



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 6, 2026 – 09:24 PM UTC

PDB ID : 1OZ0 / pdb_00001oz0
Title : CRYSTAL STRUCTURE OF THE HOMODIMERIC BIFUNCTIONAL TRANSFORMYLASE AND CYCLOHYDROLASE ENZYME AVIAN ATIC IN COMPLEX WITH A MULTISUBSTRATE ADDUCT INHIBITOR BETA-DADF.
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Deposited on : 2003-04-07
Resolution : 2.50 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](#) ①) 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

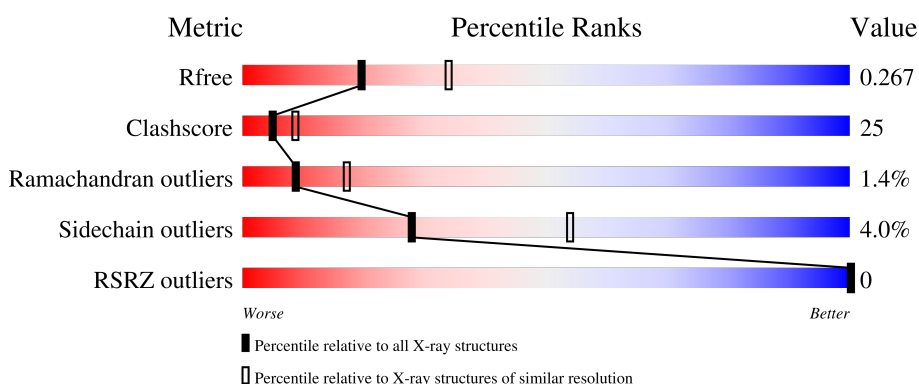
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	613	
1	B	613	

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 9526 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 Bifunctional purine biosynthesis protein PURH.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	590	Total	C	N	O	S	0	0	0
			4511	2843	800	849	19			
1	B	590	Total	C	N	O	S	4	0	0
			4511	2843	800	849	19			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP P31335
A	-18	GLY	-	expression tag	UNP P31335
A	-17	SER	-	expression tag	UNP P31335
A	-16	SER	-	expression tag	UNP P31335
A	-15	HIS	-	expression tag	UNP P31335
A	-14	HIS	-	expression tag	UNP P31335
A	-13	HIS	-	expression tag	UNP P31335
A	-12	HIS	-	expression tag	UNP P31335
A	-11	HIS	-	expression tag	UNP P31335
A	-10	HIS	-	expression tag	UNP P31335
A	-9	SER	-	expression tag	UNP P31335
A	-8	SER	-	expression tag	UNP P31335
A	-7	GLY	-	expression tag	UNP P31335
A	-6	LEU	-	expression tag	UNP P31335
A	-5	VAL	-	expression tag	UNP P31335
A	-4	PRO	-	expression tag	UNP P31335
A	-3	ARG	-	expression tag	UNP P31335
A	-2	GLY	-	expression tag	UNP P31335
A	-1	SER	-	expression tag	UNP P31335
A	0	HIS	-	expression tag	UNP P31335
B	-19	MET	-	expression tag	UNP P31335
B	-18	GLY	-	expression tag	UNP P31335
B	-17	SER	-	expression tag	UNP P31335
B	-16	SER	-	expression tag	UNP P31335
B	-15	HIS	-	expression tag	UNP P31335

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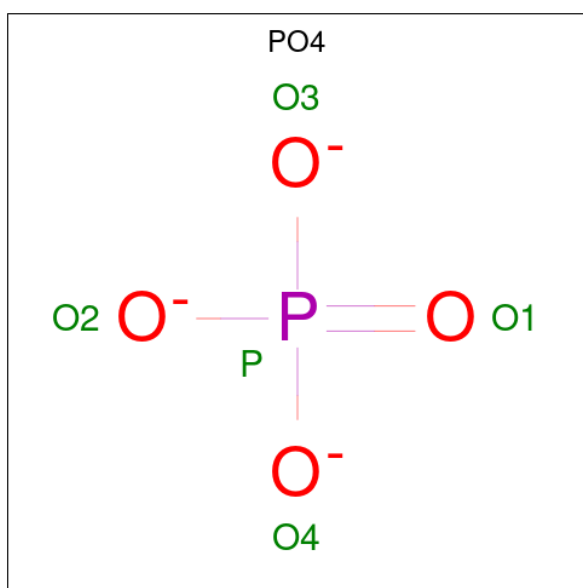
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Chain	Residue	Modelled	Actual	Comment	Reference
B	-14	HIS	-	expression tag	UNP P31335
B	-13	HIS	-	expression tag	UNP P31335
B	-12	HIS	-	expression tag	UNP P31335
B	-11	HIS	-	expression tag	UNP P31335
B	-10	HIS	-	expression tag	UNP P31335
B	-9	SER	-	expression tag	UNP P31335
B	-8	SER	-	expression tag	UNP P31335
B	-7	GLY	-	expression tag	UNP P31335
B	-6	LEU	-	expression tag	UNP P31335
B	-5	VAL	-	expression tag	UNP P31335
B	-4	PRO	-	expression tag	UNP P31335
B	-3	ARG	-	expression tag	UNP P31335
B	-2	GLY	-	expression tag	UNP P31335
B	-1	SER	-	expression tag	UNP P31335
B	0	HIS	-	expression tag	UNP P31335

- Molecule 2 is POTASSIUM ION (CCD ID: K) (formula: K).

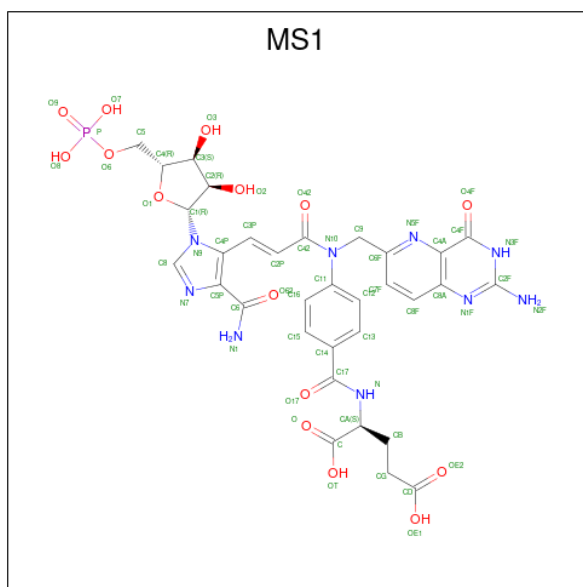
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total K 1 1	0	0
2	B	1	Total K 1 1	0	0

- Molecule 3 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



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

- Molecule 4 is 2-[4-((2-AMINO-4-OXO-3,4-DIHYDRO-PYRIDO[3,2-D]PYRIMIDIN-6-YLM ETHYL)-{3-[5-CARBAMOYL-3-(3,4- DIHYDROXY-5-PHOSPHONOOXYMETHYL-TE TRAHYDRO-FURAN-2-YL)-3H-IMIDAZOL-4-YL]-ACRYLOYL}-AMINO)-BENZOYLAMINO]- PENTANEDIOIC ACID (CCD ID: MS1) (formula: C₃₂H₃₄N₉O₁₅P).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	
			57	32	9	15	1	0
4	B	1	Total	C	N	O	P	
			57	32	9	15	1	0

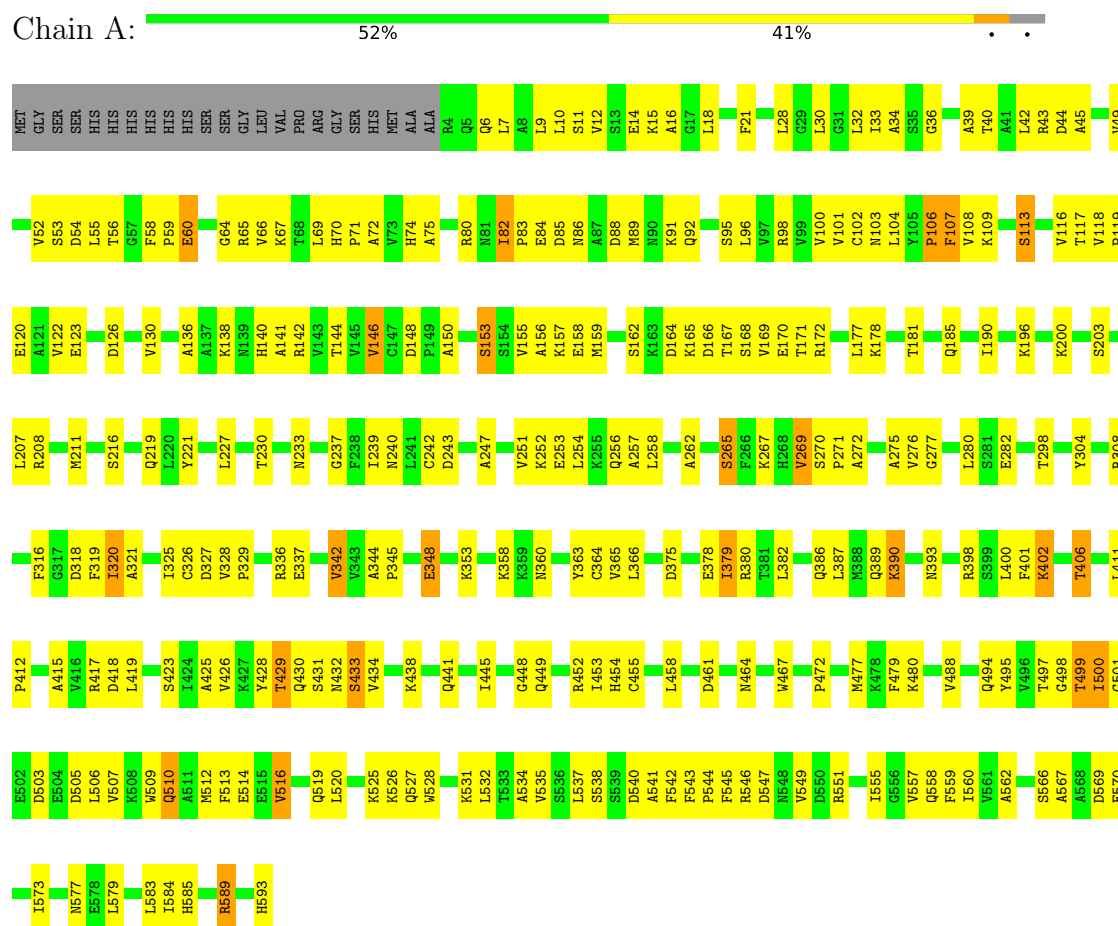
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	204	Total	O	0	0
			204	204		
5	B	174	Total	O	0	0
			174	174		

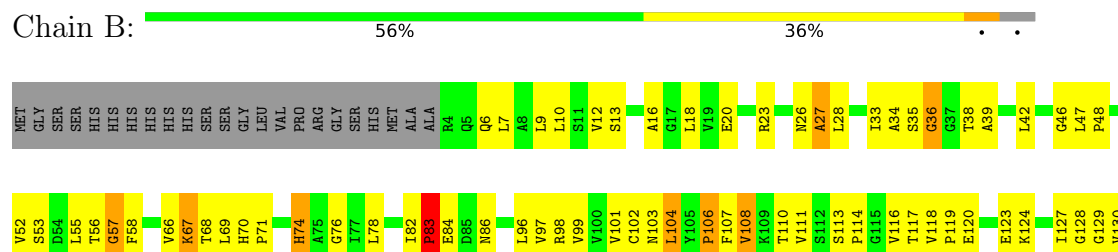
3 Residue-property plots

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: Bifunctional purine biosynthesis protein PURH



- Molecule 1: Bifunctional purine biosynthesis protein PURH



F543	G448	D327	Q219	A131
R546	Q449	V328	L220	L132
D547	Q450	P329	Y221	L133
N548	S451	V342	T222	R134
V549	R452	V343	P225	A135
D550	I453	V344	K226	A136
R551	H454	A344	L227	A137
A552	C455	P345	P228	K138
K553	K462	E349	L229	M139
R554	I555	E350	N233	V143
I555	A463	E350	G234	T144
G556	N464	M660	S235	V145
V557	R469	V365	P236	V146
Q558	V474	L366	G237	C147
F559	M477	Y372	F238	D148
I560	K478	I379	I239	Y152
V561	F479	L382	N240	V155
A562	K480	Q386	L241	A156
P563	A481	L387	C242	K157
S564	G482	L387	D243	E158
G565	V483	K388	A247	M159
S566	K484	Q390	K252	A160
A567	E487	K390	E253	A161
A568	V488	K390	L254	D164
D569	S489	N393	K255	K165
I573	N490	A394	Q256	S168
E574	Q494	F401	A262	V169
A575	Y495	T406	F266	E170
L579	T497	K407	V269	T171
L583	I500	N408	A272	R172
I584	W509	K409	G277	R173
H585	Q510	L411	I278	H174
L588	A511	A415	P279	L175
F591	M512	V416	L280	L176
H592	F513	R417	E283	L177
H593	E514	D418	H294	K178
	E515	L419	L297	T181
	V516	I420	L302	Q185
	P517	V426	A303	R195
	T521	S431	Y304	S199
	E524	N432	D311	K200
	V535	S433	F316	S203
	S536	Y436	I320	L207
	L537	A437	A321	M211
	D540	K438		N212
	A541	Q441		S216
	F542			

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	54.41Å 108.10Å 102.36Å 90.00° 91.92° 90.00°	Depositor
Resolution (Å)	50.00 – 2.50 50.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	92.1 (50.00-2.50) 92.8 (50.00-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.00 (at 2.33Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.217 , 0.269 0.215 , 0.267	Depositor DCC
R_{free} test set	3554 reflections (7.59%)	wwPDB-VP
Wilson B-factor (Å ²)	36.6	Xtriage
Anisotropy	0.507	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 47.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.060 for h,-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9526	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.70% 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: K, PO4, MS1

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	0/4595	0.98	13/6230 (0.2%)
1	B	0.49	0/4595	0.99	15/6230 (0.2%)
All	All	0.49	0/9190	0.98	28/12460 (0.2%)

There are no bond length outliers.

The worst 5 of 28 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	567	ALA	N-CA-C	-8.30	102.86	113.16
1	B	567	ALA	N-CA-C	-7.78	103.78	113.20
1	A	75	ALA	N-CA-C	-7.42	103.13	111.07
1	B	433	SER	N-CA-C	7.30	119.96	108.79
1	A	433	SER	N-CA-C	7.11	119.87	109.07

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	4511	0	4561	255	0
1	B	4511	0	4561	221	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	5	0	0	0	0
3	B	5	0	0	0	0
4	A	57	0	31	6	0
4	B	57	0	31	7	0
5	A	204	0	0	29	0
5	B	174	0	0	23	0
All	All	9526	0	9184	460	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 25.

The worst 5 of 460 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:501:GLY:HA3	1:A:505:ASP:HB2	1.45	0.97
1:A:401:PHE:HB3	1:A:584:ILE:HD13	1.48	0.94
1:B:350:GLU:HB2	5:B:1125:HOH:O	1.70	0.90
1:B:283:GLU:HG3	5:B:1067:HOH:O	1.69	0.89
1:A:64:GLY:HA3	1:A:67:LYS:HE3	1.55	0.88

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	588/613 (96%)	542 (92%)	39 (7%)	7 (1%)	10	20
1	B	588/613 (96%)	541 (92%)	38 (6%)	9 (2%)	8	16
All	All	1176/1226 (96%)	1083 (92%)	77 (6%)	16 (1%)	9	17

5 of 16 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	67	LYS
1	A	36	GLY
1	A	53	SER
1	A	499	THR
1	A	500	ILE

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	484/502 (96%)	464 (96%)	20 (4%)	27	53
1	B	484/502 (96%)	465 (96%)	19 (4%)	28	55
All	All	968/1004 (96%)	929 (96%)	39 (4%)	28	54

5 of 39 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	199	SER
1	B	379	ILE
1	B	211	MET
1	B	269	VAL
1	B	537	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 24 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	174	HIS
1	B	256	GLN
1	B	215	GLN
1	B	367	GLN
1	A	219	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 6 ligands modelled in this entry, 2 are monoatomic - leaving 4 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	MS1	B	1001	-	60,61,61	2.31	23 (38%)	78,89,89	2.93	37 (47%)
4	MS1	A	1002	-	60,61,61	3.50	21 (35%)	78,89,89	3.03	32 (41%)
3	PO4	B	1006	-	4,4,4	1.69	1 (25%)	6,6,6	0.45	0
3	PO4	A	1005	-	4,4,4	1.69	0	6,6,6	0.46	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
4	MS1	B	1001	-	-	12/48/64/64	0/5/5/5
4	MS1	A	1002	-	-	13/48/64/64	0/5/5/5

The worst 5 of 45 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1002	MS1	C42-N10	-19.44	1.11	1.36
4	A	1002	MS1	C11-N10	8.63	1.60	1.43
4	B	1001	MS1	C11-N10	7.32	1.57	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	1002	MS1	C2P-C42	5.49	1.56	1.48
4	A	1002	MS1	C13-C14	4.89	1.46	1.39

The worst 5 of 69 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1002	MS1	C8-N9-C4P	-12.66	96.51	107.30
4	A	1002	MS1	C3P-C2P-C42	11.13	151.85	121.53
4	B	1001	MS1	C3P-C2P-C42	11.03	151.57	121.53
4	B	1001	MS1	C8-N9-C4P	-8.24	100.28	107.30
4	A	1002	MS1	C11-N10-C42	-7.27	115.72	123.53

There are no chirality outliers.

5 of 25 torsion outliers are listed below:

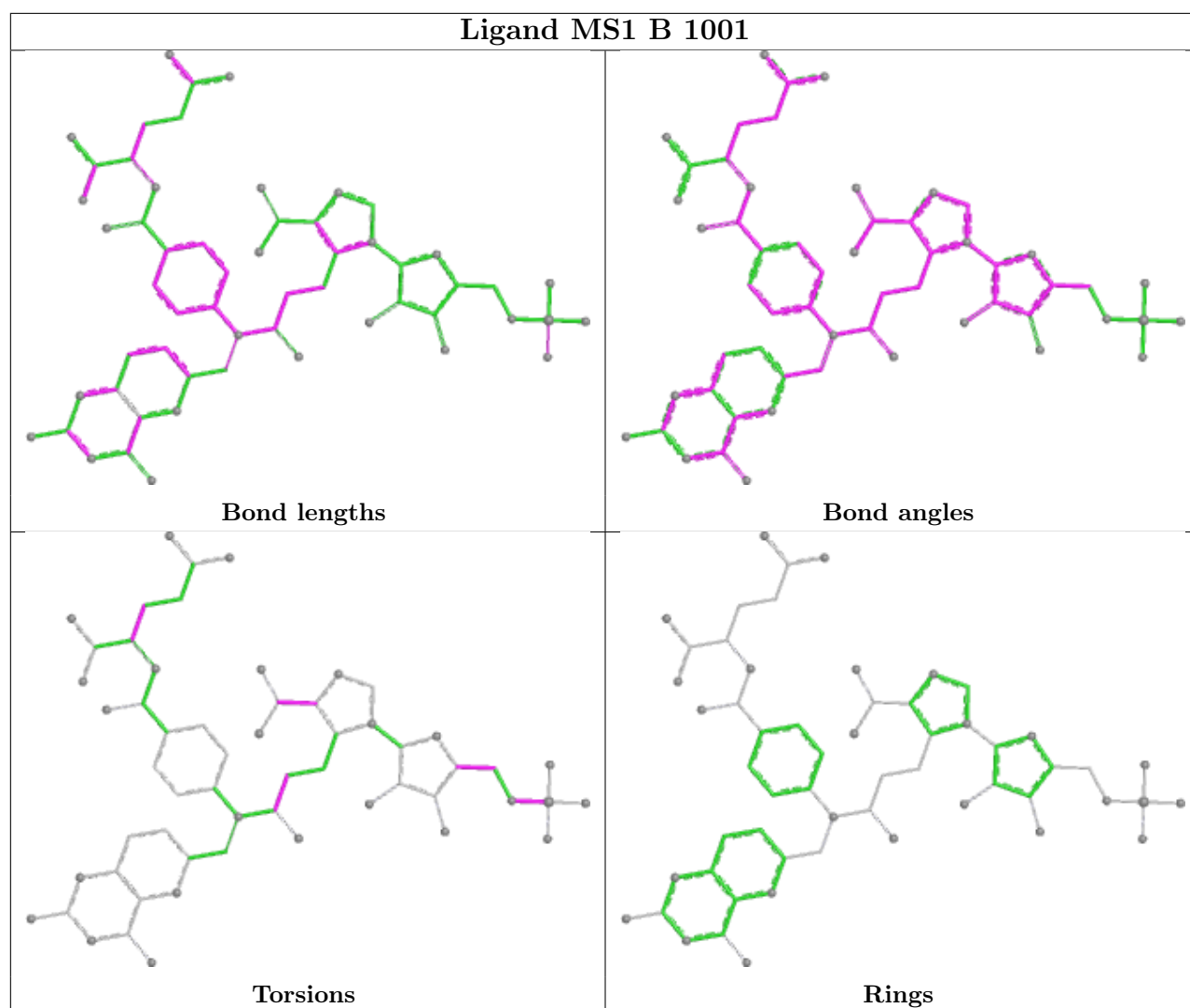
Mol	Chain	Res	Type	Atoms
4	A	1002	MS1	C4P-C5P-C6-O62
4	A	1002	MS1	N7-C5P-C6-O62
4	A	1002	MS1	N7-C5P-C6-N1
4	A	1002	MS1	C5-O6-P-O8
4	B	1001	MS1	C5-O6-P-O8

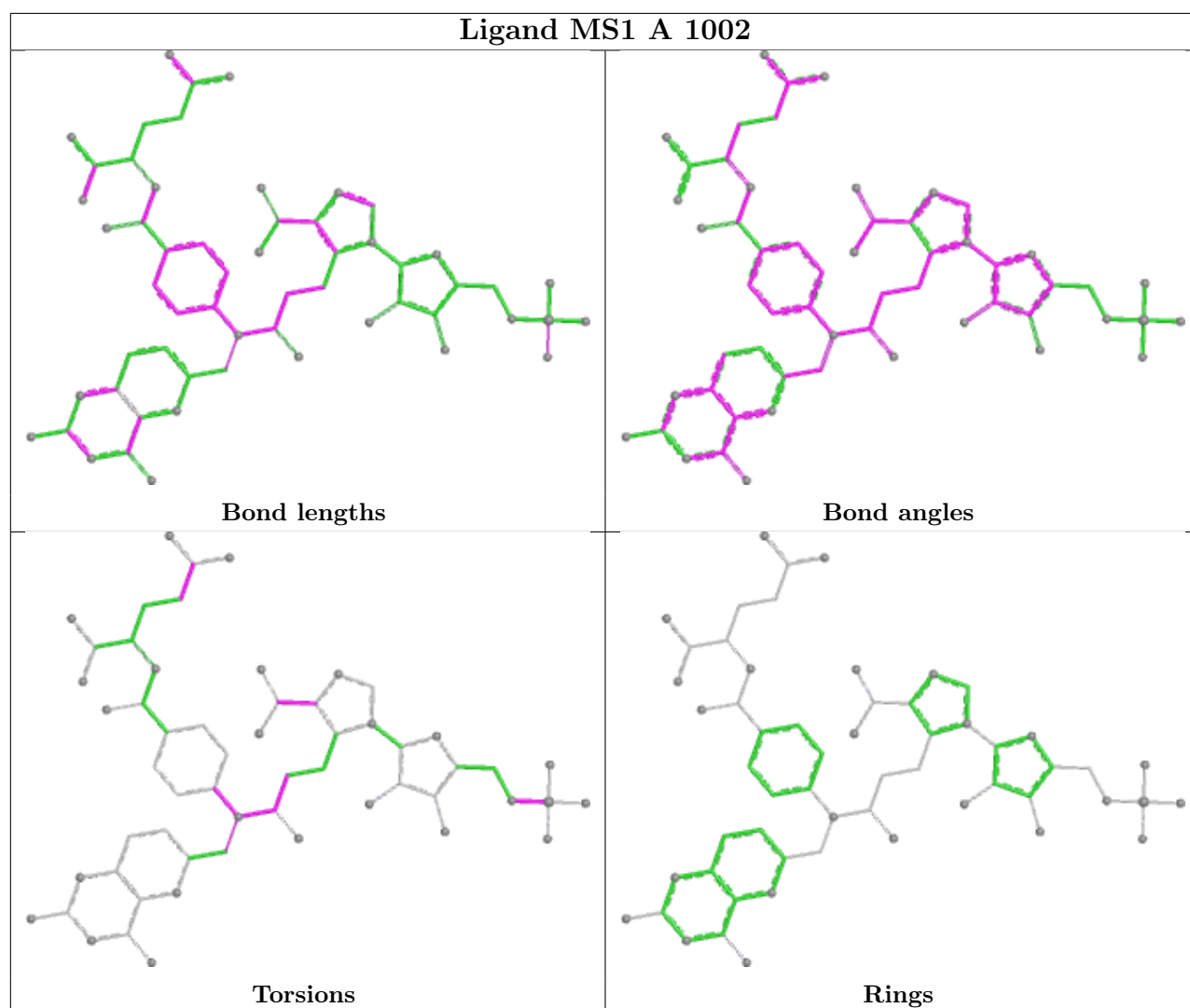
There are no ring outliers.

2 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	1001	MS1	7	0
4	A	1002	MS1	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	590/613 (96%)	-0.21	0 100 100	21, 40, 70, 87	0
1	B	590/613 (96%)	-0.18	0 100 100	21, 39, 68, 82	1 (0%)
All	All	1180/1226 (96%)	-0.20	0 100 100	21, 40, 69, 87	1 (0%)

There are no RSRZ outliers to report.

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](#)

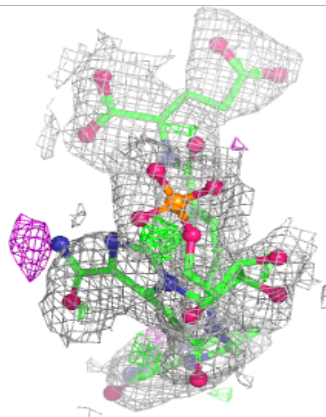
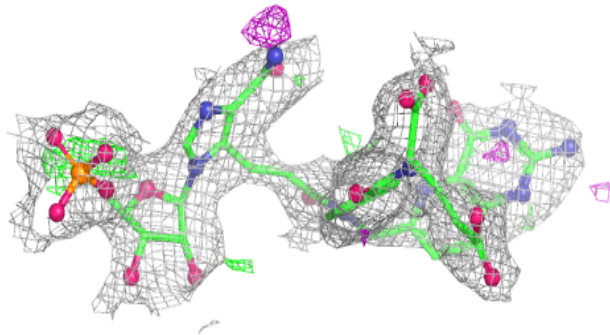
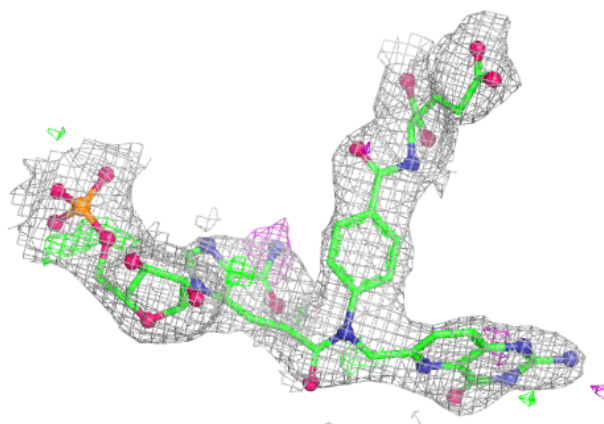
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists 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	PO4	B	1006	5/5	0.78	0.12	81,81,81,82	0
3	PO4	A	1005	5/5	0.83	0.12	97,97,97,98	0
4	MS1	B	1001	57/57	0.86	0.12	47,57,65,66	0
4	MS1	A	1002	57/57	0.92	0.09	38,50,58,58	0
2	K	A	1004	1/1	0.95	0.06	42,42,42,42	0
2	K	B	1003	1/1	0.95	0.05	35,35,35,35	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.

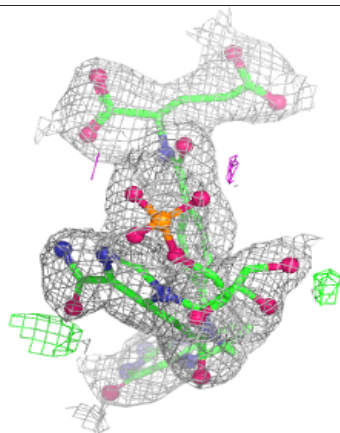
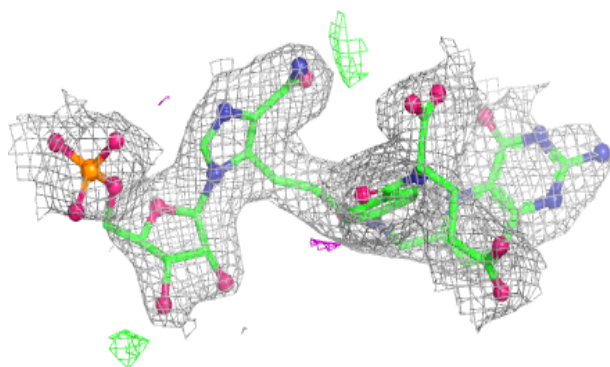
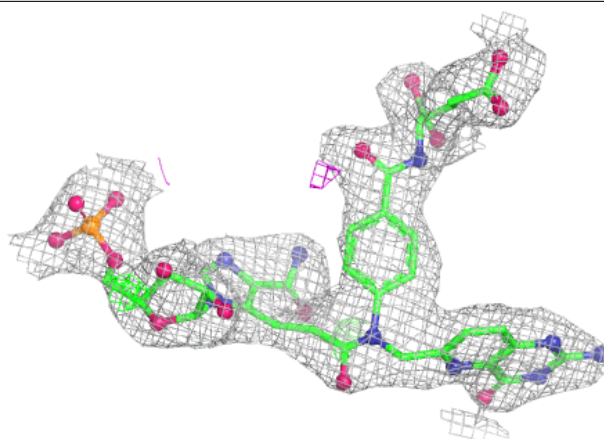
Electron density around MS1 B 1001:

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



Electron density around MS1 A 1002:

$2mF_o-DF_c$ (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.