



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 9, 2026 – 07:43 PM UTC

PDB ID : 4NXO / pdb_00004nxo
Title : Crystal Structure of Insulin Degrading Enzyme in complex with BDM44768
Authors : Liang, W.G.; Deprez, R.; Deprez, B.; Tang, W.
Deposited on : 2013-12-09
Resolution : 2.73 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 ⓘ](#)) 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

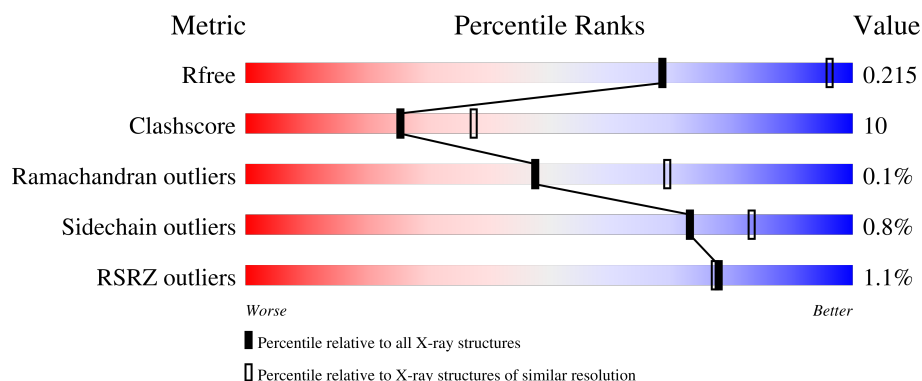
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.73 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	1819 (2.76-2.72)
Clashscore	190562	1866 (2.76-2.72)
Ramachandran outliers	187476	1830 (2.76-2.72)
Sidechain outliers	187428	1831 (2.76-2.72)
RSRZ outliers	180081	1819 (2.76-2.72)

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	990	2% 76% 20%
1	B	990	% 76% 20%

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 16088 atoms, of which 175 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 Insulin-degrading enzyme.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	958	Total	C	N	O	S	0	0	0
			7789	5020	1310	1437	22			
1	B	956	Total	C	N	O	S	0	0	0
			7775	5010	1308	1435	22			

There are 52 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	30	MET	-	expression tag	UNP P14735
A	31	HIS	-	expression tag	UNP P14735
A	32	HIS	-	expression tag	UNP P14735
A	33	HIS	-	expression tag	UNP P14735
A	34	HIS	-	expression tag	UNP P14735
A	35	HIS	-	expression tag	UNP P14735
A	36	HIS	-	expression tag	UNP P14735
A	37	ALA	-	expression tag	UNP P14735
A	38	ALA	-	expression tag	UNP P14735
A	39	GLY	-	expression tag	UNP P14735
A	40	ILE	-	expression tag	UNP P14735
A	41	PRO	-	expression tag	UNP P14735
A	110	LEU	CYS	engineered mutation	UNP P14735
A	111	GLN	GLU	engineered mutation	UNP P14735
A	171	SER	CYS	engineered mutation	UNP P14735
A	178	ALA	CYS	engineered mutation	UNP P14735
A	257	VAL	CYS	engineered mutation	UNP P14735
A	414	LEU	CYS	engineered mutation	UNP P14735
A	573	ASN	CYS	engineered mutation	UNP P14735
A	590	SER	CYS	engineered mutation	UNP P14735
A	789	SER	CYS	engineered mutation	UNP P14735
A	812	ALA	CYS	engineered mutation	UNP P14735
A	819	ALA	CYS	engineered mutation	UNP P14735
A	904	SER	CYS	engineered mutation	UNP P14735
A	966	ASN	CYS	engineered mutation	UNP P14735

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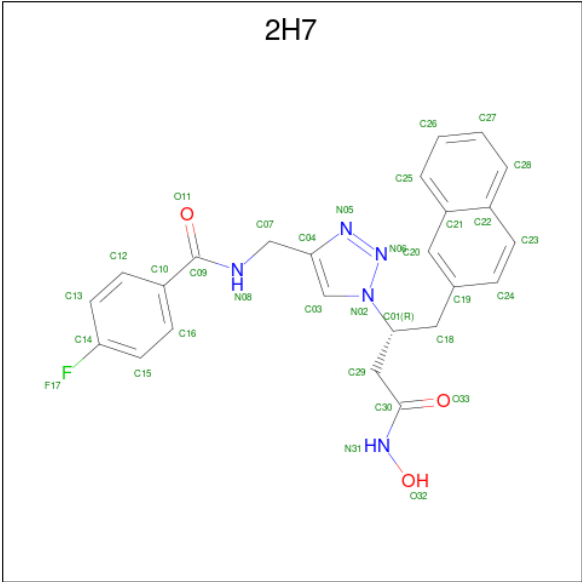
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Chain	Residue	Modelled	Actual	Comment	Reference
A	974	ALA	CYS	engineered mutation	UNP P14735
B	30	MET	-	expression tag	UNP P14735
B	31	HIS	-	expression tag	UNP P14735
B	32	HIS	-	expression tag	UNP P14735
B	33	HIS	-	expression tag	UNP P14735
B	34	HIS	-	expression tag	UNP P14735
B	35	HIS	-	expression tag	UNP P14735
B	36	HIS	-	expression tag	UNP P14735
B	37	ALA	-	expression tag	UNP P14735
B	38	ALA	-	expression tag	UNP P14735
B	39	GLY	-	expression tag	UNP P14735
B	40	ILE	-	expression tag	UNP P14735
B	41	PRO	-	expression tag	UNP P14735
B	110	LEU	CYS	engineered mutation	UNP P14735
B	111	GLN	GLU	engineered mutation	UNP P14735
B	171	SER	CYS	engineered mutation	UNP P14735
B	178	ALA	CYS	engineered mutation	UNP P14735
B	257	VAL	CYS	engineered mutation	UNP P14735
B	414	LEU	CYS	engineered mutation	UNP P14735
B	573	ASN	CYS	engineered mutation	UNP P14735
B	590	SER	CYS	engineered mutation	UNP P14735
B	789	SER	CYS	engineered mutation	UNP P14735
B	812	ALA	CYS	engineered mutation	UNP P14735
B	819	ALA	CYS	engineered mutation	UNP P14735
B	904	SER	CYS	engineered mutation	UNP P14735
B	966	ASN	CYS	engineered mutation	UNP P14735
B	974	ALA	CYS	engineered mutation	UNP P14735

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

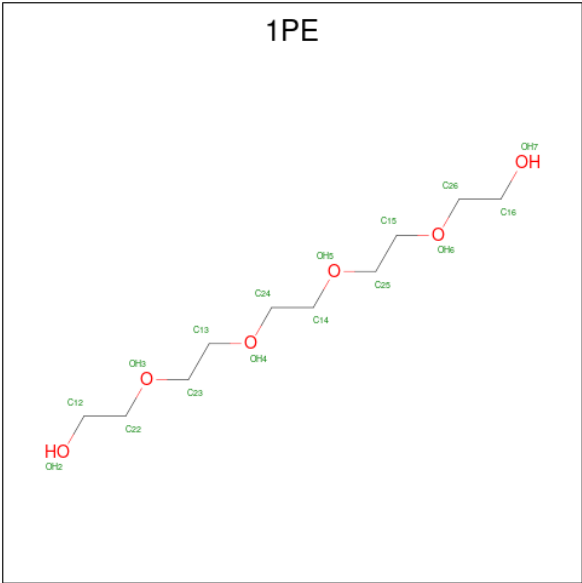
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Zn 1 1	0	0
2	B	1	Total Zn 1 1	0	0

- Molecule 3 is 4-fluoro-N-({1-[(2R)-4-(hydroxyamino)-1-(naphthalen-2-yl)-4-oxobutan-2-yl]-1H-1,2,3-triazol-4-yl}methyl)benzamide (CCD ID: 2H7) (formula: C₂₄H₂₂FN₅O₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	F	N	O	0	0
			33	24	1	5	3		
3	B	1	Total	C	F	N	O	0	0
			33	24	1	5	3		

- Molecule 4 is PENTAETHYLENE GLYCOL (CCD ID: 1PE) (formula: C₁₀H₂₂O₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	H	O	0	0
			38	10	22	6		
4	B	1	Total	C	H	O	0	0
			30	8	17	5		

- Molecule 5 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



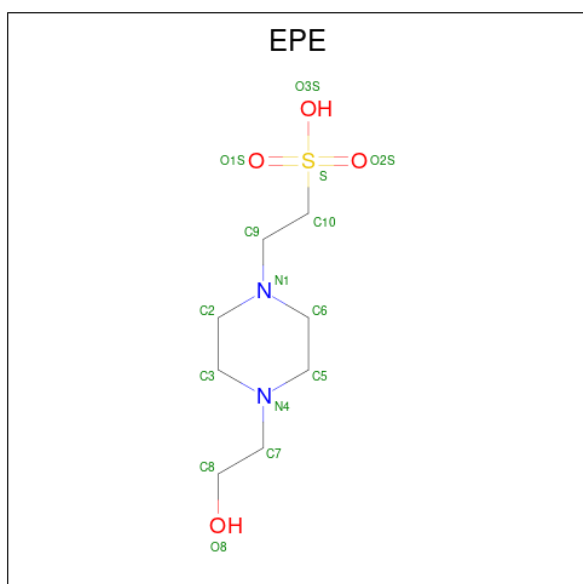
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	H	O	0	0
			17	4	10	3		
5	A	1	Total	C	H	O	0	0
			17	4	10	3		
5	A	1	Total	C	H	O	0	0
			17	4	10	3		
5	B	1	Total	C	H	O	0	0
			17	4	10	3		
5	B	1	Total	C	H	O	0	0
			17	4	10	3		
5	B	1	Total	C	H	O	0	0
			17	4	10	3		

- Molecule 6 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



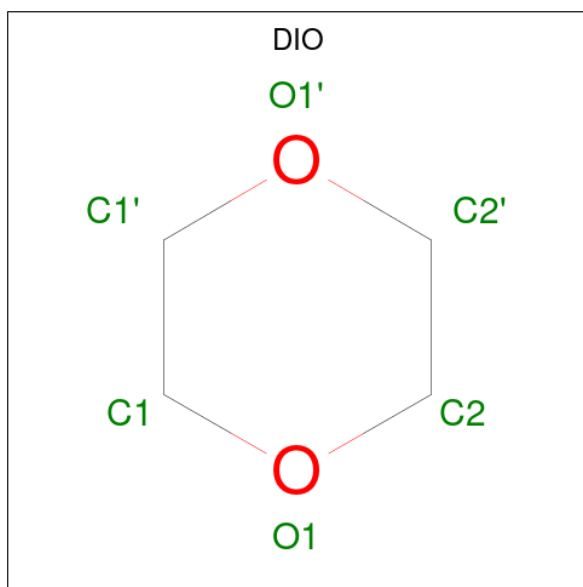
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	H	O	0	0
			10	2	6	2		
6	A	1	Total	C	H	O	0	0
			10	2	6	2		
6	B	1	Total	C	H	O	0	0
			10	2	6	2		
6	B	1	Total	C	H	O	0	0
			10	2	6	2		

- Molecule 7 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (CCD ID: EPE) (formula: $C_8H_{18}N_2O_4S$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
7	A	1	Total	C	H	N	O	S	0	0
			32	8	17	2	4	1		
7	B	1	Total	C	H	N	O	S	0	0
			32	8	17	2	4	1		

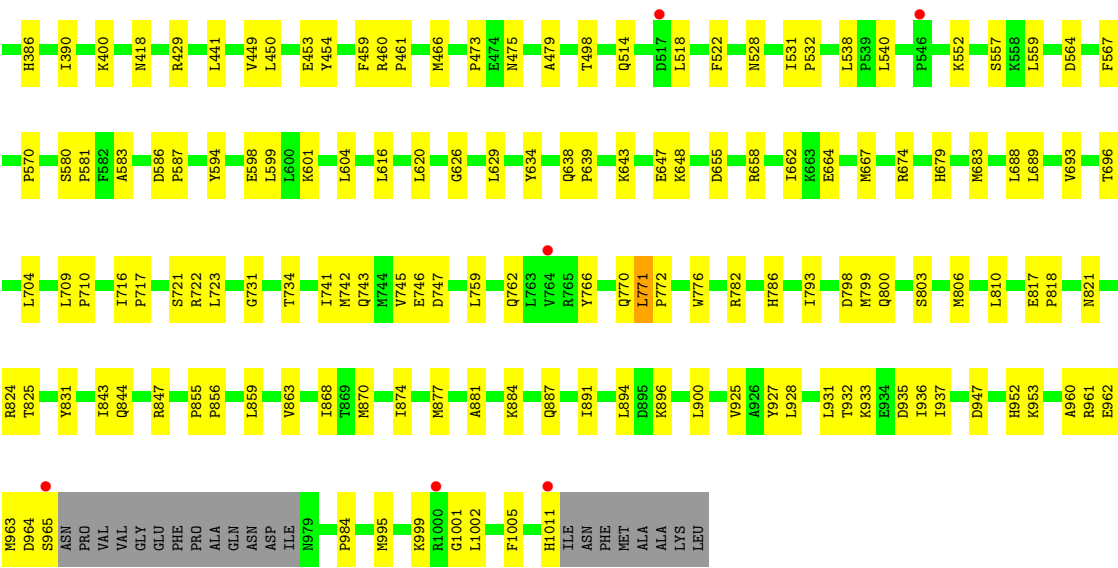
- Molecule 8 is 1,4-DIETHYLENE DIOXIDE (CCD ID: DIO) (formula: C₄H₈O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	B	1	Total	C	H	O	0	0
			14	4	8	2		

- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	62	Total	O	0	0
			62	62		
9	B	89	Total	O	0	0
			89	89		



4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α , β , γ	262.91Å 262.91Å 86.34Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	43.82 – 2.73 43.82 – 2.73	Depositor EDS
% Data completeness (in resolution range)	99.5 (43.82-2.73) 94.4 (43.82-2.73)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.07 (at 2.73Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, R_{free}	0.170 , 0.219 (Not available) , 0.215	Depositor DCC
R_{free} test set	2001 reflections (2.14%)	wwPDB-VP
Wilson B-factor (Å ²)	50.5	Xtriage
Anisotropy	1.077	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 49.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.027 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	16088	wwPDB-VP
Average B, all atoms (Å ²)	65.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.68% 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: ZN, 2H7, EDO, 1PE, PEG, EPE, DIO

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	0.31	0/7985	0.77	8/10810 (0.1%)
1	B	0.32	0/7970	0.75	4/10788 (0.0%)
All	All	0.31	0/15955	0.76	12/21598 (0.1%)

There are no bond length outliers.

The worst 5 of 12 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	756	LYS	CA-C-N	6.89	127.29	119.92
1	A	756	LYS	C-N-CA	6.89	127.29	119.92
1	B	771	LEU	CA-C-N	6.58	126.49	119.85
1	B	771	LEU	C-N-CA	6.58	126.49	119.85
1	A	210	ASN	CA-C-N	6.35	126.04	119.56

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	7789	0	7701	158	0
1	B	7775	0	7685	144	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	33	0	21	5	0
3	B	33	0	20	6	0
4	A	16	22	22	0	0
4	B	13	17	17	1	0
5	A	21	30	30	1	0
5	B	28	40	40	3	0
6	A	8	12	12	1	0
6	B	8	12	12	2	0
7	A	15	17	17	0	0
7	B	15	17	17	1	0
8	B	6	8	8	0	0
9	A	62	0	0	2	0
9	B	89	0	0	2	0
All	All	15913	175	15602	302	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 10.

The worst 5 of 302 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:688:LEU:HD13	1:B:696:THR:HG22	1.48	0.95
3:A:1102:2H7:H4	3:A:1102:2H7:H8	1.49	0.94
3:B:1102:2H7:H4	3:B:1102:2H7:C03	1.97	0.93
1:B:557:SER:HB2	1:B:742:MET:HE3	1.50	0.92
1:B:870:MET:HA	1:B:870:MET:HE2	1.53	0.90

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	954/990 (96%)	934 (98%)	19 (2%)	1 (0%)	48	69
1	B	952/990 (96%)	935 (98%)	17 (2%)	0	100	100
All	All	1906/1980 (96%)	1869 (98%)	36 (2%)	1 (0%)	48	69

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	967	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	842/879 (96%)	835 (99%)	7 (1%)	73	83
1	B	840/879 (96%)	833 (99%)	7 (1%)	73	83
All	All	1682/1758 (96%)	1668 (99%)	14 (1%)	73	83

5 of 14 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	110	LEU
1	B	111	GLN
1	B	1011	HIS
1	B	887	GLN
1	B	891	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 13 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	988	GLN
1	B	376	ASN
1	B	982	GLN
1	B	780	GLN
1	B	788	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 20 ligands modelled in this entry, 2 are monoatomic - leaving 18 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	2H7	A	1102	2	36,36,36	2.43	7 (19%)	45,49,49	1.84	10 (22%)
6	EDO	B	1107	-	3,3,3	0.42	0	2,2,2	0.35	0
8	DIO	B	1111	-	6,6,6	0.45	0	6,6,6	0.68	0
5	PEG	B	1109	-	6,6,6	0.67	0	5,5,5	0.28	0
4	1PE	B	1103	-	12,12,15	0.68	0	11,11,14	0.33	0
5	PEG	A	1104	-	6,6,6	0.67	0	5,5,5	0.20	0
5	PEG	B	1105	-	6,6,6	0.69	0	5,5,5	0.51	0
5	PEG	A	1105	-	6,6,6	0.62	0	5,5,5	0.34	0
3	2H7	B	1102	2	36,36,36	2.40	12 (33%)	45,49,49	2.54	13 (28%)
6	EDO	A	1108	-	3,3,3	0.47	0	2,2,2	0.31	0
6	EDO	B	1108	-	3,3,3	0.49	0	2,2,2	0.26	0
7	EPE	A	1109	-	15,15,15	0.86	1 (6%)	19,20,20	1.80	4 (21%)
5	PEG	B	1106	-	6,6,6	0.65	0	5,5,5	0.35	0
5	PEG	A	1106	-	6,6,6	0.66	0	5,5,5	0.34	0
4	1PE	A	1103	-	15,15,15	0.66	0	14,14,14	0.41	0
7	EPE	B	1110	-	15,15,15	0.77	1 (6%)	19,20,20	1.76	4 (21%)
6	EDO	A	1107	-	3,3,3	0.46	0	2,2,2	0.33	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	PEG	B	1104	-	6,6,6	0.65	0	5,5,5	0.24	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
3	2H7	A	1102	2	-	10/23/23/23	0/4/4/4
6	EDO	B	1107	-	-	1/1/1/1	-
8	DIO	B	1111	-	-	-	0/1/1/1
5	PEG	B	1109	-	-	2/4/4/4	-
4	1PE	B	1103	-	-	5/10/10/13	-
5	PEG	A	1104	-	-	1/4/4/4	-
5	PEG	B	1105	-	-	3/4/4/4	-
5	PEG	A	1105	-	-	2/4/4/4	-
3	2H7	B	1102	2	-	9/23/23/23	0/4/4/4
6	EDO	A	1108	-	-	0/1/1/1	-
6	EDO	B	1108	-	-	0/1/1/1	-
7	EPE	A	1109	-	-	4/9/19/19	0/1/1/1
5	PEG	B	1106	-	-	3/4/4/4	-
5	PEG	A	1106	-	-	3/4/4/4	-
4	1PE	A	1103	-	-	7/13/13/13	-
7	EPE	B	1110	-	-	4/9/19/19	0/1/1/1
6	EDO	A	1107	-	-	0/1/1/1	-
5	PEG	B	1104	-	-	1/4/4/4	-

The worst 5 of 21 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1102	2H7	C30-N31	9.68	1.42	1.32
3	B	1102	2H7	C30-N31	7.97	1.41	1.32
3	A	1102	2H7	C09-N08	5.77	1.46	1.33
3	B	1102	2H7	C09-N08	4.93	1.44	1.33
3	A	1102	2H7	C03-C04	4.40	1.43	1.36

The worst 5 of 31 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1102	2H7	C19-C18-C01	-8.37	97.05	113.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1102	2H7	C01-N02-N06	5.61	128.55	120.94
3	B	1102	2H7	C29-C30-N31	5.47	123.42	115.14
3	B	1102	2H7	C01-N02-N06	5.33	128.16	120.94
3	B	1102	2H7	C07-C04-N05	5.15	131.74	121.61

There are no chirality outliers.

5 of 55 torsion outliers are listed below:

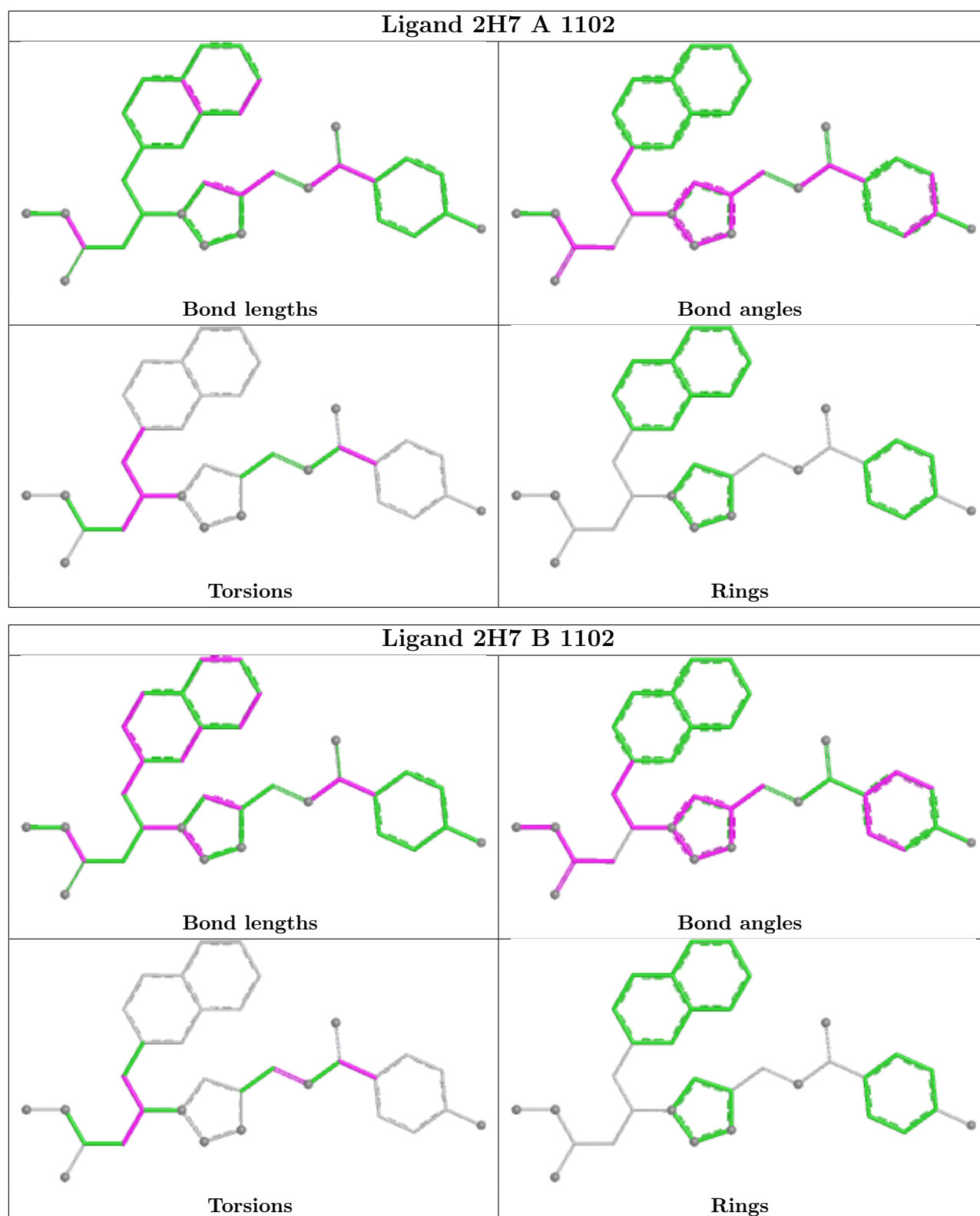
Mol	Chain	Res	Type	Atoms
3	A	1102	2H7	N02-C01-C18-C19
3	B	1102	2H7	N02-C01-C18-C19
3	B	1102	2H7	N02-C01-C29-C30
3	B	1102	2H7	C18-C01-C29-C30
7	A	1109	EPE	C10-C9-N1-C6

There are no ring outliers.

9 monomers are involved in 20 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1102	2H7	5	0
4	B	1103	1PE	1	0
5	B	1105	PEG	1	0
5	A	1105	PEG	1	0
3	B	1102	2H7	6	0
6	A	1108	EDO	1	0
6	B	1108	EDO	2	0
7	B	1110	EPE	1	0
5	B	1104	PEG	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	958/990 (96%)	0.04	15 (1%) 70 68	45, 66, 90, 131	0
1	B	956/990 (96%)	-0.15	6 (0%) 85 85	41, 59, 82, 120	0
All	All	1914/1980 (96%)	-0.06	21 (1%) 78 77	41, 62, 88, 131	0

The worst 5 of 21 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	967	PRO	5.1
1	A	968	VAL	4.3
1	A	764	VAL	4.3
1	B	965	SER	3.9
1	A	582	PHE	3.0

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](#)

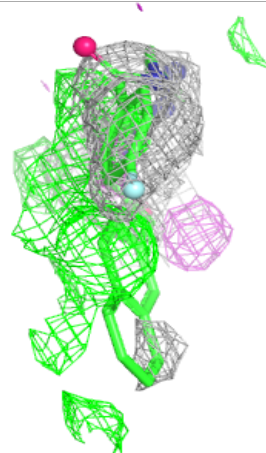
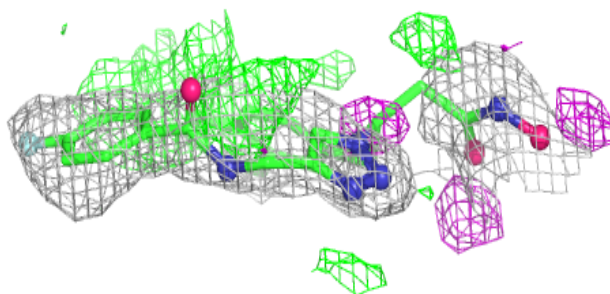
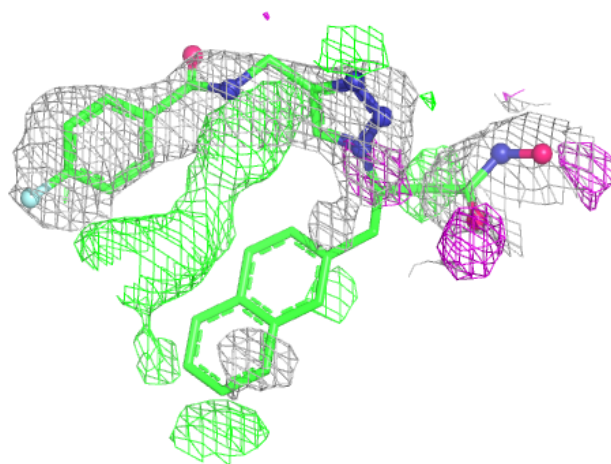
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists 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
6	EDO	A	1108	4/4	0.79	0.18	82,99,109,123	0
4	1PE	A	1103	16/16	0.80	0.18	61,99,112,119	0
3	2H7	A	1102	33/33	0.80	0.27	51,83,103,116	7
8	DIO	B	1111	6/6	0.83	0.20	78,95,126,126	0
3	2H7	B	1102	33/33	0.84	0.23	57,85,99,104	8
6	EDO	B	1108	4/4	0.84	0.19	60,80,102,102	0
4	1PE	B	1103	13/16	0.84	0.15	72,98,118,124	0
6	EDO	A	1107	4/4	0.85	0.17	70,94,113,113	0
5	PEG	B	1106	7/7	0.85	0.19	74,101,127,127	0
5	PEG	A	1106	7/7	0.86	0.16	70,88,102,108	0
5	PEG	B	1104	7/7	0.87	0.14	77,97,123,123	0
5	PEG	A	1105	7/7	0.87	0.17	73,90,108,116	0
7	EPE	B	1110	15/15	0.92	0.11	67,87,111,119	0
6	EDO	B	1107	4/4	0.93	0.14	64,77,87,89	0
5	PEG	B	1105	7/7	0.93	0.17	56,82,102,107	0
5	PEG	A	1104	7/7	0.94	0.10	60,84,110,110	0
5	PEG	B	1109	7/7	0.94	0.10	56,77,104,104	0
7	EPE	A	1109	15/15	0.95	0.10	67,87,105,113	0
2	ZN	B	1101	1/1	0.99	0.04	55,55,55,55	0
2	ZN	A	1101	1/1	1.00	0.02	58,58,58,58	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.

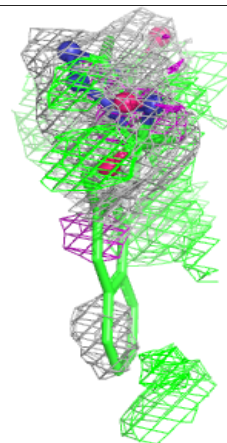
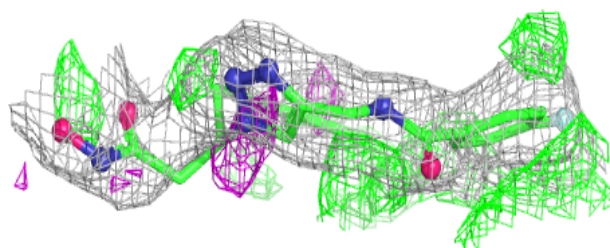
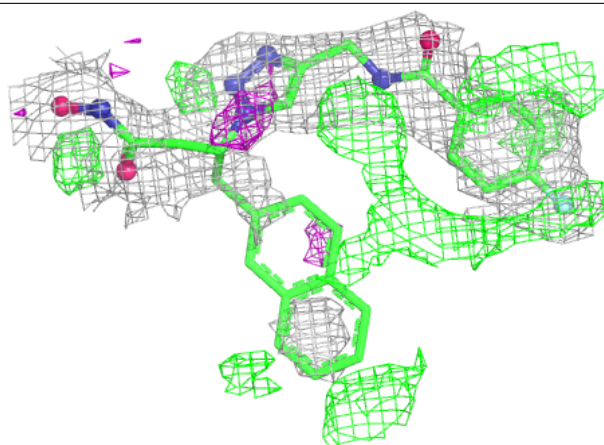
Electron density around 2H7 A 1102:

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



Electron density around 2H7 B 1102:

$2mF_o-DF_c$ (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.