



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 02:39 PM UTC

PDB ID : 6ZQZ / pdb_00006zqz
Title : [1,2,4]Triazolo[1,5-a]pyrimidine Phosphodiesterase 2 Inhibitors
Authors : Tresadern, G.; Leonard, P.M.
Deposited on : 2020-07-10
Resolution : 1.88 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

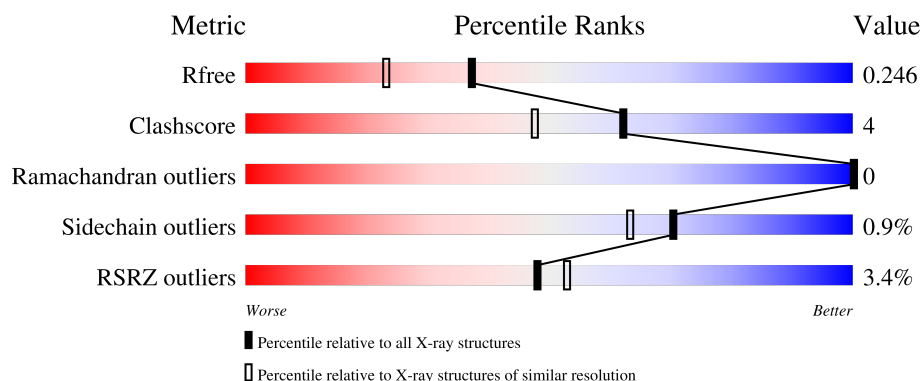
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.88 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	1220 (1.88-1.88)
Clashscore	190562	1234 (1.88-1.88)
Ramachandran outliers	187476	1222 (1.88-1.88)
Sidechain outliers	187428	1222 (1.88-1.88)
RSRZ outliers	180081	1220 (1.88-1.88)

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	353	2% 87% 9% .
1	B	353	3% 88% 7% 5%
1	C	353	7% 84% 9% 7%
1	D	353	% 85% 10% 5%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 11809 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 cGMP-dependent 3',5'-cyclic phosphodiesterase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	339	Total	C	N	O	S	0	1	0
			2738	1747	462	504	25			
1	B	337	Total	C	N	O	S	0	1	0
			2707	1728	460	494	25			
1	C	327	Total	C	N	O	S	0	1	0
			2610	1668	441	476	25			
1	D	336	Total	C	N	O	S	0	0	0
			2727	1737	466	499	25			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	576	MET	-	initiating methionine	UNP O00408
A	577	GLY	-	expression tag	UNP O00408
A	922	ARG	-	expression tag	UNP O00408
A	923	HIS	-	expression tag	UNP O00408
A	924	HIS	-	expression tag	UNP O00408
A	925	HIS	-	expression tag	UNP O00408
A	926	HIS	-	expression tag	UNP O00408
A	927	HIS	-	expression tag	UNP O00408
A	928	HIS	-	expression tag	UNP O00408
B	576	MET	-	initiating methionine	UNP O00408
B	577	GLY	-	expression tag	UNP O00408
B	922	ARG	-	expression tag	UNP O00408
B	923	HIS	-	expression tag	UNP O00408
B	924	HIS	-	expression tag	UNP O00408
B	925	HIS	-	expression tag	UNP O00408
B	926	HIS	-	expression tag	UNP O00408
B	927	HIS	-	expression tag	UNP O00408
B	928	HIS	-	expression tag	UNP O00408
C	576	MET	-	initiating methionine	UNP O00408
C	577	GLY	-	expression tag	UNP O00408
C	922	ARG	-	expression tag	UNP O00408

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Chain	Residue	Modelled	Actual	Comment	Reference
C	923	HIS	-	expression tag	UNP O00408
C	924	HIS	-	expression tag	UNP O00408
C	925	HIS	-	expression tag	UNP O00408
C	926	HIS	-	expression tag	UNP O00408
C	927	HIS	-	expression tag	UNP O00408
C	928	HIS	-	expression tag	UNP O00408
D	576	MET	-	initiating methionine	UNP O00408
D	577	GLY	-	expression tag	UNP O00408
D	922	ARG	-	expression tag	UNP O00408
D	923	HIS	-	expression tag	UNP O00408
D	924	HIS	-	expression tag	UNP O00408
D	925	HIS	-	expression tag	UNP O00408
D	926	HIS	-	expression tag	UNP O00408
D	927	HIS	-	expression tag	UNP O00408
D	928	HIS	-	expression tag	UNP O00408

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

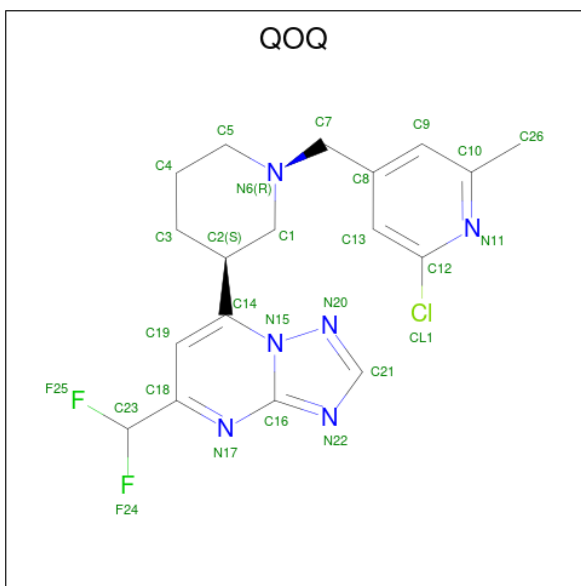
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Zn 1 1	0	0
2	B	1	Total Zn 1 1	0	0
2	C	1	Total Zn 1 1	0	0
2	D	1	Total Zn 1 1	0	0

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

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Mg 1 1	0	0
3	B	2	Total Mg 2 2	0	0
3	C	1	Total Mg 1 1	0	0
3	D	1	Total Mg 1 1	0	0

- Molecule 4 is 5-[bis(fluoranyl)methyl]-7-[(3 {S})-1-[(2-chloranyl-6-methyl-pyridin-4-yl)methyl]piperidin-3-yl]-[1,2,4]triazolo[1,5-a]pyrimidine (CCD ID: QOQ) (formula: C₁₈H₁₉ClF₂N₆)

(labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total 27	C 18	Cl 1	F 2	N 6	0	0
4	B	1	Total 27	C 18	Cl 1	F 2	N 6	0	0
4	C	1	Total 27	C 18	Cl 1	F 2	N 6	0	0
4	D	1	Total 27	C 18	Cl 1	F 2	N 6	0	0

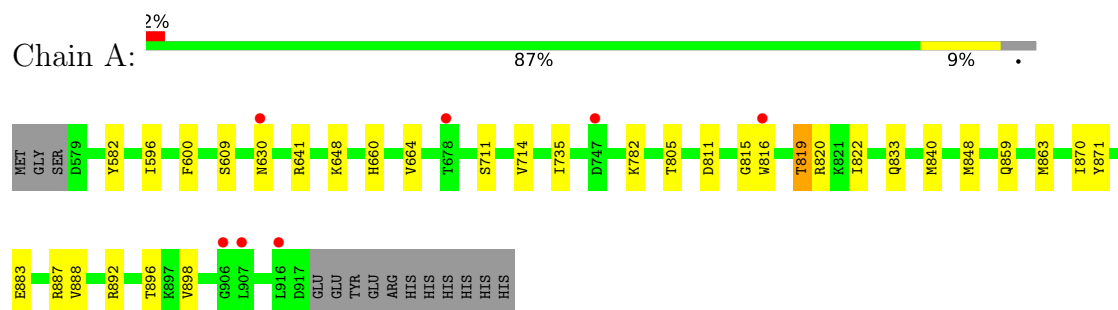
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	249	Total	O	0	0
			249	249		
5	B	240	Total	O	0	0
			240	240		
5	C	173	Total	O	0	0
			173	173		
5	D	248	Total	O	0	0
			248	248		

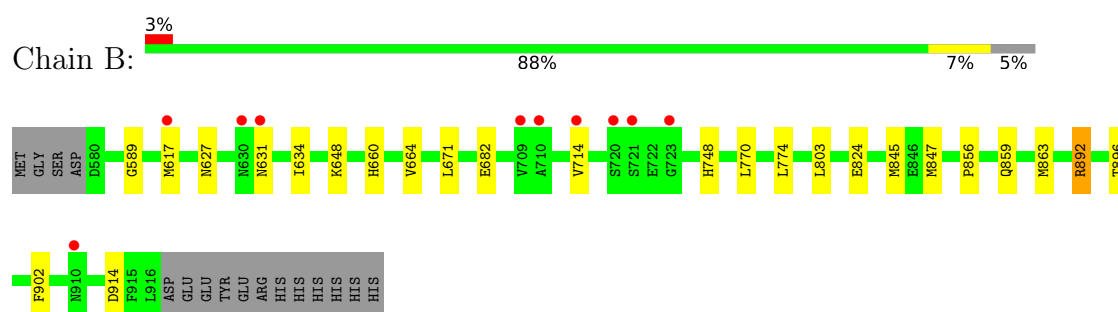
3 Residue-property plots [i](#)

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.

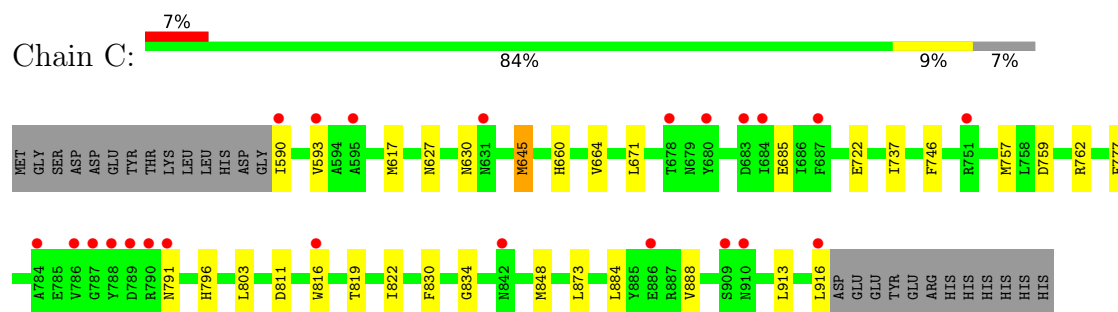
- Molecule 1: cGMP-dependent 3',5'-cyclic phosphodiesterase



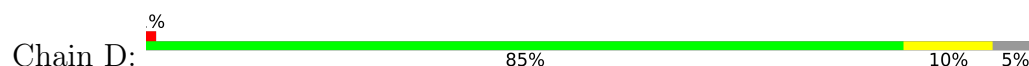
- Molecule 1: cGMP-dependent 3',5'-cyclic phosphodiesterase

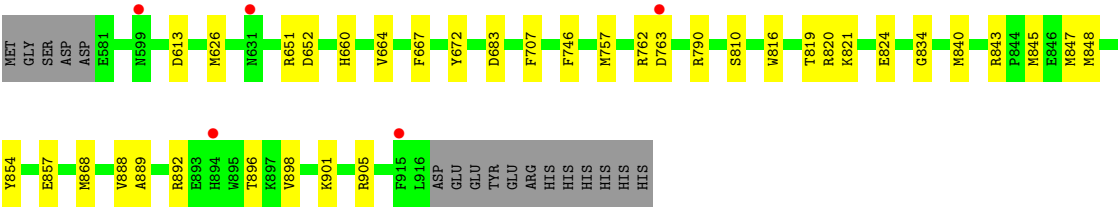


- Molecule 1: cGMP-dependent 3',5'-cyclic phosphodiesterase



- Molecule 1: cGMP-dependent 3',5'-cyclic phosphodiesterase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	55.96Å 73.46Å 92.23Å 109.16° 91.37° 91.26°	Depositor
Resolution (Å)	87.07 – 1.88 87.07 – 1.88	Depositor EDS
% Data completeness (in resolution range)	93.4 (87.07-1.88) 93.5 (87.07-1.88)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.06 (at 1.88Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.196 , 0.244 0.201 , 0.246	Depositor DCC
R_{free} test set	5262 reflections (4.65%)	wwPDB-VP
Wilson B-factor (Å ²)	19.9	Xtriage
Anisotropy	0.072	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 33.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.021 for h,-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11809	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.12% 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: QOQ, MG, ZN

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	0.78	0/2808	0.76	0/3798
1	B	0.78	0/2778	0.76	0/3762
1	C	0.75	0/2678	0.76	0/3630
1	D	0.77	0/2794	0.75	2/3778 (0.1%)
All	All	0.77	0/11058	0.76	2/14968 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	652	ASP	CA-C-N	-5.03	115.20	120.38
1	D	652	ASP	C-N-CA	-5.03	115.20	120.38

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	2738	0	2642	25	0
1	B	2707	0	2591	19	0
1	C	2610	0	2491	19	0
1	D	2727	0	2635	27	0
2	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	1	0	0	0	0
3	B	2	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	27	0	0	0	0
4	B	27	0	0	0	0
4	C	27	0	0	0	0
4	D	27	0	0	0	0
5	A	249	0	0	5	0
5	B	240	0	0	4	0
5	C	173	0	0	1	0
5	D	248	0	0	14	0
All	All	11809	0	10359	87	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 4.

All (87) 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:714:VAL:HG12	5:A:1305:HOH:O	1.52	1.10
1:D:763:ASP:HB3	5:D:1103:HOH:O	1.70	0.91
1:D:868:MET:HG2	5:D:1101:HOH:O	1.77	0.84
1:B:892:ARG:NH2	1:B:896:THR:OG1	2.24	0.70
1:D:845:MET:HB3	5:D:1110:HOH:O	1.93	0.69
1:A:883:GLU:OE2	1:A:887:ARG:NH2	2.27	0.68
1:B:682:GLU:HG3	5:B:1136:HOH:O	1.93	0.68
1:D:613:ASP:HB2	5:D:1314:HOH:O	1.94	0.68
1:C:590:ILE:N	1:C:617:MET:HE1	2.09	0.68
1:D:847:MET:HG2	5:D:1110:HOH:O	1.95	0.65
1:D:845:MET:HA	1:D:845:MET:HE2	1.78	0.64
1:C:759:ASP:OD1	1:C:762:ARG:NH2	2.30	0.64
1:B:617:MET:HE2	5:B:1311:HOH:O	1.99	0.63
1:B:770:LEU:O	1:B:774:LEU:CD2	2.48	0.62
1:C:645:MET:HE3	1:C:737:ILE:HG23	1.82	0.62
1:D:898:VAL:HG23	1:D:901:LYS:HD2	1.83	0.61
1:A:630:ASN:HB3	5:A:1322:HOH:O	2.01	0.60
1:C:627:ASN:OD1	1:C:630:ASN:ND2	2.28	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:816:TRP:NE1	5:A:1103:HOH:O	2.35	0.59
1:B:770:LEU:O	1:B:774:LEU:HD23	2.02	0.59
1:C:830:PHE:O	1:C:848:MET:HE2	2.03	0.58
1:C:830:PHE:HA	1:C:848:MET:CE	2.34	0.58
1:D:889:ALA:N	5:D:1101:HOH:O	2.38	0.57
1:A:711:SER:HA	1:A:840:MET:HE1	1.86	0.57
1:C:777:PHE:CE1	1:C:873:LEU:HD11	2.41	0.56
1:D:707:PHE:CZ	1:D:840:MET:HE1	2.41	0.55
1:D:888:VAL:HB	5:D:1101:HOH:O	2.06	0.55
1:A:833:GLN:NE2	1:A:848:MET:HE2	2.22	0.54
1:B:627:ASN:ND2	1:B:631:ASN:OD1	2.40	0.54
1:A:805:THR:HG22	1:A:870:ILE:HD13	1.89	0.54
1:A:848:MET:HG3	5:A:1111:HOH:O	2.08	0.53
1:D:660:HIS:O	1:D:664:VAL:HG23	2.09	0.53
1:B:682:GLU:CG	5:B:1136:HOH:O	2.55	0.52
1:D:845:MET:CB	5:D:1110:HOH:O	2.56	0.52
1:B:634:ILE:HD11	1:B:748[B]:HIS:CE1	2.44	0.52
1:A:820:ARG:NH2	5:A:1105:HOH:O	2.42	0.52
1:A:892:ARG:HH22	1:A:896[B]:THR:CG2	2.23	0.52
1:D:905:ARG:NE	5:D:1102:HOH:O	2.44	0.50
1:C:671:LEU:HD13	1:C:803:LEU:HD22	1.94	0.50
1:C:816:TRP:O	1:C:819[B]:THR:HG22	2.12	0.50
1:A:892:ARG:HH22	1:A:896[B]:THR:HG23	1.76	0.49
1:B:770:LEU:O	1:B:774:LEU:HD22	2.12	0.49
1:A:816:TRP:HA	1:A:819:THR:CG2	2.43	0.47
1:A:833:GLN:CD	1:A:848:MET:HE2	2.40	0.47
1:C:913:LEU:HB3	1:C:916:LEU:HD12	1.96	0.47
1:B:589:GLY:O	1:B:617:MET:HE1	2.14	0.47
1:A:582:TYR:CE1	1:A:641:ARG:HG3	2.50	0.46
1:C:811:ASP:HB3	1:C:822:ILE:HG13	1.97	0.46
1:A:596:ILE:HG21	1:A:600:PHE:CE1	2.51	0.46
1:A:815:GLY:O	1:A:819:THR:HG22	2.15	0.45
1:C:617:MET:O	1:C:617:MET:HE3	2.16	0.45
1:D:834:GLY:HA3	1:D:848:MET:O	2.17	0.45
1:A:811:ASP:HB3	1:A:822:ILE:HG13	1.99	0.45
1:C:685:GLU:OE1	1:C:796:HIS:ND1	2.48	0.45
1:B:845:MET:HE2	1:B:847:MET:HE3	1.99	0.44
1:A:833:GLN:CD	1:A:848:MET:CE	2.91	0.44
1:C:884:LEU:O	1:C:888:VAL:HG23	2.18	0.44
1:B:914:ASP:HB2	5:B:1115:HOH:O	2.17	0.44
1:A:859:GLN:O	1:A:863:MET:HG2	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:854:TYR:HD2	1:D:857:GLU:HB2	1.83	0.43
1:D:746:PHE:CE1	1:D:757:MET:HG2	2.54	0.43
1:A:660:HIS:O	1:A:664:VAL:HG23	2.17	0.43
1:C:746:PHE:CZ	1:C:757:MET:HG2	2.54	0.43
1:A:735:ILE:HG21	1:B:714:VAL:HG21	2.00	0.43
1:D:816:TRP:HA	1:D:819:THR:HG22	2.01	0.43
1:A:871:TYR:CD1	1:A:888:VAL:HG21	2.54	0.42
1:B:648:LYS:HG2	1:C:791:ASN:HD21	1.84	0.42
1:D:821:LYS:O	1:D:824:GLU:HB3	2.19	0.42
1:D:626:MET:HG2	1:D:672:TYR:CD2	2.54	0.42
1:A:892:ARG:NH2	1:A:896[A]:THR:OG1	2.43	0.42
1:B:660:HIS:O	1:B:664:VAL:HG23	2.18	0.42
1:B:671:LEU:HD13	1:B:803:LEU:HD22	2.02	0.42
1:D:845:MET:HE3	5:D:1329:HOH:O	2.19	0.42
1:B:856:PRO:HG3	1:B:902:PHE:CD2	2.54	0.42
1:C:722:GLU:HB2	5:C:1193:HOH:O	2.19	0.42
1:D:892:ARG:NH1	1:D:896:THR:OG1	2.49	0.42
1:D:667:PHE:CD1	1:D:810:SER:HB2	2.56	0.41
1:A:648:LYS:HG3	1:D:790:ARG:HD2	2.02	0.41
1:C:660:HIS:O	1:C:664:VAL:HG23	2.21	0.41
1:C:834:GLY:HA3	1:C:848:MET:O	2.20	0.41
1:D:905:ARG:CZ	5:D:1102:HOH:O	2.69	0.41
1:D:762:ARG:NH1	5:D:1112:HOH:O	2.53	0.41
1:D:843:ARG:NH2	5:D:1114:HOH:O	2.54	0.41
1:B:859:GLN:O	1:B:863:MET:HG2	2.21	0.41
1:D:651:ARG:NH1	5:D:1115:HOH:O	2.54	0.41
1:A:805:THR:CG2	1:A:870:ILE:HD13	2.49	0.40
1:B:634:ILE:CD1	1:B:748[B]:HIS:CE1	3.04	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	338/353 (96%)	334 (99%)	4 (1%)	0	100	100
1	B	336/353 (95%)	331 (98%)	5 (2%)	0	100	100
1	C	326/353 (92%)	322 (99%)	4 (1%)	0	100	100
1	D	334/353 (95%)	331 (99%)	3 (1%)	0	100	100
All	All	1334/1412 (94%)	1318 (99%)	16 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	298/319 (93%)	294 (99%)	4 (1%)	61	50
1	B	291/319 (91%)	289 (99%)	2 (1%)	76	69
1	C	279/319 (88%)	277 (99%)	2 (1%)	76	69
1	D	297/319 (93%)	295 (99%)	2 (1%)	76	69
All	All	1165/1276 (91%)	1155 (99%)	10 (1%)	70	63

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

Mol	Chain	Res	Type
1	A	609	SER
1	A	782	LYS
1	A	819	THR
1	A	898	VAL
1	B	824	GLU
1	B	892	ARG
1	C	593	VAL
1	C	645	MET
1	D	683	ASP
1	D	820	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	624	GLN
1	A	741	HIS
1	A	812	GLN
1	B	627	ASN
1	B	679	ASN
1	B	708	GLN
1	B	733	GLN
1	B	842	ASN
1	C	624	GLN
1	C	791	ASN
1	D	812	GLN
1	D	842	ASN
1	D	900	HIS

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 13 ligands modelled in this entry, 9 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	QOQ	B	1003	-	28,30,30	1.44	5 (17%)	36,43,43	2.00	9 (25%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	QOQ	D	1003	-	28,30,30	1.58	5 (17%)	36,43,43	2.21	9 (25%)
4	QOQ	C	1003	-	28,30,30	1.42	5 (17%)	36,43,43	1.97	7 (19%)
4	QOQ	A	1003	-	28,30,30	1.45	5 (17%)	36,43,43	2.01	7 (19%)

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	QOQ	B	1003	-	-	0/8/22/22	0/4/4/4
4	QOQ	D	1003	-	-	0/8/22/22	0/4/4/4
4	QOQ	C	1003	-	-	0/8/22/22	0/4/4/4
4	QOQ	A	1003	-	-	0/8/22/22	0/4/4/4

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	1003	QOQ	C16-N15	-4.80	1.31	1.38
4	A	1003	QOQ	C16-N15	-4.50	1.32	1.38
4	B	1003	QOQ	C16-N15	-4.07	1.32	1.38
4	C	1003	QOQ	C16-N15	-4.07	1.32	1.38
4	B	1003	QOQ	C19-C18	-3.45	1.39	1.45
4	D	1003	QOQ	C19-C18	-3.02	1.40	1.45
4	C	1003	QOQ	C18-N17	3.01	1.35	1.32
4	C	1003	QOQ	C19-C18	-2.94	1.40	1.45
4	D	1003	QOQ	C12-CL1	2.86	1.80	1.74
4	D	1003	QOQ	C18-N17	2.71	1.35	1.32
4	A	1003	QOQ	C19-C18	-2.60	1.41	1.45
4	B	1003	QOQ	C18-N17	2.60	1.35	1.32
4	B	1003	QOQ	C21-N20	2.43	1.36	1.32
4	B	1003	QOQ	C12-CL1	2.38	1.79	1.74
4	A	1003	QOQ	C18-N17	2.38	1.35	1.32
4	C	1003	QOQ	C12-CL1	2.34	1.79	1.74
4	C	1003	QOQ	C21-N20	2.34	1.36	1.32
4	A	1003	QOQ	C12-CL1	2.33	1.79	1.74
4	D	1003	QOQ	C21-N20	2.32	1.36	1.32
4	A	1003	QOQ	C21-N20	2.14	1.35	1.32

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	1003	QOQ	C21-N22-C16	7.86	107.61	102.36
4	C	1003	QOQ	C21-N22-C16	7.39	107.30	102.36
4	A	1003	QOQ	C21-N22-C16	7.27	107.22	102.36
4	B	1003	QOQ	C21-N22-C16	7.26	107.21	102.36
4	D	1003	QOQ	C13-C12-N11	-4.36	120.52	125.47
4	D	1003	QOQ	N20-C21-N22	-4.26	110.03	116.78
4	C	1003	QOQ	N20-C21-N22	-4.13	110.24	116.78
4	B	1003	QOQ	C13-C12-N11	-3.85	121.11	125.47
4	B	1003	QOQ	N20-C21-N22	-3.84	110.69	116.78
4	A	1003	QOQ	C16-N15-N20	3.81	112.68	110.11
4	A	1003	QOQ	N20-C21-N22	-3.62	111.04	116.78
4	D	1003	QOQ	C16-N15-N20	3.53	112.49	110.11
4	A	1003	QOQ	C13-C12-N11	-3.41	121.60	125.47
4	C	1003	QOQ	C13-C12-N11	-2.92	122.16	125.47
4	D	1003	QOQ	C2-C1-N6	-2.87	106.65	110.22
4	C	1003	QOQ	C21-N20-N15	2.84	103.18	100.86
4	C	1003	QOQ	C2-C1-N6	-2.79	106.75	110.22
4	D	1003	QOQ	C21-N20-N15	2.76	103.11	100.86
4	B	1003	QOQ	C16-N15-N20	2.73	111.95	110.11
4	D	1003	QOQ	N15-C16-N22	-2.73	106.75	109.19
4	B	1003	QOQ	C2-C1-N6	-2.71	106.85	110.22
4	A	1003	QOQ	C2-C1-N6	-2.69	106.87	110.22
4	B	1003	QOQ	C3-C2-C14	2.69	116.39	111.78
4	A	1003	QOQ	N15-C16-N22	-2.65	106.82	109.19
4	B	1003	QOQ	C21-N20-N15	2.48	102.88	100.86
4	C	1003	QOQ	C16-N15-N20	2.45	111.77	110.11
4	D	1003	QOQ	C13-C12-CL1	2.38	121.99	118.92
4	A	1003	QOQ	C3-C2-C14	2.26	115.65	111.78
4	B	1003	QOQ	CL1-C12-N11	2.21	120.15	115.45
4	B	1003	QOQ	N15-C16-N22	-2.16	107.26	109.19
4	C	1003	QOQ	C14-C19-C18	-2.10	119.30	121.94
4	D	1003	QOQ	C14-C19-C18	-2.03	119.39	121.94

There are no chirality outliers.

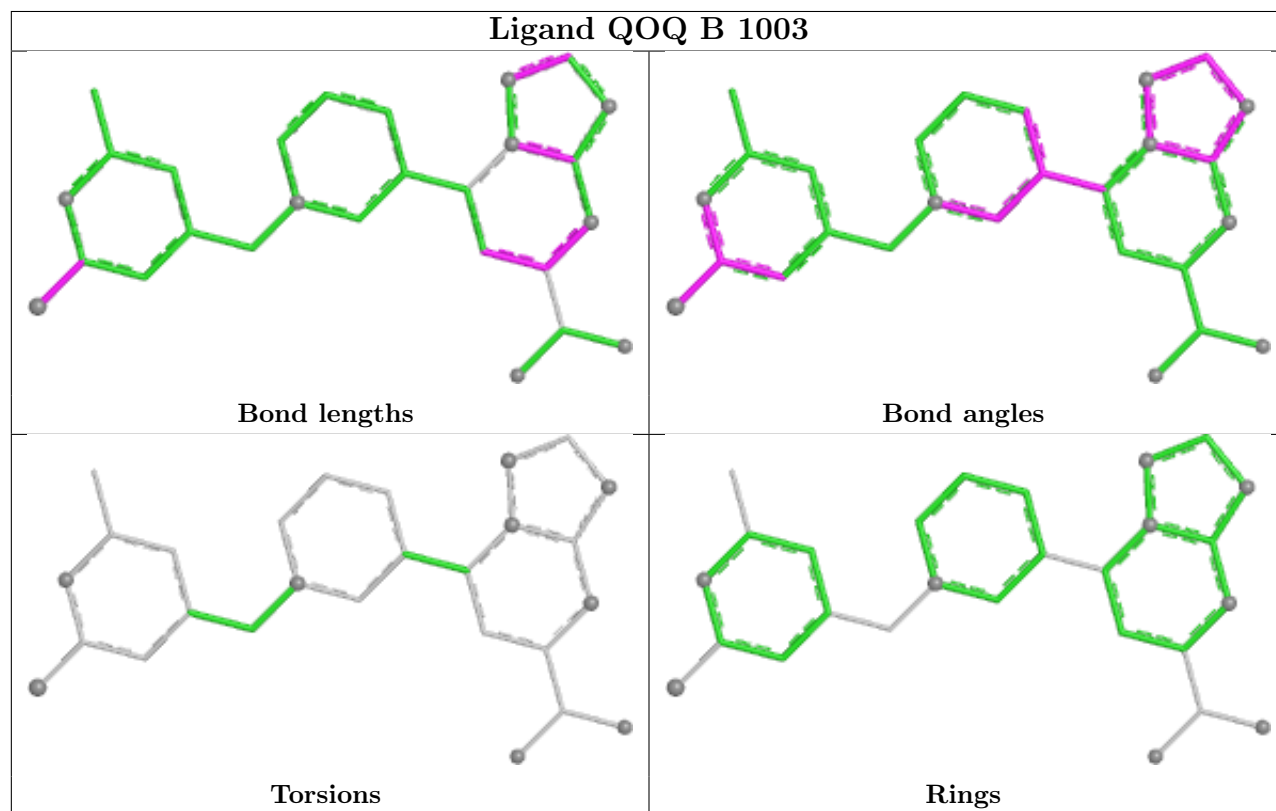
There are no torsion outliers.

There are no ring outliers.

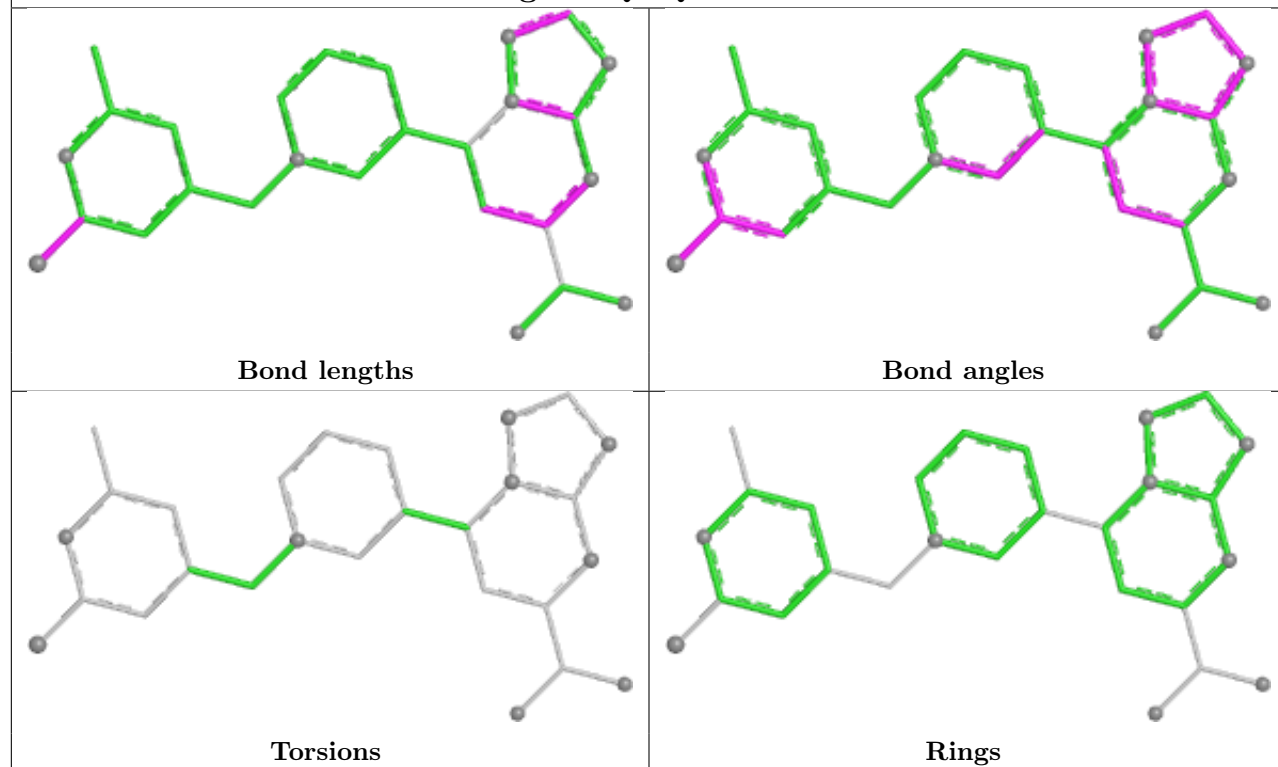
No monomer is involved in short contacts.

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.

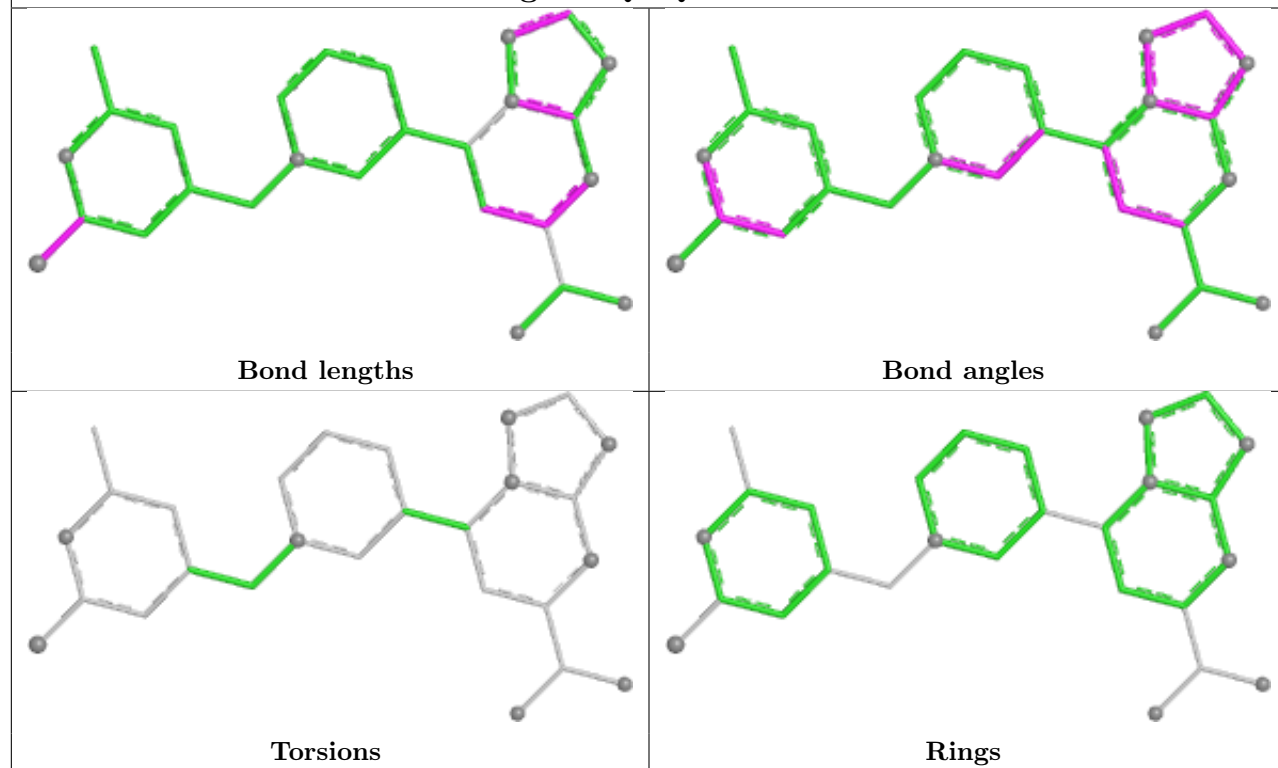
Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD 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.

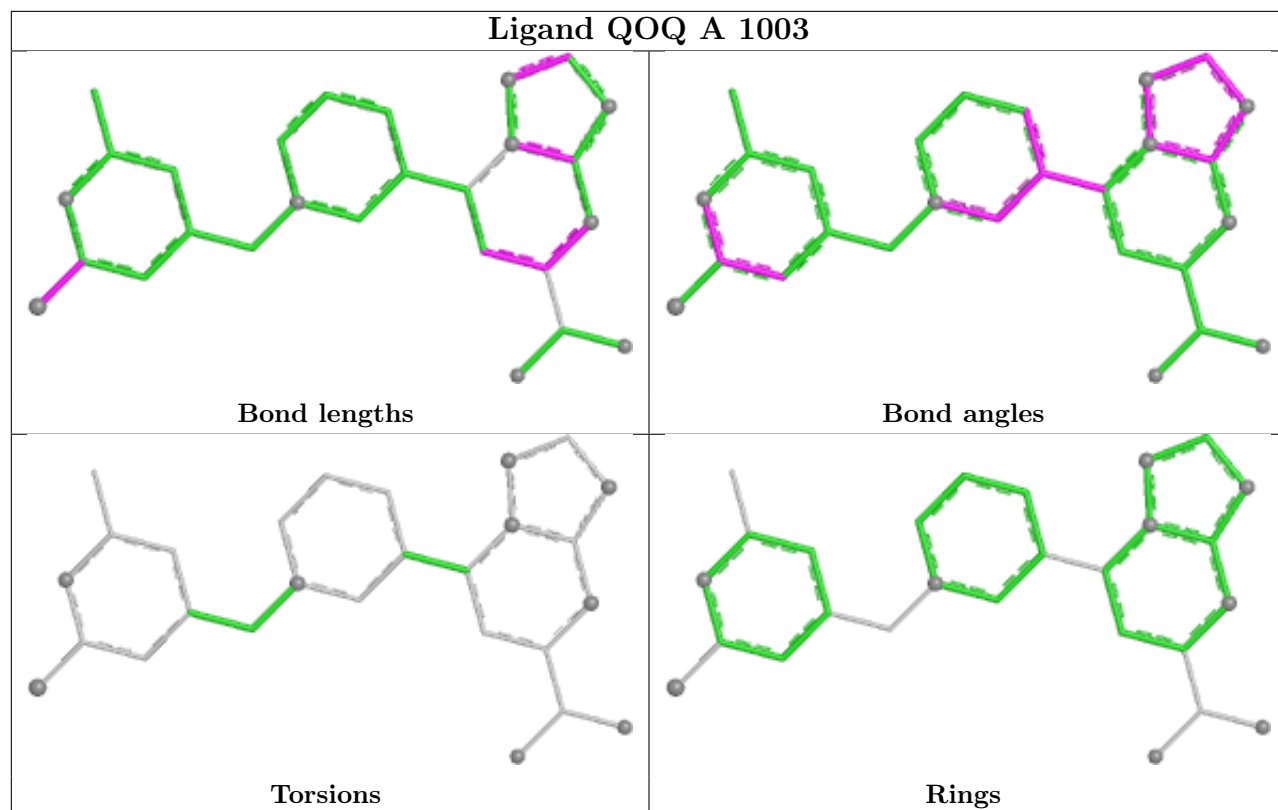


Ligand QOQ D 1003



Ligand QOQ C 1003





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	339/353 (96%)	0.20	7 (2%) 63 70	8, 18, 32, 53	1 (0%)
1	B	337/353 (95%)	0.36	10 (2%) 52 57	10, 21, 34, 47	1 (0%)
1	C	327/353 (92%)	0.63	23 (7%) 22 25	11, 25, 39, 54	1 (0%)
1	D	336/353 (95%)	0.27	5 (1%) 72 77	10, 20, 33, 44	0
All	All	1339/1412 (94%)	0.37	45 (3%) 48 53	8, 21, 36, 54	3 (0%)

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	630	ASN	5.7
1	B	723	GLY	4.0
1	C	787	GLY	3.9
1	C	678	THR	3.9
1	C	593	VAL	3.6
1	C	790	ARG	3.6
1	C	680	TYR	3.5
1	A	747	ASP	3.4
1	B	710	ALA	3.3
1	C	590	ILE	3.2
1	C	786	VAL	3.2
1	C	816	TRP	2.9
1	C	789	ASP	2.8
1	B	721	SER	2.7
1	C	791	ASN	2.7
1	B	910	ASN	2.7
1	C	784	ALA	2.7
1	C	683	ASP	2.7
1	A	678	THR	2.5
1	C	751	ARG	2.5
1	B	714	VAL	2.5

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Mol	Chain	Res	Type	RSRZ
1	C	886	GLU	2.5
1	A	906	GLY	2.4
1	D	763	ASP	2.3
1	A	816	TRP	2.3
1	D	599	ASN	2.3
1	D	894	HIS	2.3
1	B	617	MET	2.3
1	A	916	LEU	2.2
1	B	720	SER	2.2
1	C	595	ALA	2.2
1	C	684	ILE	2.2
1	B	631	ASN	2.2
1	C	910	ASN	2.2
1	B	709	VAL	2.2
1	A	630	ASN	2.1
1	A	907	LEU	2.1
1	D	915	PHE	2.1
1	C	631	ASN	2.1
1	C	687	PHE	2.0
1	C	842	ASN	2.0
1	C	916	LEU	2.0
1	C	909	SER	2.0
1	C	788	TYR	2.0
1	D	631	ASN	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](#)

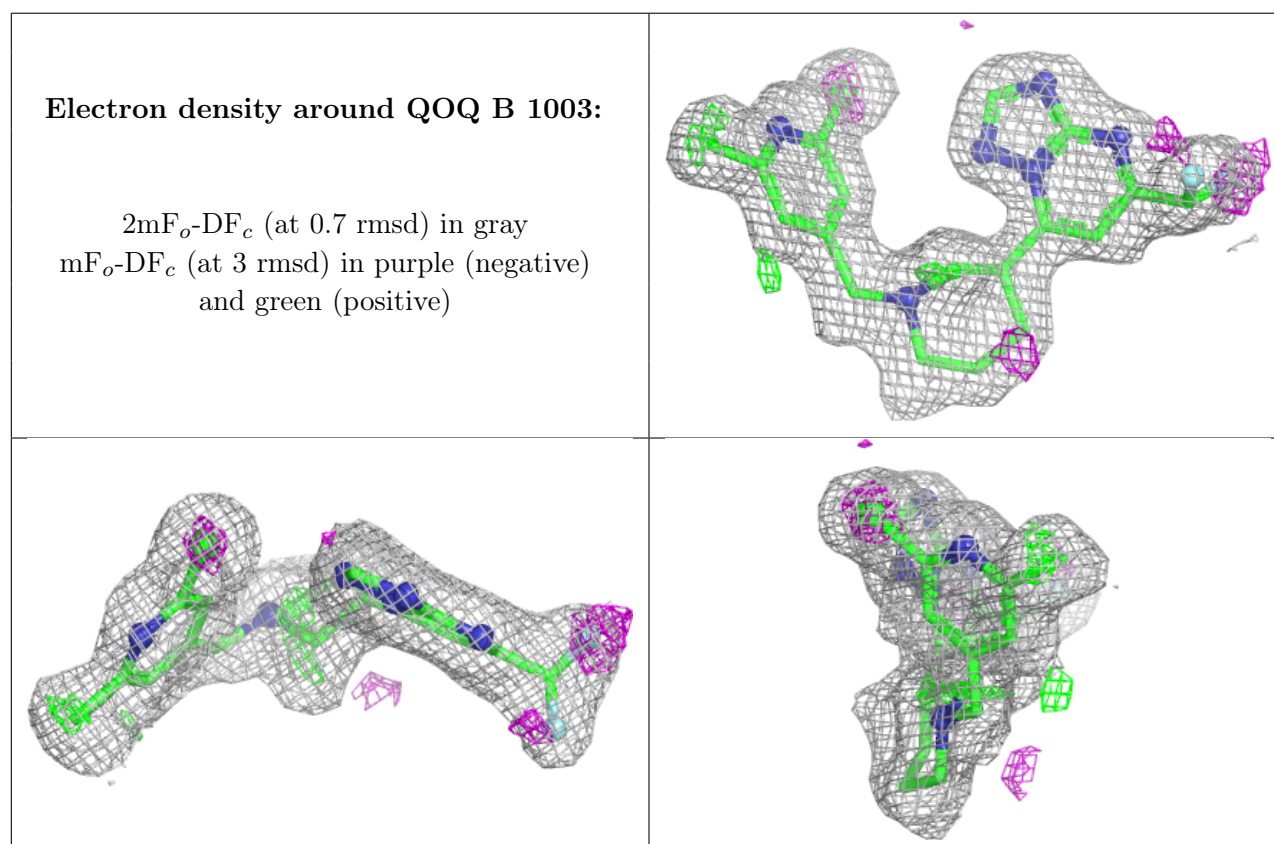
There are no oligosaccharides in this entry.

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.

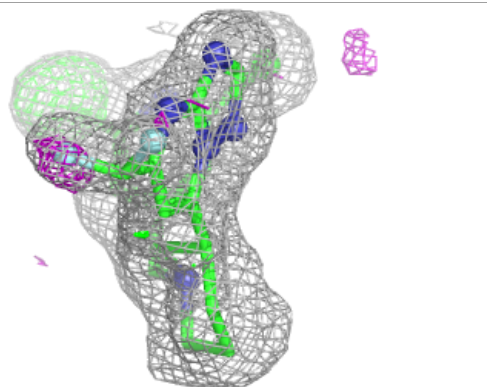
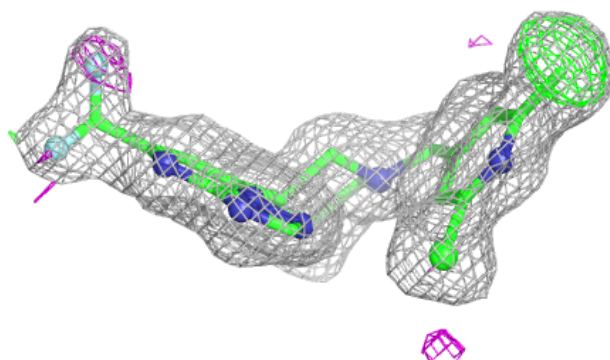
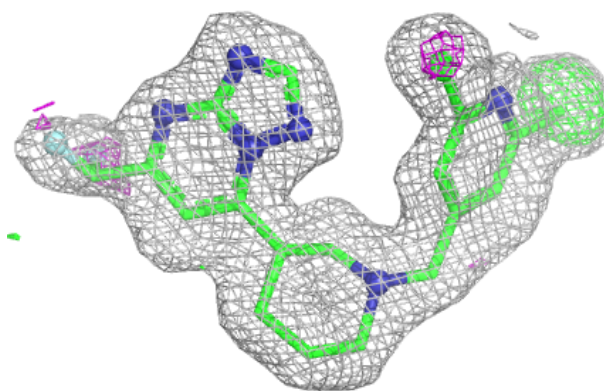
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	QOQ	B	1003	27/27	0.87	0.10	18,20,25,27	0
4	QOQ	A	1003	27/27	0.88	0.10	14,16,19,22	0
4	QOQ	D	1003	27/27	0.88	0.11	20,21,25,28	0
4	QOQ	C	1003	27/27	0.89	0.10	20,21,25,28	0
3	MG	B	1004	1/1	0.94	0.05	27,27,27,27	0
3	MG	B	1002	1/1	0.95	0.03	12,12,12,12	0
3	MG	C	1002	1/1	0.98	0.03	10,10,10,10	0
3	MG	D	1002	1/1	0.98	0.06	11,11,11,11	0
3	MG	A	1002	1/1	0.98	0.05	11,11,11,11	0
2	ZN	B	1001	1/1	1.00	0.01	15,15,15,15	0
2	ZN	C	1001	1/1	1.00	0.01	15,15,15,15	0
2	ZN	D	1001	1/1	1.00	0.01	14,14,14,14	0
2	ZN	A	1001	1/1	1.00	0.02	13,13,13,13	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.

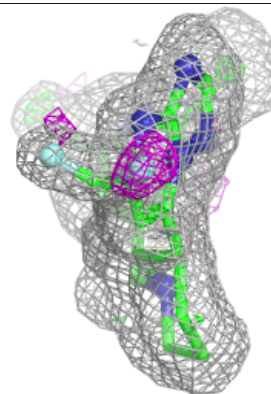
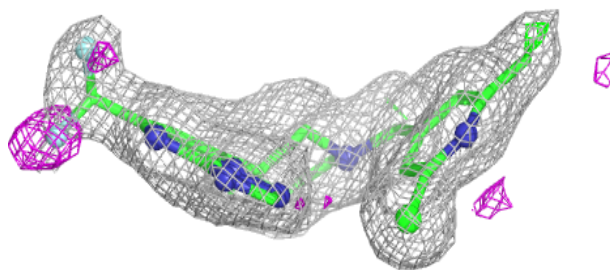
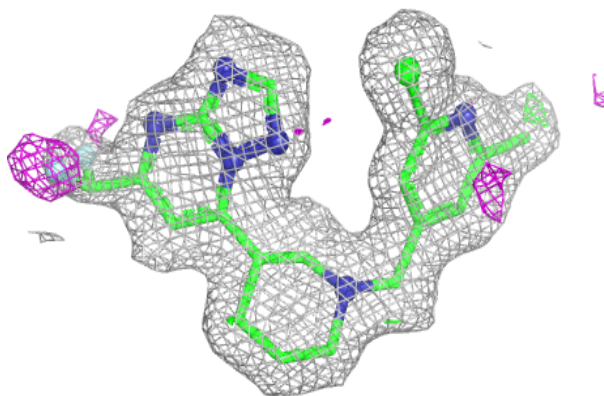


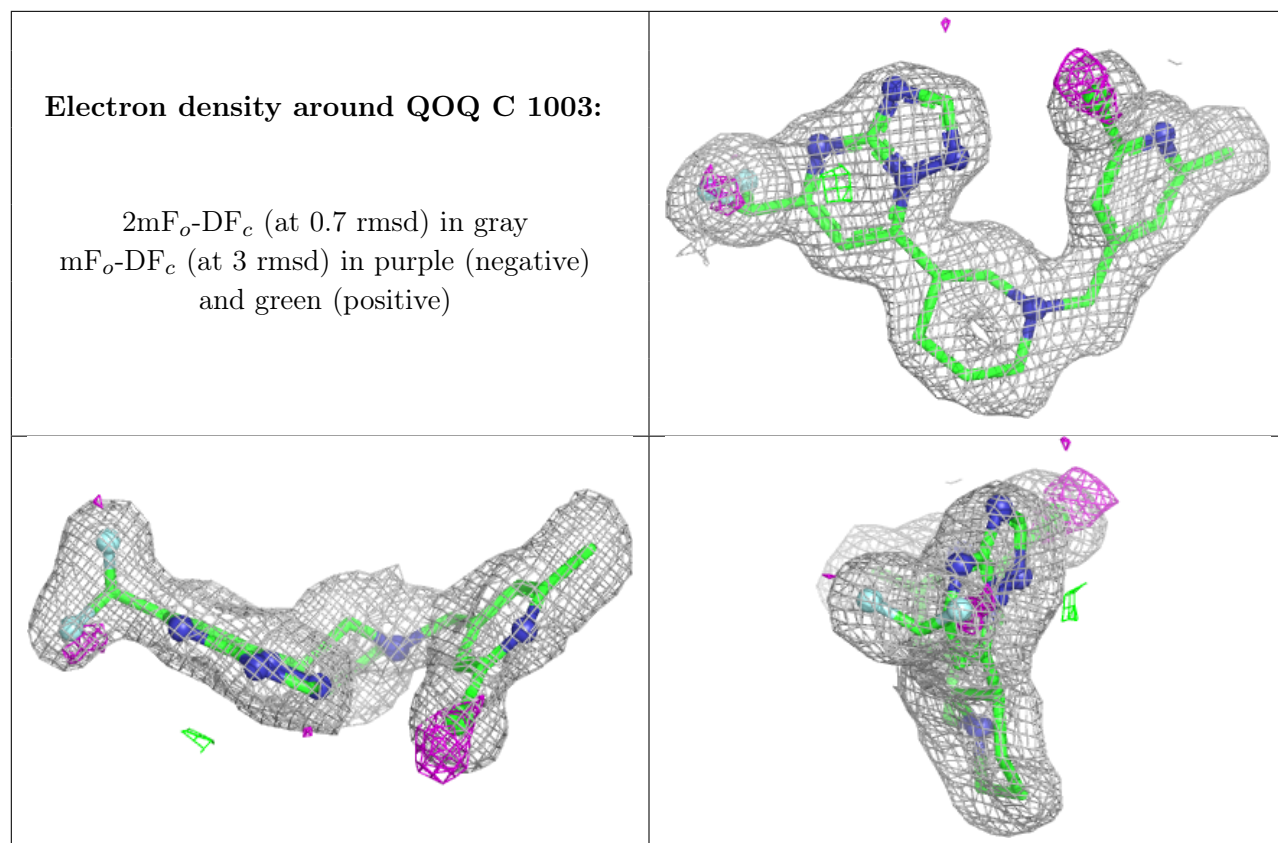
Electron density around QOQ A 1003:

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

**Electron density around QOQ D 1003:**

$2mF_o-DF_c$ (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.