



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 01:51 PM UTC

PDB ID : 5U7K / pdb_00005u7k
Title : PDE2 catalytic domain complexed with inhibitors
Authors : Pandit, J.; Parris, K.
Deposited on : 2016-12-12
Resolution : 2.06 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

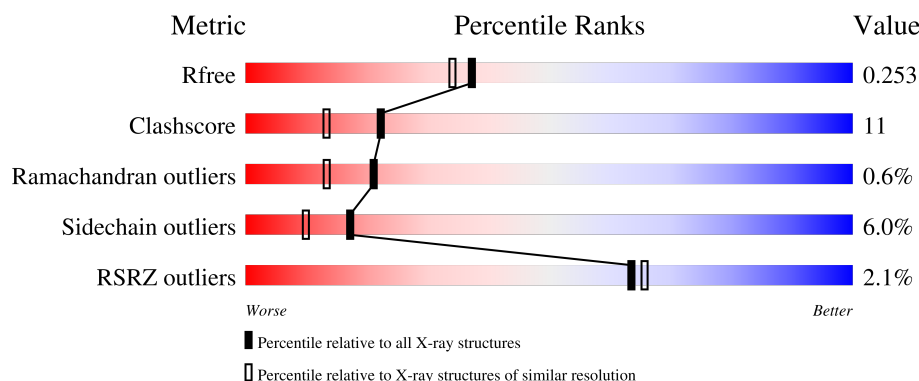
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.06 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	3774 (2.08-2.04)
Clashscore	190562	3883 (2.08-2.04)
Ramachandran outliers	187476	3860 (2.08-2.04)
Sidechain outliers	187428	3860 (2.08-2.04)
RSRZ outliers	180081	3775 (2.08-2.04)

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	345	% 74% 22% ..
1	B	345	% 77% 20% ..
1	C	345	4% 74% 21% ..
1	D	345	% 70% 22% 5%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 11869 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	340	Total	C	N	O	S	0	5	0
			2810	1793	480	510	27			
1	B	343	Total	C	N	O	S	0	2	0
			2810	1789	479	516	26			
1	C	339	Total	C	N	O	S	0	6	0
			2807	1791	477	513	26			
1	D	327	Total	C	N	O	S	0	4	0
			2693	1720	461	486	26			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	575	GLY	-	expression tag	UNP O00408
A	576	SER	-	expression tag	UNP O00408
A	577	ALA	-	expression tag	UNP O00408
A	578	MET	-	expression tag	UNP O00408
B	575	GLY	-	expression tag	UNP O00408
B	576	SER	-	expression tag	UNP O00408
B	577	ALA	-	expression tag	UNP O00408
B	578	MET	-	expression tag	UNP O00408
C	575	GLY	-	expression tag	UNP O00408
C	576	SER	-	expression tag	UNP O00408
C	577	ALA	-	expression tag	UNP O00408
C	578	MET	-	expression tag	UNP O00408
D	575	GLY	-	expression tag	UNP O00408
D	576	SER	-	expression tag	UNP O00408
D	577	ALA	-	expression tag	UNP O00408
D	578	MET	-	expression tag	UNP O00408

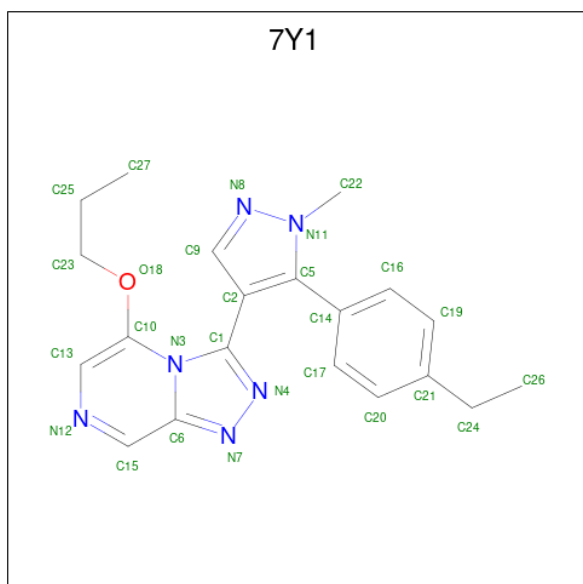
- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

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

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

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

- Molecule 4 is 3-[5-(4-ethylphenyl)-1-methyl-1H-pyrazol-4-yl]-5-propoxy[1,2,4]triazolo[4,3-a]pyrazine (CCD ID: 7Y1) (formula: C₂₀H₂₂N₆O).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			27	20	6	1		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total	C	N	O	0	0
			27	20	6	1		
4	C	1	Total	C	N	O	0	0
			27	20	6	1		
4	D	1	Total	C	N	O	0	0
			27	20	6	1		

- Molecule 5 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	1	Total	Cl	0	0
			1	1		

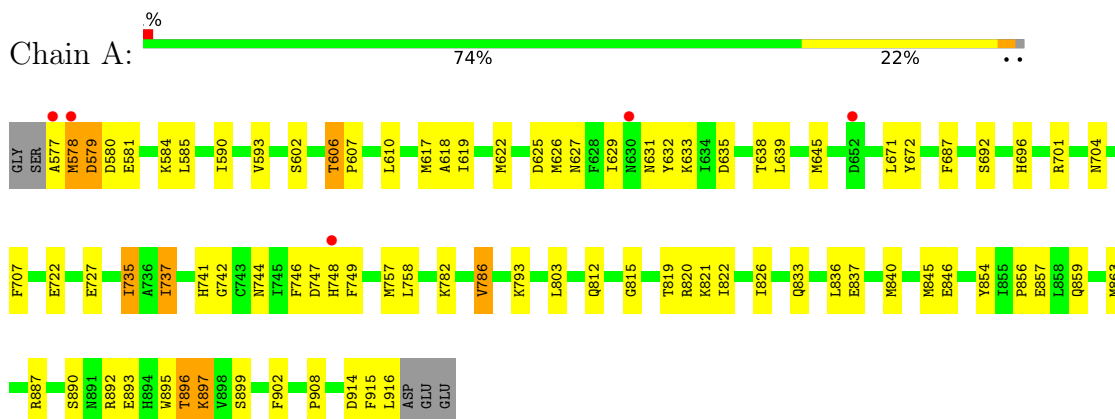
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	159	Total	O	0	0
			159	159		
6	B	192	Total	O	0	0
			192	192		
6	C	150	Total	O	0	0
			150	150		
6	D	131	Total	O	0	0
			131	131		

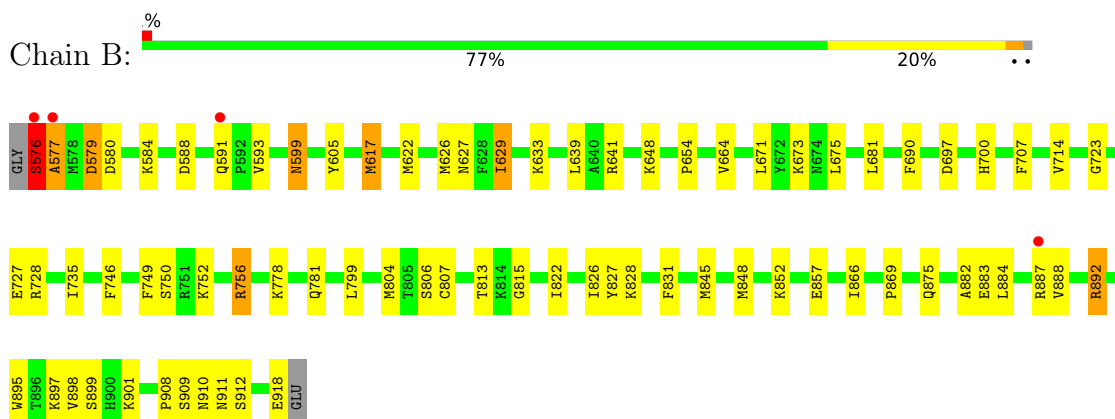
3 Residue-property plots [i](#)

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.

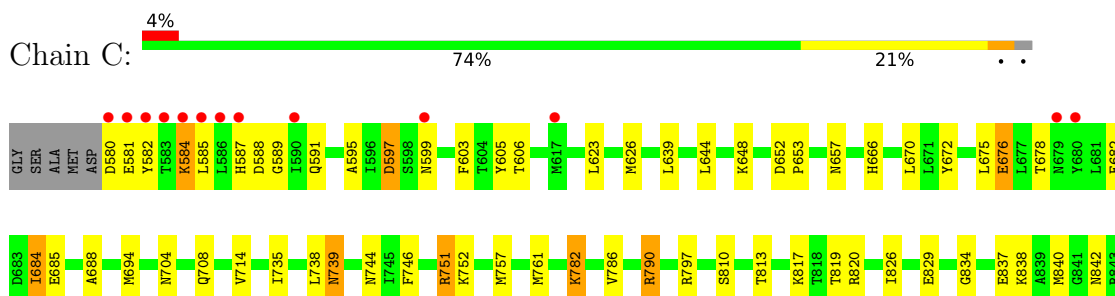
- Molecule 1: cGMP-dependent 3',5'-cyclic phosphodiesterase

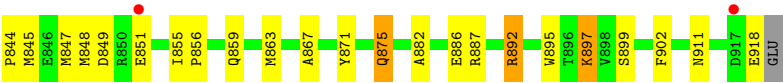


- Molecule 1: cGMP-dependent 3',5'-cyclic phosphodiesterase



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● 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	55.87Å 73.51Å 91.49Å 109.38° 91.23° 90.67°	Depositor
Resolution (Å)	36.00 – 2.06 36.00 – 2.06	Depositor EDS
% Data completeness (in resolution range)	96.2 (36.00-2.06) 96.4 (36.00-2.06)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.59 (at 2.06Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.192 , 0.255 0.193 , 0.253	Depositor DCC
R_{free} test set	4077 reflections (4.79%)	wwPDB-VP
Wilson B-factor (Å ²)	28.6	Xtriage
Anisotropy	0.068	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 46.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.019 for h,-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	11869	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.53% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, MG, 7Y1, CL

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	0.79	0/2893	0.99	5/3901 (0.1%)
1	B	0.83	2/2883 (0.1%)	1.01	3/3890 (0.1%)
1	C	0.79	0/2893	1.03	9/3903 (0.2%)
1	D	0.76	0/2770	0.94	3/3735 (0.1%)
All	All	0.79	2/11439 (0.0%)	0.99	20/15429 (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	B	0	2
1	C	0	1
All	All	0	3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	806	SER	C-O	-6.63	1.16	1.24
1	B	804	MET	C-O	-6.48	1.16	1.24

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	810	SER	N-CA-C	6.75	119.47	111.71
1	C	855	ILE	CA-C-N	-6.60	112.19	119.19
1	C	855	ILE	C-N-CA	-6.60	112.19	119.19
1	A	606	THR	CA-C-N	6.04	126.48	119.47
1	A	606	THR	C-N-CA	6.04	126.48	119.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	625	ASP	N-CA-C	5.94	118.55	111.71
1	C	606	THR	CA-C-N	5.89	125.57	119.56
1	C	606	THR	C-N-CA	5.89	125.57	119.56
1	B	727	GLU	N-CA-C	5.79	117.60	111.28
1	A	744	ASN	N-CA-C	5.70	118.72	108.17
1	C	652	ASP	CA-C-N	-5.38	114.84	120.38
1	C	652	ASP	C-N-CA	-5.38	114.84	120.38
1	D	707	PHE	N-CA-C	5.36	117.82	111.33
1	D	627	ASN	N-CA-C	5.32	118.78	111.54
1	B	735	ILE	N-CA-C	5.31	115.52	110.42
1	C	670	LEU	N-CA-C	5.30	117.74	111.33
1	C	694	MET	N-CA-C	-5.18	105.81	111.82
1	A	610	LEU	N-CA-C	5.15	116.90	109.48
1	A	737	ILE	N-CA-C	5.11	115.33	110.42
1	B	707	PHE	N-CA-C	5.02	116.75	111.28

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	576	SER	Peptide
1	B	577	ALA	Peptide
1	C	589	GLY	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	2810	0	2766	60	0
1	B	2810	0	2752	53	0
1	C	2807	0	2757	68	0
1	D	2693	0	2662	63	0
2	A	1	0	0	0	0
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

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	1	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	1	0
4	D	27	0	0	0	0
5	C	1	0	0	0	0
6	A	159	0	0	6	0
6	B	192	0	0	6	0
6	C	150	0	0	5	0
6	D	131	0	0	6	0
All	All	11869	0	10937	236	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 11.

All (236) 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:D:887:ARG:HG2	6:D:1214:HOH:O	1.43	1.18
1:B:813:THR:O	1:B:887:ARG:HD2	1.44	1.14
1:C:751:ARG:HG2	1:C:751:ARG:HH11	1.20	1.06
1:C:892:ARG:HG2	1:C:892:ARG:HH11	1.24	1.03
1:B:591:GLN:H	1:B:617:MET:CE	1.78	0.96
1:B:591:GLN:H	1:B:617:MET:HE1	1.34	0.93
1:C:751:ARG:HH11	1:C:751:ARG:CG	1.82	0.90
1:B:892:ARG:HH11	1:B:892:ARG:HG2	1.38	0.88
1:B:591:GLN:N	1:B:617:MET:HE1	1.90	0.86
1:C:751:ARG:HG2	1:C:751:ARG:NH1	1.84	0.85
1:A:854:TYR:HD2	1:A:857:GLU:HG3	1.42	0.84
1:A:836:LEU:HG	1:A:840:MET:HE3	1.59	0.83
1:D:883:GLU:OE1	1:D:887:ARG:NH2	2.14	0.81
1:B:845:MET:HE2	6:B:1214:HOH:O	1.79	0.80
1:A:581:GLU:CG	6:A:1171:HOH:O	2.28	0.80
1:B:756:ARG:HH11	1:B:756:ARG:HG2	1.45	0.80
1:C:684:ILE:HG12	1:C:757[A]:MET:HE1	1.63	0.80
1:D:617:MET:HG3	6:D:1204:HOH:O	1.82	0.79
1:C:819:THR:CG2	1:C:895:TRP:HE1	1.94	0.79
1:A:627:ASN:O	1:A:631:ASN:HB2	1.82	0.78
1:B:813:THR:O	1:B:887:ARG:CD	2.28	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:840:MET:HE3	1:C:842:ASN:ND2	1.98	0.78
1:B:591:GLN:HB2	1:B:617:MET:HE2	1.66	0.78
1:C:840:MET:HE3	1:C:842:ASN:HD22	1.50	0.76
1:C:892:ARG:HG2	1:C:892:ARG:NH1	2.00	0.76
1:D:671:LEU:HD13	1:D:803:LEU:HD22	1.68	0.76
1:B:883:GLU:OE1	1:B:887:ARG:NH2	2.21	0.74
1:A:854:TYR:CD2	1:A:857:GLU:HG3	2.21	0.74
1:D:859:GLN:O	1:D:863:MET:HG2	1.88	0.73
1:A:856:PRO:HG3	1:A:902:PHE:CD2	2.24	0.73
1:C:819:THR:HG21	1:C:895:TRP:HE1	1.52	0.72
1:C:587:HIS:NE2	6:C:1101:HOH:O	1.88	0.72
1:C:817:LYS:HE2	1:C:820:ARG:HH22	1.55	0.71
1:D:892:ARG:HH12	1:D:893:GLU:HG3	1.57	0.70
1:D:813:THR:O	1:D:887:ARG:HD2	1.91	0.70
1:D:910:ASN:ND2	1:D:910:ASN:H	1.89	0.70
1:B:884:LEU:O	1:B:888:VAL:HG23	1.94	0.68
1:B:591:GLN:HB2	1:B:617:MET:CE	2.23	0.68
1:A:854:TYR:HD2	1:A:857:GLU:CG	2.07	0.68
1:A:782:LYS:O	1:A:786:VAL:HG13	1.96	0.66
1:A:581:GLU:HG2	6:A:1171:HOH:O	1.91	0.66
1:C:675:LEU:O	1:C:676:GLU:HB2	1.95	0.66
1:C:892:ARG:HH11	1:C:892:ARG:CG	2.04	0.65
1:A:625:ASP:O	1:D:778:LYS:HE2	1.96	0.65
1:C:591:GLN:HG2	1:C:595:ALA:HB3	1.80	0.63
1:B:591:GLN:H	1:B:617:MET:HE2	1.63	0.62
1:C:845:MET:CG	1:C:848:MET:HG3	2.30	0.62
1:B:845:MET:HE3	1:B:848:MET:HG3	1.82	0.61
1:D:819:THR:CG2	1:D:895:TRP:HE1	2.12	0.61
1:B:892:ARG:HG2	1:B:892:ARG:NH1	2.11	0.61
1:D:838:LYS:NZ	6:D:1102:HOH:O	2.25	0.60
1:B:654:PRO:HG2	1:B:828:LYS:HG2	1.83	0.60
1:C:819:THR:HG22	1:C:895:TRP:HE1	1.67	0.60
1:A:692:SER:O	1:A:696:HIS:HB3	2.01	0.60
1:C:837:GLU:OE1	1:C:844:PRO:HA	2.02	0.59
1:B:756:ARG:HH11	1:B:756:ARG:CG	2.14	0.59
1:D:910:ASN:H	1:D:910:ASN:HD22	1.50	0.59
1:A:735:ILE:HD11	1:A:758:LEU:HD22	1.83	0.59
1:A:581:GLU:HG3	6:A:1171:HOH:O	1.96	0.58
1:A:833:GLN:O	1:A:837:GLU:HG3	2.04	0.58
1:D:681:LEU:HD21	1:D:799:LEU:HD23	1.83	0.58
1:A:812:GLN:OE1	1:A:819:THR:HG23	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:817:LYS:CE	1:C:820:ARG:HH22	2.17	0.58
1:A:893:GLU:O	1:A:896:THR:HB	2.03	0.58
1:D:590:ILE:HG22	1:D:590:ILE:O	2.03	0.57
1:D:654:PRO:CG	1:D:828[A]:LYS:HD3	2.34	0.57
1:B:875:GLN:NE2	1:B:882:ALA:CB	2.67	0.57
1:D:748:HIS:H	1:D:748:HIS:CD2	2.21	0.57
1:B:629:ILE:HD13	1:B:639:LEU:HD23	1.85	0.57
1:C:746:PHE:CZ	1:C:757[A]:MET:HG2	2.39	0.57
1:C:840:MET:CE	1:C:842:ASN:HD22	2.18	0.57
1:B:910:ASN:O	1:B:911:ASN:HB2	2.05	0.57
1:A:578:MET:O	1:A:581:GLU:HB3	2.05	0.56
1:B:875:GLN:HE21	1:B:882:ALA:CB	2.17	0.56
1:C:580:ASP:O	1:C:584:LYS:HG3	2.04	0.56
1:D:819:THR:HG22	1:D:895:TRP:HE1	1.69	0.56
1:A:707:PHE:CZ	1:A:840:MET:HE1	2.40	0.56
1:D:622:MET:O	1:D:626:MET:HG3	2.06	0.56
1:B:875:GLN:HE21	1:B:882:ALA:HB2	1.69	0.56
1:C:837:GLU:HA	1:C:840:MET:HE2	1.87	0.56
1:D:782:LYS:HE2	1:D:786:VAL:CG2	2.36	0.56
1:A:632:TYR:HB3	1:A:748[A]:HIS:CE1	2.41	0.55
1:A:704:ASN:HD21	1:A:845:MET:HE1	1.71	0.55
1:B:599:ASN:O	1:B:605:TYR:HB2	2.06	0.55
1:D:875:GLN:HG3	1:D:882:ALA:HA	1.87	0.55
1:A:645[B]:MET:HG2	1:A:737:ILE:HG23	1.87	0.55
1:A:629:ILE:HG12	1:A:639:LEU:CD2	2.36	0.55
1:C:623:LEU:HD13	1:C:639:LEU:HD21	1.89	0.55
1:D:746:PHE:CZ	1:D:757[A]:MET:HG2	2.42	0.55
1:A:635:ASP:HB3	1:A:638:THR:HB	1.88	0.55
1:B:815:GLY:HA3	6:B:1155:HOH:O	2.06	0.54
1:C:845:MET:HG2	1:C:848:MET:HG3	1.90	0.54
1:C:845:MET:HE2	1:C:847:MET:HG2	1.89	0.54
1:D:746:PHE:CE1	1:D:757[B]:MET:HE2	2.43	0.53
1:A:617:MET:HG3	6:A:1204:HOH:O	2.08	0.53
1:B:577:ALA:HB1	1:B:580:ASP:H	1.72	0.53
1:D:884:LEU:O	1:D:888:VAL:HG23	2.08	0.53
1:D:910:ASN:ND2	1:D:910:ASN:N	2.57	0.53
1:C:897:LYS:NZ	1:C:897:LYS:CB	2.70	0.53
1:A:892:ARG:HG2	1:A:892:ARG:HH11	1.74	0.53
1:D:680:TYR:HB3	1:D:788:TYR:CE2	2.43	0.53
1:C:790:ARG:HD2	6:C:1205:HOH:O	2.06	0.52
1:C:892:ARG:NH1	1:C:892:ARG:CG	2.66	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:855:ILE:N	1:D:856:PRO:HD2	2.25	0.52
1:B:664:VAL:HG13	1:B:807:CYS:HB3	1.91	0.52
1:B:626:MET:HE1	1:B:690:PHE:CD1	2.45	0.52
1:C:684:ILE:HG12	1:C:757[A]:MET:CE	2.37	0.52
1:D:664:VAL:HG13	1:D:807:CYS:HB3	1.92	0.51
1:D:875:GLN:HE21	1:D:882:ALA:HA	1.76	0.51
1:C:626:MET:HG2	1:C:672:TYR:CD2	2.45	0.51
1:A:746:PHE:CD1	1:A:749:PHE:HE2	2.29	0.51
1:B:591:GLN:N	1:B:617:MET:CE	2.55	0.50
1:B:895:TRP:O	1:B:899:SER:HB2	2.12	0.50
1:C:813:THR:O	1:C:887:ARG:HD2	2.11	0.50
1:B:576:SER:HB3	1:C:678:THR:HB	1.92	0.50
1:D:910:ASN:HD22	1:D:910:ASN:N	2.09	0.50
1:A:635:ASP:HB3	1:A:638:THR:CB	2.42	0.49
1:B:591:GLN:HG2	6:B:1102:HOH:O	2.11	0.49
1:D:797:ARG:HG2	1:D:797:ARG:HH11	1.77	0.49
1:C:782:LYS:O	1:C:786:VAL:HG22	2.13	0.49
1:D:628:PHE:CD1	1:D:632:TYR:HE2	2.31	0.49
1:C:746:PHE:CE1	1:C:757[B]:MET:HE2	2.48	0.48
1:A:687:PHE:CE2	1:A:746:PHE:HE1	2.31	0.48
1:A:749:PHE:CZ	1:A:757:MET:HE3	2.48	0.48
1:D:892:ARG:HH12	1:D:893:GLU:CG	2.25	0.48
1:D:692:SER:O	1:D:696:HIS:HB3	2.13	0.48
1:C:581:GLU:HA	1:C:584:LYS:HD2	1.94	0.48
1:B:866:ILE:O	1:B:869:PRO:HD2	2.14	0.48
1:C:739:ASN:OD1	1:D:714:VAL:HG21	2.13	0.48
1:B:622:MET:O	1:B:626:MET:HG3	2.13	0.48
1:A:914:ASP:C	1:A:916:LEU:N	2.72	0.48
1:C:597:ASP:C	1:C:599:ASN:H	2.20	0.48
1:C:856:PRO:HG3	1:C:902:PHE:CE2	2.48	0.48
1:C:859:GLN:O	1:C:863:MET:HG2	2.13	0.47
1:D:711:SER:HA	1:D:840:MET:HE1	1.96	0.47
1:B:898:VAL:O	1:B:901:LYS:HB2	2.15	0.47
1:C:714:VAL:HG21	1:D:739:ASN:OD1	2.15	0.47
1:D:593:VAL:HA	1:D:596:ILE:HD12	1.97	0.47
1:D:646:VAL:HG13	1:D:698:LEU:HD21	1.95	0.47
1:D:680:TYR:O	1:D:788:TYR:OH	2.17	0.47
1:A:579:ASP:C	1:A:581:GLU:N	2.69	0.47
1:B:714[B]:VAL:HG22	6:B:1249:HOH:O	2.14	0.47
1:C:688:ALA:HB2	1:C:757[A]:MET:SD	2.55	0.47
1:D:627:ASN:ND2	1:D:630:ASN:HB2	2.29	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:626:MET:HG2	1:C:672:TYR:CE2	2.50	0.46
1:A:707:PHE:HB2	1:A:833:GLN:NE2	2.30	0.46
1:A:707:PHE:HZ	1:A:840:MET:HE1	1.77	0.46
1:A:859:GLN:O	1:A:863:MET:HG2	2.15	0.46
1:C:587:HIS:CD2	6:C:1101:HOH:O	2.56	0.46
1:A:856:PRO:HG3	1:A:902:PHE:CE2	2.50	0.46
1:B:673:LYS:HA	1:B:673:LYS:HD2	1.67	0.46
1:C:875:GLN:OE1	1:C:882:ALA:HA	2.15	0.46
1:C:897:LYS:NZ	1:C:897:LYS:HB3	2.31	0.46
1:A:577:ALA:O	1:A:581:GLU:HB2	2.16	0.46
1:B:671:LEU:O	1:B:675:LEU:HB2	2.15	0.46
1:D:592:PRO:HA	6:D:1208:HOH:O	2.16	0.46
1:C:582:TYR:HA	1:C:644:LEU:HD11	1.98	0.46
1:D:603:PHE:HA	1:D:670:LEU:HD13	1.96	0.45
1:D:632:TYR:HB2	1:D:634:ILE:HD12	1.98	0.45
1:D:652:ASP:OD1	6:D:1101:HOH:O	2.21	0.45
1:A:741:HIS:NE2	1:C:797:ARG:HD3	2.32	0.45
1:D:777:PHE:O	1:D:780:LEU:HB2	2.15	0.45
1:D:819:THR:HG21	1:D:895:TRP:HE1	1.82	0.45
1:B:584:LYS:O	1:B:588:ASP:HB2	2.17	0.45
1:B:908:PRO:HG2	1:B:912:SER:O	2.17	0.45
1:A:618:ALA:O	1:A:622:MET:HG3	2.16	0.45
1:C:738:LEU:O	1:C:744:ASN:HB2	2.16	0.45
1:B:633:LYS:NZ	6:B:1119:HOH:O	2.50	0.45
1:B:648:LYS:HG3	1:C:790:ARG:HD3	1.99	0.45
1:C:682:GLU:O	1:C:685:GLU:HB2	2.16	0.45
1:A:747:ASP:C	1:A:747:ASP:OD1	2.60	0.45
1:B:746:PHE:O	1:B:749:PHE:HB2	2.16	0.45
1:D:625:ASP:OD1	1:D:673:LYS:NZ	2.48	0.45
1:D:824:GLU:O	1:D:828[B]:LYS:HD2	2.17	0.45
1:B:577:ALA:HB1	1:B:580:ASP:HB2	1.98	0.45
1:A:746:PHE:C	1:A:748[A]:HIS:H	2.25	0.44
1:B:723:GLY:O	1:B:728:ARG:NH1	2.50	0.44
1:D:780:LEU:HD23	1:D:780:LEU:HA	1.84	0.44
1:A:584:LYS:HB2	6:A:1226:HOH:O	2.17	0.44
1:B:697:ASP:O	1:B:700:HIS:HB2	2.17	0.44
1:C:751:ARG:CG	1:C:751:ARG:NH1	2.51	0.44
1:A:645[B]:MET:HG2	1:A:737:ILE:CG2	2.48	0.44
1:A:914:ASP:C	1:A:916:LEU:H	2.26	0.44
1:A:897:LYS:HE3	1:A:897:LYS:HB3	1.51	0.43
1:C:826:ILE:HD13	4:C:1004:7Y1:C15	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:619:ILE:HD13	1:A:622:MET:HE3	2.00	0.43
1:B:629:ILE:HD13	1:B:639:LEU:CD2	2.48	0.43
1:C:657:ASN:HB2	6:C:1135:HOH:O	2.19	0.43
1:C:653:PRO:HB2	1:C:829:GLU:OE1	2.18	0.43
1:D:685:GLU:CD	1:D:796:HIS:HD1	2.27	0.43
1:D:880:LYS:HB2	6:D:1182:HOH:O	2.18	0.42
1:A:671:LEU:HD13	1:A:803:LEU:HD22	2.01	0.42
1:A:914:ASP:O	1:A:916:LEU:N	2.51	0.42
1:C:704:ASN:O	1:C:708:GLN:HG2	2.19	0.42
1:C:746:PHE:CZ	1:C:757[B]:MET:HE2	2.53	0.42
1:D:755:GLN:NE2	1:D:759:ASP:OD1	2.51	0.42
1:A:626:MET:HG2	1:A:672:TYR:CD2	2.54	0.42
1:A:645[B]:MET:HG2	1:A:737:ILE:HG12	2.01	0.42
1:A:895:TRP:O	1:A:899:SER:HB2	2.19	0.42
1:B:827:TYR:HB3	1:B:831:PHE:CE2	2.54	0.42
1:C:882:ALA:O	1:C:886[B]:GLU:HG3	2.20	0.42
1:D:628:PHE:HD1	1:D:632:TYR:HE2	1.67	0.42
1:D:654:PRO:HG2	1:D:828[A]:LYS:HD3	2.02	0.42
1:C:834:GLY:HA3	1:C:848:MET:O	2.19	0.42
1:A:638:THR:HA	1:A:742:GLY:O	2.20	0.42
1:B:675:LEU:O	6:B:1101:HOH:O	2.21	0.42
1:A:645[B]:MET:HB2	1:A:645[B]:MET:HE3	1.54	0.41
1:A:815:GLY:HA3	6:A:1203:HOH:O	2.19	0.41
1:B:577:ALA:C	1:B:579:ASP:N	2.76	0.41
1:A:821:LYS:O	1:A:821:LYS:HG3	2.19	0.41
1:C:797:ARG:HE	1:C:797:ARG:HB3	1.68	0.41
1:D:827:TYR:HB3	1:D:831:PHE:CE2	2.54	0.41
1:A:606:THR:HA	1:A:607:PRO:HD2	1.74	0.41
1:C:838:LYS:NZ	1:C:849:ASP:OD1	2.42	0.41
1:D:746:PHE:CZ	1:D:757[A]:MET:CG	3.03	0.41
1:C:867:ALA:O	1:C:871:TYR:HD2	2.03	0.41
1:D:627:ASN:HD21	1:D:630:ASN:HB2	1.85	0.41
1:B:892:ARG:NH1	1:B:892:ARG:CG	2.76	0.41
1:A:746:PHE:HB3	1:A:749:PHE:HD2	1.85	0.41
1:B:892:ARG:HG2	1:B:892:ARG:O	2.20	0.41
1:C:735:ILE:O	1:C:739:ASN:HB2	2.21	0.41
1:C:790:ARG:NH2	6:C:1116:HOH:O	2.54	0.41
1:D:674:ASN:O	1:D:880:LYS:HD3	2.21	0.41
1:A:590:ILE:O	1:D:774:LEU:HD13	2.21	0.41
1:C:605:TYR:HD2	1:C:666:HIS:CE1	2.39	0.41
1:D:650:TYR:CE1	1:D:698:LEU:HD23	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:762:ARG:HG2	1:D:766:LEU:HD12	2.03	0.41
1:B:681:LEU:HD21	1:B:799:LEU:HD23	2.03	0.40
1:C:597:ASP:C	1:C:599:ASN:N	2.78	0.40
1:A:585:LEU:HD12	1:A:585:LEU:HA	1.91	0.40
1:A:727:GLU:H	1:A:727:GLU:CD	2.29	0.40
1:D:664:VAL:HG22	1:D:807:CYS:O	2.21	0.40
1:D:632:TYR:OH	1:D:683:ASP:HB2	2.21	0.40
1:A:741:HIS:CD2	1:C:797:ARG:HD3	2.56	0.40
1:B:875:GLN:NE2	1:B:882:ALA:HB1	2.37	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	343/345 (99%)	324 (94%)	16 (5%)	3 (1%)	14	7
1	B	343/345 (99%)	326 (95%)	17 (5%)	0	100	100
1	C	343/345 (99%)	328 (96%)	12 (4%)	3 (1%)	14	7
1	D	329/345 (95%)	313 (95%)	14 (4%)	2 (1%)	21	13
All	All	1358/1380 (98%)	1291 (95%)	59 (4%)	8 (1%)	21	13

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	676	GLU
1	D	786	VAL
1	A	915	PHE
1	A	580	ASP
1	C	584	LYS
1	C	603	PHE

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Mol	Chain	Res	Type
1	D	676	GLU
1	A	908	PRO

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	311/310 (100%)	293 (94%)	18 (6%)	18	11
1	B	311/310 (100%)	290 (93%)	21 (7%)	14	8
1	C	312/310 (101%)	295 (95%)	17 (5%)	20	12
1	D	299/310 (96%)	282 (94%)	17 (6%)	18	11
All	All	1233/1240 (99%)	1160 (94%)	73 (6%)	17	10

All (73) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	578	MET
1	A	579	ASP
1	A	593	VAL
1	A	602	SER
1	A	633	LYS
1	A	701	ARG
1	A	722	GLU
1	A	735	ILE
1	A	786	VAL
1	A	793	LYS
1	A	820	ARG
1	A	822	ILE
1	A	826	ILE
1	A	846	GLU
1	A	887	ARG
1	A	890	SER
1	A	896	THR
1	A	897	LYS
1	B	576	SER

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Mol	Chain	Res	Type
1	B	579	ASP
1	B	593	VAL
1	B	599	ASN
1	B	617	MET
1	B	627	ASN
1	B	629	ILE
1	B	641	ARG
1	B	750	SER
1	B	752	LYS
1	B	756	ARG
1	B	778	LYS
1	B	781	GLN
1	B	822	ILE
1	B	826	ILE
1	B	852	LYS
1	B	857	GLU
1	B	892	ARG
1	B	897	LYS
1	B	909	SER
1	B	918	GLU
1	C	585	LEU
1	C	588	ASP
1	C	597	ASP
1	C	648	LYS
1	C	684	ILE
1	C	739	ASN
1	C	751	ARG
1	C	752	LYS
1	C	782	LYS
1	C	790	ARG
1	C	851	GLU
1	C	875	GLN
1	C	892	ARG
1	C	897	LYS
1	C	899	SER
1	C	911	ASN
1	C	918	GLU
1	D	597	ASP
1	D	617	MET
1	D	634	ILE
1	D	663	SER
1	D	680	TYR

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Mol	Chain	Res	Type
1	D	684	ILE
1	D	739	ASN
1	D	752	LYS
1	D	782	LYS
1	D	790	ARG
1	D	817	LYS
1	D	820	ARG
1	D	822	ILE
1	D	846	GLU
1	D	892	ARG
1	D	909	SER
1	D	910	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	627	ASN
1	A	739	ASN
1	A	755	GLN
1	A	781	GLN
1	A	865	HIS
1	B	630	ASN
1	B	657	ASN
1	B	875	GLN
1	B	894	HIS
1	B	900	HIS
1	C	627	ASN
1	C	781	GLN
1	C	842	ASN
1	C	859	GLN
1	C	911	ASN
1	D	624	GLN
1	D	748	HIS
1	D	859	GLN
1	D	865	HIS
1	D	875	GLN
1	D	894	HIS
1	D	910	ASN

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 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	7Y1	C	1004	-	29,30,30	1.51	6 (20%)	35,42,42	2.56	9 (25%)
4	7Y1	A	1003	-	29,30,30	1.37	4 (13%)	35,42,42	2.25	11 (31%)
4	7Y1	B	1003	-	29,30,30	1.27	2 (6%)	35,42,42	2.73	14 (40%)
4	7Y1	D	1003	-	29,30,30	1.50	6 (20%)	35,42,42	2.56	9 (25%)

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	7Y1	C	1004	-	-	2/14/14/14	0/4/4/4
4	7Y1	A	1003	-	-	5/14/14/14	0/4/4/4
4	7Y1	B	1003	-	-	4/14/14/14	0/4/4/4
4	7Y1	D	1003	-	-	2/14/14/14	0/4/4/4

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	1003	7Y1	C9-N8	3.53	1.38	1.32
4	D	1003	7Y1	N11-N8	3.17	1.41	1.36
4	A	1003	7Y1	C2-C1	3.14	1.53	1.47
4	C	1004	7Y1	C6-N3	-3.06	1.33	1.39
4	C	1004	7Y1	C1-N3	-2.86	1.33	1.39
4	C	1004	7Y1	C9-N8	2.85	1.37	1.32
4	D	1003	7Y1	C15-N12	2.75	1.34	1.31
4	C	1004	7Y1	C15-N12	2.74	1.34	1.31
4	A	1003	7Y1	C15-N12	2.70	1.34	1.31
4	D	1003	7Y1	C6-N3	-2.57	1.34	1.39
4	A	1003	7Y1	C9-N8	2.45	1.36	1.32
4	B	1003	7Y1	C6-N3	-2.44	1.35	1.39
4	B	1003	7Y1	N11-N8	2.42	1.40	1.36
4	C	1004	7Y1	N11-N8	2.34	1.40	1.36
4	D	1003	7Y1	C1-N3	-2.14	1.35	1.39
4	D	1003	7Y1	C22-N11	2.13	1.50	1.46
4	C	1004	7Y1	C22-N11	2.05	1.49	1.46
4	A	1003	7Y1	N11-N8	2.04	1.39	1.36

All (43) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1003	7Y1	C9-C2-C5	10.84	112.36	104.76
4	C	1004	7Y1	C9-C2-C5	9.88	111.68	104.76
4	D	1003	7Y1	C9-C2-C5	9.59	111.48	104.76
4	A	1003	7Y1	C9-C2-C5	7.24	109.84	104.76
4	B	1003	7Y1	C2-C9-N8	-5.69	105.39	112.45
4	D	1003	7Y1	C2-C1-N4	-5.45	117.71	125.25
4	C	1004	7Y1	C2-C9-N8	-5.33	105.83	112.45
4	D	1003	7Y1	C2-C9-N8	-5.13	106.08	112.45
4	C	1004	7Y1	C2-C1-N4	-4.53	118.98	125.25
4	D	1003	7Y1	O18-C10-C13	-4.10	117.86	128.49
4	B	1003	7Y1	O18-C10-C13	-3.92	118.33	128.49
4	A	1003	7Y1	C2-C9-N8	-3.92	107.59	112.45
4	A	1003	7Y1	O18-C10-C13	-3.90	118.39	128.49
4	A	1003	7Y1	C13-N12-C15	3.87	122.47	117.30
4	A	1003	7Y1	C2-C1-N4	-3.56	120.33	125.25
4	A	1003	7Y1	C9-C2-C1	-3.50	116.86	126.82
4	C	1004	7Y1	C9-C2-C1	-3.24	117.59	126.82
4	A	1003	7Y1	C6-C15-N12	-3.11	118.78	123.17
4	C	1004	7Y1	O18-C10-C13	-3.11	120.44	128.49
4	B	1003	7Y1	C6-C15-N12	-3.10	118.80	123.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1003	7Y1	C2-C1-N4	-2.96	121.15	125.25
4	B	1003	7Y1	C9-C2-C1	-2.95	118.41	126.82
4	B	1003	7Y1	C1-N4-N7	-2.87	102.68	106.98
4	D	1003	7Y1	C15-C6-N3	2.82	121.38	117.91
4	A	1003	7Y1	C22-N11-C5	2.78	133.10	129.03
4	D	1003	7Y1	C9-C2-C1	-2.78	118.91	126.82
4	C	1004	7Y1	C6-C15-N12	-2.68	119.39	123.17
4	B	1003	7Y1	C13-N12-C15	2.67	120.87	117.30
4	A	1003	7Y1	C22-N11-N8	-2.62	113.99	119.22
4	C	1004	7Y1	C17-C14-C16	2.56	121.83	118.57
4	A	1003	7Y1	C20-C17-C14	-2.47	118.16	120.80
4	D	1003	7Y1	C6-C15-N12	-2.42	119.75	123.17
4	B	1003	7Y1	N3-C1-N4	2.40	113.70	110.17
4	C	1004	7Y1	C1-N4-N7	-2.40	103.40	106.98
4	D	1003	7Y1	C22-N11-C5	2.37	132.50	129.03
4	B	1003	7Y1	C15-C6-N3	2.30	120.74	117.91
4	C	1004	7Y1	C15-C6-N3	2.27	120.71	117.91
4	B	1003	7Y1	C6-N7-N4	2.26	109.61	106.97
4	A	1003	7Y1	C17-C14-C16	2.25	121.43	118.57
4	B	1003	7Y1	C2-C5-N11	-2.24	102.83	106.63
4	B	1003	7Y1	C22-N11-C5	2.15	132.19	129.03
4	D	1003	7Y1	C23-O18-C10	2.09	119.18	115.94
4	B	1003	7Y1	C22-N11-N8	-2.03	115.17	119.22

There are no chirality outliers.

All (13) torsion outliers are listed below:

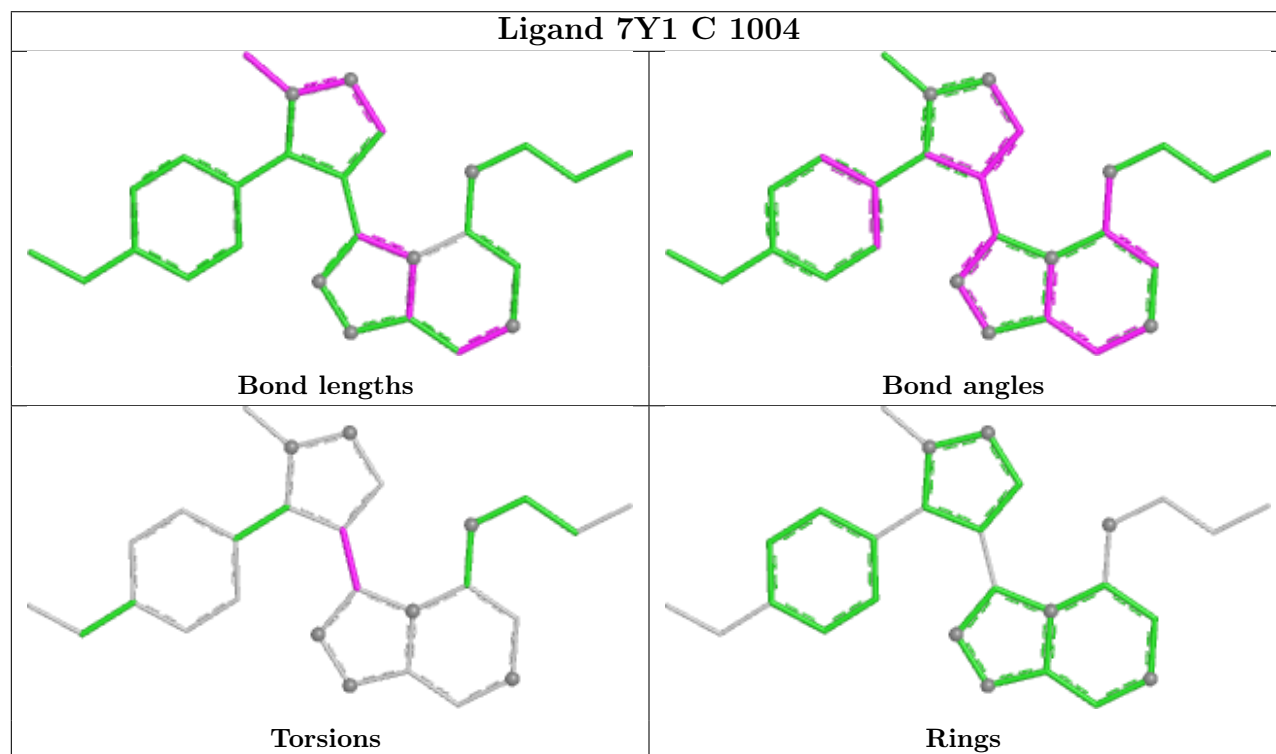
Mol	Chain	Res	Type	Atoms
4	A	1003	7Y1	N3-C10-O18-C23
4	A	1003	7Y1	C13-C10-O18-C23
4	B	1003	7Y1	N3-C10-O18-C23
4	B	1003	7Y1	C13-C10-O18-C23
4	A	1003	7Y1	O18-C23-C25-C27
4	B	1003	7Y1	O18-C23-C25-C27
4	A	1003	7Y1	N3-C1-C2-C9
4	B	1003	7Y1	N3-C1-C2-C9
4	C	1004	7Y1	N3-C1-C2-C9
4	D	1003	7Y1	N3-C1-C2-C9
4	A	1003	7Y1	N4-C1-C2-C9
4	C	1004	7Y1	N4-C1-C2-C9
4	D	1003	7Y1	N4-C1-C2-C9

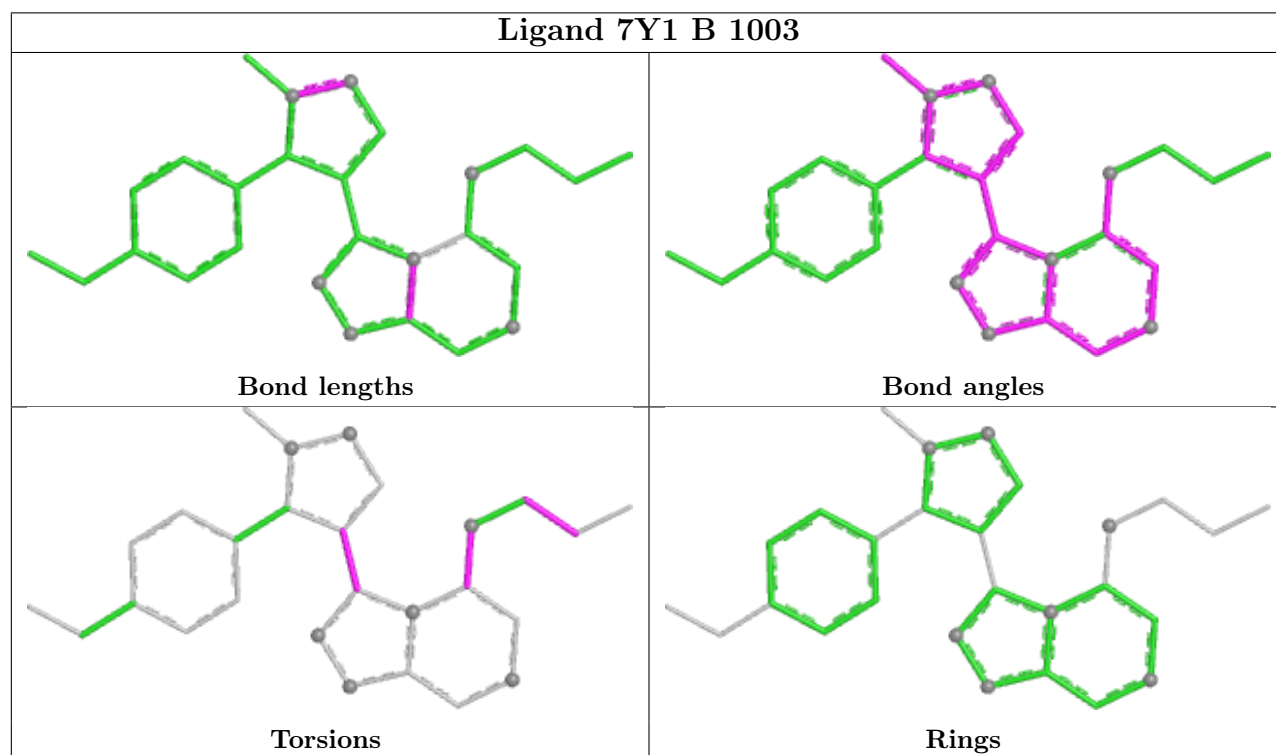
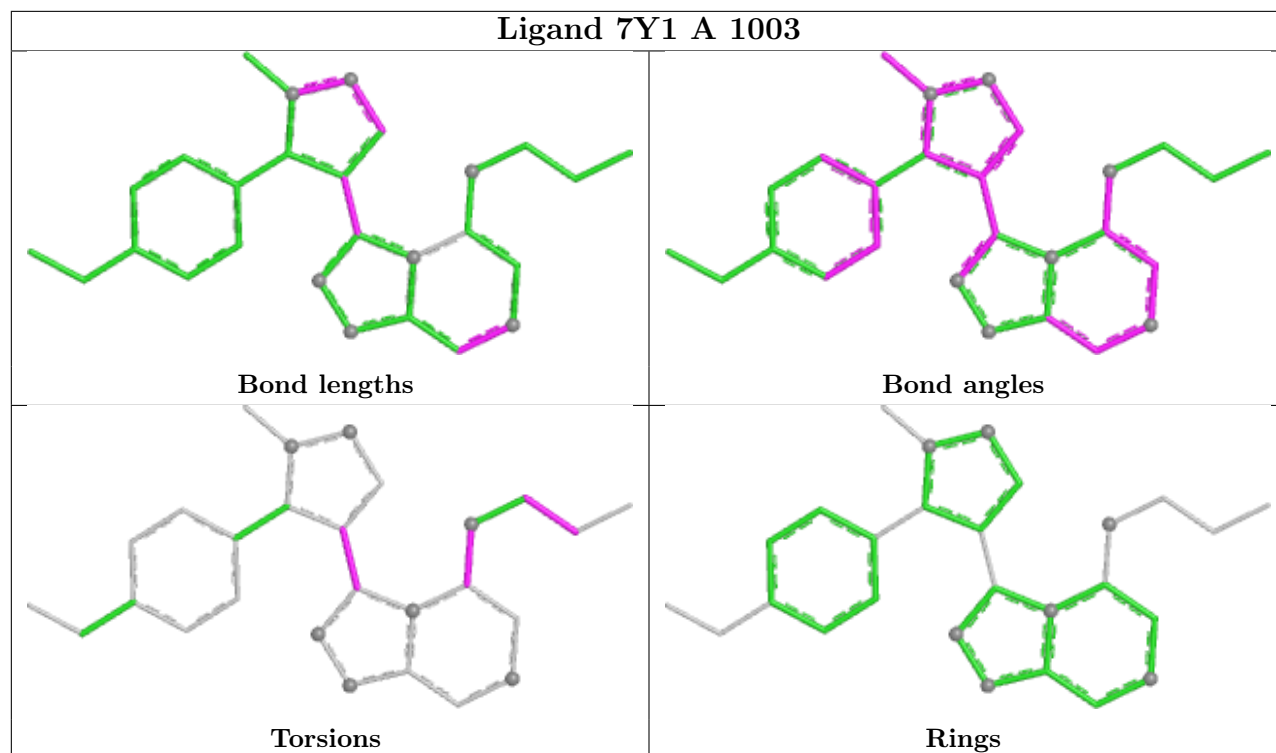
There are no ring outliers.

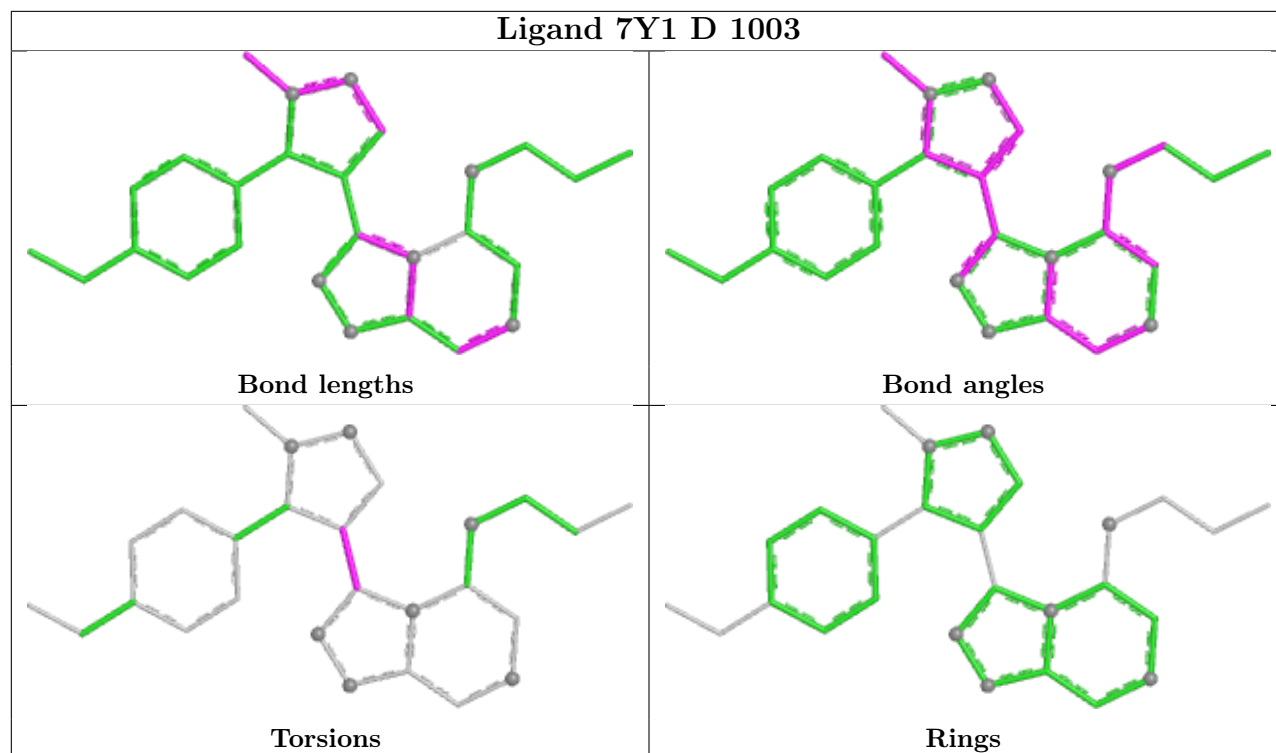
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	1004	7Y1	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 ⓘ

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	340/345 (98%)	-0.01	5 (1%) 72 74	16, 34, 57, 92	7 (2%)
1	B	343/345 (99%)	-0.08	4 (1%) 76 78	17, 32, 53, 85	8 (2%)
1	C	339/345 (98%)	0.04	15 (4%) 39 39	15, 34, 60, 116	13 (3%)
1	D	327/345 (94%)	0.14	5 (1%) 72 74	16, 38, 59, 77	7 (2%)
All	All	1349/1380 (97%)	0.02	29 (2%) 63 66	15, 34, 59, 116	35 (2%)

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	577	ALA	5.4
1	B	577	ALA	4.6
1	C	587	HIS	4.2
1	C	586	LEU	4.1
1	C	582	TYR	3.6
1	C	585	LEU	3.3
1	C	679	ASN	3.3
1	C	680	TYR	3.2
1	C	580	ASP	3.1
1	D	680	TYR	3.0
1	D	593	VAL	2.9
1	D	590	ILE	2.8
1	C	584	LYS	2.8
1	C	851	GLU	2.7
1	C	583	THR	2.7
1	C	581	GLU	2.7
1	A	578	MET	2.5
1	C	617	MET	2.5
1	B	576	SER	2.5
1	B	887	ARG	2.4
1	A	630	ASN	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	652	ASP	2.3
1	C	917	ASP	2.2
1	D	678	THR	2.1
1	C	599	ASN	2.1
1	A	748[A]	HIS	2.1
1	B	591	GLN	2.1
1	C	590	ILE	2.1
1	D	775	ARG	2.0

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

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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.

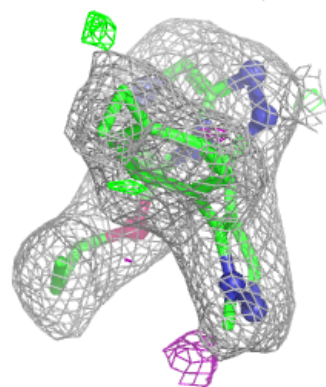
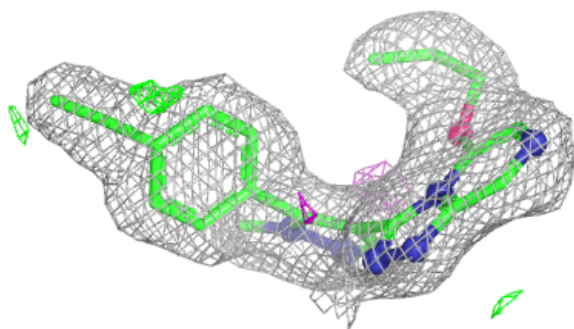
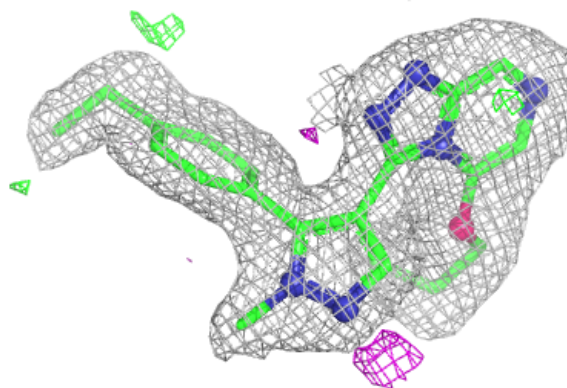
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	7Y1	D	1003	27/27	0.93	0.07	25,30,35,36	0
4	7Y1	C	1004	27/27	0.94	0.07	20,26,30,32	0
4	7Y1	B	1003	27/27	0.94	0.07	18,26,32,32	0
5	CL	C	1003	1/1	0.94	0.11	61,61,61,61	0
4	7Y1	A	1003	27/27	0.95	0.07	23,28,32,38	0
3	MG	B	1002	1/1	0.97	0.03	20,20,20,20	0
3	MG	D	1002	1/1	0.98	0.02	19,19,19,19	0
3	MG	C	1002	1/1	0.98	0.03	15,15,15,15	0
2	ZN	D	1001	1/1	0.99	0.02	29,29,29,29	0
3	MG	A	1002	1/1	0.99	0.02	17,17,17,17	0
2	ZN	B	1001	1/1	1.00	0.01	29,29,29,29	0
2	ZN	C	1001	1/1	1.00	0.01	29,29,29,29	0
2	ZN	A	1001	1/1	1.00	0.01	28,28,28,28	0

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

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

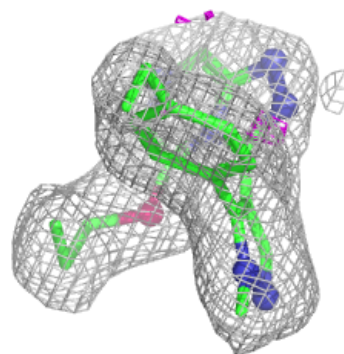
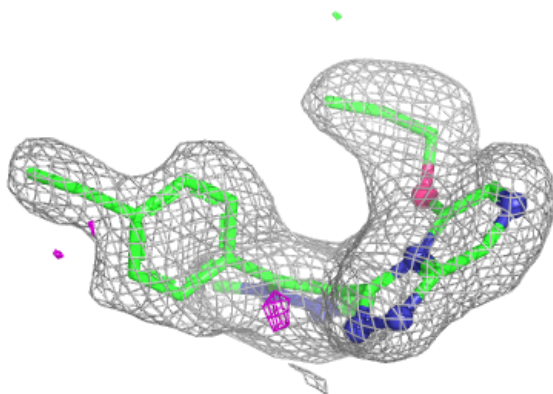
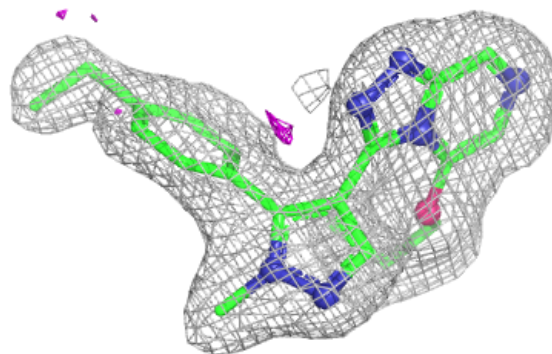
Electron density around 7Y1 D 1003:

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

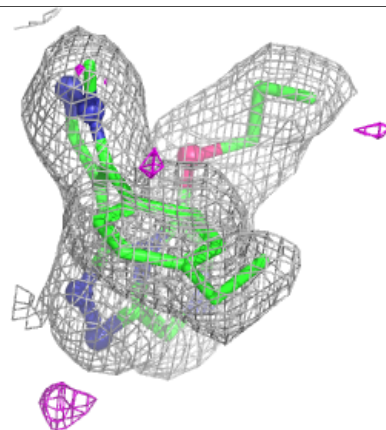
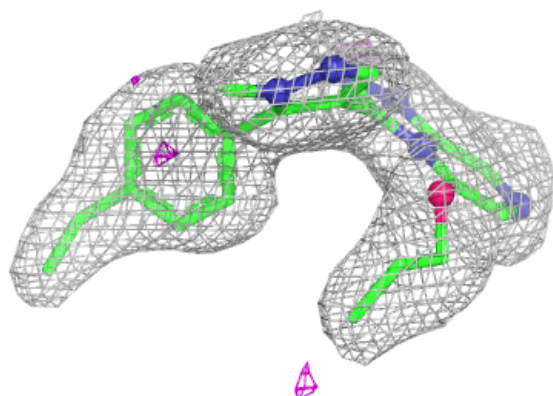
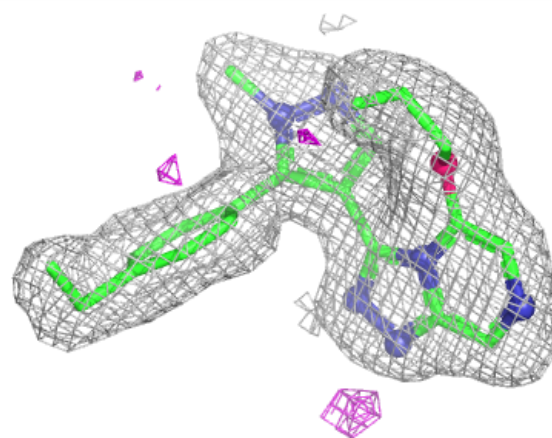


Electron density around 7Y1 C 1004:

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

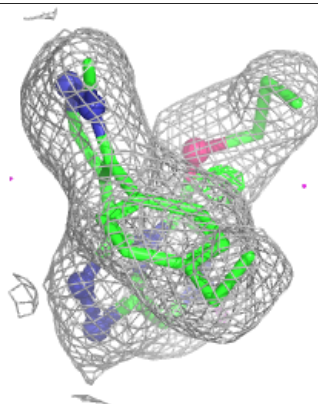
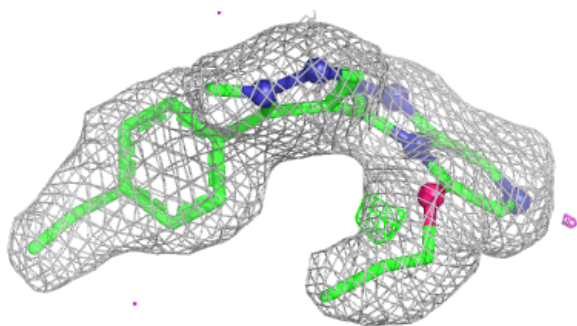
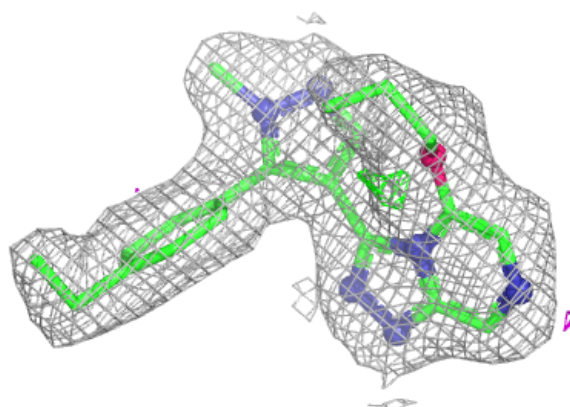
**Electron density around 7Y1 B 1003:**

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



Electron density around 7Y1 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)



6.5 Other polymers [i](#)

There are no such residues in this entry.