



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 05:11 AM UTC

PDB ID : 3DT7 / pdb_00003dt7
Title : The structure of rat cytosolic PEPCK in complex with beta-sulfoxyruvate and GTP
Authors : Sullivan, S.M.; Holyoak, T.
Deposited on : 2008-07-14
Resolution : 1.50 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

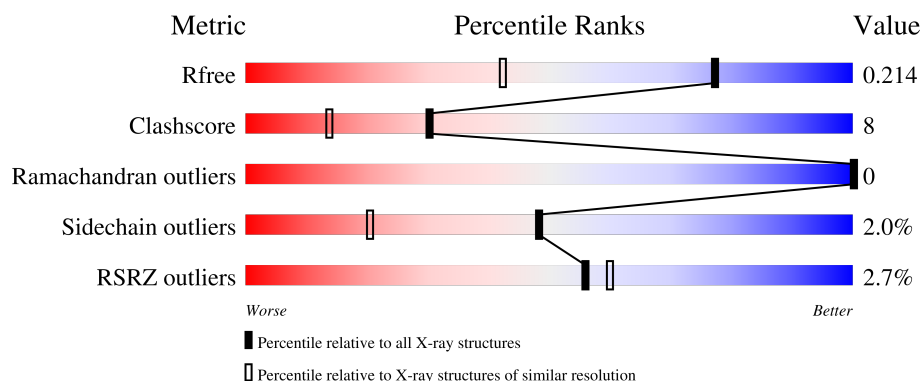
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.50 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	4037 (1.50-1.50)
Clashscore	190562	4235 (1.50-1.50)
Ramachandran outliers	187476	4153 (1.50-1.50)
Sidechain outliers	187428	4150 (1.50-1.50)
RSRZ outliers	180081	4039 (1.50-1.50)

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	

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 11144 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	624	Total	C	N	O	S	0	31	0
			5068	3255	858	920	35			
1	B	611	Total	C	N	O	S	0	27	0
			4964	3184	842	905	33			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP P07379
A	0	SER	-	expression tag	UNP P07379
B	-1	GLY	-	expression tag	UNP P07379
B	0	SER	-	expression tag	UNP P07379

- Molecule 2 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

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

- Molecule 3 is SODIUM ION (CCD ID: NA) (formula: Na).

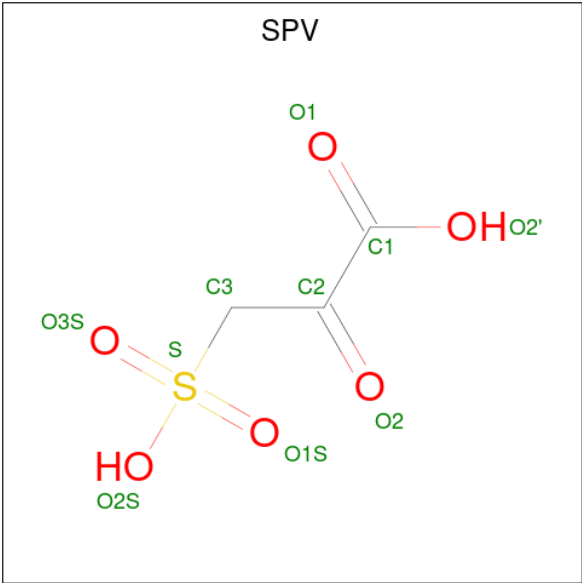
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Na	0	0
			1	1		
3	B	1	Total	Na	0	0
			1	1		

- Molecule 4 is GUANOSINE-5'-TRIPHOSPHATE (CCD ID: GTP) (formula: C₁₀H₁₆N₅O₁₄P₃).



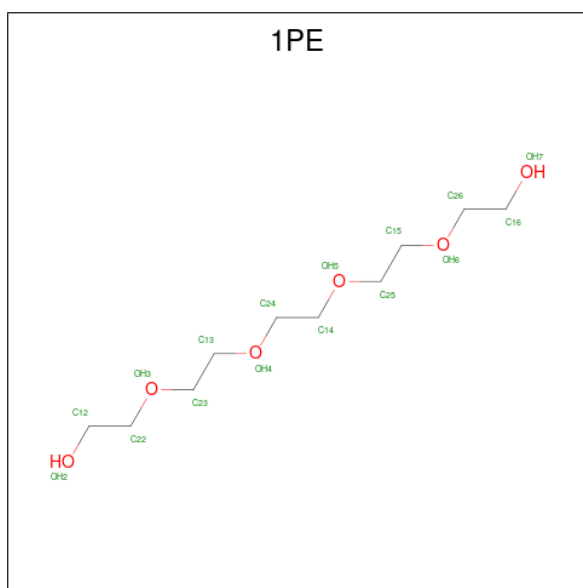
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total 32	C 10	N 5	O 14	P 3	0	0
4	B	1	Total 32	C 10	N 5	O 14	P 3	0	0

- Molecule 5 is SULFOPYRUVATE (CCD ID: SPV) (formula: C₃H₄O₆S).



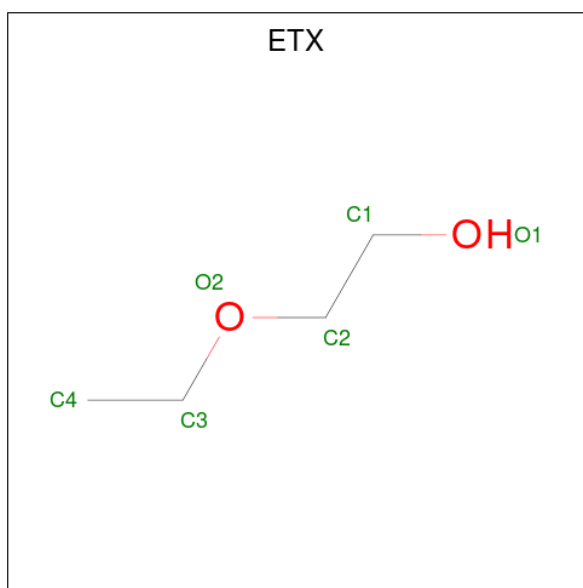
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	O	S	0	0
			10	3	6	1		
5	B	1	Total	C	O	S	0	0
			10	3	6	1		

- Molecule 6 is PENTAETHYLENE GLYCOL (CCD ID: 1PE) (formula: $C_{10}H_{22}O_6$).



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

- Molecule 7 is 2-ETHOXYETHANOL (CCD ID: ETX) (formula: $C_4H_{10}O_2$).



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

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	502	Total 502	O 502	0	0
8	B	497	Total 497	O 497	0	0

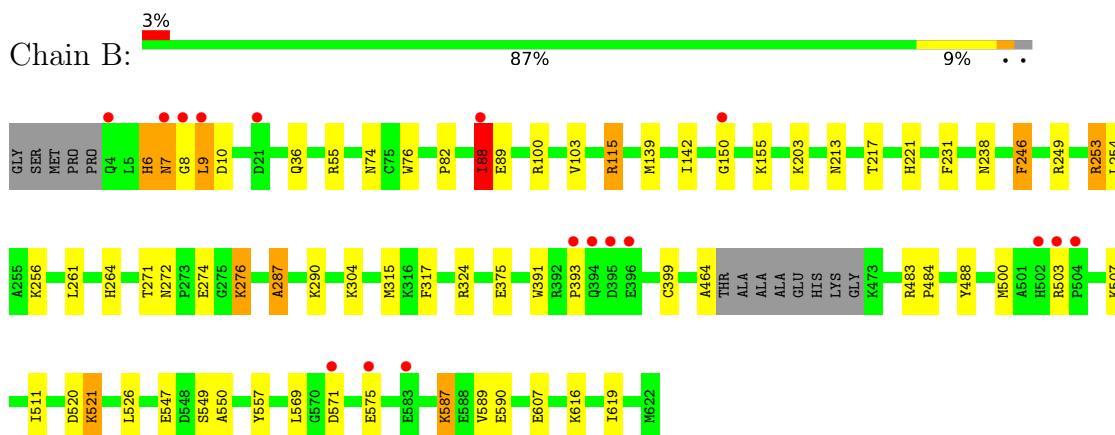
3 Residue-property plots [i](#)

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: Phosphoenolpyruvate carboxykinase, cytosolic [GTP]



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4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	62.02Å 119.50Å 87.28Å 90.00° 107.07° 90.00°	Depositor
Resolution (Å)	29.15 – 1.50 29.15 – 1.50	Depositor EDS
% Data completeness (in resolution range)	97.3 (29.15-1.50) 97.3 (29.15-1.50)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.90 (at 1.50Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.174 , 0.209 0.182 , 0.214	Depositor DCC
R_{free} test set	9440 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å ²)	13.9	Xtriage
Anisotropy	0.136	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 38.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	11144	wwPDB-VP
Average B, all atoms (Å ²)	8.0	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 67.64 % 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 4.9764e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 1PE, SPV, ETX, NA, GTP, MN

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	1.02	1/5290 (0.0%)	1.01	2/7157 (0.0%)
1	B	1.02	3/5172 (0.1%)	0.99	6/6996 (0.1%)
All	All	1.02	4/10462 (0.0%)	1.00	8/14153 (0.1%)

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	A	0	1
1	B	0	1
All	All	0	2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	139	MET	SD-CE	-5.81	1.65	1.79
1	B	142	ILE	CA-CB	5.70	1.59	1.53
1	B	287	ALA	CA-CB	5.19	1.61	1.53
1	A	444	VAL	CA-CB	5.12	1.61	1.55

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	150	GLY	N-CA-C	-13.38	95.69	115.63
1	B	88	ILE	CB-CA-C	-8.94	100.51	111.32
1	B	150	GLY	N-CA-C	7.28	120.95	112.50
1	B	7	ASN	N-CA-C	6.04	123.66	110.80
1	B	103	VAL	N-CA-C	-5.62	104.58	109.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	81	ASP	N-CA-C	-5.40	102.97	109.83
1	B	503	ARG	CA-C-N	5.27	126.09	119.98
1	B	503	ARG	C-N-CA	5.27	126.09	119.98

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	149	LEU	Peptide
1	B	6	HIS	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	5068	0	5099	91	0
1	B	4964	0	4977	79	0
2	A	3	0	0	0	0
2	B	2	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	32	0	12	4	0
4	B	32	0	12	1	0
5	A	10	0	2	0	0
5	B	10	0	2	0	0
6	A	10	0	13	6	0
7	A	6	0	10	1	0
7	B	6	0	10	3	0
8	A	502	0	0	23	1
8	B	497	0	0	21	0
All	All	11144	0	10137	171	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 8.

All (171) 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:7:ASN:OD1	1:B:8:GLY:HA3	1.19	1.30
1:B:115[A]:ARG:HB3	1:B:115[A]:ARG:NH1	1.49	1.28
1:A:324:ARG:NH2	8:A:3913:HOH:O	1.75	1.15
1:A:601[B]:ASN:OD1	8:A:3457:HOH:O	1.64	1.15
1:A:190[B]:ILE:HG21	1:A:246:PHE:CE1	1.83	1.14
1:B:511:ILE:H	7:B:3430:ETX:H11	1.10	1.11
1:B:115[A]:ARG:HH11	1:B:115[A]:ARG:CG	1.63	1.10
1:B:7:ASN:CG	1:B:8:GLY:HA3	1.76	1.09
1:A:153:LEU:HA	8:A:3817:HOH:O	1.52	1.08
1:B:115[A]:ARG:NH1	1:B:115[A]:ARG:CB	2.19	1.06
1:B:115[A]:ARG:HH11	1:B:115[A]:ARG:CB	1.68	1.06
1:A:93:VAL:HG12	1:A:115[B]:ARG:HG2	1.41	1.02
1:B:7:ASN:OD1	1:B:8:GLY:CA	2.06	1.02
1:B:375[A]:GLU:CG	8:B:4189:HOH:O	2.08	1.00
1:B:115[A]:ARG:HH11	1:B:115[A]:ARG:HG2	1.28	0.97
1:B:375[A]:GLU:HG3	8:B:4189:HOH:O	1.63	0.96
1:A:578[B]:PHE:HD1	8:A:3918:HOH:O	1.48	0.95
1:A:316:LYS:HZ3	6:A:1000:1PE:H231	1.33	0.93
1:B:256[B]:LYS:HE3	8:B:4240:HOH:O	1.68	0.92
1:B:590[A]:GLU:CD	8:B:3953:HOH:O	2.14	0.92
1:B:274:GLU:HG3	8:B:3983:HOH:O	1.71	0.90
1:A:112[B]:GLN:NE2	8:A:3776:HOH:O	2.03	0.90
1:A:317:PHE:HE2	1:A:500[B]:MET:HE3	1.39	0.86
1:B:511:ILE:N	7:B:3430:ETX:H11	1.91	0.86
1:A:268:LEU:HD13	1:A:270[B]:ILE:HG22	1.59	0.85
1:B:82[B]:PRO:HG3	8:B:4053:HOH:O	1.77	0.83
1:A:42:TYR:CE2	8:A:3743:HOH:O	2.30	0.83
1:A:42:TYR:HE2	8:A:3743:HOH:O	1.60	0.82
1:A:292[B]:ASN:OD1	4:A:800:GTP:H8	1.63	0.81
1:B:9:LEU:HG	1:B:36:GLN:NE2	1.98	0.79
1:A:292[B]:ASN:OD1	4:A:800:GTP:C8	2.37	0.78
1:A:224[B]:ASP:OD2	8:A:3906:HOH:O	2.02	0.77
1:A:317:PHE:CE2	1:A:500[B]:MET:HE3	2.20	0.77
1:A:253:ARG:NH1	8:A:3800:HOH:O	2.16	0.76
1:A:268:LEU:HD13	1:A:270[B]:ILE:CG2	2.16	0.75
1:A:574:VAL:O	1:A:578[A]:PHE:HD2	1.69	0.75
1:B:375[A]:GLU:HG2	8:B:4189:HOH:O	1.76	0.75
1:B:7:ASN:CG	1:B:8:GLY:CA	2.56	0.74
1:A:317:PHE:HE2	1:A:500[B]:MET:CE	2.01	0.73
1:A:0:SER:H	1:A:25:GLN:HE22	1.34	0.73
1:A:152:PRO:HG2	1:A:153:LEU:HD13	1.71	0.73
1:A:316:LYS:NZ	6:A:1000:1PE:H122	2.04	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:375[B]:GLU:OE2	8:B:4008:HOH:O	2.07	0.71
1:B:464:ALA:HB3	8:B:4049:HOH:O	1.89	0.70
1:A:153:LEU:CD2	8:A:3813:HOH:O	2.39	0.70
1:A:268:LEU:CD1	1:A:270[B]:ILE:CG2	2.69	0.70
1:A:152:PRO:HG2	1:A:153:LEU:CD1	2.22	0.69
1:A:342:LYS:HG2	8:A:3662:HOH:O	1.92	0.68
1:A:413:CYS:SG	1:A:415[B]:ILE:HD13	2.34	0.68
1:A:224[B]:ASP:OD1	1:A:253:ARG:NH2	2.27	0.68
1:A:190[B]:ILE:CG2	1:A:246:PHE:CE1	2.71	0.66
1:B:317:PHE:CE2	1:B:500[B]:MET:HE3	2.31	0.66
1:A:153:LEU:HD22	8:A:3813:HOH:O	1.95	0.66
1:B:115[A]:ARG:NH1	1:B:115[A]:ARG:HG2	2.03	0.66
1:A:316:LYS:HZ1	6:A:1000:1PE:H122	1.61	0.65
1:A:151:SER:OG	1:A:153:LEU:HD22	1.97	0.64
1:B:7:ASN:HD21	1:B:10:ASP:H	1.44	0.64
1:B:256[B]:LYS:HE2	8:B:4218:HOH:O	1.96	0.64
1:B:317:PHE:HE2	1:B:500[B]:MET:HE3	1.64	0.63
1:A:578[B]:PHE:CD1	8:A:3918:HOH:O	2.33	0.63
1:A:60[B]:MET:HE2	1:A:66:ILE:HG23	1.79	0.63
1:A:520:ASP:OD1	1:A:524:LYS:HG2	1.99	0.63
1:B:547[A]:GLU:HG2	1:B:549:SER:HB3	1.80	0.62
1:B:7:ASN:ND2	1:B:10:ASP:H	1.96	0.62
1:B:607:GLU:HG3	8:B:4007:HOH:O	1.98	0.62
1:A:74:ASN:ND2	1:A:76:TRP:HE1	1.98	0.61
1:B:55[A]:ARG:NH2	8:B:4227:HOH:O	2.32	0.61
1:B:6:HIS:O	1:B:7:ASN:HB2	2.01	0.61
1:A:7:ASN:HB3	8:A:3814:HOH:O	2.00	0.61
1:B:115[A]:ARG:HD3	8:B:3913:HOH:O	1.99	0.61
1:B:587:LYS:O	1:B:587:LYS:HD3	2.00	0.60
7:B:3430:ETX:H32	8:B:4330:HOH:O	2.02	0.59
1:A:292[B]:ASN:CG	4:A:800:GTP:C8	2.80	0.59
1:B:520:ASP:HB3	1:B:526[B]:LEU:HD11	1.84	0.59
1:A:293:LEU:HD12	1:A:296:MET:HE3	1.83	0.59
1:A:268:LEU:CD1	1:A:270[B]:ILE:HG23	2.32	0.59
1:A:317:PHE:CE2	1:A:500[B]:MET:CE	2.84	0.59
1:B:155:LYS:HE2	1:B:254:LEU:CD1	2.33	0.59
1:B:587:LYS:HD3	1:B:587:LYS:C	2.28	0.59
1:B:203:LYS:HE3	8:B:3957:HOH:O	2.02	0.58
1:A:601[B]:ASN:OD1	1:A:602:ALA:N	2.30	0.58
1:B:115[A]:ARG:CB	1:B:115[A]:ARG:CZ	2.81	0.58
1:A:153:LEU:HD13	1:A:153:LEU:N	2.21	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:105:ILE:O	1:A:601[B]:ASN:ND2	2.39	0.56
1:A:573:ASN:ND2	1:A:576:GLU:HB2	2.20	0.56
1:B:115[A]:ARG:HB3	1:B:115[A]:ARG:CZ	2.30	0.55
1:B:272:ASN:HD21	1:B:276:LYS:NZ	2.05	0.54
1:B:88:ILE:H	1:B:238:ASN:HD21	1.54	0.54
1:B:89:GLU:OE1	1:B:483[B]:ARG:NH1	2.40	0.54
1:B:74:ASN:ND2	1:B:76:TRP:HE1	2.05	0.54
1:A:574:VAL:O	1:A:578[A]:PHE:CD2	2.56	0.54
8:A:3836:HOH:O	1:B:276:LYS:HE2	2.08	0.54
7:A:3430:ETX:H31	8:A:3929:HOH:O	2.07	0.53
1:B:589:VAL:HG11	1:B:619[A]:ILE:HD12	1.90	0.53
1:B:9:LEU:HG	1:B:36:GLN:CD	2.34	0.53
1:A:149:LEU:HD21	1:A:179[A]:SER:HB2	1.90	0.53
1:A:151:SER:HG	1:A:153:LEU:HD22	1.73	0.52
1:B:375[B]:GLU:HG2	8:B:4189:HOH:O	2.09	0.51
1:A:292[B]:ASN:OD1	4:A:800:GTP:H2'	2.10	0.51
1:A:316:LYS:NZ	6:A:1000:1PE:H242	2.26	0.51
1:A:107:LYS:O	1:B:507:LYS:NZ	2.44	0.50
1:A:153:LEU:HD23	1:A:326:ILE:HD12	1.93	0.50
1:A:149:LEU:HD21	1:A:179[B]:SER:HB3	1.92	0.50
1:B:155:LYS:HE2	1:B:254:LEU:HD13	1.94	0.50
1:A:435:GLY:CA	8:A:3702:HOH:O	2.60	0.50
1:B:88:ILE:HD11	1:B:217:THR:HG21	1.94	0.50
1:A:153:LEU:HD21	1:A:423[A]:PRO:HA	1.92	0.50
1:A:190[B]:ILE:HG21	1:A:246:PHE:CZ	2.41	0.50
1:B:317:PHE:CE2	1:B:500[B]:MET:CE	2.96	0.49
1:B:55[A]:ARG:HH11	1:B:55[A]:ARG:HG2	1.77	0.49
1:B:7:ASN:OD1	1:B:9:LEU:N	2.45	0.49
1:B:317:PHE:HE2	1:B:500[B]:MET:CE	2.25	0.49
1:A:316:LYS:HZ3	6:A:1000:1PE:H242	1.77	0.49
1:A:190[B]:ILE:CG2	1:A:246:PHE:CZ	2.95	0.49
1:B:253:ARG:HG2	1:B:488:TYR:HB3	1.94	0.48
1:A:415[B]:ILE:HD13	1:A:415[B]:ILE:H	1.78	0.48
1:A:292[B]:ASN:ND2	1:A:530:PHE:CE1	2.82	0.47
1:B:521:LYS:HD2	8:B:4020:HOH:O	2.15	0.47
1:A:607[A]:GLU:HG2	1:A:610:ARG:NH2	2.29	0.47
1:A:153:LEU:HD21	1:A:423[B]:PRO:HA	1.96	0.47
1:A:435:GLY:N	8:A:3702:HOH:O	2.48	0.47
1:B:391:TRP:CH2	1:B:393:PRO:HA	2.50	0.47
1:A:71[A]:LYS:NZ	8:A:3796:HOH:O	2.47	0.46
1:B:483[A]:ARG:HB3	1:B:484:PRO:HD3	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:483[B]:ARG:CD	8:B:3911:HOH:O	2.64	0.46
1:B:287:ALA:HA	4:B:800:GTP:O2A	2.15	0.46
1:A:483:ARG:HB3	1:A:484:PRO:HD3	1.98	0.46
1:B:271:THR:HB	1:B:304:LYS:HD3	1.98	0.46
1:A:97[A]:GLN:HE21	1:A:97[A]:GLN:HB2	1.56	0.45
1:A:268:LEU:C	1:A:268:LEU:HD12	2.41	0.45
1:A:342:LYS:HE3	8:A:3662:HOH:O	2.15	0.45
1:A:56:LEU:O	1:A:60[B]:MET:HG3	2.17	0.45
1:A:213:ASN:C	1:A:213:ASN:HD22	2.24	0.45
1:A:589[B]:VAL:HG11	1:A:615:LEU:HG	1.99	0.45
1:B:89:GLU:CD	1:B:483[B]:ARG:HH22	2.24	0.45
1:A:159:GLU:HA	1:A:192:CYS:HB2	1.99	0.44
1:A:520:ASP:HB2	8:A:3740:HOH:O	2.18	0.44
1:A:148:PRO:HD2	8:A:3813:HOH:O	2.17	0.44
1:B:616[B]:LYS:HD2	8:B:4329:HOH:O	2.15	0.44
1:B:590[A]:GLU:CG	8:B:3953:HOH:O	2.57	0.44
1:B:315:MET:HA	1:B:324:ARG:O	2.18	0.44
1:A:292[B]:ASN:ND2	1:A:530:PHE:HE1	2.16	0.44
1:A:15:VAL:HG21	1:A:20:LEU:HG	2.01	0.43
1:B:203:LYS:HD2	8:B:4086:HOH:O	2.18	0.43
1:A:45[B]:ILE:CD1	1:A:141:VAL:HG21	2.48	0.43
1:B:272:ASN:HD21	1:B:276:LYS:HZ1	1.66	0.43
1:A:502:HIS:CD2	8:A:3673:HOH:O	2.71	0.42
1:B:7:ASN:HD22	1:B:10:ASP:HB3	1.83	0.42
1:A:477:HIS:HE2	1:A:591:GLU:CD	2.27	0.42
1:B:246:PHE:C	1:B:246:PHE:CD1	2.98	0.41
1:B:261:LEU:HD11	1:B:500[B]:MET:SD	2.60	0.41
1:A:190[B]:ILE:HD11	1:A:228:ILE:HG13	2.01	0.41
1:A:266:LEU:C	1:A:266:LEU:HD12	2.45	0.41
1:A:589[B]:VAL:HG11	1:A:619:ILE:HD12	2.02	0.41
1:A:56:LEU:HD11	1:A:164:PRO:HB3	2.02	0.41
1:B:217:THR:HA	1:B:231:PHE:O	2.20	0.41
1:A:45[B]:ILE:HD13	1:A:141:VAL:CG2	2.50	0.41
1:B:391:TRP:CD2	1:B:399:CYS:HB3	2.55	0.41
1:B:569:LEU:HD23	1:B:569:LEU:HA	1.81	0.41
1:B:213:ASN:HD22	1:B:213:ASN:C	2.29	0.41
1:B:221:HIS:CE1	1:B:249:ARG:HD2	2.56	0.41
1:B:264:HIS:CD2	1:B:290:LYS:HE3	2.56	0.41
1:B:520:ASP:HB3	1:B:526[B]:LEU:CD1	2.50	0.41
1:A:316:LYS:NZ	6:A:1000:1PE:H231	2.18	0.41
1:A:0:SER:N	1:A:25:GLN:HE22	2.12	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:550:ALA:HB1	1:B:557:TYR:HB3	2.02	0.40
1:A:389:LYS:HE2	1:A:389:LYS:HA	2.03	0.40
1:A:95:ILE:HD12	1:A:95:ILE:N	2.36	0.40
1:A:413:CYS:SG	1:A:415[B]:ILE:CD1	3.09	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:A:3626:HOH:O	8:A:3743:HOH:O[2_645]	2.14	0.06

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	653/624 (105%)	631 (97%)	22 (3%)	0	100	100
1	B	634/624 (102%)	616 (97%)	18 (3%)	0	100	100
All	All	1287/1248 (103%)	1247 (97%)	40 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	551/520 (106%)	539 (98%)	12 (2%)	45	17
1	B	539/520 (104%)	528 (98%)	11 (2%)	48	20
All	All	1090/1040 (105%)	1067 (98%)	23 (2%)	48	19

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	GLU
1	A	97[A]	GLN
1	A	97[B]	GLN
1	A	107	LYS
1	A	153	LEU
1	A	213	ASN
1	A	326	ILE
1	A	394	GLN
1	A	471	LYS
1	A	521	LYS
1	A	577	LEU
1	A	587	LYS
1	B	9	LEU
1	B	88	ILE
1	B	115[A]	ARG
1	B	115[B]	ARG
1	B	246	PHE
1	B	253	ARG
1	B	276	LYS
1	B	521	LYS
1	B	571	ASP
1	B	575	GLU
1	B	587	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	25	GLN
1	A	74	ASN
1	A	213	ASN
1	A	320	GLN
1	A	322	ASN
1	A	352	GLN
1	A	515	ASN

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Mol	Chain	Res	Type
1	B	4	GLN
1	B	25	GLN
1	B	36	GLN
1	B	59	HIS
1	B	74	ASN
1	B	213	ASN
1	B	238	ASN
1	B	272	ASN
1	B	292	ASN
1	B	320	GLN
1	B	515	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 14 ligands modelled in this entry, 7 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
5	SPV	A	900	2	9,9,9	2.58	3 (33%)	12,13,13	1.83	1 (8%)
5	SPV	B	900	2	9,9,9	3.19	3 (33%)	12,13,13	1.20	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	1PE	A	1000	-	9,9,15	0.67	0	8,8,14	0.63	0
7	ETX	A	3430	-	5,5,5	0.58	0	4,4,4	0.57	0
7	ETX	B	3430	-	5,5,5	0.62	0	4,4,4	0.86	0
4	GTP	A	800	2	33,34,34	1.25	3 (9%)	50,54,54	1.73	12 (24%)
4	GTP	B	800	2	33,34,34	1.40	2 (6%)	50,54,54	1.44	8 (16%)

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	SPV	A	900	2	-	0/8/9/9	-
5	SPV	B	900	2	-	0/8/9/9	-
6	1PE	A	1000	-	-	3/7/7/13	-
7	ETX	A	3430	-	-	1/3/3/3	-
7	ETX	B	3430	-	-	0/3/3/3	-
4	GTP	A	800	2	-	2/22/38/38	0/3/3/3
4	GTP	B	800	2	-	1/22/38/38	0/3/3/3

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	900	SPV	C3-S	-6.78	1.71	1.78
5	B	900	SPV	C2-C1	-5.68	1.45	1.53
5	A	900	SPV	C2-C1	-5.14	1.45	1.53
5	A	900	SPV	C3-S	-4.89	1.73	1.78
4	B	800	GTP	PA-O3A	4.64	1.64	1.59
4	B	800	GTP	PB-O3A	3.70	1.63	1.59
5	B	900	SPV	O1-C1	2.64	1.29	1.22
5	A	900	SPV	O1-C1	2.62	1.29	1.22
4	A	800	GTP	PG-O2G	-2.60	1.45	1.54
4	A	800	GTP	C4-N9	-2.30	1.32	1.38
4	A	800	GTP	O4'-C4'	-2.19	1.40	1.45

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	800	GTP	C5-C4-N3	-5.30	119.95	128.39
5	A	900	SPV	O1-C1-C2	-5.16	115.39	121.81
4	A	800	GTP	C2-N3-C4	4.09	119.34	112.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	800	GTP	C2-N3-C4	3.37	118.11	112.30
4	A	800	GTP	C2-N1-C6	-3.28	119.17	125.11
4	A	800	GTP	C4-C5-N7	-3.10	105.76	110.67
4	B	800	GTP	C5-C4-N3	-3.05	123.53	128.39
4	B	800	GTP	C6-C5-N7	2.86	135.50	130.29
4	A	800	GTP	C5-C6-N1	2.70	120.12	113.25
4	A	800	GTP	O6-C6-N1	-2.66	115.12	120.11
4	A	800	GTP	O2B-PB-O3B	2.57	114.23	107.27
4	A	800	GTP	O2G-PG-O3B	-2.56	96.06	104.64
4	A	800	GTP	C6-C5-N7	2.47	134.78	130.29
4	B	800	GTP	N2-C2-N1	2.45	121.94	116.76
4	A	800	GTP	C5-C4-N9	2.23	109.65	105.66
4	A	800	GTP	N9-C4-N3	2.21	130.37	125.95
4	B	800	GTP	N1-C2-N3	-2.10	119.48	123.32
4	B	800	GTP	O2B-PB-O3B	2.06	112.85	107.27
4	B	800	GTP	N9-C8-N7	-2.06	109.58	113.40
4	B	800	GTP	O4'-C1'-C2'	-2.04	102.25	106.62
4	A	800	GTP	N2-C2-N1	2.03	121.04	116.76

There are no chirality outliers.

All (7) torsion outliers are listed below:

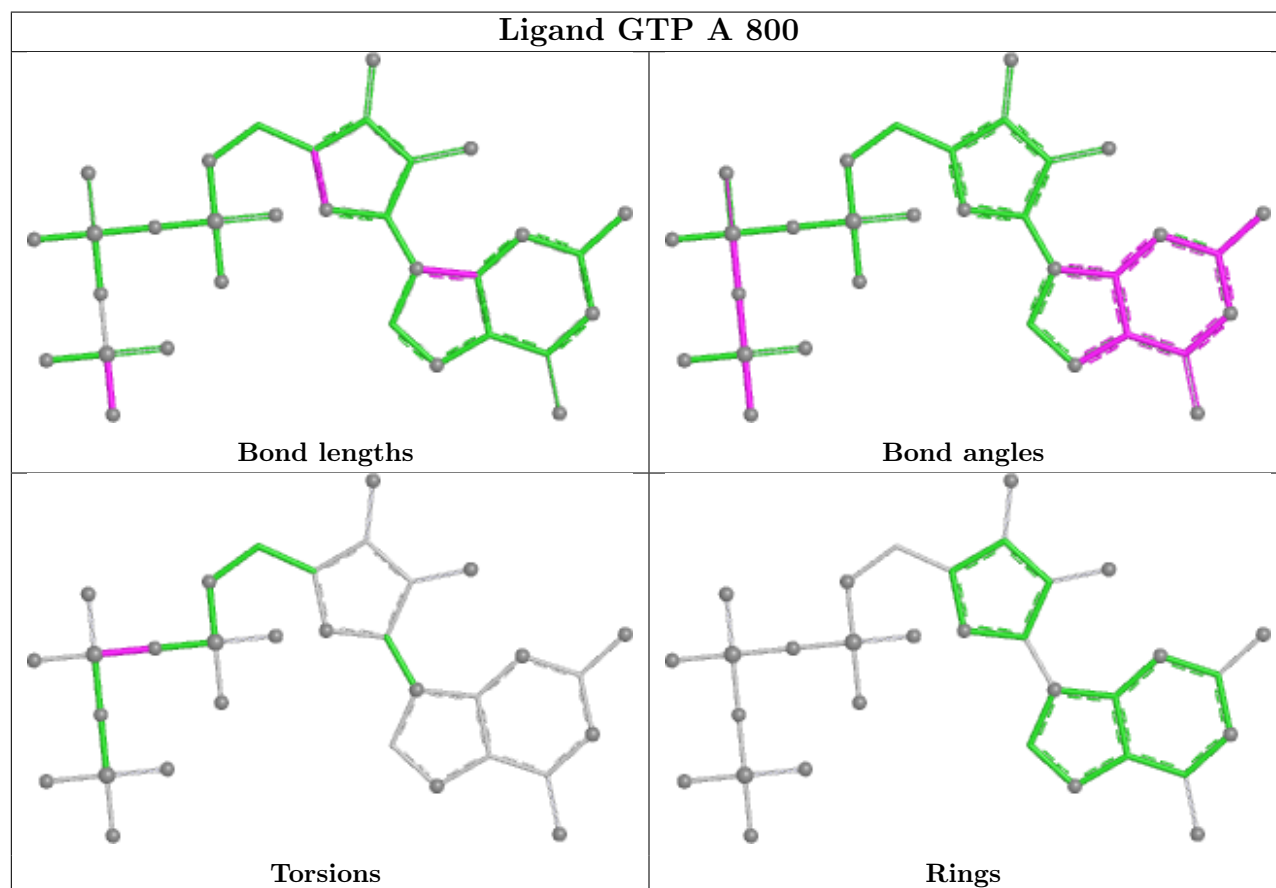
Mol	Chain	Res	Type	Atoms
4	B	800	GTP	C5'-O5'-PA-O1A
6	A	1000	1PE	OH5-C14-C24-OH4
7	A	3430	ETX	C4-C3-O2-C2
4	A	800	GTP	PA-O3A-PB-O2B
4	A	800	GTP	PA-O3A-PB-O1B
6	A	1000	1PE	OH4-C13-C23-OH3
6	A	1000	1PE	C23-C13-OH4-C24

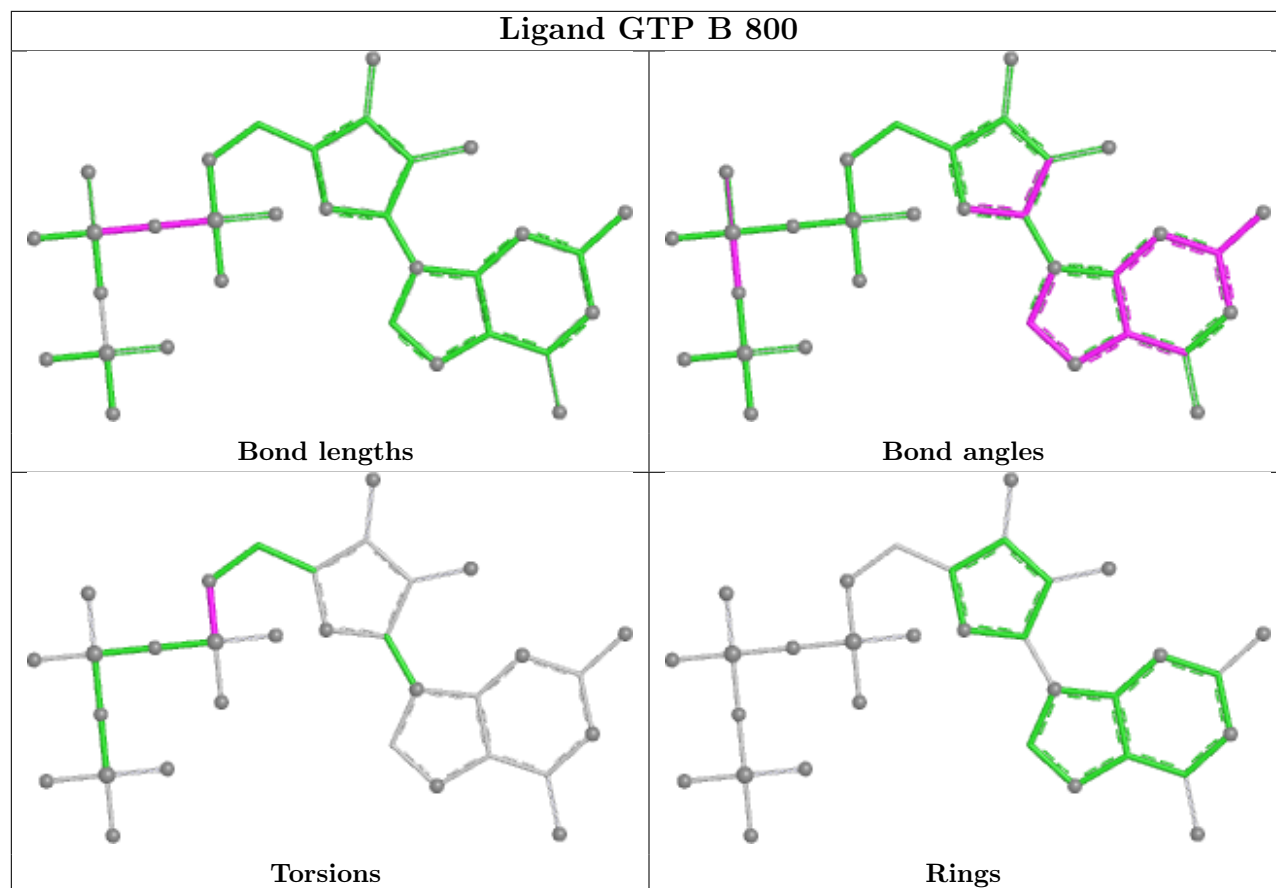
There are no ring outliers.

5 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	1000	1PE	6	0
7	A	3430	ETX	1	0
7	B	3430	ETX	3	0
4	A	800	GTP	4	0
4	B	800	GTP	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	624/624 (100%)	-0.11	16 (2%) 57 61	2, 6, 18, 35	42 (6%)
1	B	611/624 (97%)	-0.08	17 (2%) 55 59	2, 6, 19, 34	29 (4%)
All	All	1235/1248 (98%)	-0.10	33 (2%) 56 60	2, 6, 19, 35	71 (5%)

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	7	ASN	5.9
1	B	8	GLY	5.9
1	A	504	PRO	4.7
1	B	504	PRO	4.2
1	A	152	PRO	3.8
1	A	153	LEU	3.6
1	B	396	GLU	3.5
1	B	575	GLU	3.4
1	A	150	GLY	3.3
1	A	149	LEU	3.0
1	B	150	GLY	2.9
1	A	394	GLN	2.8
1	A	575	GLU	2.8
1	A	396	GLU	2.6
1	B	394	GLN	2.6
1	B	571	ASP	2.6
1	A	503	ARG	2.6
1	B	88	ILE	2.6
1	B	393	PRO	2.6
1	A	55	ARG	2.6
1	A	393	PRO	2.5
1	A	472	GLY	2.4
1	B	4	GLN	2.4
1	B	395	ASP	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	583	GLU	2.3
1	B	9	LEU	2.2
1	B	502	HIS	2.2
1	A	186	ASP	2.2
1	A	1	MET	2.1
1	B	21	ASP	2.1
1	A	522	ASN	2.1
1	B	503	ARG	2.0
1	A	120	GLU	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](#)

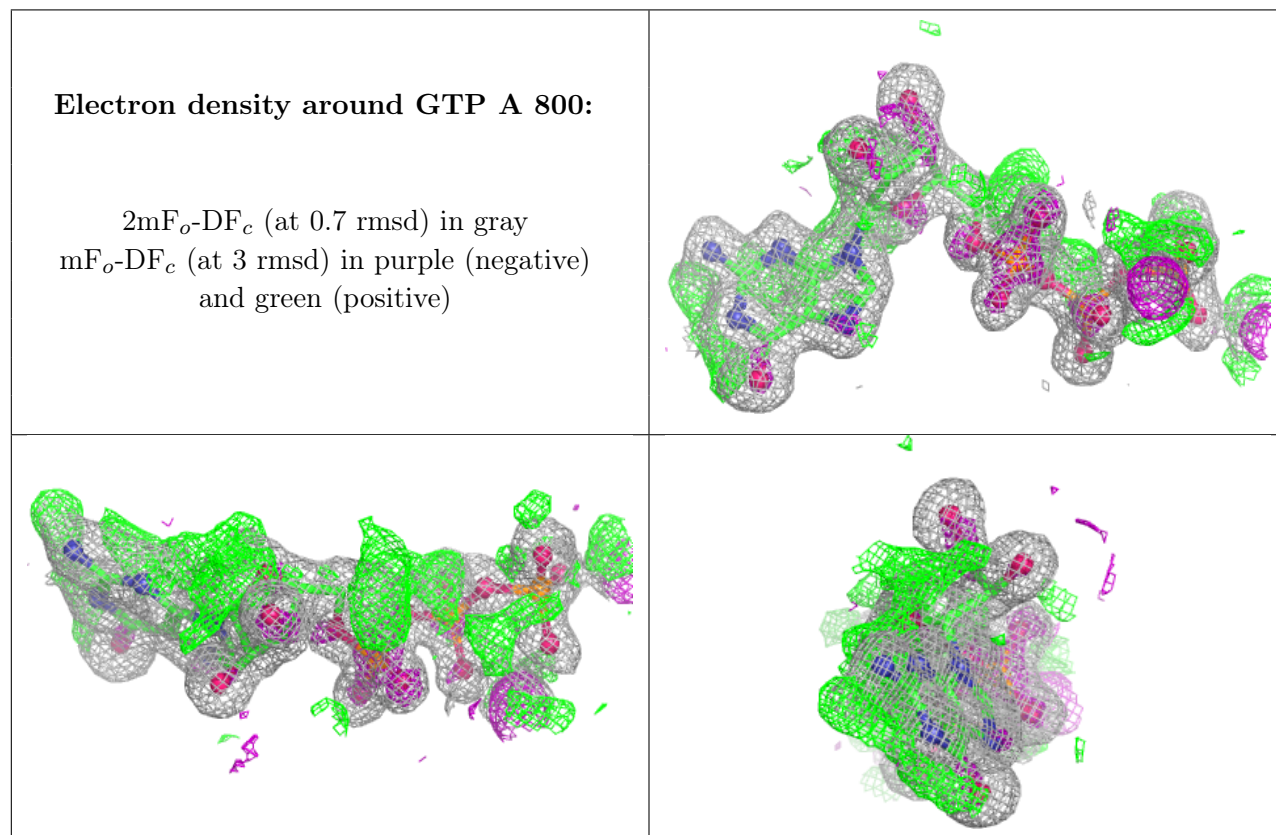
There are no oligosaccharides in this entry.

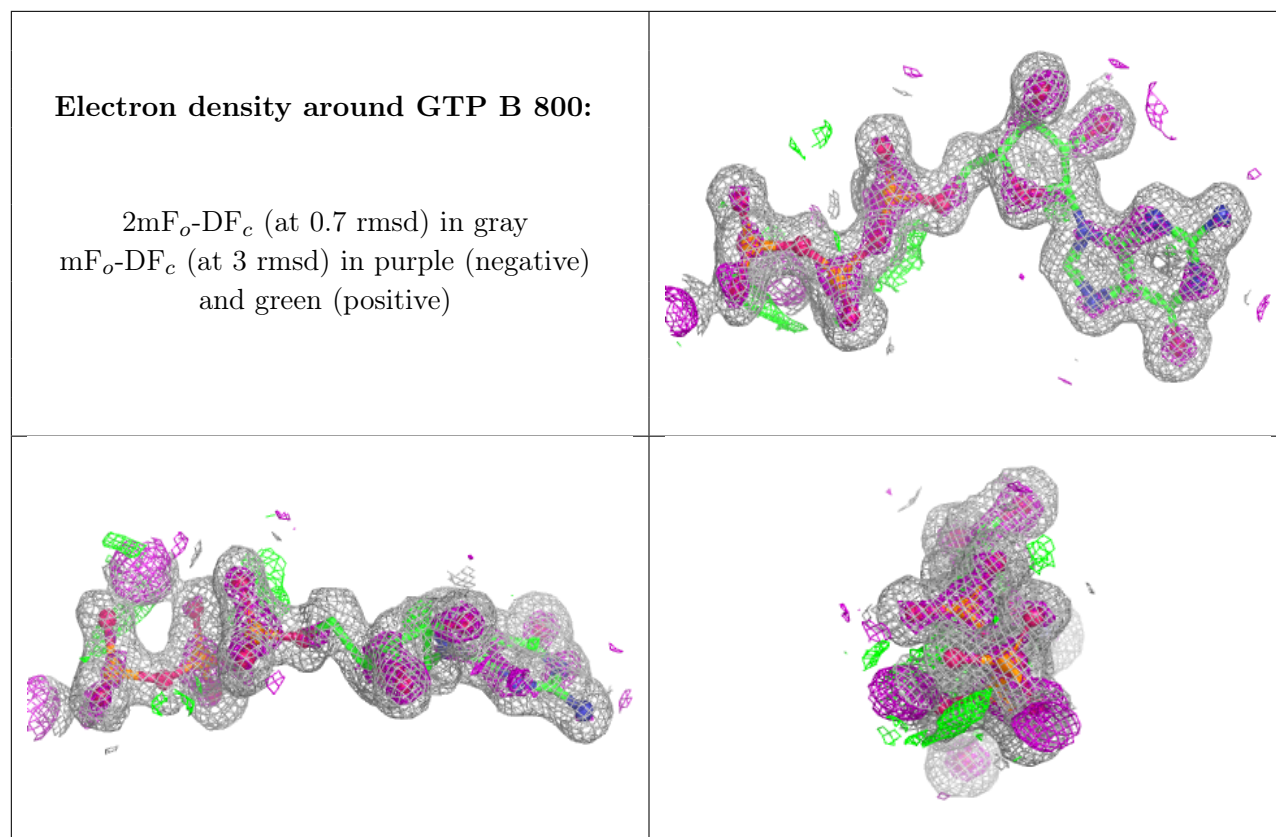
6.4 Ligands [i](#)

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($\text{\AA}^2$)	Q<0.9
7	ETX	B	3430	6/6	0.80	0.19	4,16,18,25	0
6	1PE	A	1000	10/16	0.82	0.11	13,19,24,26	0
7	ETX	A	3430	6/6	0.94	0.15	6,14,21,23	0
3	NA	A	1100	1/1	0.95	0.12	22,22,22,22	0
4	GTP	A	800	32/32	0.96	0.07	2,2,6,7	32
3	NA	B	1100	1/1	0.98	0.13	16,16,16,16	0
4	GTP	B	800	32/32	0.98	0.05	2,4,6,7	0
5	SPV	A	900	10/10	0.98	0.07	2,4,7,7	0
2	MN	A	700	1/1	0.99	0.09	3,3,3,3	0
5	SPV	B	900	10/10	0.99	0.05	2,3,5,6	0
2	MN	A	703	1/1	0.99	0.09	2,2,2,2	1
2	MN	B	700	1/1	0.99	0.07	4,4,4,4	0
2	MN	B	701	1/1	0.99	0.11	4,4,4,4	0
2	MN	A	701	1/1	1.00	0.13	2,2,2,2	1

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





6.5 Other polymers [i](#)

There are no such residues in this entry.