



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 6, 2026 – 04:56 AM UTC

PDB ID : 2I47 / pdb_00002i47
Title : Crystal structure of catalytic domain of TACE with inhibitor
Authors : Xu, W.; Condon, J.S.; Lovering, F.E.
Deposited on : 2006-08-21
Resolution : 1.90 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

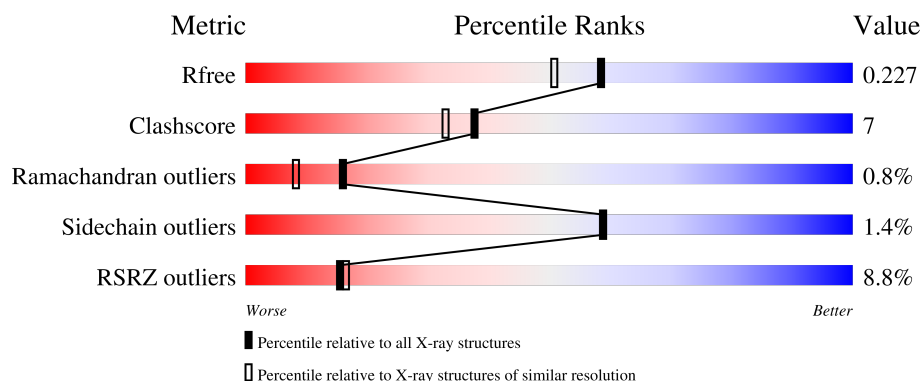
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	288	9% 70% 16% • 11%
1	B	288	6% 73% 14% • 12%
1	C	288	9% 69% 16% • 12%
1	D	288	7% 72% 16% • 11%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 8926 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 ADAM 17.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	255	Total 2023	C 1273	N 342	O 395	S 13	0	0	0
1	B	254	Total 2016	C 1268	N 341	O 394	S 13	0	0	0
1	C	254	Total 2016	C 1268	N 341	O 394	S 13	0	0	0
1	D	255	Total 2023	C 1273	N 342	O 395	S 13	0	0	0

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	266	ALA	SER	engineered mutation	UNP P78536
A	452	GLN	ASN	engineered mutation	UNP P78536
A	493	SER	-	cloning artifact	UNP P78536
A	494	HIS	-	expression tag	UNP P78536
A	495	HIS	-	expression tag	UNP P78536
A	496	HIS	-	expression tag	UNP P78536
A	497	HIS	-	expression tag	UNP P78536
A	498	HIS	-	expression tag	UNP P78536
A	499	HIS	-	expression tag	UNP P78536
B	266	ALA	SER	engineered mutation	UNP P78536
B	452	GLN	ASN	engineered mutation	UNP P78536
B	493	SER	-	cloning artifact	UNP P78536
B	494	HIS	-	expression tag	UNP P78536
B	495	HIS	-	expression tag	UNP P78536
B	496	HIS	-	expression tag	UNP P78536
B	497	HIS	-	expression tag	UNP P78536
B	498	HIS	-	expression tag	UNP P78536
B	499	HIS	-	expression tag	UNP P78536
C	266	ALA	SER	engineered mutation	UNP P78536
C	452	GLN	ASN	engineered mutation	UNP P78536
C	493	SER	-	cloning artifact	UNP P78536

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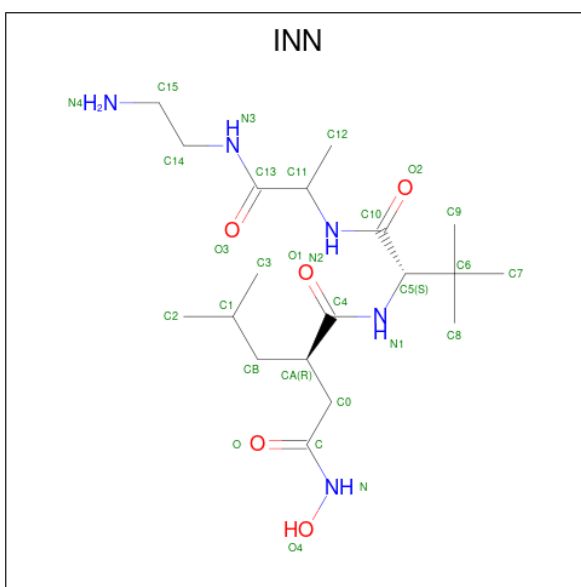
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Chain	Residue	Modelled	Actual	Comment	Reference
C	494	HIS	-	expression tag	UNP P78536
C	495	HIS	-	expression tag	UNP P78536
C	496	HIS	-	expression tag	UNP P78536
C	497	HIS	-	expression tag	UNP P78536
C	498	HIS	-	expression tag	UNP P78536
C	499	HIS	-	expression tag	UNP P78536
D	266	ALA	SER	engineered mutation	UNP P78536
D	452	GLN	ASN	engineered mutation	UNP P78536
D	493	SER	-	cloning artifact	UNP P78536
D	494	HIS	-	expression tag	UNP P78536
D	495	HIS	-	expression tag	UNP P78536
D	496	HIS	-	expression tag	UNP P78536
D	497	HIS	-	expression tag	UNP P78536
D	498	HIS	-	expression tag	UNP P78536
D	499	HIS	-	expression tag	UNP P78536

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

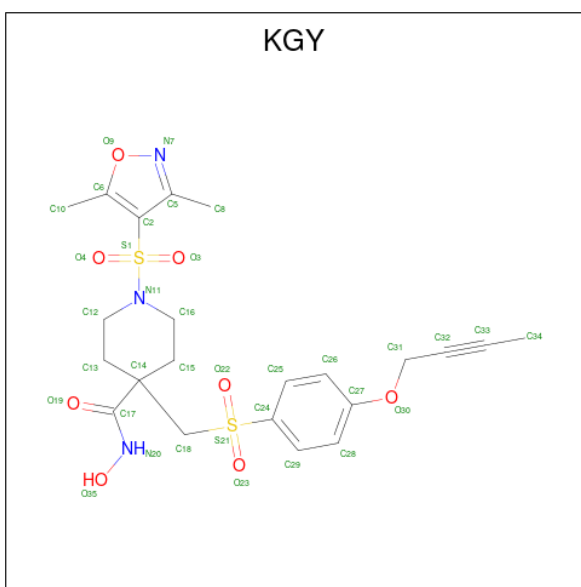
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Zn 1 1	0	0
2	B	1	Total Zn 1 1	0	0
2	C	1	Total Zn 1 1	0	0
2	D	1	Total Zn 1 1	0	0

- Molecule 3 is N-{(2R)-2-[2-(hydroxyamino)-2-oxoethyl]-4-methylpentanoyl}-3-methyl-L-valyl-N-(2-aminoethyl)-L-alaninamide (CCD ID: INN) (formula: C₁₉H₃₇N₅O₅).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			29	19	5	5		
3	B	1	Total	C	N	O	0	0
			29	19	5	5		

- Molecule 4 is 4-({[4-(BUT-2-YN-1-YLOXY)PHENYL]SULFONYL}METHYL)-1-[(3,5-DIMETHYLISOXAZOL-4-YL)SULFONYL]-N-HYDROXYPIPERIDINE-4-CARBOXAMIDE (CCD ID: KGY) (formula: C₂₂H₂₇N₃O₈S₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	C	1	Total	C	N	O	S	0	0
			35	22	3	8	2		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	D	1	Total	C	N	O	S	0	0
			35	22	3	8	2		

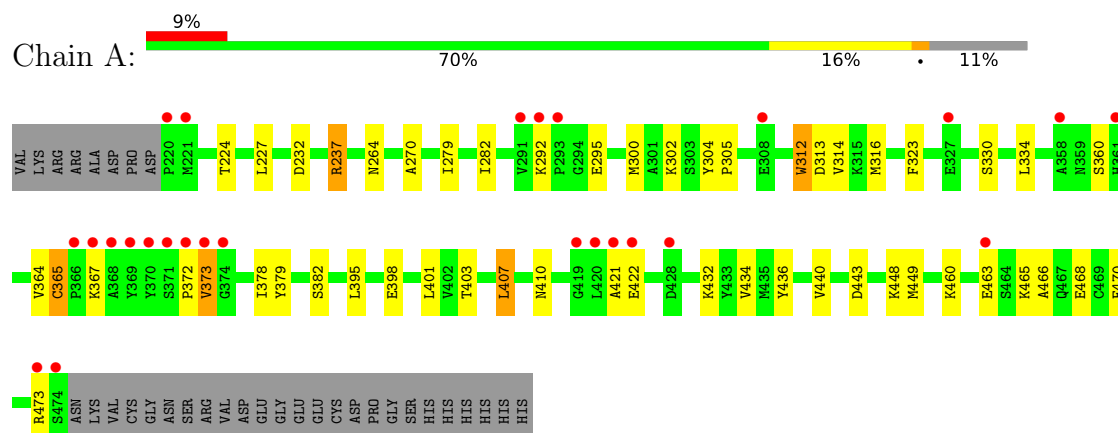
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	213	Total	O	0	0
			213	213		
5	B	165	Total	O	0	0
			165	165		
5	C	169	Total	O	0	0
			169	169		
5	D	169	Total	O	0	0
			169	169		

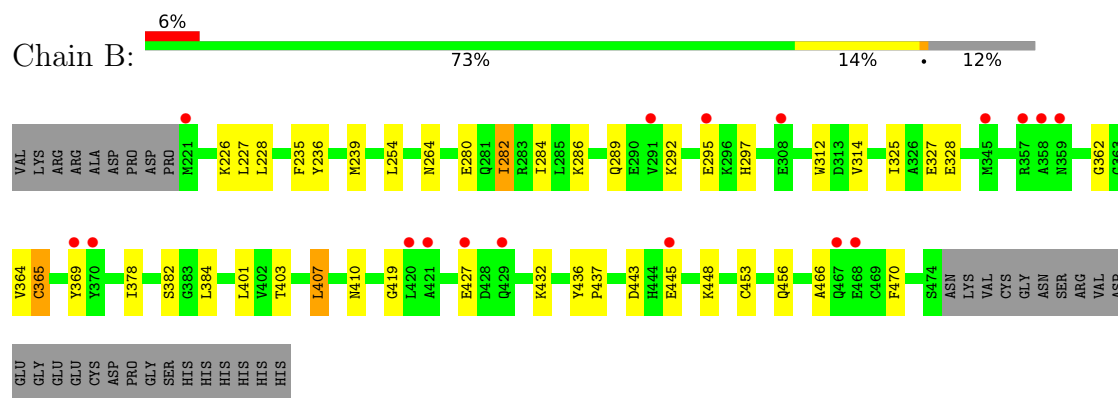
3 Residue-property plots [i](#)

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.

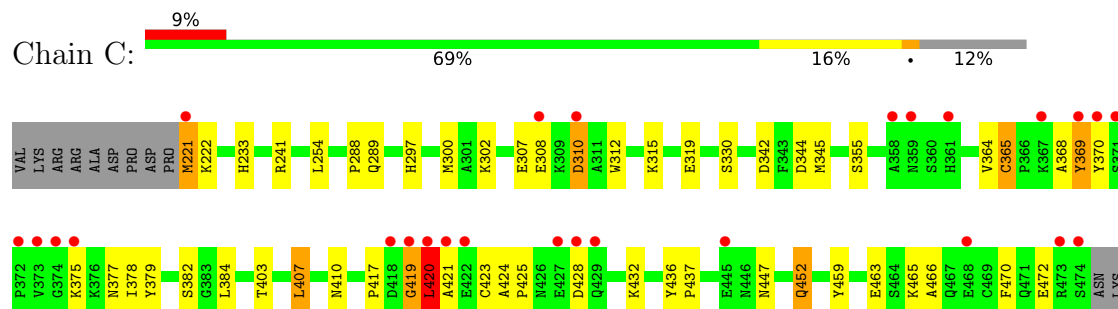
• Molecule 1: ADAM 17



• Molecule 1: ADAM 17



• Molecule 1: ADAM 17



VAL
CYS
GLY
ASN
ARG
SER
VAL
ASP
GLU
GLY
GLU
GLY
CYS
ASP
PRO
GLY
SER
HIS
HIS
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HIS
HIS

● Molecule 1: ADAM 17



VAL	LYS	ARG	ALA	ASP	PRO	ASP	P220	M221	K226	R234	G240	T247	L254	G271	I279	E280	Q281	I282	Q289	H297	K300	Y304	P305	N306	W312	D313	V314	K315	K316	E319	E327	Q341	V353	R357	A358	N359	S360	H361	V364	C365
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P366	K367	P372	V373	G374	L396	T396	K397	E398	L401	V402	T403	L407	E414	D418	G419	L420	A421	E422	G423	A424	P425	N426	E427	D428	Y436	P437	V440	D443	H444	E445	N446	N447	Q456	K460	A466	Q467	E468	C469	F470	Q471	E472	R473	S474	ASN	LYS	VAL	CYS
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GLY
ASN
SER
ARG
VAL
ASP
GLU
GLY
GLU
GLU
CYS
ASP
PRO
GLY
SER
HIS
HIS
HIS
HIS

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	61.77Å 126.19Å 81.22Å 90.00° 107.41° 90.00°	Depositor
Resolution (Å)	20.00 – 1.90 20.00 – 1.90	Depositor EDS
% Data completeness (in resolution range)	(Not available) (20.00-1.90) 96.0 (20.00-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.21 (at 1.81Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.198 , 0.227 0.199 , 0.227	Depositor DCC
R_{free} test set	5164 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	15.3	Xtriage
Anisotropy	0.290	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 41.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8926	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 3.73% 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: KGY, INN, 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	0.64	0/2069	0.98	8/2791 (0.3%)
1	B	0.64	0/2061	0.96	8/2780 (0.3%)
1	C	0.61	1/2061 (0.0%)	0.94	5/2780 (0.2%)
1	D	0.63	0/2069	0.94	5/2791 (0.2%)
All	All	0.63	1/8260 (0.0%)	0.95	26/11142 (0.2%)

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	A	0	1
1	B	0	2
1	C	0	2
1	D	0	1
All	All	0	6

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	417	PRO	CA-C	-5.28	1.47	1.52

The worst 5 of 26 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	221	MET	N-CA-C	-7.29	104.41	112.72
1	C	410	ASN	CA-CB-CG	6.85	119.45	112.60
1	B	314	VAL	N-CA-C	6.72	117.48	110.62
1	B	410	ASN	CA-CB-CG	6.49	119.09	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	428	ASP	CA-CB-CG	-6.29	106.31	112.60

There are no chirality outliers.

5 of 6 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	312	TRP	Peptide
1	B	227	LEU	Peptide
1	B	312	TRP	Peptide
1	C	312	TRP	Peptide
1	C	419	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	2023	0	1940	32	0
1	B	2016	0	1932	20	0
1	C	2016	0	1932	34	0
1	D	2023	0	1940	29	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	29	0	36	0	0
3	B	29	0	36	0	0
4	C	35	0	26	0	0
4	D	35	0	26	1	0
5	A	213	0	0	3	0
5	B	165	0	0	0	0
5	C	169	0	0	2	0
5	D	169	0	0	3	0
All	All	8926	0	7868	113	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 7.

The worst 5 of 113 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:372:PRO:O	1:A:373:VAL:HG23	1.64	0.97
1:B:282:ILE:HD11	1:B:284:ILE:HD11	1.62	0.79
1:A:372:PRO:O	1:A:373:VAL:CG2	2.32	0.78
1:B:466:ALA:O	1:B:470:PHE:HB2	1.97	0.65
1:A:372:PRO:O	1:A:373:VAL:CB	2.47	0.61

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	253/288 (88%)	239 (94%)	11 (4%)	3 (1%)	10	4
1	B	252/288 (88%)	241 (96%)	10 (4%)	1 (0%)	30	22
1	C	252/288 (88%)	243 (96%)	7 (3%)	2 (1%)	16	8
1	D	253/288 (88%)	243 (96%)	8 (3%)	2 (1%)	16	8
All	All	1010/1152 (88%)	966 (96%)	36 (4%)	8 (1%)	16	8

5 of 8 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	373	VAL
1	C	420	LEU
1	A	312	TRP
1	A	365	CYS
1	B	365	CYS

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	218/247 (88%)	217 (100%)	1 (0%)	81	84
1	B	217/247 (88%)	215 (99%)	2 (1%)	70	73
1	C	217/247 (88%)	210 (97%)	7 (3%)	34	27
1	D	218/247 (88%)	216 (99%)	2 (1%)	70	73
All	All	870/988 (88%)	858 (99%)	12 (1%)	59	59

5 of 12 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	407	LEU
1	C	420	LEU
1	D	422	GLU
1	C	452	GLN
1	C	221	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 16 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	452	GLN
1	D	361	HIS
1	C	269	ASN
1	D	320	GLN
1	B	471	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 4 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	KGY	D	1002	2	34,37,37	1.85	8 (23%)	45,55,55	2.18	16 (35%)
3	INN	A	901	2	28,28,28	0.55	0	35,38,38	1.34	4 (11%)
4	KGY	C	1001	2	34,37,37	1.88	9 (26%)	45,55,55	2.23	13 (28%)
3	INN	B	902	2	28,28,28	0.75	1 (3%)	35,38,38	1.32	4 (11%)

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	KGY	D	1002	2	-	8/34/49/49	0/3/3/3
3	INN	A	901	2	-	0/40/40/40	-
4	KGY	C	1001	2	-	10/34/49/49	0/3/3/3
3	INN	B	902	2	-	0/40/40/40	-

The worst 5 of 18 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	1001	KGY	O9-N7	-5.83	1.31	1.42
4	D	1002	KGY	O9-N7	-5.81	1.31	1.42
4	D	1002	KGY	C28-C27	3.72	1.45	1.38
4	C	1001	KGY	C28-C27	3.52	1.45	1.38
4	C	1001	KGY	C25-C24	3.13	1.43	1.38

The worst 5 of 37 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	1002	KGY	C16-N11-C12	7.09	120.27	112.12
4	C	1001	KGY	C16-N11-C12	5.22	118.12	112.12
4	C	1001	KGY	C13-C14-C18	-5.04	102.46	110.50
3	A	901	INN	C14-N3-C13	-4.97	113.61	122.55
4	C	1001	KGY	C12-C13-C14	4.84	114.87	109.17

There are no chirality outliers.

5 of 18 torsion outliers are listed below:

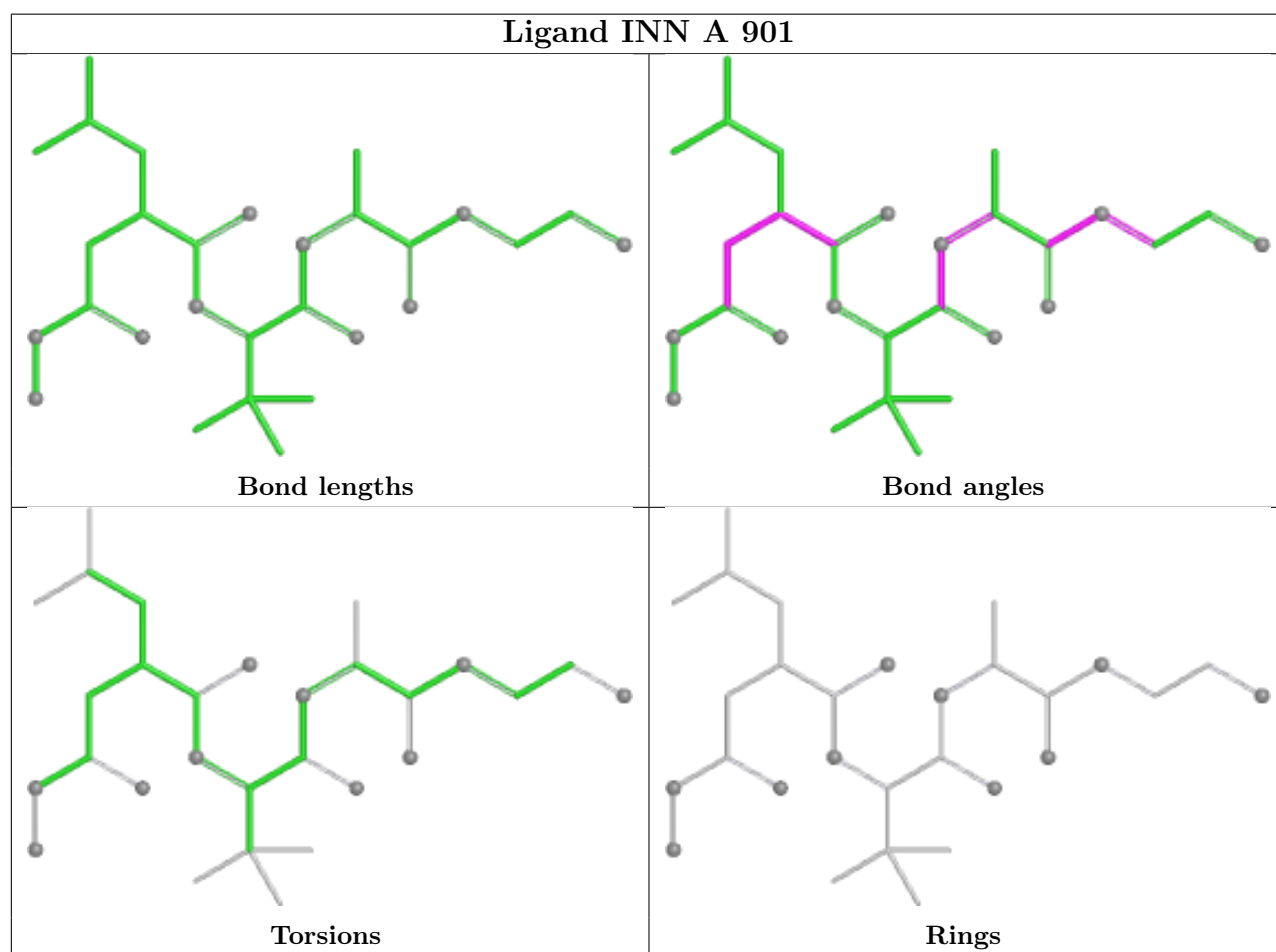
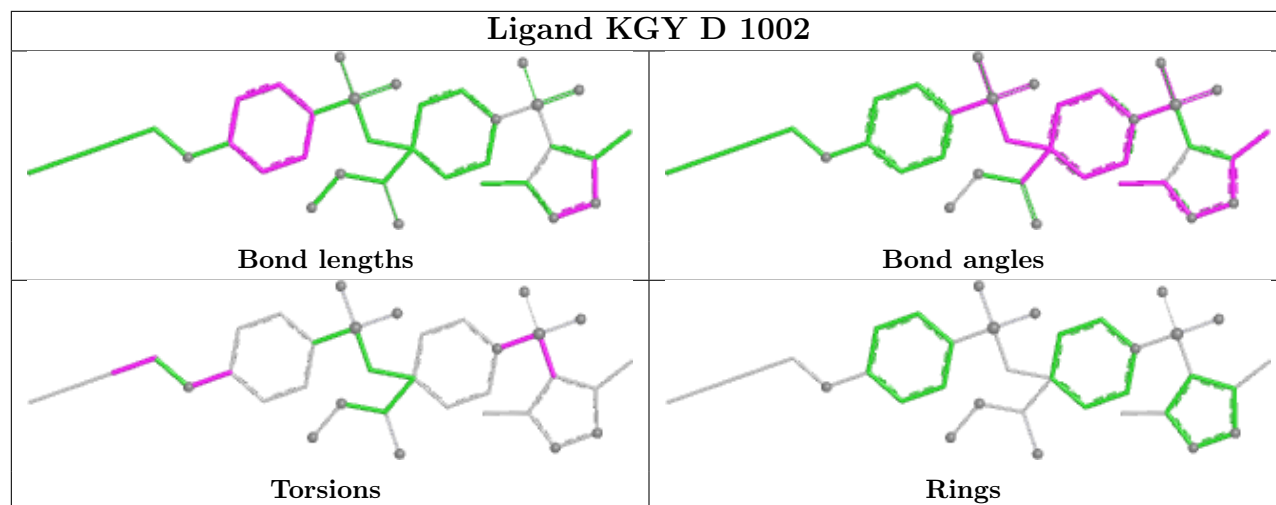
Mol	Chain	Res	Type	Atoms
4	C	1001	KGY	C6-C2-S1-O3
4	C	1001	KGY	C6-C2-S1-O4
4	C	1001	KGY	O30-C31-C32-C33
4	D	1002	KGY	C6-C2-S1-O3
4	D	1002	KGY	C12-N11-S1-O4

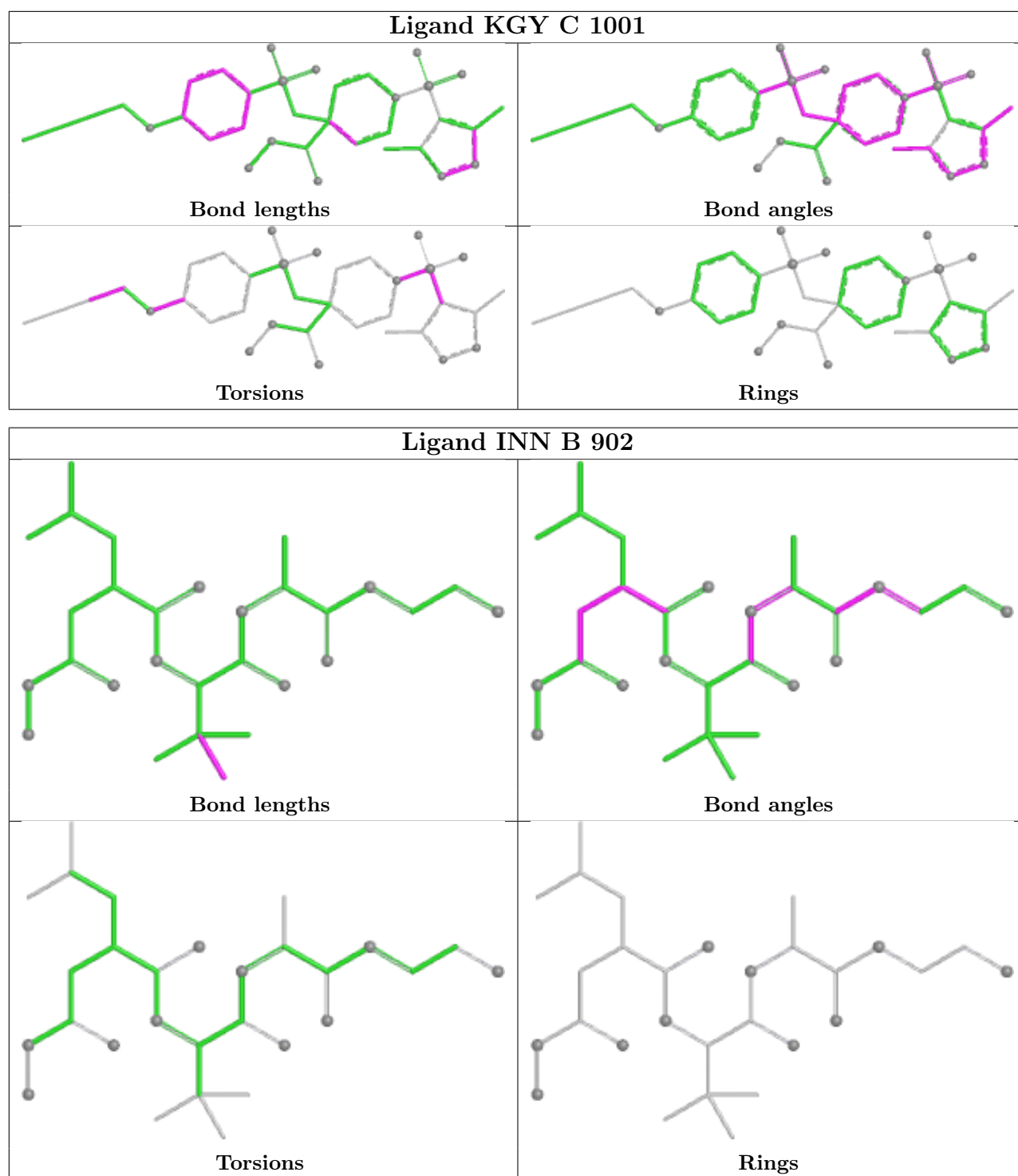
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	1002	KGY	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 ⓘ

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	255/288 (88%)	0.35	26 (10%)	12 12	7, 19, 44, 54	0
1	B	254/288 (88%)	0.17	17 (6%)	24 25	7, 18, 40, 53	0
1	C	254/288 (88%)	0.38	26 (10%)	12 12	7, 20, 48, 54	0
1	D	255/288 (88%)	0.21	21 (8%)	17 18	7, 18, 47, 55	0
All	All	1018/1152 (88%)	0.28	90 (8%)	15 16	7, 19, 45, 55	0

The worst 5 of 90 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	420	LEU	9.6
1	D	220	PRO	6.9
1	A	421	ALA	6.4
1	C	421	ALA	5.9
1	A	369	TYR	5.8

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](#)

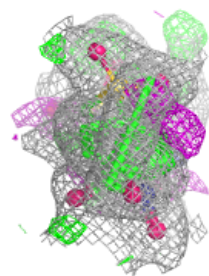
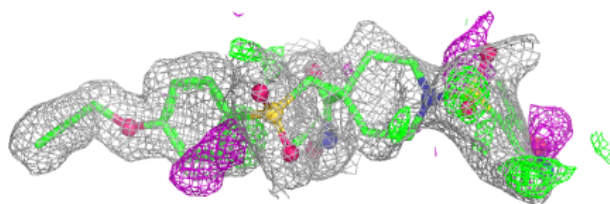
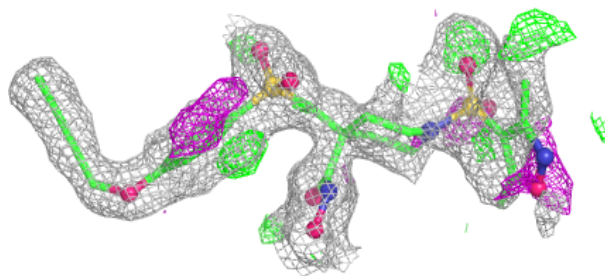
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists 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	KGY	C	1001	35/35	0.88	0.16	17,24,61,62	0
4	KGY	D	1002	35/35	0.93	0.10	13,21,42,43	0
3	INN	A	901	29/29	0.96	0.06	6,11,27,34	0
3	INN	B	902	29/29	0.96	0.06	8,11,26,28	0
2	ZN	C	804	1/1	0.99	0.02	14,14,14,14	0
2	ZN	D	803	1/1	0.99	0.03	14,14,14,14	0
2	ZN	A	802	1/1	1.00	0.01	11,11,11,11	0
2	ZN	B	801	1/1	1.00	0.03	12,12,12,12	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.

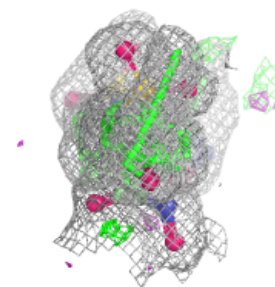
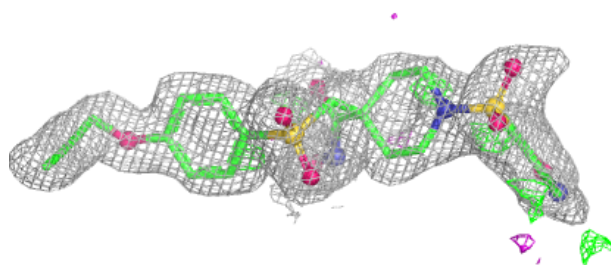
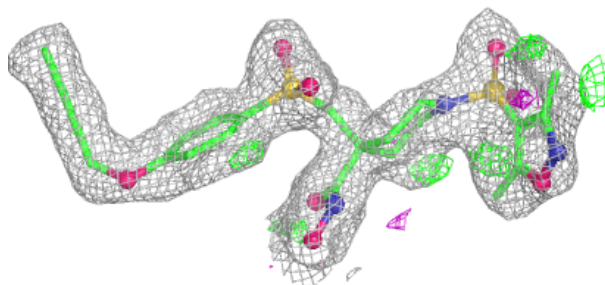
Electron density around KGY C 1001:

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

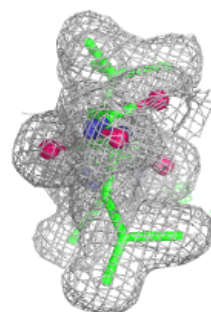
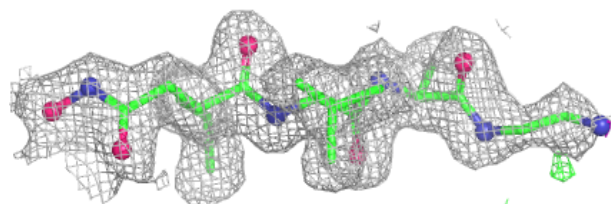
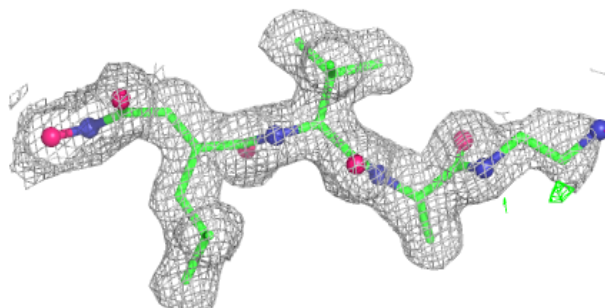


Electron density around KGY D 1002:

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

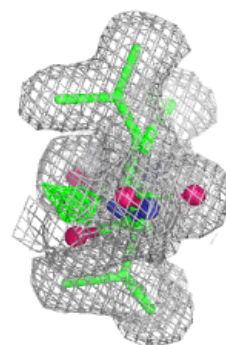
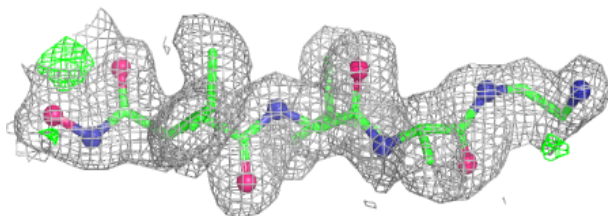
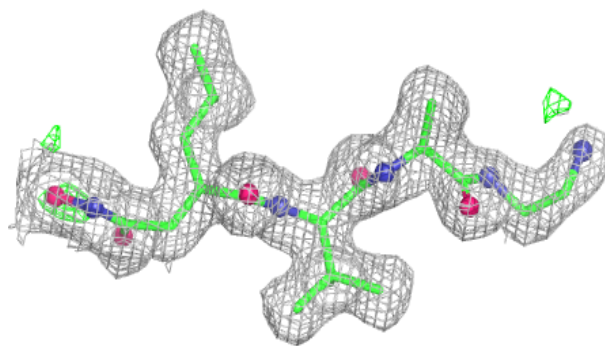
**Electron density around INN A 901:**

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



Electron density around INN B 902:

$2mF_o-DF_c$ (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.