



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 11:01 PM UTC

PDB ID : 1KHE / pdb_00001khe
Title : PEPCK complex with nonhydrolyzable GTP analog, MAD data
Authors : Dunten, P.; Belunis, C.; Crowther, R.; Hollfelder, K.; Kammlott, U.; Levin, W.; Michel, H.; Ramsey, G.B.; Swain, A.; Weber, D.; Wertheimer, S.J.
Deposited on : 2001-11-29
Resolution : 2.40 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

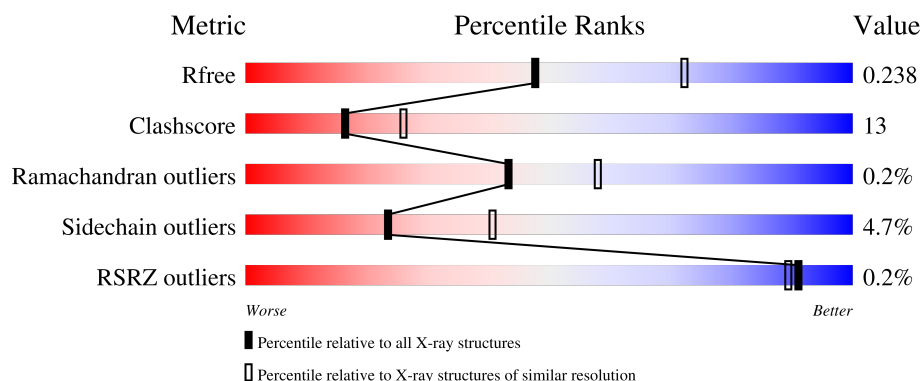
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	625	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4904 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	603	Total	C	N	O	S	Se	0	1	0
			4725	3015	807	867	14	22			

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	cloning artifact	UNP P35558
A	-1	GLU	-	cloning artifact	UNP P35558
A	0	LEU	-	cloning artifact	UNP P35558
A	1	MSE	MET	modified residue	UNP P35558
A	60	MSE	MET	modified residue	UNP P35558
A	117	MSE	MET	modified residue	UNP P35558
A	134	MSE	MET	modified residue	UNP P35558
A	139	MSE	MET	modified residue	UNP P35558
A	146	MSE	MET	modified residue	UNP P35558
A	170	MSE	MET	modified residue	UNP P35558
A	173	MSE	MET	modified residue	UNP P35558
A	176	MSE	MET	modified residue	UNP P35558
A	250	MSE	MET	modified residue	UNP P35558
A	265	MSE	MET	modified residue	UNP P35558
A	267	VAL	ILE	variant	UNP P35558
A	295	MSE	MET	modified residue	UNP P35558
A	296	MSE	MET	modified residue	UNP P35558
A	315	MSE	MET	modified residue	UNP P35558
A	460	MSE	MET	modified residue	UNP P35558
A	476	MSE	MET	modified residue	UNP P35558
A	482	MSE	MET	modified residue	UNP P35558
A	500	MSE	MET	modified residue	UNP P35558
A	540	MSE	MET	modified residue	UNP P35558
A	574	MSE	MET	modified residue	UNP P35558
A	575	MSE	MET	modified residue	UNP P35558
A	586	ASP	GLU	variant	UNP P35558
A	597	VAL	GLU	variant	UNP P35558

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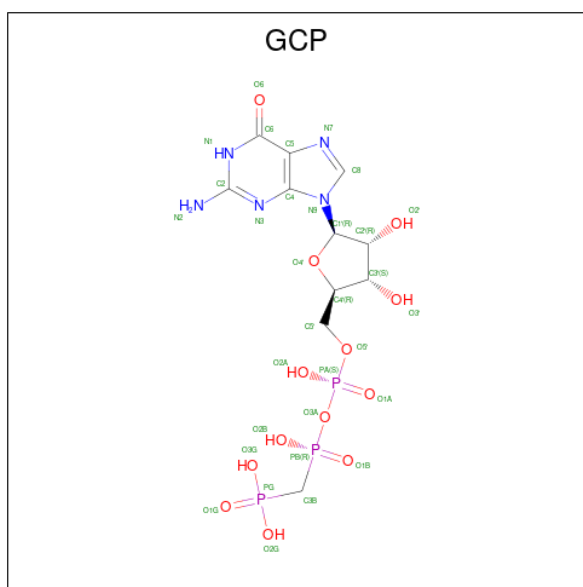
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Chain	Residue	Modelled	Actual	Comment	Reference
A	622	MSE	MET	modified residue	UNP P35558

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

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

- Molecule 3 is PHOSPHOMETHYLPHOSPHONIC ACID GUANYLATE ESTER (CCD ID: GCP) (formula: C₁₁H₁₈N₅O₁₃P₃).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C N O P 32 11 5 13 3	0	0

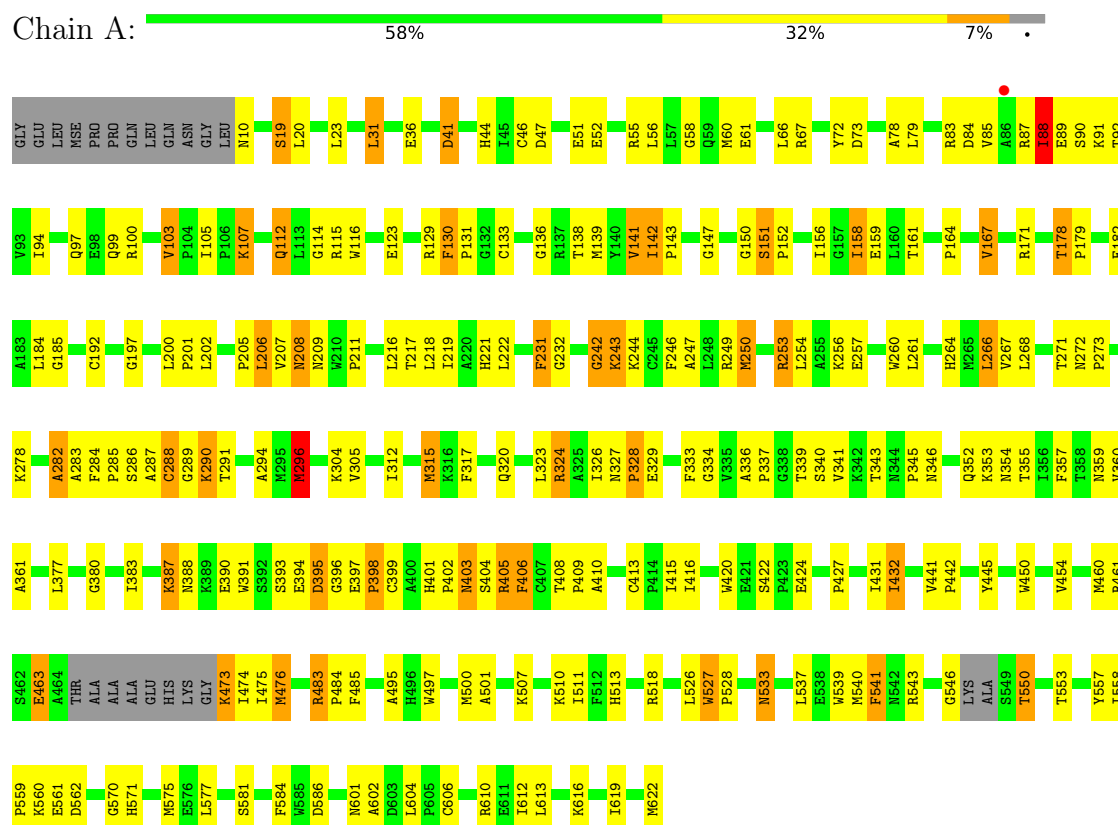
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	145	Total O 145 145	0	0

3 Residue-property plots

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

- Molecule 1: Phosphoenolpyruvate Carboxykinase, cytosolic (GTP)



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	45.39Å 60.54Å 61.83Å 88.36° 70.39° 72.37°	Depositor
Resolution (Å)	20.00 – 2.40 20.00 – 2.40	Depositor EDS
% Data completeness (in resolution range)	97.9 (20.00-2.40) 97.8 (20.00-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.02	Depositor
$\langle I/\sigma(I) \rangle$ ¹	15.12 (at 2.41Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.170 , 0.246 0.169 , 0.238	Depositor DCC
R_{free} test set	1125 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	14.0	Xtriage
Anisotropy	0.164	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 28.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.009 for -h,-l,-k	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4904	wwPDB-VP
Average B, all atoms (Å ²)	12.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.02% 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: MN, GCP

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	1.89	88/4829 (1.8%)	1.69	62/6504 (1.0%)

All (88) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	139	MSE	SE-CE	-18.43	1.40	1.95
1	A	296	MSE	SE-CE	-17.84	1.42	1.95
1	A	250	MSE	SE-CE	11.05	2.28	1.95
1	A	305	VAL	C-O	-9.50	1.14	1.24
1	A	327	ASN	C-O	-8.98	1.13	1.24
1	A	206	LEU	N-CA	-8.54	1.35	1.46
1	A	142	ILE	CA-CB	-7.53	1.44	1.54
1	A	242	GLY	C-O	-7.47	1.15	1.24
1	A	339	THR	C-O	7.42	1.32	1.23
1	A	460	MSE	SE-CE	-7.28	1.73	1.95
1	A	107	LYS	C-O	-7.24	1.15	1.24
1	A	340	SER	CA-C	-7.17	1.44	1.52
1	A	352	GLN	C-O	-7.11	1.15	1.24
1	A	540	MSE	SE-CE	-7.11	1.74	1.95
1	A	283	ALA	CA-CB	6.99	1.62	1.53
1	A	247	ALA	CA-CB	6.89	1.65	1.53
1	A	282	ALA	CA-CB	6.85	1.67	1.53
1	A	360	VAL	CA-CB	-6.81	1.42	1.55
1	A	161	THR	CA-CB	6.77	1.65	1.53
1	A	290	LYS	N-CA	-6.75	1.38	1.46
1	A	355	THR	N-CA	-6.71	1.37	1.46
1	A	139	MSE	C-O	6.70	1.32	1.24
1	A	410	ALA	CA-CB	6.69	1.64	1.53
1	A	404	SER	CA-C	6.59	1.60	1.52
1	A	156	ILE	N-CA	-6.56	1.38	1.46
1	A	291	THR	CA-CB	6.55	1.63	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	511	ILE	CA-CB	-6.54	1.46	1.54
1	A	158	ILE	CA-CB	-6.50	1.46	1.54
1	A	510	LYS	CA-C	6.48	1.61	1.52
1	A	256	LYS	C-O	-6.43	1.16	1.24
1	A	454	VAL	CB-CG2	6.39	1.73	1.52
1	A	105	ILE	CA-C	-6.36	1.48	1.53
1	A	612	ILE	C-O	-6.31	1.17	1.24
1	A	219	ILE	CA-C	-6.25	1.44	1.52
1	A	284	PHE	CA-C	-6.17	1.46	1.52
1	A	602	ALA	CA-CB	6.16	1.64	1.53
1	A	422	SER	CA-C	6.16	1.59	1.52
1	A	476	MSE	C-O	-6.15	1.16	1.23
1	A	178	THR	CA-C	6.12	1.59	1.52
1	A	586	ASP	CA-C	-6.03	1.45	1.52
1	A	357	PHE	C-O	-5.99	1.16	1.23
1	A	116	TRP	C-O	5.95	1.30	1.23
1	A	253	ARG	CZ-NH1	5.92	1.41	1.32
1	A	294	ALA	CA-CB	-5.91	1.43	1.53
1	A	431	ILE	CA-CB	-5.89	1.47	1.54
1	A	341	VAL	CA-CB	5.86	1.61	1.54
1	A	527	TRP	C-O	-5.85	1.16	1.24
1	A	413	CYS	C-O	-5.80	1.16	1.24
1	A	83	ARG	CZ-NH1	5.80	1.40	1.32
1	A	501	ALA	CA-CB	-5.79	1.44	1.53
1	A	315	MSE	CA-C	-5.79	1.45	1.52
1	A	405	ARG	N-CA	-5.78	1.38	1.45
1	A	483	ARG	NE-CZ	5.73	1.39	1.33
1	A	66	LEU	CA-C	-5.69	1.45	1.52
1	A	405	ARG	CZ-NH1	5.67	1.40	1.32
1	A	539	TRP	N-CA	-5.61	1.39	1.46
1	A	211	PRO	CG-CD	5.59	1.69	1.50
1	A	94	ILE	CA-CB	-5.55	1.47	1.54
1	A	78	ALA	CA-CB	-5.54	1.45	1.53
1	A	463	GLU	CA-C	-5.53	1.46	1.52
1	A	445	TYR	CA-C	-5.49	1.45	1.52
1	A	420	TRP	C-O	-5.41	1.17	1.24
1	A	495	ALA	CA-CB	5.37	1.61	1.53
1	A	403	ASN	CA-C	5.36	1.59	1.52
1	A	288	CYS	C-O	5.35	1.31	1.24
1	A	61	GLU	N-CA	-5.33	1.40	1.46
1	A	361	ALA	CA-CB	5.30	1.62	1.53
1	A	222	LEU	CA-C	5.28	1.58	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	208	ASN	CA-C	-5.26	1.45	1.52
1	A	133	CYS	C-O	-5.25	1.17	1.24
1	A	47	ASP	CA-C	-5.23	1.45	1.52
1	A	167	VAL	CA-CB	5.21	1.60	1.54
1	A	267	VAL	CA-C	-5.20	1.46	1.52
1	A	408	THR	C-O	-5.19	1.18	1.24
1	A	92	THR	CA-C	-5.19	1.46	1.52
1	A	84	ASP	C-O	-5.18	1.17	1.24
1	A	99	GLN	CA-C	5.14	1.59	1.52
1	A	571	HIS	N-CA	-5.12	1.39	1.46
1	A	604	LEU	CA-C	5.12	1.58	1.52
1	A	533	ASN	CG-OD1	5.11	1.33	1.23
1	A	141	VAL	C-O	-5.10	1.18	1.24
1	A	216	LEU	CG-CD2	5.09	1.69	1.52
1	A	268	LEU	CA-C	-5.09	1.46	1.52
1	A	527	TRP	CA-C	-5.07	1.48	1.52
1	A	403	ASN	CG-ND2	5.03	1.43	1.33
1	A	218	LEU	CA-C	5.02	1.58	1.52
1	A	476	MSE	SE-CE	5.01	2.10	1.95
1	A	513	HIS	C-O	5.01	1.29	1.24

All (62) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	543	ARG	N-CA-C	-10.12	99.24	111.69
1	A	85	VAL	CB-CA-C	-9.88	98.49	110.84
1	A	586	ASP	N-CA-C	-7.87	102.64	111.14
1	A	85	VAL	N-CA-C	7.64	121.76	113.43
1	A	289	GLY	N-CA-C	7.47	127.00	115.72
1	A	550	THR	N-CA-C	7.47	121.71	110.14
1	A	333	PHE	CA-C-N	7.41	129.44	120.92
1	A	333	PHE	C-N-CA	7.41	129.44	120.92
1	A	405	ARG	NE-CZ-NH2	-7.09	112.81	119.20
1	A	58	GLY	N-CA-C	-7.04	104.33	112.50
1	A	343	THR	CA-C-N	-6.93	115.92	122.37
1	A	343	THR	C-N-CA	-6.93	115.92	122.37
1	A	103	VAL	CB-CA-C	-6.84	102.50	109.89
1	A	19	SER	N-CA-C	6.78	119.46	108.34
1	A	231	PHE	CA-C-N	-6.65	114.10	123.08
1	A	231	PHE	C-N-CA	-6.65	114.10	123.08
1	A	88	ILE	CB-CA-C	-6.52	101.44	110.96
1	A	359	ASN	CA-C-N	-6.41	115.41	122.27

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	359	ASN	C-N-CA	-6.41	115.41	122.27
1	A	406	PHE	N-CA-C	-6.28	99.92	109.85
1	A	243	LYS	N-CA-C	6.10	118.90	111.82
1	A	197	GLY	N-CA-C	6.05	124.52	114.48
1	A	158	ILE	CA-C-N	-6.01	114.70	123.00
1	A	158	ILE	C-N-CA	-6.01	114.70	123.00
1	A	540	MSE	CA-C-O	5.99	126.77	120.42
1	A	427	PRO	CB-CA-C	5.98	116.84	111.40
1	A	112	GLN	N-CA-C	-5.84	105.34	112.88
1	A	575	MSE	N-CA-C	-5.75	104.62	111.69
1	A	409	PRO	N-CD-CG	-5.65	94.72	103.20
1	A	327	ASN	N-CA-C	-5.60	100.53	109.04
1	A	432	ILE	N-CA-C	5.45	115.80	108.17
1	A	41	ASP	CA-C-N	-5.44	113.68	122.82
1	A	41	ASP	C-N-CA	-5.44	113.68	122.82
1	A	398	PRO	N-CA-C	-5.40	102.67	111.68
1	A	511	ILE	N-CA-C	5.37	116.31	108.53
1	A	380	GLY	CA-C-N	-5.37	116.17	122.93
1	A	380	GLY	C-N-CA	-5.37	116.17	122.93
1	A	253	ARG	NE-CZ-NH2	-5.34	114.39	119.20
1	A	284	PHE	N-CA-C	-5.34	99.79	108.82
1	A	476	MSE	CA-CB-CG	-5.29	103.53	114.10
1	A	484	PRO	CB-CA-C	-5.27	103.45	110.88
1	A	55	ARG	CA-C-N	-5.26	112.82	120.29
1	A	55	ARG	C-N-CA	-5.26	112.82	120.29
1	A	114	GLY	O-C-N	-5.25	117.45	123.23
1	A	55	ARG	N-CA-C	-5.24	105.65	111.36
1	A	405	ARG	N-CA-CB	-5.21	102.48	111.55
1	A	94	ILE	CA-C-N	-5.20	116.36	123.12
1	A	94	ILE	C-N-CA	-5.20	116.36	123.12
1	A	266	LEU	CA-C-N	-5.19	114.03	122.33
1	A	266	LEU	C-N-CA	-5.19	114.03	122.33
1	A	171	ARG	NE-CZ-NH1	-5.18	116.31	121.50
1	A	570	GLY	N-CA-C	-5.18	104.64	112.41
1	A	72	TYR	N-CA-C	5.13	117.69	110.14
1	A	461	ARG	NE-CZ-NH2	5.13	123.82	119.20
1	A	290	LYS	N-CA-C	-5.10	105.61	111.07
1	A	205	PRO	CA-C-O	5.08	127.45	121.56
1	A	395	ASP	N-CA-CB	5.07	117.69	110.44
1	A	130	PHE	N-CA-C	5.05	120.97	109.81
1	A	83	ARG	NH1-CZ-NH2	5.04	125.86	119.30
1	A	324	ARG	NE-CZ-NH1	-5.04	116.46	121.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	232	GLY	N-CA-C	5.00	120.67	114.37
1	A	546	GLY	CA-C-O	-5.00	110.30	120.80

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	4725	0	4667	122	0
2	A	2	0	0	0	0
3	A	32	0	14	2	0
4	A	145	0	0	2	0
All	All	4904	0	4681	123	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 13.

All (123) 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:A:250:MSE:CE	1:A:250:MSE:SE	2.28	1.31
1:A:497:TRP:HA	1:A:500[A]:MSE:HE2	1.50	0.94
1:A:90:SER:O	1:A:115:ARG:NH1	2.02	0.93
1:A:395:ASP:OD1	1:A:396:GLY:N	2.05	0.88
1:A:497:TRP:CA	1:A:500[A]:MSE:HE2	2.03	0.87
1:A:90:SER:C	1:A:115:ARG:HH11	1.83	0.86
1:A:497:TRP:CE3	1:A:500[A]:MSE:HE1	2.12	0.84
1:A:398:PRO:HD2	4:A:716:HOH:O	1.80	0.80
1:A:497:TRP:O	1:A:500[A]:MSE:HE2	1.84	0.78
1:A:497:TRP:O	1:A:500[A]:MSE:CE	2.32	0.76
1:A:90:SER:C	1:A:115:ARG:NH1	2.43	0.76
1:A:315:MSE:HA	1:A:324:ARG:O	1.87	0.75
1:A:619:ILE:O	1:A:622:MSE:HG2	1.87	0.73
1:A:473:LYS:O	1:A:474:ILE:HD13	1.88	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:346:ASN:ND2	1:A:415:ILE:HD13	2.05	0.71
1:A:613:LEU:HD12	1:A:616:LYS:HD2	1.73	0.70
1:A:142:ILE:HD11	1:A:167:VAL:HG22	1.75	0.68
1:A:497:TRP:C	1:A:500[A]:MSE:HE2	2.20	0.67
1:A:242:GLY:HA2	1:A:246:PHE:HB3	1.80	0.63
1:A:41:ASP:OD2	1:A:136:GLY:HA2	2.01	0.61
1:A:31:LEU:HD12	1:A:31:LEU:C	2.26	0.60
1:A:450:TRP:HZ2	1:A:500[A]:MSE:CE	2.15	0.60
1:A:264:HIS:CD2	1:A:290:LYS:HE3	2.35	0.60
1:A:377:LEU:HD11	1:A:383:ILE:HD11	1.83	0.60
1:A:424:GLU:H	1:A:424:GLU:CD	2.09	0.60
1:A:317:PHE:CE1	1:A:500[B]:MSE:HE3	2.37	0.59
1:A:19:SER:O	1:A:20:LEU:C	2.44	0.59
1:A:89:GLU:O	1:A:90:SER:C	2.46	0.58
1:A:90:SER:CA	1:A:115:ARG:HH11	2.17	0.57
1:A:67:ARG:HH21	1:A:79:LEU:HD12	1.71	0.56
1:A:178:THR:N	1:A:179:PRO:CD	2.68	0.56
1:A:497:TRP:O	1:A:500[A]:MSE:HE3	2.04	0.56
1:A:336:ALA:HB3	1:A:337:PRO:HD3	1.88	0.56
1:A:250:MSE:CE	1:A:250:MSE:HB2	2.37	0.55
1:A:272:ASN:HB2	1:A:273:PRO:CD	2.36	0.55
1:A:178:THR:HB	1:A:179:PRO:HD3	1.89	0.55
1:A:282:ALA:HB2	1:A:432:ILE:HB	1.88	0.54
1:A:67:ARG:HH22	1:A:388:ASN:ND2	2.05	0.54
1:A:23:LEU:CD1	1:A:31:LEU:HD23	2.39	0.53
1:A:200:LEU:HA	1:A:201:PRO:C	2.33	0.53
1:A:202:LEU:HD11	1:A:206:LEU:HD22	1.91	0.53
1:A:260:TRP:HB2	1:A:315:MSE:O	2.09	0.53
1:A:296:MSE:HE1	1:A:533:ASN:HB2	1.92	0.52
1:A:31:LEU:HD22	1:A:141:VAL:HG21	1.91	0.51
1:A:450:TRP:HZ2	1:A:500[A]:MSE:HE3	1.76	0.51
1:A:88:ILE:HD13	1:A:91:LYS:HD3	1.91	0.51
1:A:320:GLN:O	1:A:507:LYS:NZ	2.43	0.51
1:A:474:ILE:HG22	1:A:476:MSE:HG3	1.92	0.51
1:A:463:GLU:HA	1:A:475:ILE:HA	1.91	0.51
1:A:450:TRP:CZ2	1:A:500[A]:MSE:HE3	2.46	0.50
1:A:326:ILE:O	1:A:326:ILE:HG13	2.10	0.50
1:A:261:LEU:HD11	1:A:500[A]:MSE:SE	2.62	0.50
1:A:278:LYS:HG3	1:A:541:PHE:CE1	2.47	0.49
1:A:73:ASP:HB3	1:A:353:LYS:HG2	1.95	0.48
1:A:334:GLY:HA3	1:A:406:PHE:CZ	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:51:GLU:H	1:A:51:GLU:CD	2.20	0.48
1:A:353:LYS:HE2	1:A:353:LYS:HB2	1.68	0.48
1:A:441:VAL:HA	1:A:442:PRO:HD3	1.57	0.48
1:A:254:LEU:O	1:A:257:GLU:HB2	2.14	0.47
1:A:334:GLY:O	1:A:405:ARG:HA	2.13	0.47
1:A:345:PRO:HD2	4:A:761:HOH:O	2.15	0.47
1:A:129:ARG:HB3	1:A:231:PHE:CZ	2.50	0.47
1:A:221:HIS:CE1	1:A:249:ARG:HD2	2.50	0.47
1:A:184:LEU:O	1:A:185:GLY:C	2.56	0.47
1:A:473:LYS:C	1:A:474:ILE:HD13	2.39	0.47
1:A:558:ILE:HB	1:A:559:PRO:CD	2.45	0.47
1:A:56:LEU:HA	1:A:56:LEU:HD23	1.52	0.46
1:A:90:SER:O	1:A:115:ARG:HD3	2.15	0.46
1:A:150:GLY:O	1:A:151:SER:C	2.58	0.46
1:A:221:HIS:NE2	1:A:249:ARG:HD2	2.30	0.46
1:A:208:ASN:O	1:A:209:ASN:C	2.57	0.46
1:A:606:CYS:O	1:A:610:ARG:HG3	2.16	0.46
1:A:527:TRP:CD2	1:A:528:PRO:HD2	2.51	0.45
1:A:550:THR:HB	1:A:557:TYR:HB3	1.98	0.45
1:A:130:PHE:N	1:A:131:PRO:CD	2.79	0.45
1:A:282:ALA:CB	1:A:432:ILE:HB	2.46	0.45
1:A:553:THR:HG21	1:A:558:ILE:HD13	1.99	0.45
1:A:159:GLU:HA	1:A:192:CYS:HB2	1.98	0.45
1:A:207:VAL:O	1:A:208:ASN:C	2.60	0.45
1:A:271:THR:HB	1:A:304:LYS:HB3	2.00	0.44
1:A:518:ARG:HG2	1:A:526:LEU:HD12	2.00	0.44
1:A:112:GLN:HB3	1:A:483:ARG:HH21	1.82	0.44
1:A:244:LYS:HD2	1:A:485:PHE:CE2	2.53	0.44
1:A:286:SER:O	1:A:287:ALA:HB3	2.18	0.44
1:A:387:LYS:O	1:A:388:ASN:CB	2.65	0.44
1:A:424:GLU:CD	1:A:424:GLU:N	2.74	0.44
1:A:20:LEU:HD23	1:A:20:LEU:HA	1.76	0.43
1:A:243:LYS:NZ	1:A:329:GLU:OE2	2.41	0.43
1:A:107:LYS:NZ	1:A:601:ASN:OD1	2.45	0.43
1:A:67:ARG:HH22	1:A:388:ASN:HD21	1.66	0.43
1:A:88:ILE:HD12	1:A:88:ILE:H	1.84	0.42
1:A:147:GLY:HA2	1:A:328:PRO:HA	2.01	0.42
1:A:323:LEU:HD12	1:A:323:LEU:HA	1.77	0.42
1:A:390:GLU:HG3	1:A:391:TRP:N	2.32	0.42
3:A:703:GCP:H3B2	3:A:703:GCP:O2A	2.18	0.42
1:A:67:ARG:HH22	1:A:388:ASN:CG	2.26	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:143:PRO:HA	1:A:158:ILE:HD13	2.01	0.42
1:A:60:MSE:SE	1:A:164:PRO:HB2	2.70	0.42
1:A:266:LEU:C	1:A:266:LEU:HD12	2.44	0.42
1:A:395:ASP:CG	1:A:396:GLY:N	2.75	0.42
1:A:217:THR:HA	1:A:231:PHE:O	2.19	0.42
1:A:272:ASN:CB	1:A:273:PRO:CD	2.96	0.42
1:A:391:TRP:CD2	1:A:399:CYS:HB3	2.54	0.42
1:A:129:ARG:HH11	1:A:129:ARG:HG3	1.85	0.42
1:A:151:SER:HA	1:A:152:PRO:HD3	1.77	0.42
1:A:130:PHE:N	1:A:131:PRO:HD3	2.35	0.41
1:A:401:HIS:CE1	1:A:402:PRO:HD2	2.55	0.41
1:A:246:PHE:HD1	1:A:246:PHE:C	2.28	0.41
1:A:353:LYS:O	1:A:354:ASN:C	2.63	0.41
1:A:474:ILE:HD13	1:A:474:ILE:HA	1.72	0.41
1:A:285:PRO:O	1:A:288:CYS:HB2	2.20	0.41
1:A:397:GLU:HB2	1:A:398:PRO:HD2	2.03	0.41
1:A:584:PHE:CD2	1:A:584:PHE:C	2.99	0.41
1:A:243:LYS:O	1:A:243:LYS:HG3	2.20	0.41
1:A:44:HIS:CD2	1:A:44:HIS:C	2.98	0.41
1:A:46:CYS:HA	1:A:52:GLU:OE2	2.20	0.41
1:A:561:GLU:O	1:A:562:ASP:HB2	2.21	0.41
1:A:158:ILE:HD12	1:A:158:ILE:HG23	1.82	0.40
1:A:290:LYS:HB2	3:A:703:GCP:O1B	2.20	0.40
1:A:246:PHE:C	1:A:246:PHE:CD1	2.99	0.40
1:A:442:PRO:HA	1:A:577:LEU:O	2.21	0.40
1:A:416:ILE:HD12	1:A:416:ILE:HA	1.99	0.40
1:A:500[A]:MSE:HE3	1:A:500[A]:MSE:HB2	1.97	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	598/625 (96%)	571 (96%)	26 (4%)	1 (0%)	43 58

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	151	SER

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	509/501 (102%)	485 (95%)	24 (5%)	23 41

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	ASN
1	A	31	LEU
1	A	36	GLU
1	A	87	ARG
1	A	88	ILE
1	A	97	GLN
1	A	100	ARG
1	A	103	VAL
1	A	123	GLU
1	A	138	THR
1	A	182	GLU
1	A	253	ARG
1	A	296	MSE
1	A	312	ILE
1	A	328	PRO
1	A	387	LYS
1	A	393	SER
1	A	394	GLU
1	A	403	ASN
1	A	473	LYS

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Mol	Chain	Res	Type
1	A	537	LEU
1	A	541	PHE
1	A	560	LYS
1	A	581	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	53	ASN
1	A	112	GLN
1	A	320	GLN
1	A	330	ASN
1	A	346	ASN
1	A	388	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 3 ligands modelled in this entry, 2 are monoatomic - leaving 1 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
3	GCP	A	703	2	32,34,34	1.60	7 (21%)	49,54,54	1.98	14 (28%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GCP	A	703	2	-	1/19/38/38	0/3/3/3

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	703	GCP	PB-O2B	-3.54	1.47	1.56
3	A	703	GCP	PA-O3A	3.48	1.63	1.59
3	A	703	GCP	PG-O2G	-3.24	1.47	1.55
3	A	703	GCP	C4-N3	-2.49	1.28	1.34
3	A	703	GCP	PB-O3A	2.46	1.61	1.58
3	A	703	GCP	C6-N1	2.30	1.43	1.38
3	A	703	GCP	O3'-C3'	2.19	1.48	1.43

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	703	GCP	O1G-PG-C3B	6.43	125.39	111.37
3	A	703	GCP	O2G-PG-O1G	-4.90	99.72	112.39
3	A	703	GCP	O1B-PB-C3B	4.02	119.64	109.05
3	A	703	GCP	O3G-PG-O1G	-3.59	103.09	112.39
3	A	703	GCP	O2'-C2'-C1'	-3.10	99.42	110.10
3	A	703	GCP	O2A-PA-O1A	3.01	126.43	112.44
3	A	703	GCP	C5'-C4'-C3'	-2.67	105.61	115.21
3	A	703	GCP	O2B-PB-C3B	2.47	116.97	106.73
3	A	703	GCP	O3A-PA-O1A	-2.44	103.36	110.70
3	A	703	GCP	O2G-PG-C3B	2.39	112.20	106.40
3	A	703	GCP	C4'-O4'-C1'	-2.23	104.54	109.47
3	A	703	GCP	C3'-C2'-C1'	-2.17	97.36	101.46
3	A	703	GCP	C2-N3-C4	2.13	115.97	112.30
3	A	703	GCP	C5-C4-N9	2.04	109.31	105.66

There are no chirality outliers.

All (1) torsion outliers are listed below:

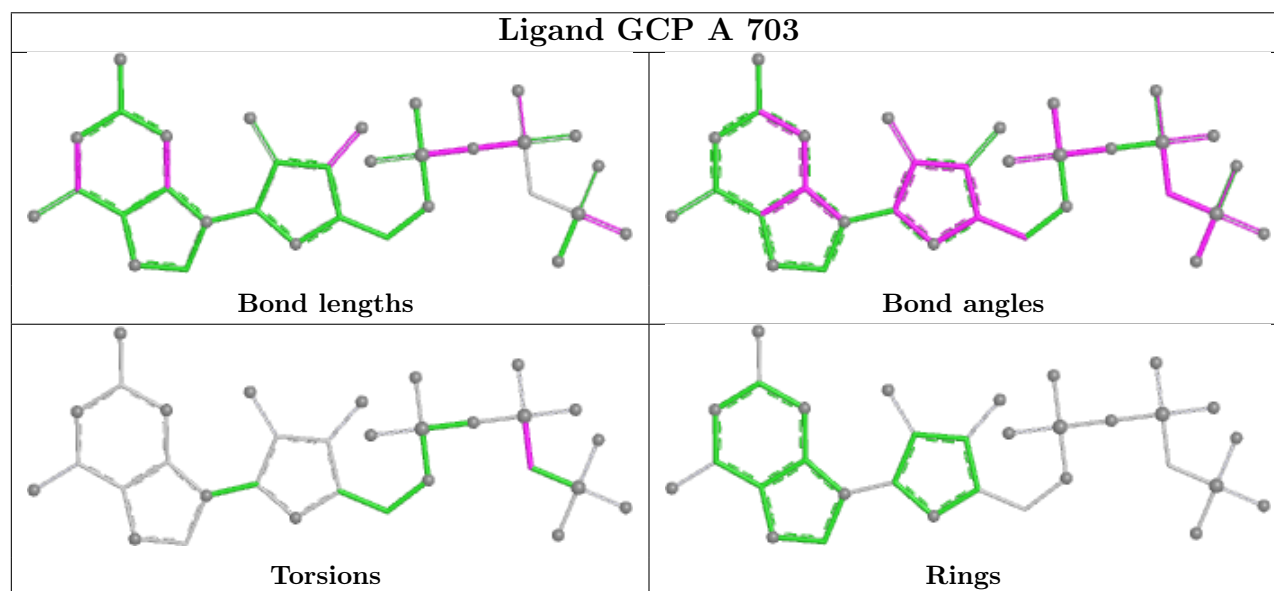
Mol	Chain	Res	Type	Atoms
3	A	703	GCP	PG-C3B-PB-O1B

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	703	GCP	2	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 ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	582/625 (93%)	-0.74	1 (0%) 91 89	3, 10, 28, 48	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	86	ALA	2.2

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 oligosaccharides in this entry.

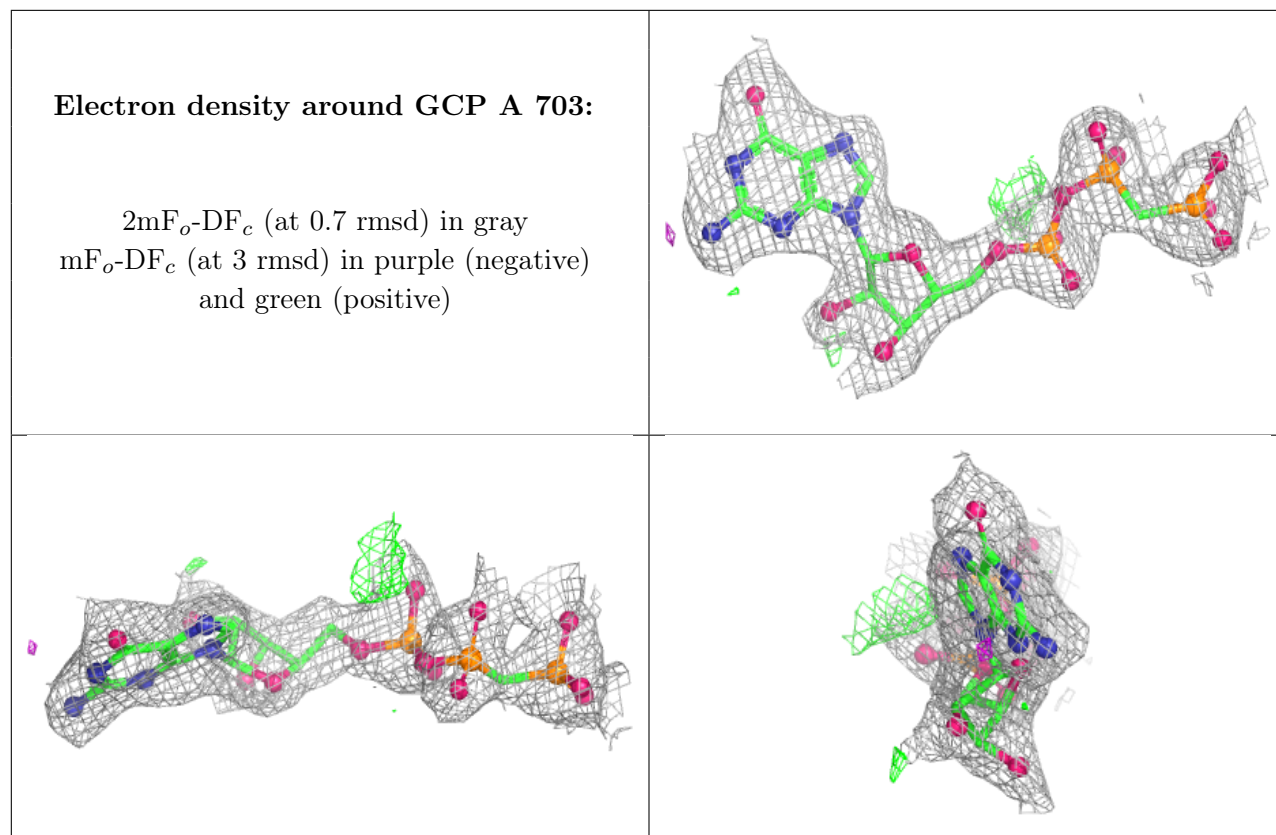
6.4 Ligands [i](#)

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(Å ²)	Q<0.9
3	GCP	A	703	32/32	0.98	0.06	2,8,21,23	0
2	MN	A	702	1/1	1.00	0.03	6,6,6,6	0
2	MN	A	701	1/1	1.00	0.02	6,6,6,6	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 [i](#)

There are no such residues in this entry.