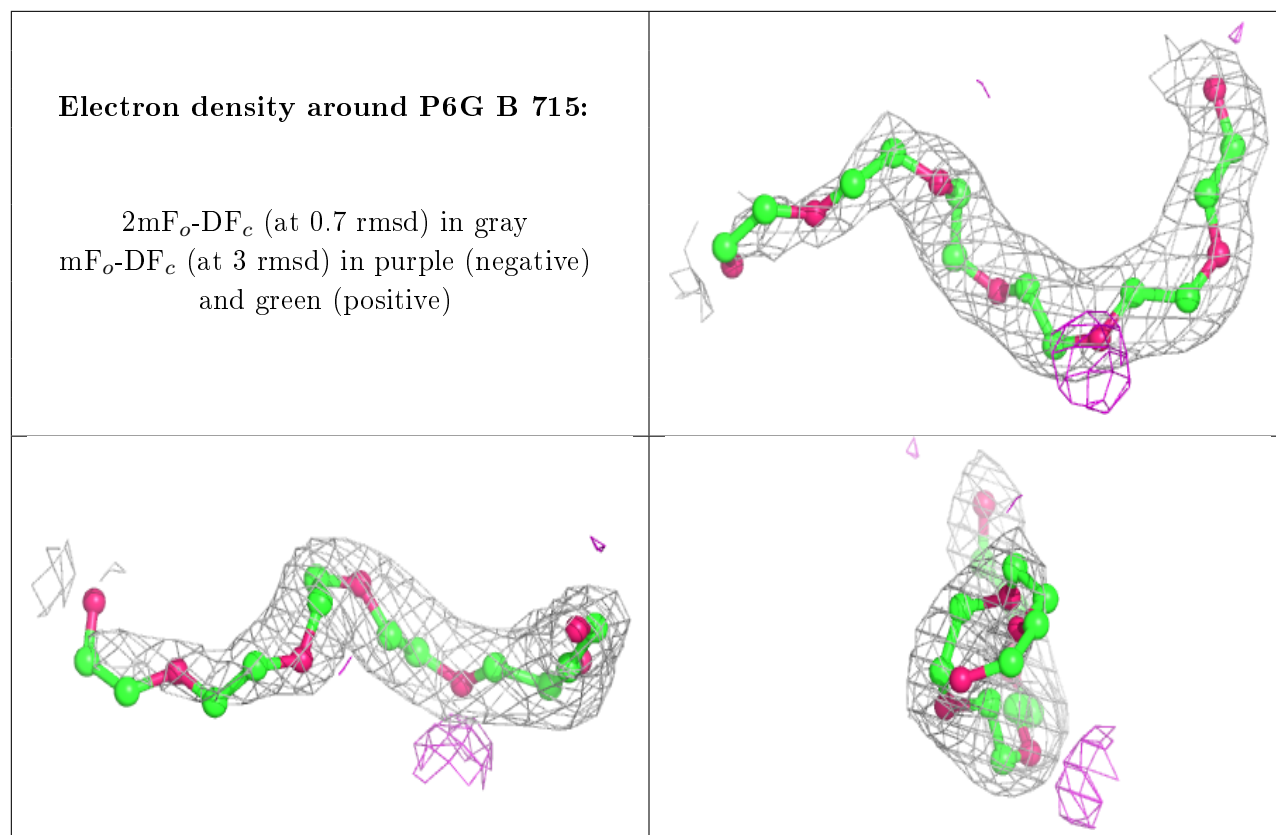
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around P6G C 712:

$2mF_o-DF_c$ (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.