



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 5, 2026 – 11:18 AM UTC

PDB ID : 6PD2 / pdb_00006pd2
Title : PntC-AEPT: fusion protein of phosphonate-specific cytidyltransferase and 2-aminoethylphosphonate (AEP) transaminase from *Treponema denticola* in complex with cytidine monophosphate-AEP
Authors : Suits, M.D.L.; Whiteside, J.
Deposited on : 2019-06-18
Resolution : 1.95 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 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)
Xtrriage (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

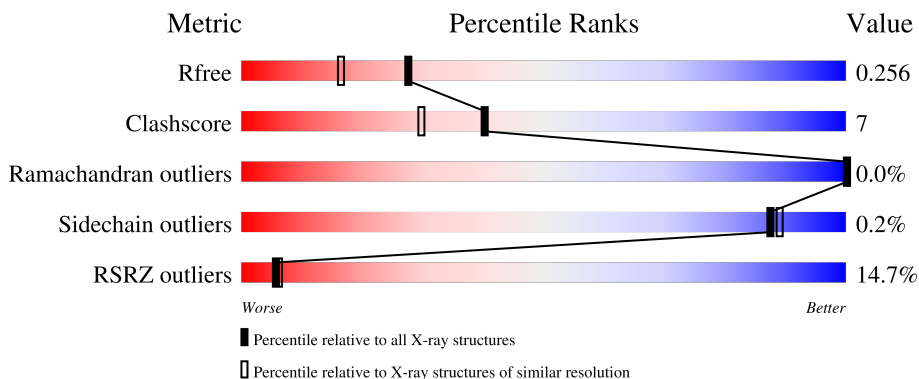
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

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	624	
1	B	624	
1	C	624	
1	D	624	

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
4	PO4	D	709	-	-	X	-
6	EDO	A	709	-	-	X	-
6	EDO	A	716	-	-	X	-
6	EDO	B	706	-	-	X	-

2 Entry composition [i](#)

There are 9 unique types of molecules in this entry. The entry contains 21090 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 Nucleotidyl transferase/aminotransferase, class V.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	615	4831	3084	798	914	35	0	0	0
1	B	615	4831	3084	798	914	35	0	0	0
1	C	616	4838	3089	799	915	35	0	0	0
1	D	608	4776	3051	786	904	35	0	0	0

There are 32 discrepancies between the modelled and reference sequences:

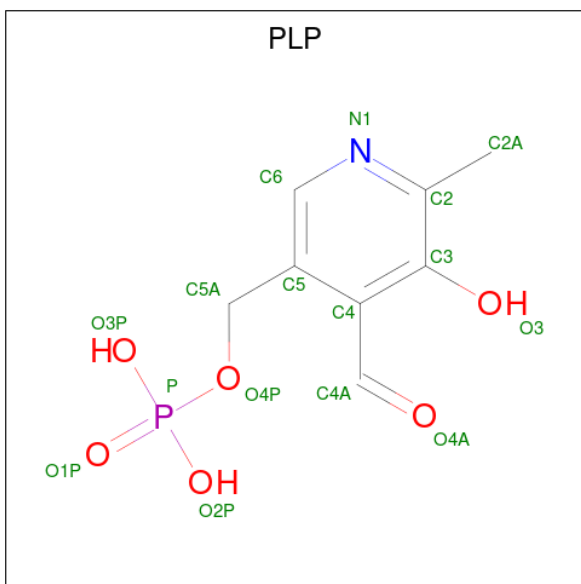
Chain	Residue	Modelled	Actual	Comment	Reference
A	617	LEU	-	expression tag	UNP Q73MU2
A	618	GLU	-	expression tag	UNP Q73MU2
A	619	HIS	-	expression tag	UNP Q73MU2
A	620	HIS	-	expression tag	UNP Q73MU2
A	621	HIS	-	expression tag	UNP Q73MU2
A	622	HIS	-	expression tag	UNP Q73MU2
A	623	HIS	-	expression tag	UNP Q73MU2
A	624	HIS	-	expression tag	UNP Q73MU2
B	617	LEU	-	expression tag	UNP Q73MU2
B	618	GLU	-	expression tag	UNP Q73MU2
B	619	HIS	-	expression tag	UNP Q73MU2
B	620	HIS	-	expression tag	UNP Q73MU2
B	621	HIS	-	expression tag	UNP Q73MU2
B	622	HIS	-	expression tag	UNP Q73MU2
B	623	HIS	-	expression tag	UNP Q73MU2
B	624	HIS	-	expression tag	UNP Q73MU2
C	617	LEU	-	expression tag	UNP Q73MU2
C	618	GLU	-	expression tag	UNP Q73MU2
C	619	HIS	-	expression tag	UNP Q73MU2
C	620	HIS	-	expression tag	UNP Q73MU2
C	621	HIS	-	expression tag	UNP Q73MU2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	622	HIS	-	expression tag	UNP Q73MU2
C	623	HIS	-	expression tag	UNP Q73MU2
C	624	HIS	-	expression tag	UNP Q73MU2
D	617	LEU	-	expression tag	UNP Q73MU2
D	618	GLU	-	expression tag	UNP Q73MU2
D	619	HIS	-	expression tag	UNP Q73MU2
D	620	HIS	-	expression tag	UNP Q73MU2
D	621	HIS	-	expression tag	UNP Q73MU2
D	622	HIS	-	expression tag	UNP Q73MU2
D	623	HIS	-	expression tag	UNP Q73MU2
D	624	HIS	-	expression tag	UNP Q73MU2

- Molecule 2 is PYRIDOXAL-5'-PHOSPHATE (CCD ID: PLP) (formula: $C_8H_{10}NO_6P$).

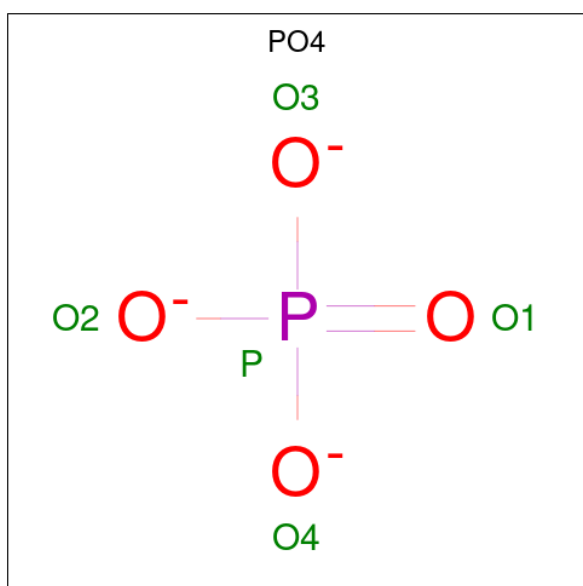


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
			Total	C	N	O	P			
2	A	1	Total	16	8	1	6	1	0	0
2	B	1	Total	16	8	1	6	1	0	0
2	C	1	Total	16	8	1	6	1	0	0
2	D	1	Total	16	8	1	6	1	0	0

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	2	Total Mg 2 2	0	0
3	B	2	Total Mg 2 2	0	0
3	C	2	Total Mg 2 2	0	0
3	D	2	Total Mg 2 2	0	0

- Molecule 4 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



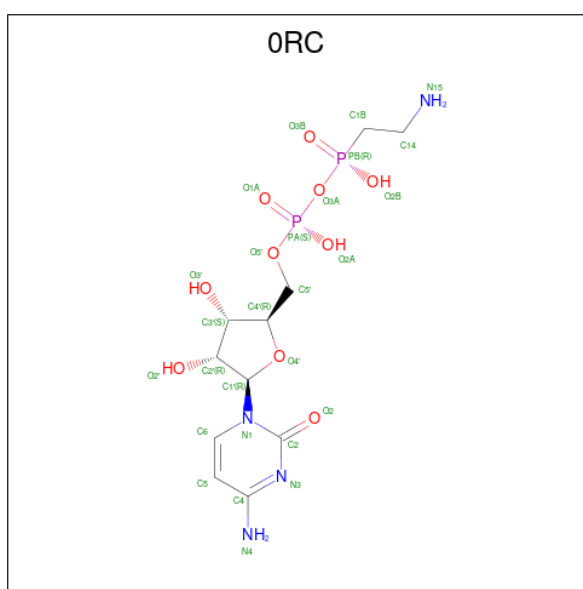
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total O P 5 4 1	0	0
4	A	1	Total O P 5 4 1	0	0
4	B	1	Total O P 5 4 1	0	0
4	B	1	Total O P 5 4 1	0	0
4	B	1	Total O P 5 4 1	0	0
4	C	1	Total O P 5 4 1	0	0
4	C	1	Total O P 5 4 1	0	0
4	C	1	Total O P 5 4 1	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	D	1	Total O P 5 4 1	0	0
4	D	1	Total O P 5 4 1	0	0
4	D	1	Total O P 5 4 1	0	0

- Molecule 5 is 5'-O-[(S)-{[(R)-(2-aminoethyl)(hydroxy)phosphoryl]oxy}(hydroxy)phosphoryl]cytidine (CCD ID: 0RC) (formula: C₁₁H₂₀N₄O₁₀P₂) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C N O P 27 11 4 10 2	0	0
5	B	1	Total C N O P 27 11 4 10 2	0	0
5	C	1	Total C N O P 27 11 4 10 2	0	0
5	D	1	Total C N O P 27 11 4 10 2	0	0

- Molecule 6 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



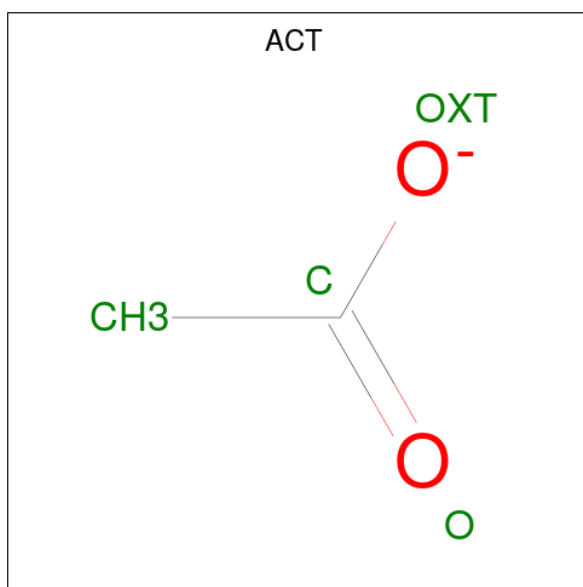
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total C O 4 2 2	0	0
6	A	1	Total C O 4 2 2	0	0
6	A	1	Total C O 4 2 2	0	0
6	A	1	Total C O 4 2 2	0	0
6	A	1	Total C O 4 2 2	0	0
6	A	1	Total C O 4 2 2	0	0
6	A	1	Total C O 4 2 2	0	0
6	A	1	Total C O 4 2 2	0	0
6	A	1	Total C O 4 2 2	0	0
6	A	1	Total C O 4 2 2	0	0
6	B	1	Total C O 4 2 2	0	0
6	B	1	Total C O 4 2 2	0	0
6	B	1	Total C O 4 2 2	0	0
6	C	1	Total C O 4 2 2	0	0

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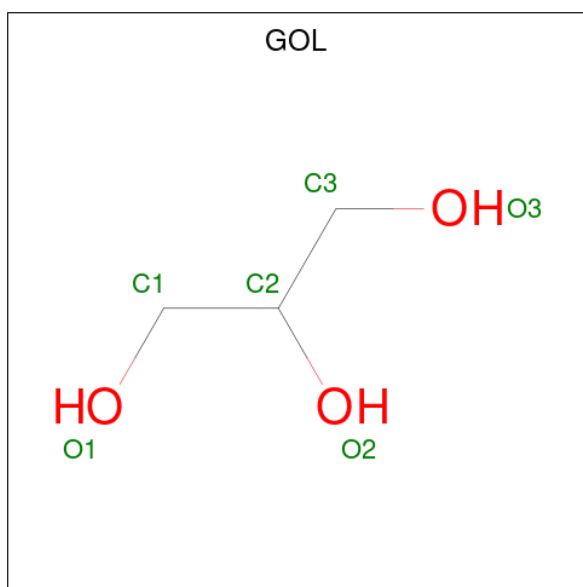
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	C	1	Total C O 4 2 2	0	0
6	C	1	Total C O 4 2 2	0	0
6	C	1	Total C O 4 2 2	0	0
6	C	1	Total C O 4 2 2	0	0
6	C	1	Total C O 4 2 2	0	0
6	C	1	Total C O 4 2 2	0	0
6	D	1	Total C O 4 2 2	0	0
6	D	1	Total C O 4 2 2	0	0
6	D	1	Total C O 4 2 2	0	0
6	D	1	Total C O 4 2 2	0	0
6	D	1	Total C O 4 2 2	0	0
6	D	1	Total C O 4 2 2	0	0
6	D	1	Total C O 4 2 2	0	0

- Molecule 7 is ACETATE ION (CCD ID: ACT) (formula: C₂H₃O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			4	2	2		
7	A	1	Total	C	O	0	0
			4	2	2		
7	C	1	Total	C	O	0	0
			4	2	2		

- Molecule 8 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

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

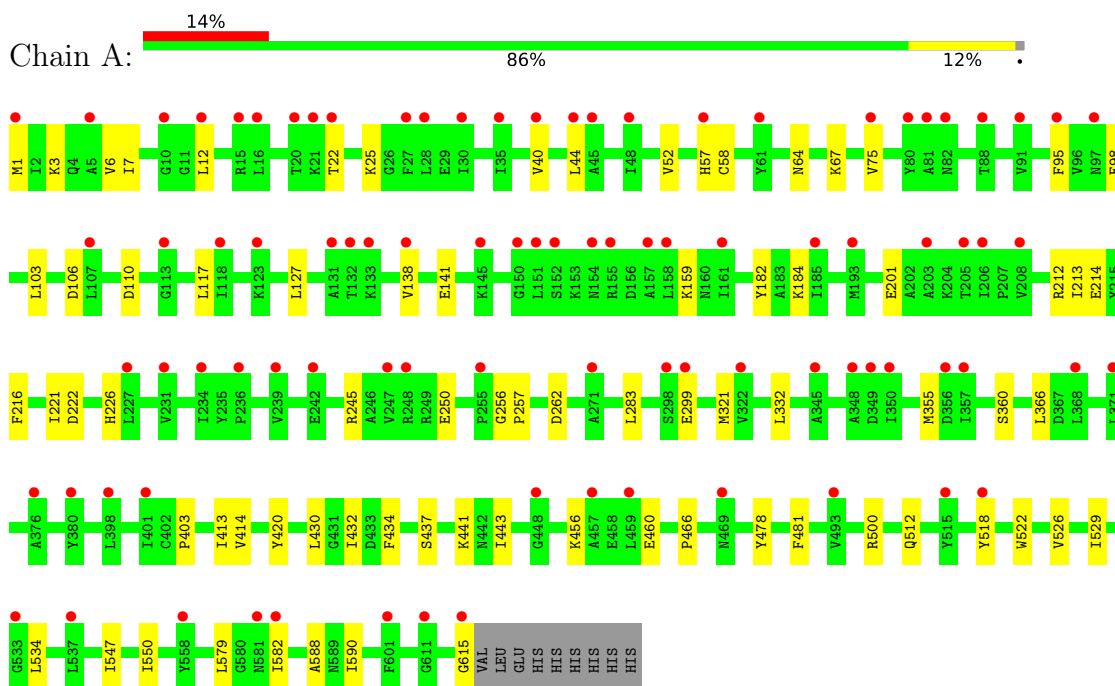
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	354	Total	O	0	0
			354	354		
9	B	352	Total	O	0	0
			352	352		
9	C	411	Total	O	0	0
			411	411		
9	D	330	Total	O	0	0
			330	330		

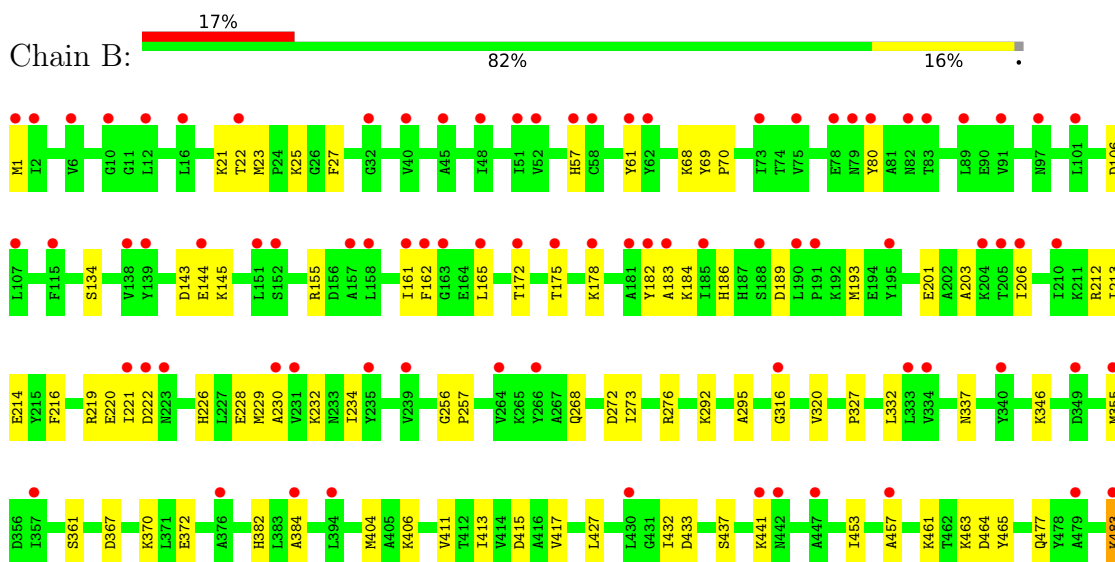
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: Nucleotidyl transferase/aminotransferase, class V



- Molecule 1: Nucleotidyl transferase/aminotransferase, class V



4 Data and refinement statistics i

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	76.45Å 154.05Å 134.58Å 90.00° 90.90° 90.00°	Depositor
Resolution (Å)	47.98 – 1.95 47.98 – 1.95	Depositor EDS
% Data completeness (in resolution range)	98.0 (47.98-1.95) 98.1 (47.98-1.95)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.01 (at 1.95Å)	Xtrriage
Refinement program	PHENIX (1.13_2998: ???)	Depositor
R, R_{free}	0.220 , 0.254 0.222 , 0.256	Depositor DCC
R_{free} test set	10992 reflections (4.87%)	wwPDB-VP
Wilson B-factor (Å ²)	23.1	Xtrriage
Anisotropy	0.768	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 35.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtrriage
Estimated twinning fraction	0.039 for h,-k,-l	Xtrriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	21090	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.21% 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: 0RC, MG, GOL, ACT, EDO, PLP, PO4

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	0.49	0/4925	0.64	0/6655
1	B	0.48	0/4925	0.63	0/6655
1	C	0.53	0/4932	0.66	4/6665 (0.1%)
1	D	0.49	0/4869	0.64	0/6581
All	All	0.50	0/19651	0.64	4/26556 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	276	ARG	NE-CZ-NH1	-5.93	115.57	121.50
1	C	276	ARG	NE-CZ-NH2	5.71	124.34	119.20
1	C	509	GLU	CA-C-N	5.09	131.25	121.54
1	C	509	GLU	C-N-CA	5.09	131.25	121.54

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	4831	0	4866	68	0
1	B	4831	0	4866	80	0
1	C	4838	0	4875	63	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	4776	0	4803	43	0
2	A	16	0	7	1	0
2	B	16	0	8	0	0
2	C	16	0	7	1	0
2	D	16	0	7	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
4	A	10	0	0	0	0
4	B	15	0	0	0	0
4	C	15	0	0	0	0
4	D	15	0	0	3	0
5	A	27	0	0	2	0
5	B	27	0	0	2	0
5	C	27	0	0	0	0
5	D	27	0	0	0	0
6	A	40	0	58	16	0
6	B	12	0	18	9	0
6	C	28	0	42	9	0
6	D	28	0	42	2	0
7	A	8	0	6	0	0
7	C	4	0	3	0	0
8	B	12	0	14	3	0
9	A	354	0	0	7	1
9	B	352	0	0	13	1
9	C	411	0	0	15	0
9	D	330	0	0	9	0
All	All	21090	0	19622	259	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 7.

The worst 5 of 259 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:337:ASN:ND2	1:C:366:LEU:HD11	1.77	0.99
1:A:321:MET:HE2	1:A:413:ILE:HG23	1.45	0.99
1:B:1:MET:SD	9:B:1144:HOH:O	2.30	0.89
1:C:83:THR:O	1:C:192:LYS:NZ	2.05	0.88
1:D:417:VAL:HG12	1:D:441:LYS:HD2	1.54	0.88

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)
9:A:813:HOH:O	9:B:832:HOH:O[1_554]	2.18	0.02

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	613/624 (98%)	589 (96%)	24 (4%)	0	100	100
1	B	613/624 (98%)	594 (97%)	19 (3%)	0	100	100
1	C	614/624 (98%)	594 (97%)	19 (3%)	1 (0%)	43	36
1	D	604/624 (97%)	584 (97%)	20 (3%)	0	100	100
All	All	2444/2496 (98%)	2361 (97%)	82 (3%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	56	GLY

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	524/533 (98%)	524 (100%)	0	100	100
1	B	524/533 (98%)	522 (100%)	2 (0%)	84	85

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	525/533 (98%)	524 (100%)	1 (0%)	87	89
1	D	518/533 (97%)	516 (100%)	2 (0%)	84	85
All	All	2091/2132 (98%)	2086 (100%)	5 (0%)	87	89

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	165	LEU
1	B	483	LYS
1	C	589	ASN
1	D	355	MET
1	D	589	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 11 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	285	GLN
1	C	369	GLN
1	D	589	ASN
1	C	589	ASN
1	B	523	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 59 ligands modelled in this entry, 8 are monoatomic - leaving 51 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	PO4	C	2215	-	4,4,4	0.99	0	6,6,6	0.87	0
6	EDO	B	707	-	3,3,3	0.37	0	2,2,2	0.46	0
8	GOL	B	709	-	5,5,5	0.55	0	5,5,5	1.13	0
4	PO4	A	718	-	4,4,4	1.10	0	6,6,6	0.65	0
2	PLP	C	2202	-	16,16,16	1.07	1 (6%)	20,23,23	1.10	1 (5%)
6	EDO	D	711	-	3,3,3	0.28	0	2,2,2	0.84	0
6	EDO	A	717	-	3,3,3	0.24	0	2,2,2	0.36	0
6	EDO	C	2209	-	3,3,3	0.77	0	2,2,2	0.10	0
4	PO4	D	714	-	4,4,4	0.45	0	6,6,6	0.74	0
6	EDO	A	706	-	3,3,3	0.70	0	2,2,2	0.17	0
6	EDO	D	707	-	3,3,3	0.49	0	2,2,2	0.57	0
4	PO4	D	709	-	4,4,4	1.11	0	6,6,6	0.82	0
2	PLP	A	701	-	16,16,16	1.13	2 (12%)	20,23,23	1.49	4 (20%)
7	ACT	A	707	-	3,3,3	2.12	1 (33%)	3,3,3	1.36	0
6	EDO	A	710	-	3,3,3	0.33	0	2,2,2	0.95	0
6	EDO	B	706	-	3,3,3	0.23	0	2,2,2	0.41	0
4	PO4	C	2204	-	4,4,4	0.65	0	6,6,6	1.30	1 (16%)
6	EDO	A	714	-	3,3,3	0.40	0	2,2,2	0.45	0
6	EDO	B	710	-	3,3,3	0.61	0	2,2,2	0.56	0
6	EDO	D	713	-	3,3,3	0.58	0	2,2,2	0.42	0
8	GOL	B	711	-	5,5,5	1.87	3 (60%)	5,5,5	1.18	0
6	EDO	D	708	-	3,3,3	0.56	0	2,2,2	0.38	0
6	EDO	C	2212	-	3,3,3	0.53	0	2,2,2	0.15	0
4	PO4	C	2207	-	4,4,4	0.77	0	6,6,6	0.72	0
7	ACT	C	2201	-	3,3,3	1.51	1 (33%)	3,3,3	1.24	0
5	ORC	A	705	3	28,28,28	4.57	13 (46%)	37,42,42	1.96	7 (18%)
5	ORC	D	705	3	28,28,28	4.34	15 (53%)	37,42,42	1.40	5 (13%)
6	EDO	A	713	-	3,3,3	0.41	0	2,2,2	0.35	0
6	EDO	C	2206	-	3,3,3	0.38	0	2,2,2	0.68	0
4	PO4	B	708	-	4,4,4	0.89	0	6,6,6	0.67	0
6	EDO	A	711	-	3,3,3	0.60	0	2,2,2	1.22	0
6	EDO	C	2208	-	3,3,3	0.35	0	2,2,2	0.97	0
5	ORC	B	705	3	28,28,28	4.41	15 (53%)	37,42,42	1.76	8 (21%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	EDO	C	2213	-	3,3,3	0.58	0	2,2,2	0.29	0
6	EDO	D	706	-	3,3,3	0.45	0	2,2,2	0.78	0
4	PO4	B	712	-	4,4,4	0.35	0	6,6,6	0.83	0
6	EDO	C	2210	-	3,3,3	0.54	0	2,2,2	0.53	0
6	EDO	C	2214	-	3,3,3	0.52	0	2,2,2	0.62	0
4	PO4	D	704	-	4,4,4	0.67	0	6,6,6	0.73	0
6	EDO	D	712	-	3,3,3	0.56	0	2,2,2	0.58	0
6	EDO	A	712	-	3,3,3	0.29	0	2,2,2	0.93	0
2	PLP	B	701	-	16,16,16	1.28	1 (6%)	20,23,23	1.23	3 (15%)
7	ACT	A	715	-	3,3,3	1.30	0	3,3,3	1.35	0
5	ORC	C	2205	3	28,28,28	4.58	14 (50%)	37,42,42	1.51	6 (16%)
6	EDO	A	708	-	3,3,3	0.53	0	2,2,2	0.30	0
4	PO4	B	704	-	4,4,4	0.91	0	6,6,6	0.88	0
6	EDO	A	709	-	3,3,3	0.49	0	2,2,2	0.76	0
2	PLP	D	702	-	16,16,16	1.21	2 (12%)	20,23,23	1.21	2 (10%)
4	PO4	A	704	-	4,4,4	0.94	0	6,6,6	1.09	0
6	EDO	D	710	-	3,3,3	0.60	0	2,2,2	0.09	0
6	EDO	A	716	-	3,3,3	0.39	0	2,2,2	0.86	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
6	EDO	B	707	-	-	1/1/1/1	-
8	GOL	B	709	-	-	2/4/4/4	-
2	PLP	C	2202	-	-	2/8/8/8	0/1/1/1
6	EDO	D	711	-	-	0/1/1/1	-
6	EDO	A	717	-	-	0/1/1/1	-
6	EDO	C	2209	-	-	0/1/1/1	-
6	EDO	A	706	-	-	0/1/1/1	-
6	EDO	D	707	-	-	0/1/1/1	-
2	PLP	A	701	-	-	1/8/8/8	0/1/1/1
6	EDO	A	710	-	-	0/1/1/1	-
6	EDO	A	714	-	-	1/1/1/1	-
6	EDO	B	710	-	-	1/1/1/1	-
6	EDO	D	713	-	-	0/1/1/1	-
8	GOL	B	711	-	-	2/4/4/4	-
6	EDO	D	708	-	-	0/1/1/1	-
6	EDO	C	2212	-	-	1/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	ORC	A	705	3	-	4/17/36/36	0/2/2/2
5	ORC	D	705	3	-	3/17/36/36	0/2/2/2
6	EDO	A	713	-	-	1/1/1/1	-
6	EDO	C	2206	-	-	1/1/1/1	-
6	EDO	A	711	-	-	0/1/1/1	-
6	EDO	C	2208	-	-	0/1/1/1	-
5	ORC	B	705	3	-	3/17/36/36	0/2/2/2
6	EDO	C	2213	-	-	0/1/1/1	-
6	EDO	D	706	-	-	1/1/1/1	-
6	EDO	C	2210	-	-	0/1/1/1	-
6	EDO	C	2214	-	-	0/1/1/1	-
6	EDO	D	712	-	-	0/1/1/1	-
6	EDO	A	712	-	-	1/1/1/1	-
2	PLP	B	701	-	-	1/8/8/8	0/1/1/1
5	ORC	C	2205	3	-	2/17/36/36	0/2/2/2
6	EDO	A	708	-	-	0/1/1/1	-
6	EDO	A	709	-	-	1/1/1/1	-
2	PLP	D	702	-	-	0/8/8/8	0/1/1/1
6	EDO	B	706	-	-	1/1/1/1	-
6	EDO	D	710	-	-	0/1/1/1	-
6	EDO	A	716	-	-	0/1/1/1	-

The worst 5 of 68 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	C	2205	ORC	PA-O3A	11.00	1.71	1.59
5	C	2205	ORC	PB-O3A	10.60	1.70	1.58
5	A	705	ORC	PB-O3A	10.39	1.70	1.58
5	A	705	ORC	PA-O3A	9.96	1.70	1.59
5	B	705	ORC	PB-O3A	9.65	1.69	1.58

The worst 5 of 37 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	705	ORC	PB-C1B-C14	-7.59	97.86	114.06
5	B	705	ORC	PB-O3A-PA	-4.23	118.57	132.37
5	B	705	ORC	C3'-C2'-C1'	4.02	109.08	101.46
5	A	705	ORC	PB-O3A-PA	-3.99	119.35	132.37
5	B	705	ORC	O4'-C1'-C2'	-3.91	98.26	106.62

There are no chirality outliers.

5 of 30 torsion outliers are listed below:

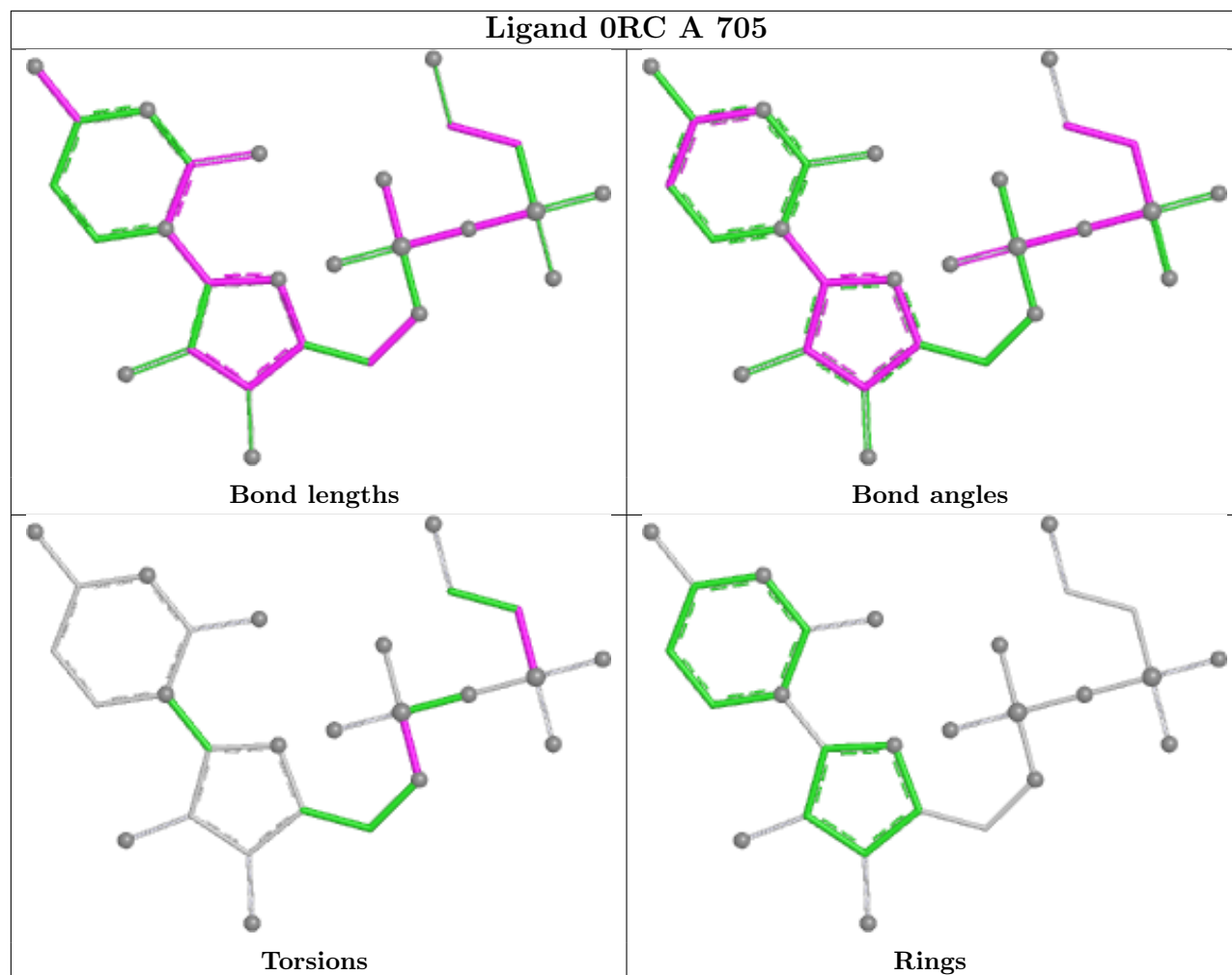
Mol	Chain	Res	Type	Atoms
5	A	705	0RC	C14-C1B-PB-O3B
5	A	705	0RC	C14-C1B-PB-O3A
5	B	705	0RC	C14-C1B-PB-O3B
5	B	705	0RC	C14-C1B-PB-O2B
5	B	705	0RC	C14-C1B-PB-O3A

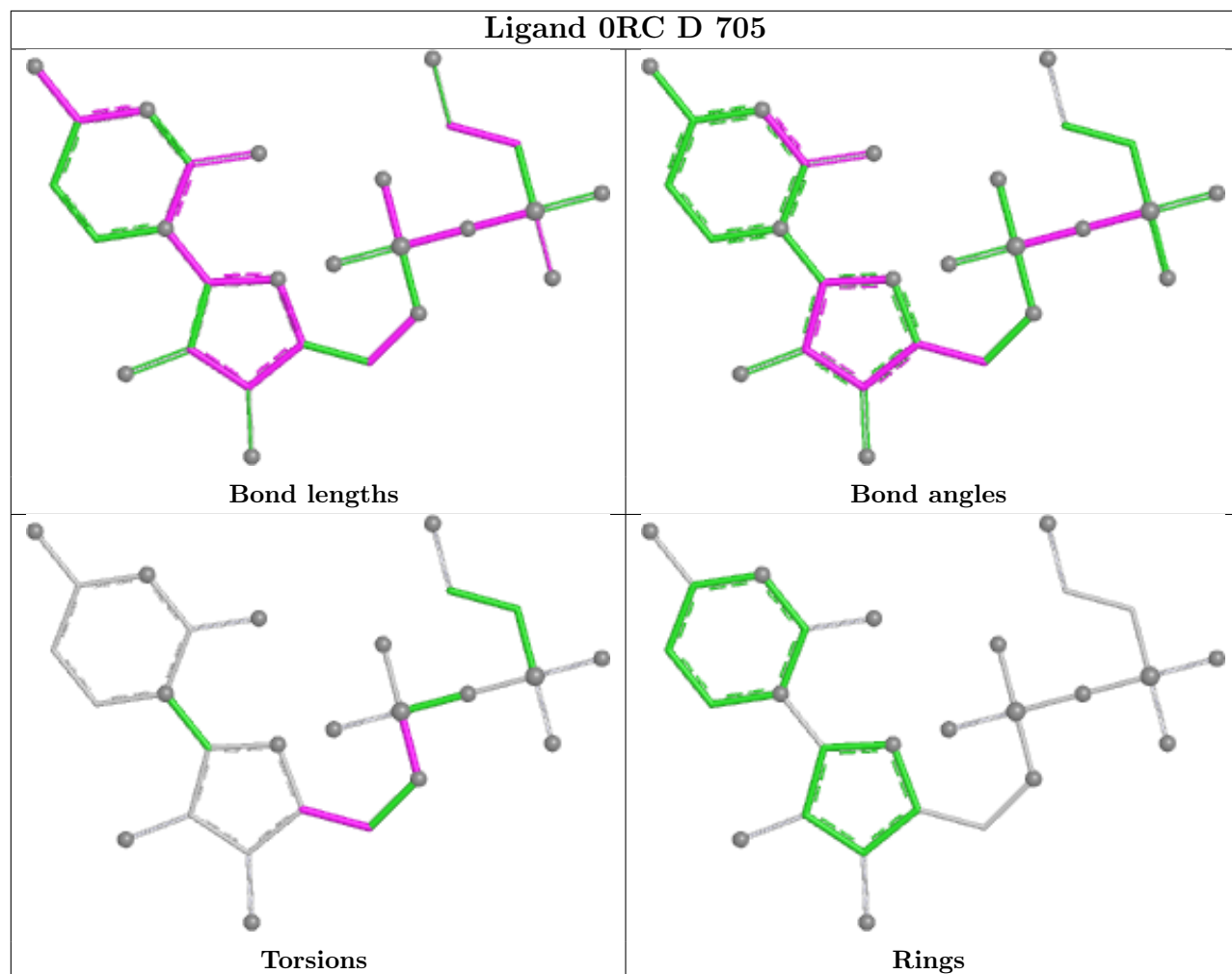
There are no ring outliers.

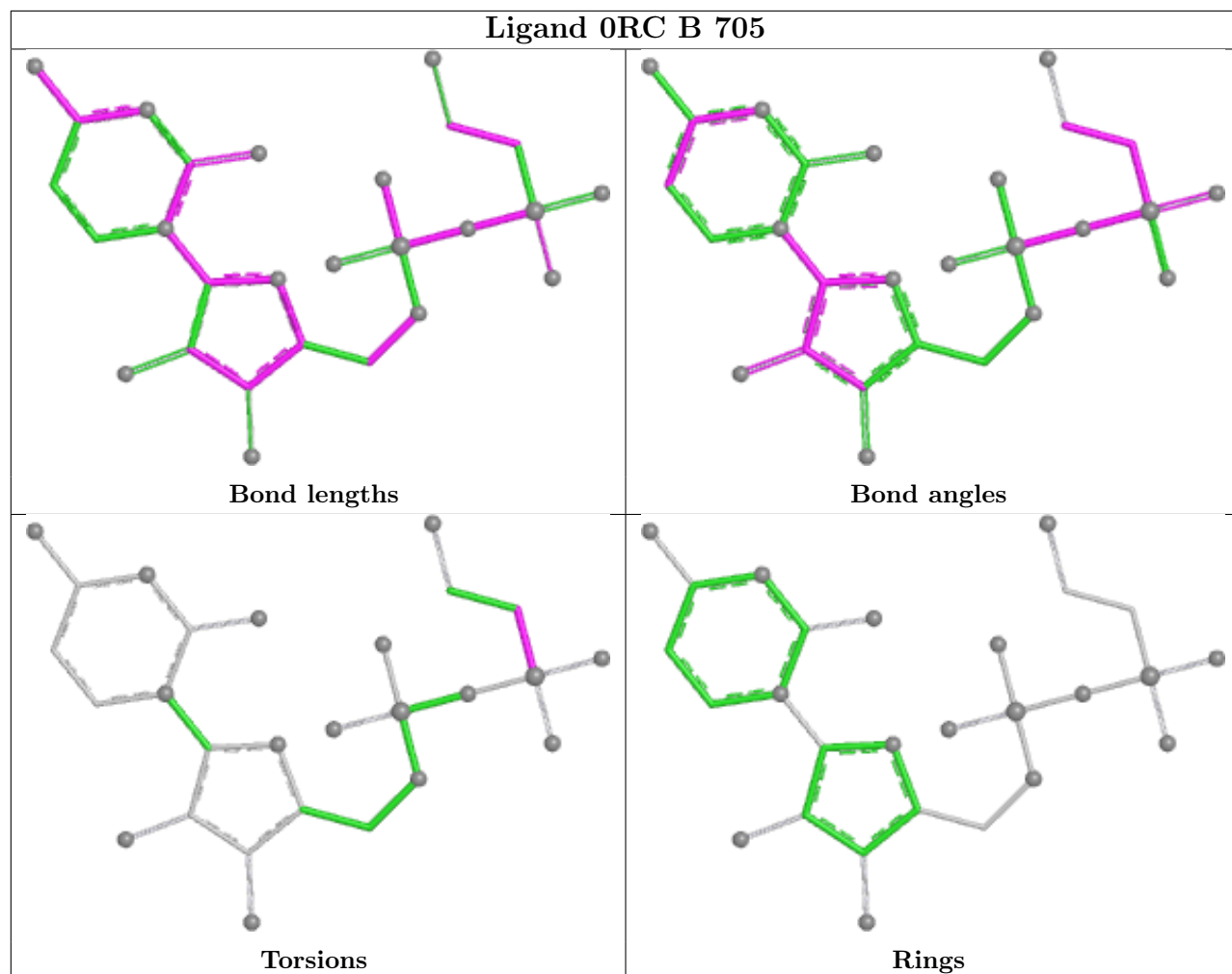
18 monomers are involved in 47 short contacts:

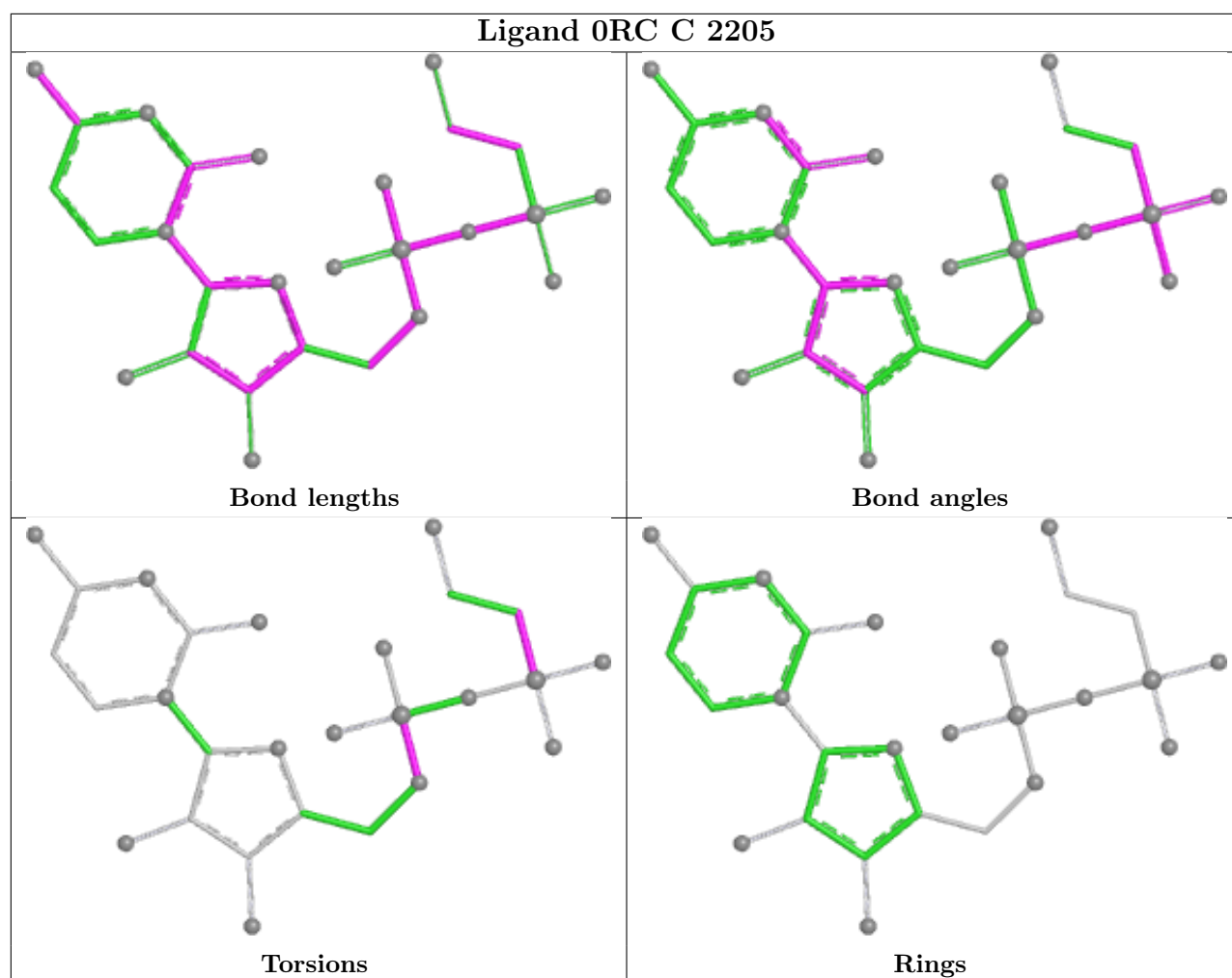
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	2202	PLP	1	0
6	A	717	EDO	3	0
6	C	2209	EDO	3	0
4	D	714	PO4	1	0
4	D	709	PO4	2	0
2	A	701	PLP	1	0
6	B	706	EDO	9	0
6	A	714	EDO	1	0
8	B	711	GOL	3	0
5	A	705	0RC	2	0
6	C	2206	EDO	1	0
6	C	2208	EDO	2	0
5	B	705	0RC	2	0
6	D	706	EDO	2	0
6	C	2210	EDO	3	0
6	A	712	EDO	1	0
6	A	709	EDO	5	0
6	A	716	EDO	6	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	615/624 (98%)	1.18	90 (14%) 6 6	15, 25, 37, 51	0
1	B	615/624 (98%)	1.27	106 (17%) 4 4	15, 26, 42, 60	0
1	C	616/624 (98%)	1.08	59 (9%) 13 15	13, 23, 36, 54	0
1	D	608/624 (97%)	1.31	106 (17%) 4 4	15, 25, 42, 51	0
All	All	2454/2496 (98%)	1.21	361 (14%) 6 6	13, 25, 40, 60	0

The worst 5 of 361 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	580	GLY	5.7
1	C	366	LEU	5.1
1	D	185	ILE	4.8
1	C	616	VAL	4.7
1	B	181	ALA	4.5

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 oligosaccharides 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.

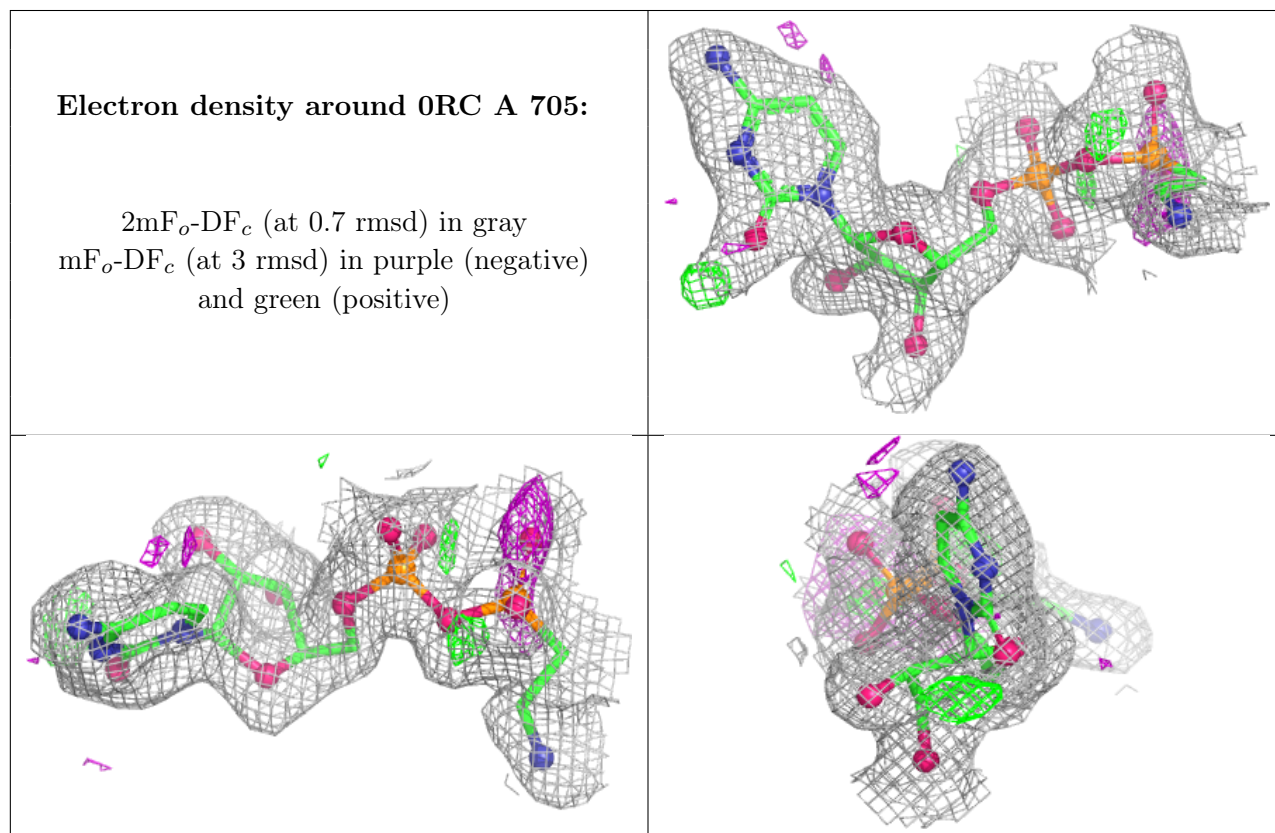
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	MG	D	703	1/1	0.55	0.19	46,46,46,46	0
4	PO4	D	704	5/5	0.58	0.17	54,60,67,78	0
4	PO4	B	712	5/5	0.66	0.19	30,31,43,51	0
6	EDO	A	711	4/4	0.67	0.24	35,40,42,42	0
7	ACT	A	707	4/4	0.68	0.19	25,29,31,33	0
6	EDO	C	2209	4/4	0.70	0.20	22,29,31,32	0
4	PO4	C	2215	5/5	0.71	0.19	30,36,40,48	0
7	ACT	A	715	4/4	0.72	0.20	27,34,37,46	0
4	PO4	A	718	5/5	0.75	0.19	30,32,43,51	0
6	EDO	D	707	4/4	0.77	0.17	30,32,38,39	0
6	EDO	B	706	4/4	0.77	0.24	20,21,22,24	0
6	EDO	A	716	4/4	0.77	0.36	21,22,23,29	0
3	MG	B	702	1/1	0.78	0.12	34,34,34,34	0
8	GOL	B	709	6/6	0.78	0.15	24,36,38,41	0
4	PO4	B	708	5/5	0.79	0.14	36,37,48,50	0
7	ACT	C	2201	4/4	0.79	0.16	21,22,29,34	0
3	MG	A	702	1/1	0.79	0.13	31,31,31,31	0
3	MG	C	2203	1/1	0.80	0.14	28,28,28,28	0
6	EDO	B	707	4/4	0.80	0.17	22,22,25,26	0
6	EDO	C	2210	4/4	0.82	0.14	23,33,33,37	0
6	EDO	A	709	4/4	0.82	0.17	20,22,25,26	0
4	PO4	D	714	5/5	0.82	0.18	28,33,38,46	0
4	PO4	C	2207	5/5	0.83	0.17	30,37,41,43	0
6	EDO	D	708	4/4	0.83	0.16	30,31,32,40	0
6	EDO	B	710	4/4	0.83	0.16	23,24,24,29	0
6	EDO	D	706	4/4	0.84	0.17	19,21,24,30	0
6	EDO	D	713	4/4	0.84	0.14	23,25,25,27	0
6	EDO	A	714	4/4	0.84	0.14	26,27,36,38	0
6	EDO	A	717	4/4	0.86	0.22	26,27,30,34	0
6	EDO	A	712	4/4	0.86	0.18	17,20,26,27	0
6	EDO	A	713	4/4	0.87	0.13	24,27,28,28	0
6	EDO	A	710	4/4	0.87	0.11	21,21,22,25	0
6	EDO	A	706	4/4	0.87	0.12	14,18,18,20	0
5	ORC	A	705	27/27	0.87	0.12	25,27,32,37	0
6	EDO	D	710	4/4	0.87	0.11	21,24,26,29	0
8	GOL	B	711	6/6	0.87	0.20	15,20,27,27	0
2	PLP	D	702	16/16	0.89	0.12	17,20,31,32	0
6	EDO	D	712	4/4	0.89	0.09	15,15,17,18	0
5	ORC	B	705	27/27	0.89	0.11	27,33,41,43	0
6	EDO	C	2212	4/4	0.89	0.12	16,17,17,22	0
5	ORC	D	705	27/27	0.90	0.11	31,35,37,40	0
2	PLP	C	2202	16/16	0.91	0.11	14,19,23,30	0
6	EDO	C	2206	4/4	0.91	0.17	24,25,28,33	0

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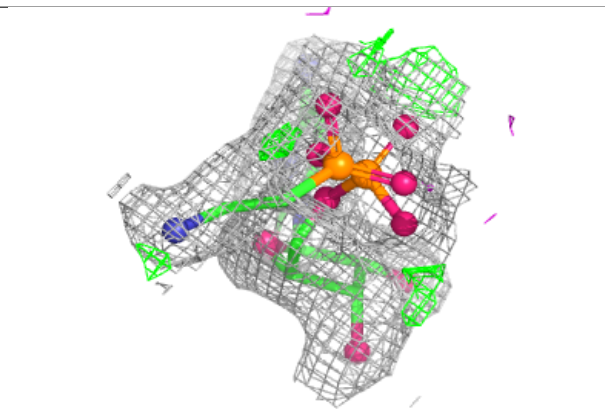
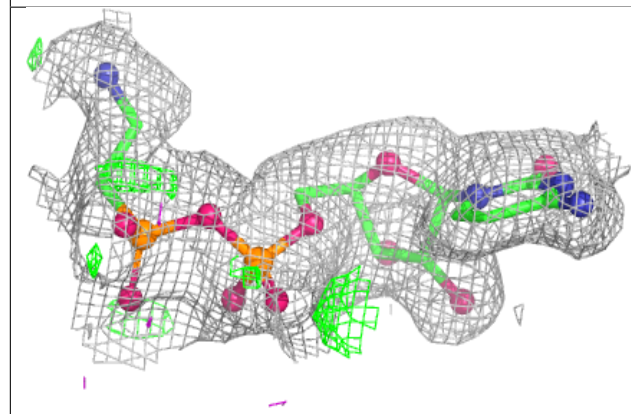
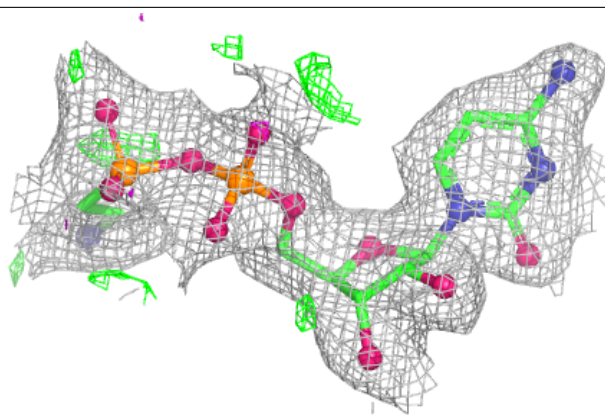
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	MG	A	703	1/1	0.91	0.06	27,27,27,27	0
4	PO4	C	2204	5/5	0.91	0.12	24,25,26,29	0
5	ORC	C	2205	27/27	0.92	0.10	22,26,31,34	0
6	EDO	C	2208	4/4	0.92	0.12	19,24,24,25	0
6	EDO	C	2214	4/4	0.93	0.10	19,23,24,26	0
4	PO4	A	704	5/5	0.93	0.10	23,24,25,35	0
6	EDO	C	2213	4/4	0.93	0.12	17,20,21,23	0
3	MG	D	701	1/1	0.94	0.07	41,41,41,41	0
4	PO4	D	709	5/5	0.94	0.12	22,24,29,31	0
2	PLP	A	701	16/16	0.94	0.09	16,20,25,32	0
3	MG	B	703	1/1	0.94	0.06	33,33,33,33	0
2	PLP	B	701	16/16	0.94	0.11	18,23,29,36	0
4	PO4	B	704	5/5	0.94	0.11	24,27,30,33	0
6	EDO	D	711	4/4	0.94	0.09	20,24,25,26	0
6	EDO	A	708	4/4	0.95	0.07	21,21,23,29	0
3	MG	C	2211	1/1	0.97	0.04	26,26,26,26	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.

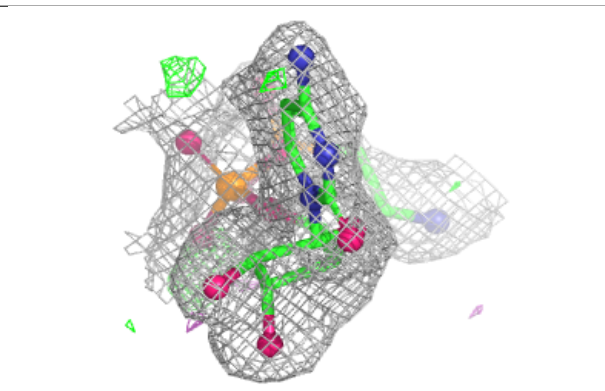
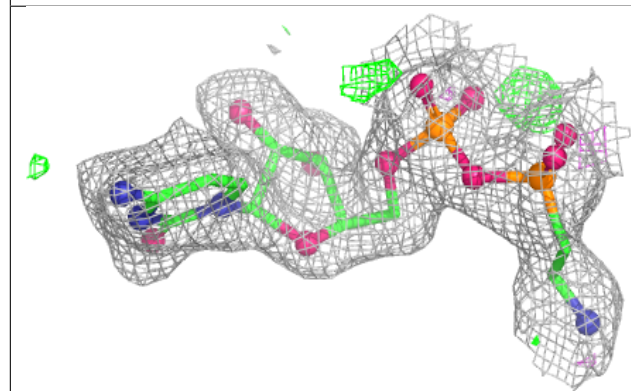
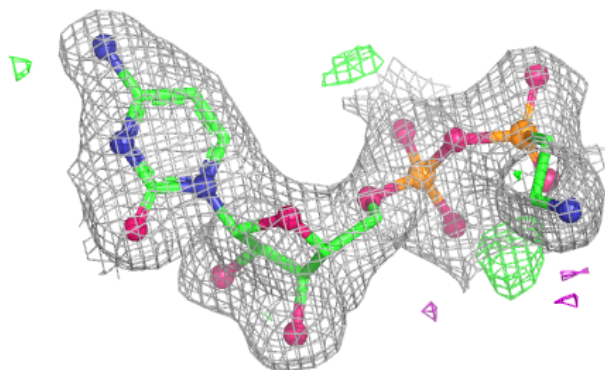


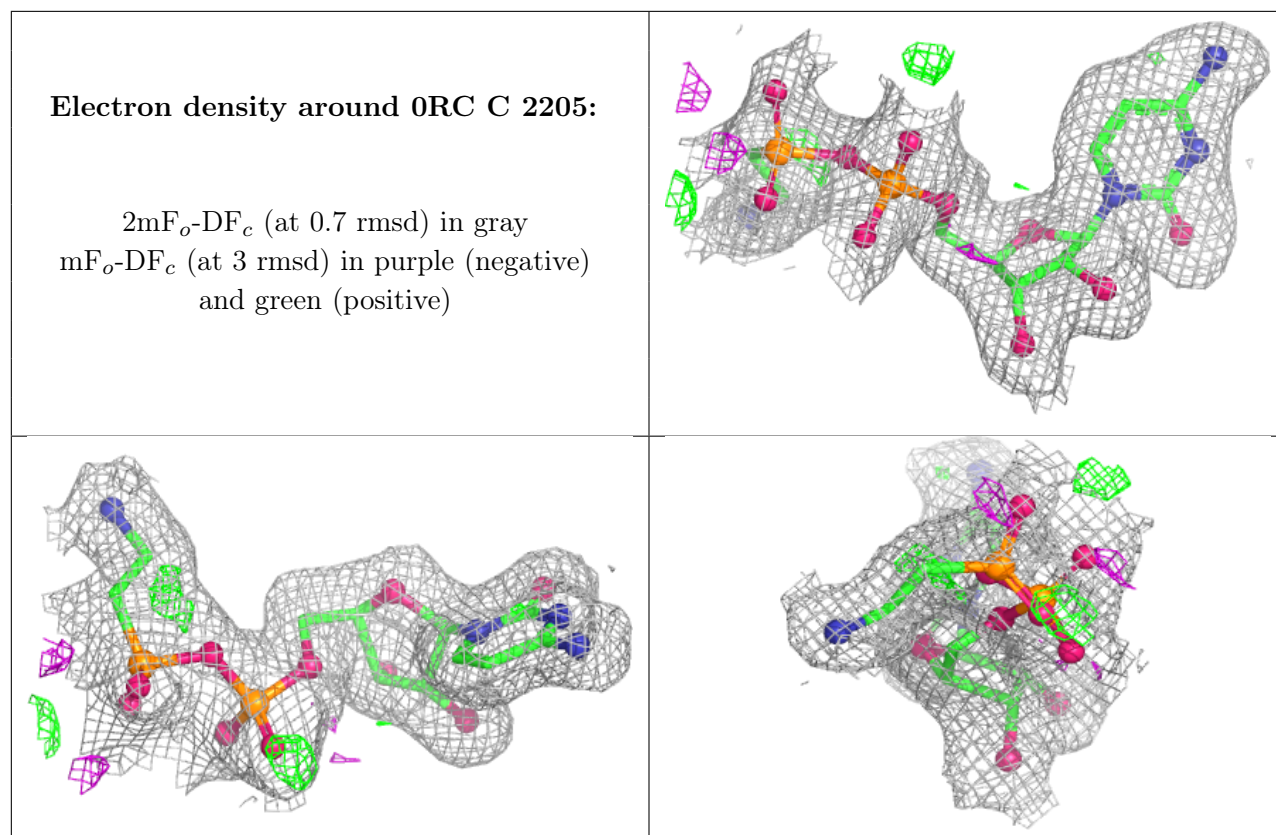
Electron density around 0RC B 705:

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

**Electron density around 0RC D 705:**

$2mF_o-DF_c$ (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.