



Full wwPDB X-ray Structure Validation Report ⓘ

May 17, 2020 – 02:55 pm BST

PDB ID : 4OX2
Title : I45T cytosolic phosphoenolpyruvate carboxykinase in complex with beta-sulfofpyruvate and GTP
Authors : Holyoak, T.
Deposited on : 2014-02-04
Resolution : 2.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The 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

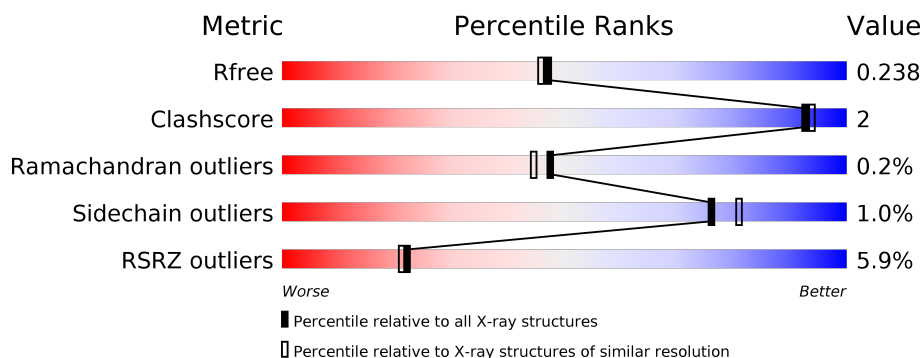
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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	622	<div> <div>5%</div> <div> <div></div> <div>92%</div> <div>5%</div> </div> <div></div> </div>
1	B	622	<div> <div>6%</div> <div> <div></div> <div>96%</div> <div></div> </div> <div></div> </div>

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 10254 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 Phosphoenolpyruvate carboxykinase, cytosolic [GTP].

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	608	Total	C	N	O	S	0	2	0
			4792	3066	820	873	33			
1	B	620	Total	C	N	O	S	0	3	0
			4890	3123	839	896	32			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	45	THR	ILE	engineered mutation	UNP P07379
B	45	THR	ILE	engineered mutation	UNP P07379

- Molecule 2 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

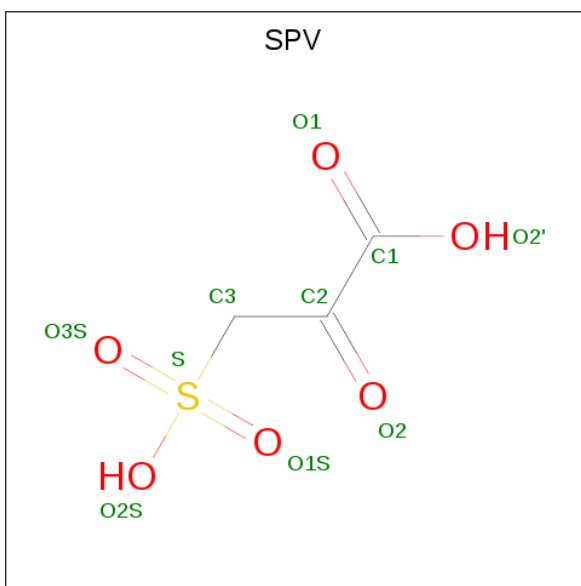
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	3	Total	Mn	0	0
			3	3		
2	A	3	Total	Mn	0	0
			3	3		

- Molecule 3 is GUANOSINE-5'-TRIPHOSPHATE (three-letter code: GTP) (formula: C₁₀H₁₆N₅O₁₄P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 32	C 10	N 5	O 14	P 3	0	0
3	B	1	Total 32	C 10	N 5	O 14	P 3	0	0

- Molecule 4 is SULFOPYRUVATE (three-letter code: SPV) (formula: $\text{C}_3\text{H}_4\text{O}_6\text{S}$).

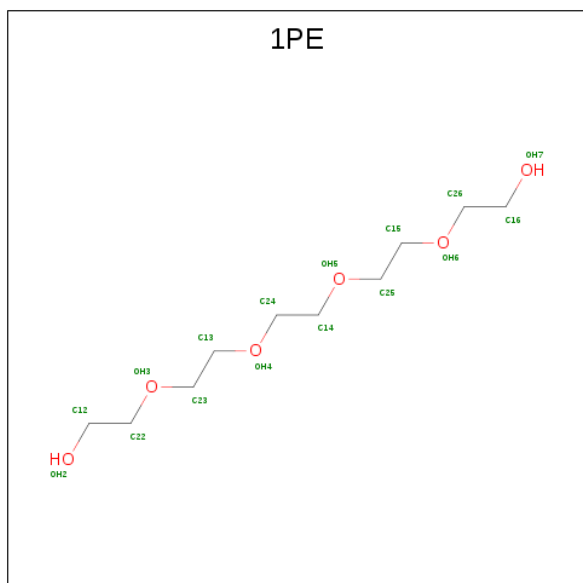


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total 10	C 3	O 6	S 1	0	0
4	B	1	Total 10	C 3	O 6	S 1	0	0

- Molecule 5 is SODIUM ION (three-letter code: NA) (formula: Na).

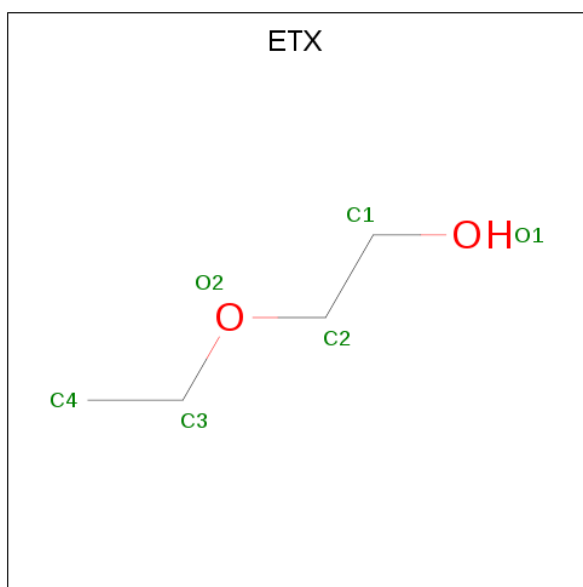
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total	Na	0	0
			1	1		
5	A	1	Total	Na	0	0
			1	1		

- Molecule 6 is PENTAETHYLENE GLYCOL (three-letter code: 1PE) (formula: C₁₀H₂₂O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			7	4	3		
6	B	1	Total	C	O	0	0
			10	6	4		

- Molecule 7 is 2-ETHOXYETHANOL (three-letter code: ETX) (formula: C₄H₁₀O₂).



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

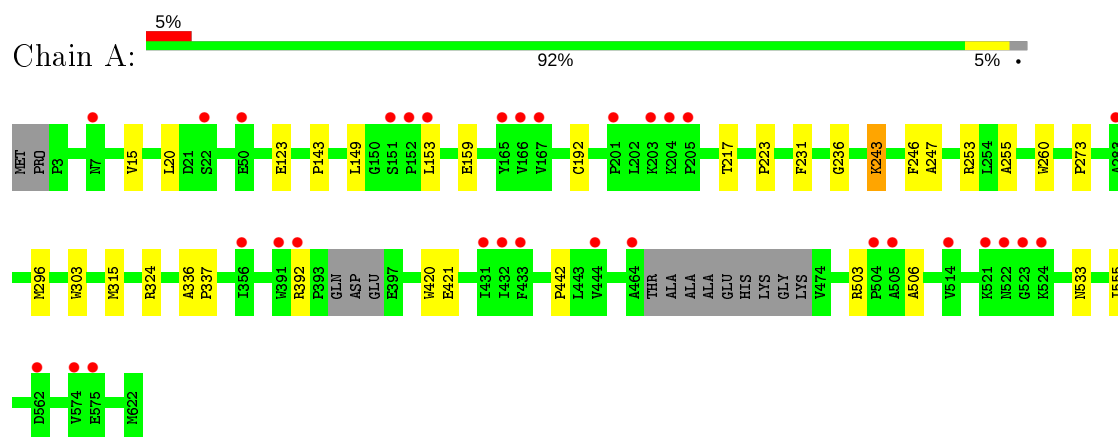
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	223	Total	O	0	0
			223	223		
8	B	234	Total	O	0	0
			234	234		

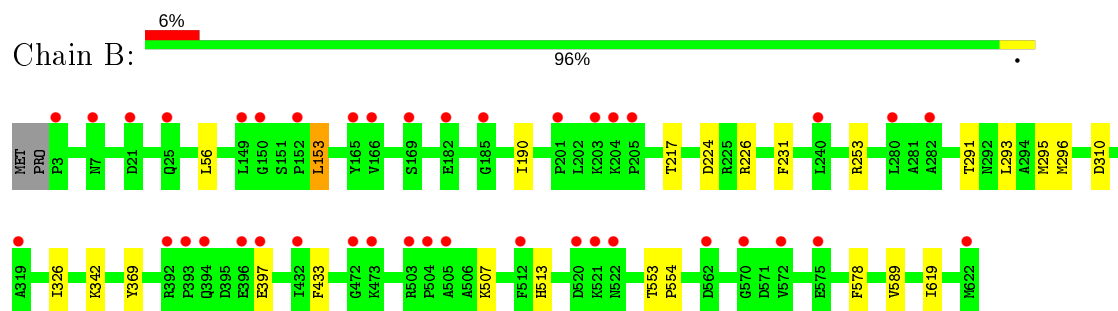
3 Residue-property plots [i](#)

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: Phosphoenolpyruvate carboxykinase, cytosolic [GTP]



- Molecule 1: Phosphoenolpyruvate carboxykinase, cytosolic [GTP]



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	61.92Å 119.25Å 87.40Å 90.00° 106.92° 90.00°	Depositor
Resolution (Å)	29.13 – 2.00 29.13 – 2.00	Depositor EDS
% Data completeness (in resolution range)	98.1 (29.13-2.00) 98.1 (29.13-2.00)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.08 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.189 , 0.233 0.195 , 0.238	Depositor DCC
R_{free} test set	4021 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	32.4	Xtriage
Anisotropy	0.160	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 50.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	10254	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 69.88 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.4354e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NA, ETX, MN, SPV, IPE, GTP

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	0.42	0/4921	0.62	0/6660
1	B	0.43	0/5022	0.60	0/6798
All	All	0.43	0/9943	0.61	0/13458

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4792	0	4743	14	0
1	B	4890	0	4834	16	0
2	A	3	0	0	0	0
2	B	3	0	0	0	0
3	A	32	0	12	0	0
3	B	32	0	12	0	0
4	A	10	0	2	0	0
4	B	10	0	2	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
6	A	7	0	9	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	10	0	12	0	0
7	B	6	0	10	0	0
8	A	223	0	0	0	0
8	B	234	0	0	4	0
All	All	10254	0	9636	30	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 2.

All (30) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:293:LEU:HD12	1:B:296:MET:HE3	1.65	0.76
1:B:224[A]:ASP:OD2	8:B:801:HOH:O	2.07	0.73
1:B:153:LEU:HD11	1:B:326:ILE:CD1	2.25	0.67
1:B:224[A]:ASP:OD1	1:B:253:ARG:NH1	2.39	0.56
1:B:153:LEU:HD11	1:B:326:ILE:HD13	1.91	0.52
1:A:296:MET:HE2	1:A:533:ASN:HB2	1.95	0.48
1:A:217:THR:HA	1:A:231:PHE:O	2.13	0.48
1:A:315:MET:HA	1:A:324:ARG:O	2.15	0.47
1:A:255:ALA:HB1	1:A:260:TRP:O	2.14	0.47
1:B:507:LYS:NZ	8:B:803:HOH:O	2.48	0.46
1:B:293:LEU:HD12	1:B:296:MET:CE	2.42	0.46
1:B:342:LYS:HD3	8:B:1028:HOH:O	2.15	0.46
1:A:336:ALA:HB3	1:A:337:PRO:HD3	1.98	0.45
1:A:243:LYS:O	1:A:247:ALA:HB3	2.17	0.44
1:B:217:THR:HA	1:B:231:PHE:O	2.17	0.44
1:A:420:TRP:CE2	1:A:421:GLU:HG3	2.53	0.43
1:A:442:PRO:HG2	1:A:555:ILE:HD11	2.01	0.43
1:A:159:GLU:HA	1:A:192:CYS:HB2	2.01	0.43
1:A:255:ALA:HB1	1:A:260:TRP:C	2.40	0.43
1:A:15:VAL:HG21	1:A:20:LEU:HG	2.00	0.42
1:B:190:ILE:HD12	1:B:226:ARG:O	2.19	0.42
1:B:291:THR:OG1	1:B:310:ASP:OD2	2.38	0.42
1:B:578:PHE:HD1	8:B:1033:HOH:O	2.02	0.42
1:B:291:THR:O	1:B:295:MET:HG2	2.20	0.41
1:A:223:PRO:HB2	1:A:253:ARG:HD3	2.02	0.41
1:B:589:VAL:HG21	1:B:619:ILE:HD12	2.02	0.41
1:A:503:ARG:HB3	1:A:506:ALA:HB2	2.02	0.40
1:B:433:PHE:O	1:B:513:HIS:HA	2.20	0.40
1:A:273:PRO:HD3	1:A:303:TRP:CE2	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:553:THR:HB	1:B:554:PRO:HD2	2.03	0.40

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	604/622 (97%)	587 (97%)	15 (2%)	2 (0%)	41	37
1	B	621/622 (100%)	606 (98%)	15 (2%)	0	100	100
All	All	1225/1244 (98%)	1193 (97%)	30 (2%)	2 (0%)	47	44

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	243	LYS
1	A	236	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	511/519 (98%)	505 (99%)	6 (1%)	71	76
1	B	520/519 (100%)	516 (99%)	4 (1%)	81	86
All	All	1031/1038 (99%)	1021 (99%)	10 (1%)	76	81

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	123	GLU
1	A	143	PRO
1	A	149	LEU
1	A	153	LEU
1	A	246	PHE
1	A	392	ARG
1	B	56	LEU
1	B	153	LEU
1	B	369	TYR
1	B	397	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	522	ASN
1	B	297	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 15 ligands modelled in this entry, 8 are monoatomic - leaving 7 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	SPV	B	704	2	6,9,9	4.37	2 (33%)	6,13,13	1.00	0
6	1PE	A	707	-	6,6,15	0.45	0	5,5,14	0.35	0
3	GTP	B	703	2	26,34,34	1.15	2 (7%)	33,54,54	1.95	7 (21%)
7	ETX	B	708	-	5,5,5	0.52	0	4,4,4	0.32	0
6	1PE	B	707	-	9,9,15	0.47	0	8,8,14	0.26	0
4	SPV	A	704	2	6,9,9	4.69	2 (33%)	6,13,13	1.79	2 (33%)
3	GTP	A	703	2	26,34,34	1.09	3 (11%)	33,54,54	2.06	8 (24%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	SPV	B	704	2	-	4/5/9/9	-
6	1PE	A	707	-	-	0/4/4/13	-
3	GTP	B	703	2	-	1/18/38/38	0/3/3/3
7	ETX	B	708	-	-	0/3/3/3	-
6	1PE	B	707	-	-	2/7/7/13	-
4	SPV	A	704	2	-	1/5/9/9	-
3	GTP	A	703	2	-	3/18/38/38	0/3/3/3

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	704	SPV	C3-S	-9.81	1.68	1.78
4	B	704	SPV	C3-S	-9.03	1.69	1.78
4	A	704	SPV	O3S-S	5.70	1.61	1.45
4	B	704	SPV	O3S-S	5.60	1.61	1.45
3	B	703	GTP	C6-C5	4.16	1.48	1.41
3	A	703	GTP	C6-C5	2.60	1.45	1.41
3	A	703	GTP	O4'-C1'	2.43	1.44	1.41
3	B	703	GTP	C5-C4	2.35	1.47	1.40
3	A	703	GTP	C5-C4	2.24	1.46	1.40

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	703	GTP	C2-N3-C4	5.91	122.11	115.36
3	A	703	GTP	C6-C5-C4	-5.11	115.92	120.80
3	A	703	GTP	C6-N1-C2	4.76	123.50	115.93
3	A	703	GTP	C2-N3-C4	4.54	120.54	115.36
3	A	703	GTP	N3-C2-N1	-4.50	121.22	127.22
3	A	703	GTP	C5-C6-N1	-3.82	118.21	123.43
3	B	703	GTP	C4-C5-N7	-3.72	105.52	109.40
3	B	703	GTP	C5-C6-N1	-3.58	118.53	123.43
3	B	703	GTP	C6-C5-C4	-3.46	117.49	120.80
3	B	703	GTP	C6-N1-C2	3.26	121.10	115.93
3	B	703	GTP	N3-C2-N1	-3.16	123.01	127.22
4	A	704	SPV	O3S-S-C3	2.56	113.66	107.44
3	A	703	GTP	N2-C2-N1	2.47	121.10	117.25
3	B	703	GTP	O3G-PG-O2G	2.37	116.70	107.64
4	A	704	SPV	O2-C2-C3	2.29	123.49	119.56
3	A	703	GTP	PA-O3A-PB	-2.24	125.15	132.83
3	A	703	GTP	O3G-PG-O2G	2.15	115.85	107.64

There are no chirality outliers.

All (11) torsion outliers are listed below:

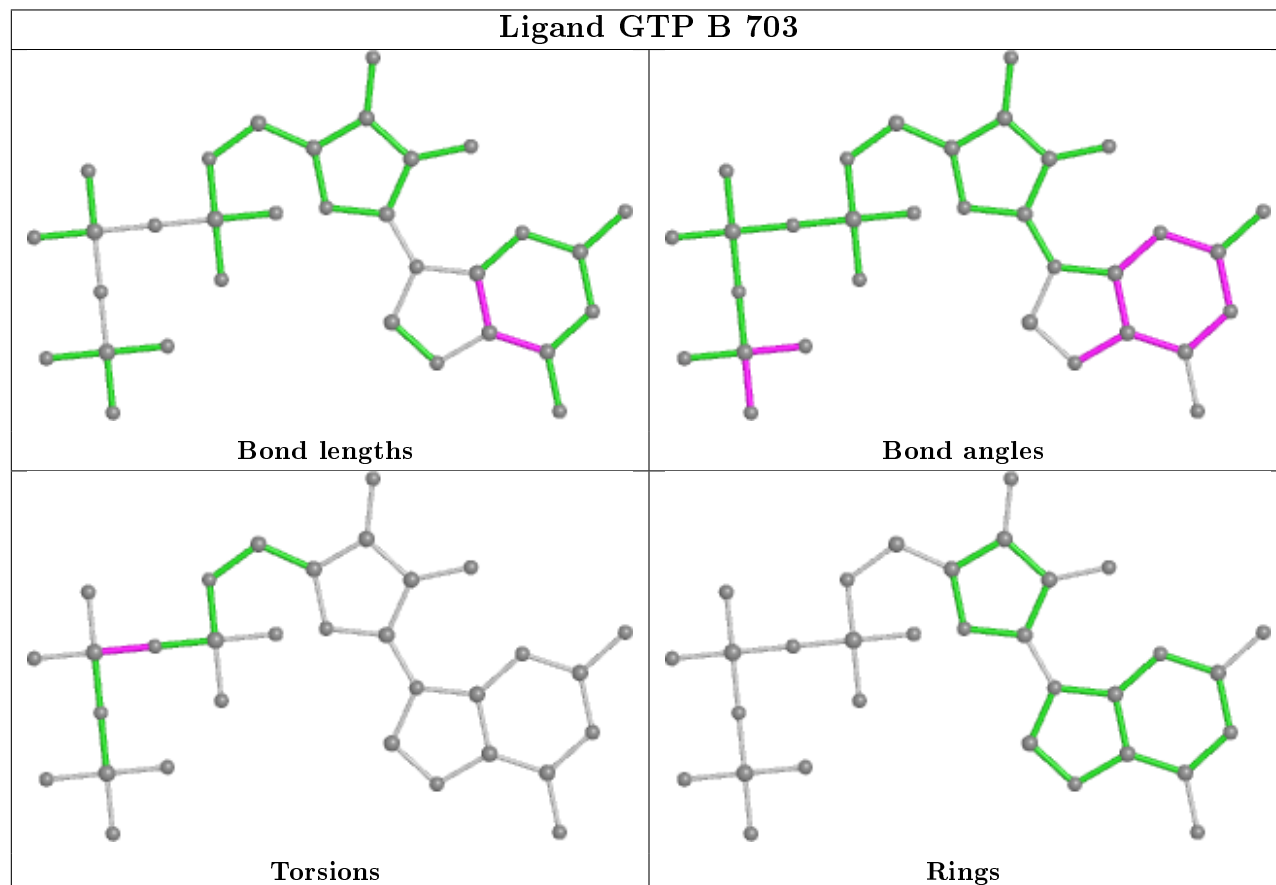
Mol	Chain	Res	Type	Atoms
4	B	704	SPV	C1-C2-C3-S
3	A	703	GTP	C5'-O5'-PA-O3A
6	B	707	1PE	OH4-C13-C23-OH3
3	A	703	GTP	PA-O3A-PB-O1B
3	A	703	GTP	C5'-O5'-PA-O1A
4	B	704	SPV	O2-C2-C3-S
4	B	704	SPV	C2-C3-S-O2S
4	A	704	SPV	C2-C3-S-O2S
3	B	703	GTP	PA-O3A-PB-O1B
6	B	707	1PE	OH6-C15-C25-OH5
4	B	704	SPV	C2-C3-S-O1S

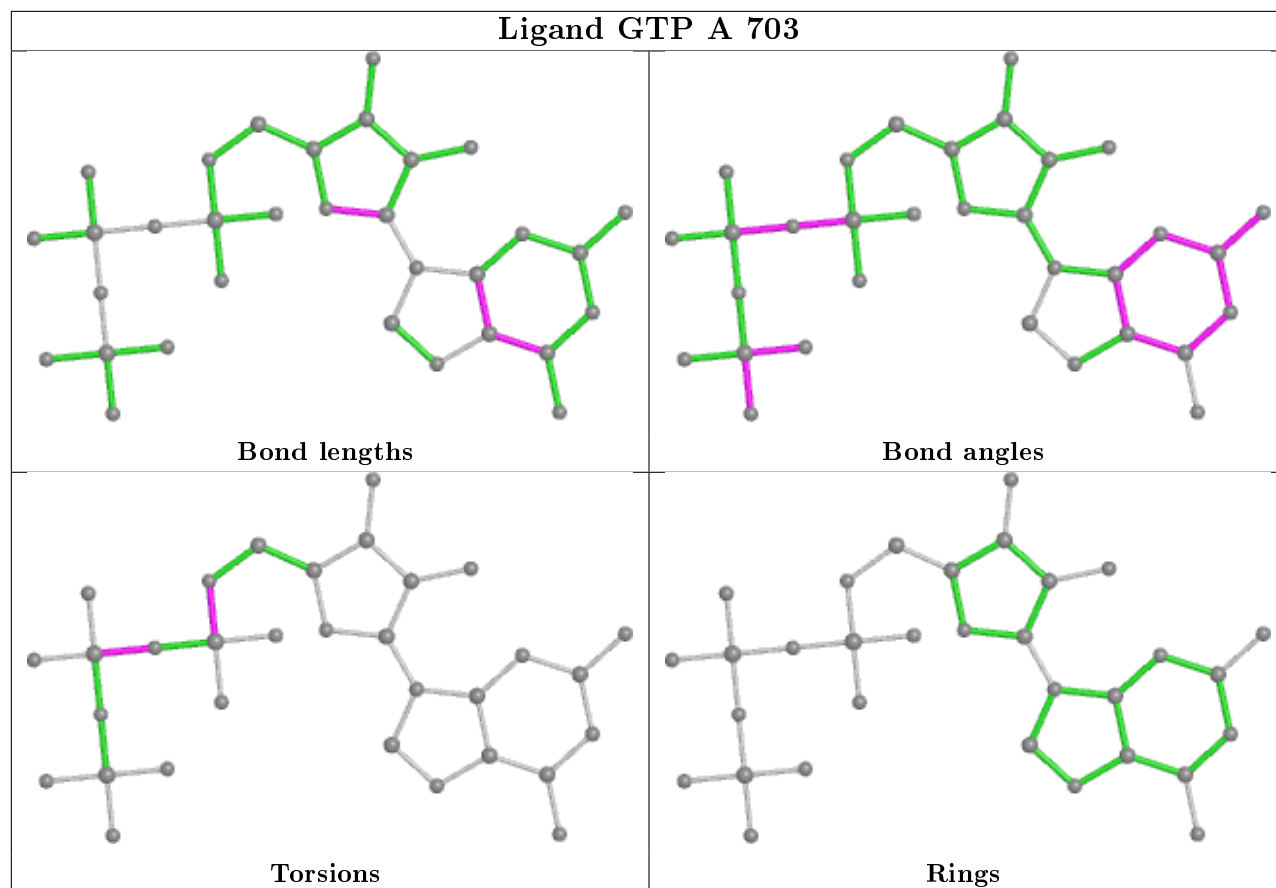
There are no ring outliers.

No monomer is involved in short contacts.

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

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	608/622 (97%)	0.15	32 (5%) 26 25	22, 32, 55, 74	0
1	B	620/622 (99%)	0.14	40 (6%) 18 18	22, 31, 53, 87	0
All	All	1228/1244 (98%)	0.15	72 (5%) 22 21	22, 32, 55, 87	0

All (72) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	504	PRO	6.8
1	B	394	GLN	6.6
1	B	393	PRO	6.1
1	B	392	ARG	6.0
1	A	152	PRO	5.5
1	B	473	LYS	5.1
1	B	396	GLU	4.6
1	A	504	PRO	4.5
1	A	153	LEU	4.3
1	A	203	LYS	4.3
1	A	523	GLY	3.8
1	B	150	GLY	3.8
1	B	203	LYS	3.7
1	A	464	ALA	3.5
1	B	204	LYS	3.5
1	A	432	ILE	3.4
1	B	205	PRO	3.4
1	B	562	ASP	3.4
1	B	521	LYS	3.3
1	A	392	ARG	3.3
1	A	575	GLU	3.3
1	B	575	GLU	3.1
1	A	7	ASN	3.0
1	A	201	PRO	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	562	ASP	3.0
1	A	522	ASN	3.0
1	B	432	ILE	2.9
1	A	431	ILE	2.9
1	A	505	ALA	2.8
1	B	397	GLU	2.8
1	B	152	PRO	2.8
1	B	472	GLY	2.7
1	B	166	VAL	2.7
1	A	205	PRO	2.7
1	B	240	LEU	2.7
1	B	572	VAL	2.7
1	A	165	TYR	2.6
1	B	165	TYR	2.6
1	A	521	LYS	2.5
1	B	185	GLY	2.5
1	B	201	PRO	2.5
1	B	505	ALA	2.5
1	A	22	SER	2.4
1	B	319	ALA	2.4
1	A	204	LYS	2.4
1	B	3	PRO	2.4
1	A	166	VAL	2.3
1	B	282	ALA	2.3
1	A	524	LYS	2.3
1	B	622	MET	2.3
1	B	169	SER	2.3
1	B	149	LEU	2.2
1	A	514	VAL	2.2
1	B	570	GLY	2.2
1	B	21	ASP	2.2
1	B	7	ASN	2.2
1	A	151	SER	2.2
1	B	280	LEU	2.2
1	B	503	ARG	2.2
1	A	167	VAL	2.1
1	A	283	ALA	2.1
1	A	433	PHE	2.1
1	B	512	PHE	2.1
1	A	391	TRP	2.1
1	A	356	ILE	2.1
1	B	182	GLU	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	50	GLU	2.0
1	B	522	ASN	2.0
1	A	444	VAL	2.0
1	A	574	VAL	2.0
1	B	25	GLN	2.0
1	B	520	ASP	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

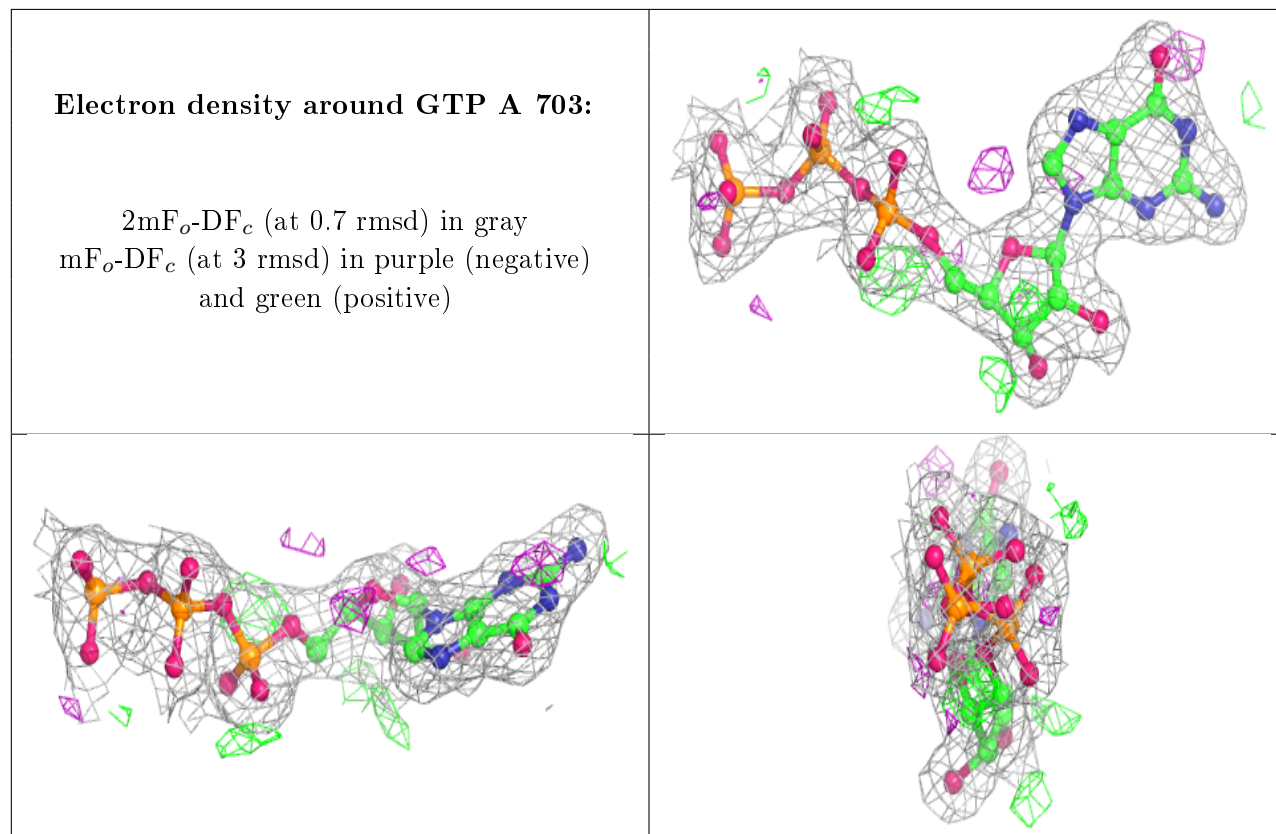
6.4 Ligands [i](#)

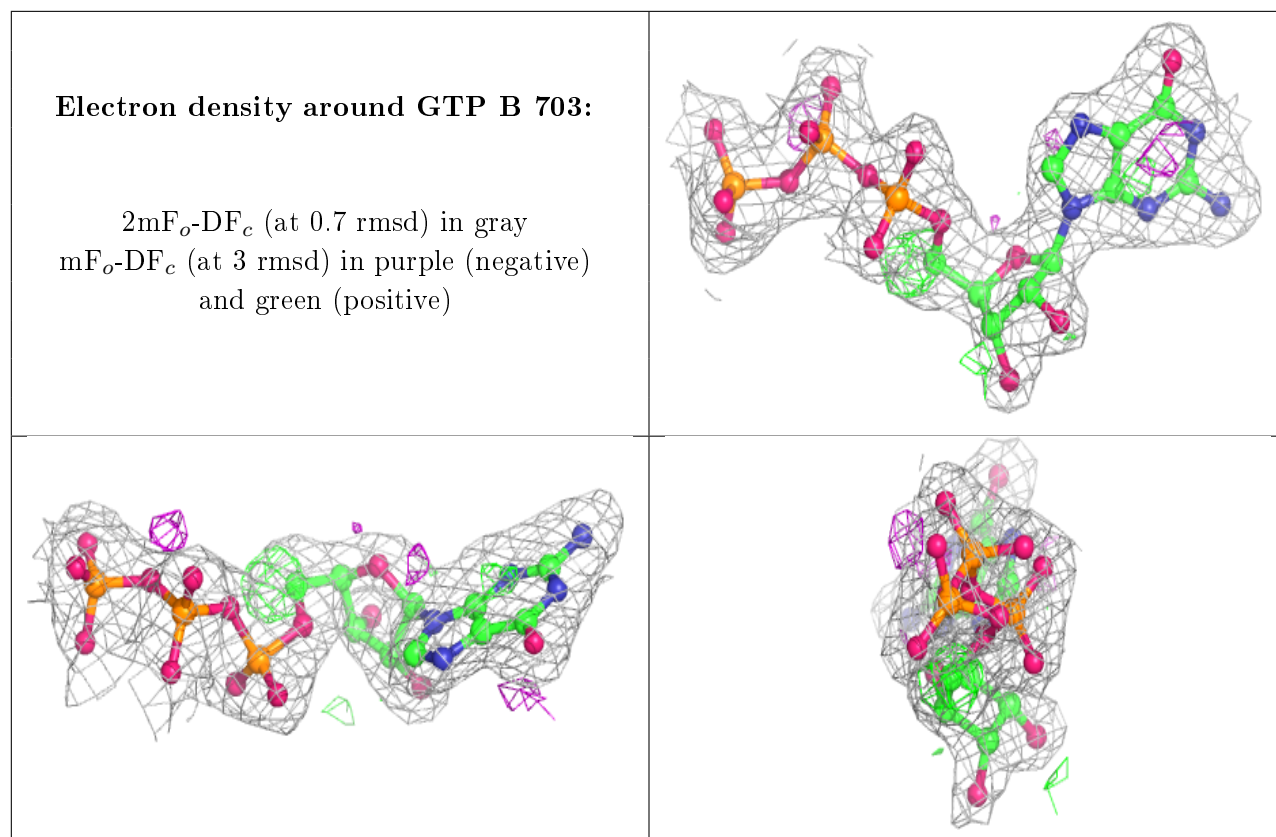
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	1PE	A	707	7/16	0.84	0.14	34,35,41,47	0
6	1PE	B	707	10/16	0.85	0.17	44,45,48,48	0
7	ETX	B	708	6/6	0.89	0.12	26,31,36,37	0
2	MN	B	705	1/1	0.91	0.06	40,40,40,40	1
5	NA	A	706	1/1	0.95	0.10	56,56,56,56	0
3	GTP	A	703	32/32	0.96	0.09	24,33,37,42	0
5	NA	B	706	1/1	0.96	0.09	55,55,55,55	0
3	GTP	B	703	32/32	0.97	0.09	26,30,34,38	0
4	SPV	B	704	10/10	0.97	0.10	19,27,32,32	0
2	MN	A	705	1/1	0.97	0.08	41,41,41,41	1
2	MN	A	702	1/1	0.99	0.06	31,31,31,31	0
4	SPV	A	704	10/10	0.99	0.07	27,27,28,30	0
2	MN	A	701	1/1	1.00	0.05	23,23,23,23	0
2	MN	B	702	1/1	1.00	0.09	23,23,23,23	0
2	MN	B	701	1/1	1.00	0.05	28,28,28,28	0

The following is a graphical depiction of the model fit to experimental electron density of all

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





6.5 Other polymers ⓘ

There are no such residues in this entry.