



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 10:36 AM UTC

PDB ID : 2YPU / pdb_00002ypu
Title : human insulin degrading enzyme E111Q in complex with inhibitor compound 41367
Authors : Guo, Q.; Deprez-Poulain, R.; Deprez, B.; Tang, W.-J.
Deposited on : 2012-11-01
Resolution : 2.80 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

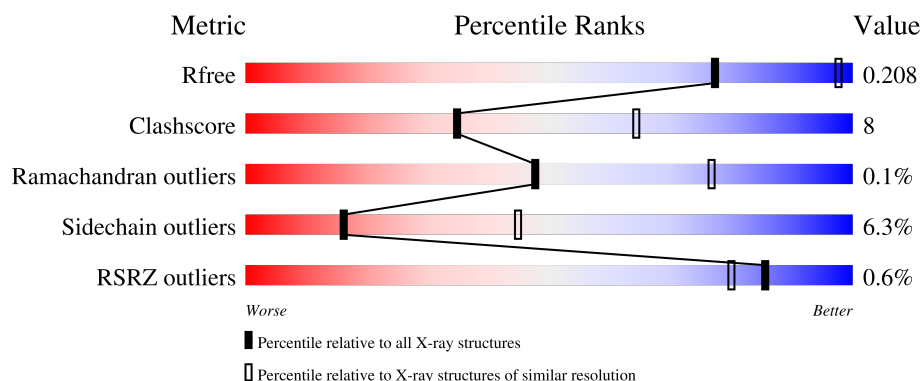
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

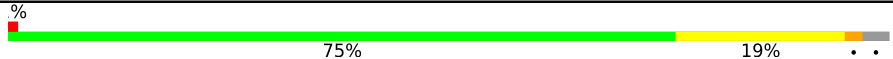
The reported resolution of this entry is 2.80 Å.

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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	
1	B	990	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	I41	A	1999	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 15874 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 INSULIN-DEGRADING ENZYME.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	956	Total	C	N	O	S	0	0	0
			7768	5008	1307	1431	22			
1	B	954	Total	C	N	O	S	0	0	0
			7758	5002	1305	1429	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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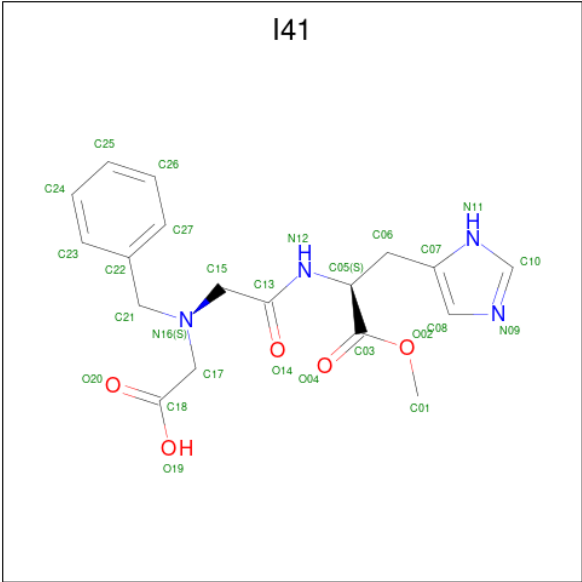
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Chain	Residue	Modelled	Actual	Comment	Reference
A	974	ALA	CYS	engineered mutation	UNP P14735
B	30	MET	-	expression tag	UNP P14735
B	31	HIS	-	expression tag	UNP P14735
B	32	HIS	-	expression tag	UNP P14735
B	33	HIS	-	expression tag	UNP P14735
B	34	HIS	-	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 2-[[2-[[[(2S)-3-(3H-IMIDAZOL-4-YL)-1-METHOXY-1-OXO-PROPAN-2-YL]AMINO]-2-OXO-ETHYL]-(PHENYLMETHYL)AMINO]ETHANOIC ACID (CCD ID: I41) (formula: C₁₈H₂₂N₄O₅).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			27	18	4	5		
3	B	1	Total	C	N	O	0	0
			27	18	4	5		

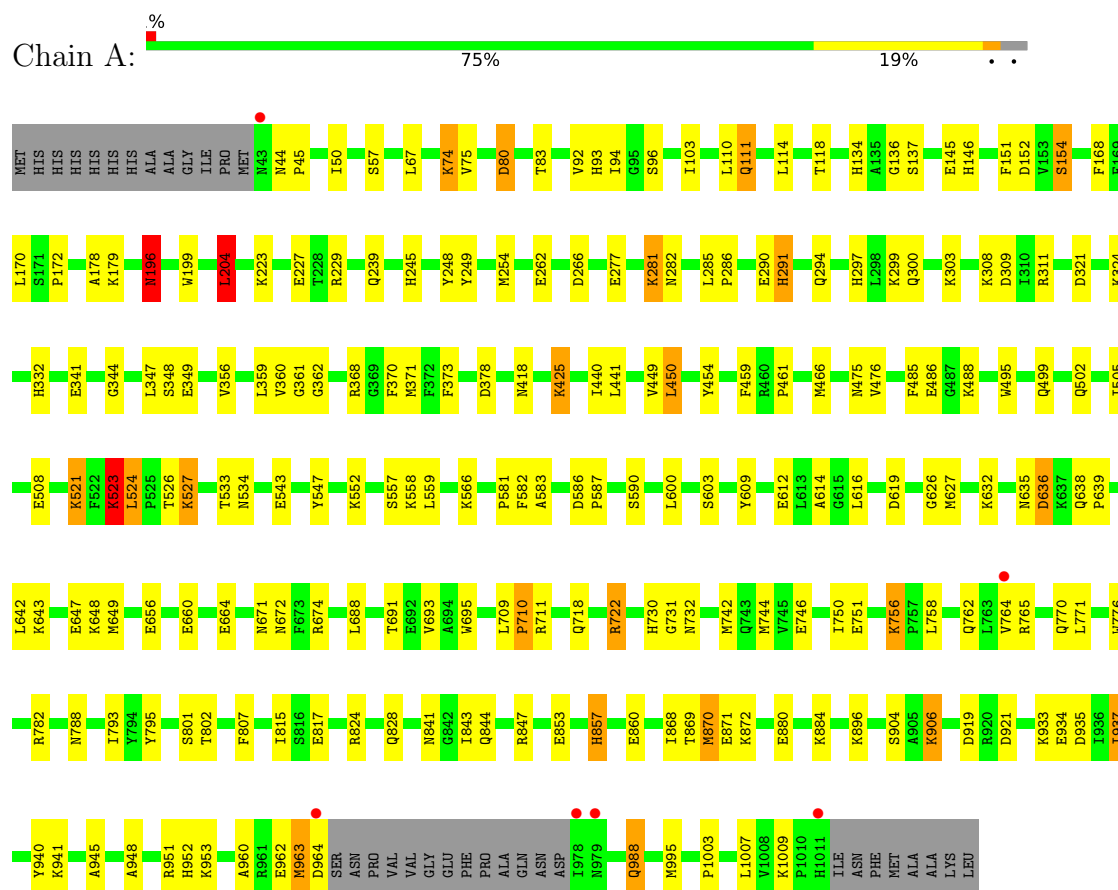
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	178	Total	O	0	0
			178	178		
4	B	114	Total	O	0	0
			114	114		

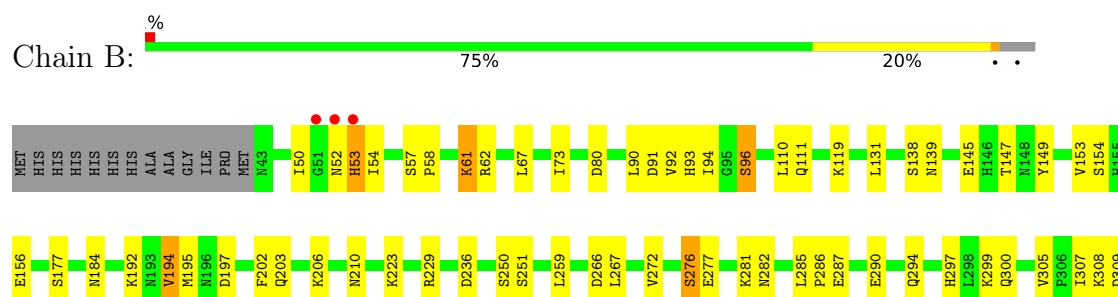
3 Residue-property plots

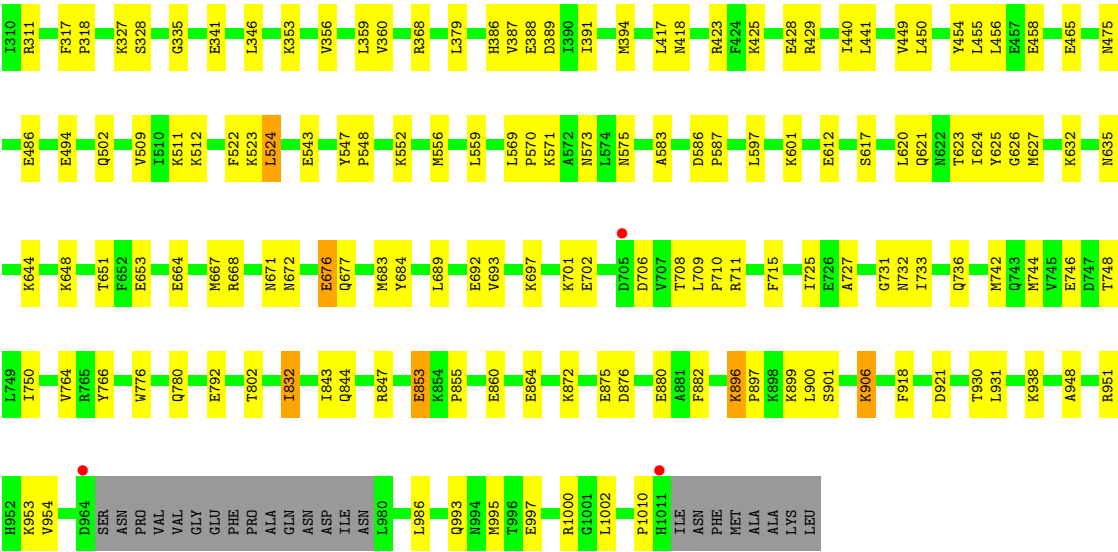
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: INSULIN-DEGRADING ENZYME



• Molecule 1: INSULIN-DEGRADING ENZYME





4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α , β , γ	262.29Å 262.29Å 90.63Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 2.80 50.00 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.9 (50.00-2.80) 99.8 (50.00-2.80)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.22 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.173 , 0.225 (Not available) , 0.208	Depositor DCC
R_{free} test set	4383 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	39.5	Xtriage
Anisotropy	0.042	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 33.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.024 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	15874	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.85% 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, I41

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	1.24	16/7963 (0.2%)	1.25	35/10779 (0.3%)
1	B	1.21	14/7953 (0.2%)	1.22	33/10765 (0.3%)
All	All	1.22	30/15916 (0.2%)	1.24	68/21544 (0.3%)

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	855	PRO	CA-C	10.21	1.57	1.51
1	A	178	ALA	CA-CB	-8.58	1.40	1.53
1	B	511	LYS	CA-CB	8.29	1.66	1.53
1	B	764	VAL	CA-CB	7.94	1.62	1.53
1	A	204	LEU	CG-CD2	7.52	1.77	1.52
1	A	614	ALA	CA-CB	-7.27	1.42	1.53
1	A	146	HIS	CA-C	-6.98	1.44	1.52
1	A	508	GLU	CA-CB	6.65	1.63	1.53
1	A	523	LYS	CG-CD	6.43	1.71	1.52
1	B	575	ASN	CA-C	-6.32	1.45	1.52
1	B	53	HIS	CA-C	6.22	1.61	1.52
1	B	543	GLU	CA-CB	6.20	1.61	1.52
1	A	693	VAL	CA-CB	6.18	1.62	1.54
1	B	305	VAL	CA-CB	6.17	1.60	1.53
1	B	997	GLU	CA-CB	5.99	1.62	1.53
1	B	465	GLU	CG-CD	5.94	1.66	1.52
1	A	963	MET	CA-C	5.90	1.60	1.52
1	B	494	GLU	CG-CD	5.50	1.65	1.52
1	A	543	GLU	CA-CB	5.48	1.61	1.53
1	A	964	ASP	CA-CB	5.46	1.64	1.53
1	B	1002	LEU	N-CA	5.32	1.50	1.45
1	A	764	VAL	CA-CB	5.30	1.60	1.53
1	A	765	ARG	CA-C	5.29	1.59	1.52
1	A	111	GLN	C-O	-5.26	1.18	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	53	HIS	CD2-NE2	5.21	1.43	1.37
1	A	750	ILE	CA-C	5.14	1.59	1.52
1	A	92	VAL	CA-C	-5.11	1.46	1.52
1	A	619	ASP	N-CA	-5.10	1.40	1.46
1	B	853	GLU	N-CA	-5.10	1.39	1.46
1	B	317	PHE	CA-C	-5.08	1.48	1.52

All (68) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	667	MET	N-CA-C	-7.87	102.65	111.07
1	B	706	ASP	N-CA-C	7.66	122.21	113.01
1	B	853	GLU	N-CA-C	-7.45	104.33	113.50
1	A	870	MET	N-CA-C	7.43	119.03	111.07
1	A	731	GLY	N-CA-C	7.38	119.83	110.38
1	A	636	ASP	CB-CA-C	-7.33	95.84	110.42
1	A	229	ARG	CA-C-N	-7.18	111.58	119.19
1	A	229	ARG	C-N-CA	-7.18	111.58	119.19
1	B	210	ASN	CA-C-N	-7.16	112.42	119.87
1	B	210	ASN	C-N-CA	-7.16	112.42	119.87
1	A	547	TYR	CA-C-N	-7.13	112.63	119.76
1	A	547	TYR	C-N-CA	-7.13	112.63	119.76
1	B	617	SER	N-CA-C	6.99	119.55	109.14
1	B	197	ASP	N-CA-C	6.91	118.81	111.28
1	B	92	VAL	CB-CA-C	-6.77	101.70	110.98
1	A	344	GLY	N-CA-C	-6.60	106.83	114.69
1	A	627	MET	N-CA-C	6.37	119.94	110.52
1	A	291	HIS	CA-C-N	-6.33	113.28	119.87
1	A	291	HIS	C-N-CA	-6.33	113.28	119.87
1	B	715	PHE	N-CA-C	6.21	118.05	111.28
1	B	387	VAL	N-CA-C	6.12	116.30	110.42
1	A	495	TRP	N-CA-C	6.11	118.91	111.82
1	B	832	ILE	CB-CA-C	-6.11	103.57	110.96
1	A	636	ASP	N-CA-C	6.05	123.68	110.80
1	B	455	LEU	CB-CA-C	6.03	119.48	109.53
1	A	919	ASP	N-CA-C	-6.01	103.39	111.87
1	B	54	ILE	CB-CA-C	-6.00	103.70	110.96
1	A	80	ASP	CA-C-N	-5.87	113.77	119.87
1	A	80	ASP	C-N-CA	-5.87	113.77	119.87
1	A	475	ASN	CB-CA-C	-5.75	99.66	110.01
1	A	505	ILE	CA-C-N	5.74	125.67	119.76
1	A	505	ILE	C-N-CA	5.74	125.67	119.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	621	GLN	N-CA-C	5.70	117.84	109.24
1	B	360	VAL	CA-C-N	-5.68	115.50	121.35
1	B	360	VAL	C-N-CA	-5.68	115.50	121.35
1	B	429	ARG	CA-C-N	5.68	125.15	119.24
1	B	429	ARG	C-N-CA	5.68	125.15	119.24
1	B	524	LEU	CA-C-N	-5.66	114.43	120.03
1	B	524	LEU	C-N-CA	-5.66	114.43	120.03
1	A	616	LEU	CA-C-N	-5.64	113.33	122.36
1	A	616	LEU	C-N-CA	-5.64	113.33	122.36
1	B	648	LYS	N-CA-C	-5.64	105.28	111.82
1	B	53	HIS	CE1-NE2-CD2	5.60	114.60	109.00
1	A	857	HIS	N-CA-C	-5.58	105.29	111.71
1	A	619	ASP	N-CA-CB	-5.55	101.73	110.77
1	A	92	VAL	CB-CA-C	-5.47	102.67	110.83
1	B	731	GLY	N-CA-C	5.45	117.36	110.38
1	B	651	THR	N-CA-C	-5.43	104.92	112.30
1	B	267	LEU	N-CA-C	-5.43	105.52	111.82
1	B	918	PHE	N-CA-C	5.38	118.64	111.75
1	B	855	PRO	N-CA-C	5.36	115.62	110.47
1	B	50	ILE	CA-C-N	-5.33	117.19	122.36
1	B	50	ILE	C-N-CA	-5.33	117.19	122.36
1	A	254	MET	N-CA-C	5.32	118.06	109.40
1	B	766	TYR	N-CA-C	5.30	118.33	110.48
1	B	335	GLY	N-CA-C	-5.26	106.04	112.77
1	B	882	PHE	N-CA-C	5.23	117.38	111.11
1	A	196	ASN	CB-CA-C	5.21	118.73	109.72
1	A	74	LYS	N-CA-C	-5.19	102.00	110.20
1	A	884	LYS	CA-CB-CG	5.15	124.40	114.10
1	A	466	MET	CG-SD-CE	-5.11	89.66	100.90
1	A	75	VAL	CB-CA-C	-5.09	102.89	110.33
1	A	170	LEU	CA-C-N	-5.07	117.66	122.37
1	A	170	LEU	C-N-CA	-5.07	117.66	122.37
1	A	817	GLU	CA-C-N	-5.06	113.71	119.28
1	A	817	GLU	C-N-CA	-5.06	113.71	119.28
1	B	938	LYS	N-CA-CB	5.03	117.52	110.12
1	A	137	SER	N-CA-C	-5.01	101.29	108.60

There are no chirality outliers.

There are no planarity outliers.

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	7768	0	7676	128	1
1	B	7758	0	7672	104	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	27	0	21	9	0
3	B	27	0	21	3	0
4	A	178	0	0	12	0
4	B	114	0	0	4	1
All	All	15874	0	15390	233	1

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 8.

All (233) 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:204:LEU:CD2	1:A:204:LEU:CG	1.77	1.60
1:B:309:ASP:H	1:B:672:ASN:HD21	1.12	0.95
1:A:332:HIS:HD2	3:A:1999:I41:C01	1.80	0.94
1:B:184:ASN:HD21	1:B:223:LYS:NZ	1.67	0.93
1:A:74:LYS:NZ	4:A:2009:HOH:O	2.01	0.92
1:A:309:ASP:H	1:A:672:ASN:HD21	1.19	0.90
1:B:597:LEU:HD11	1:B:627:MET:HG2	1.53	0.90
1:B:96:SER:HB3	4:B:2008:HOH:O	1.71	0.89
1:A:557:SER:HB2	1:A:742:MET:HE3	1.54	0.88
1:B:847:ARG:NH1	4:B:2103:HOH:O	1.83	0.84
1:A:776:TRP:CD1	1:A:953:LYS:HG2	2.15	0.82
1:A:332:HIS:CD2	3:A:1999:I41:C01	2.63	0.81
1:A:635:ASN:ND2	1:A:732:ASN:HD22	1.79	0.81
1:B:727:ALA:HB3	1:B:742:MET:HE1	1.61	0.81
1:B:93:HIS:HE1	1:B:368:ARG:HH21	1.29	0.80
1:A:988:GLN:HB2	4:A:2174:HOH:O	1.82	0.80
1:B:184:ASN:HD21	1:B:223:LYS:HZ2	1.27	0.79
1:B:309:ASP:H	1:B:672:ASN:ND2	1.82	0.77
1:A:332:HIS:CD2	3:A:1999:I41:H011	2.21	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:857:HIS:HD2	4:A:2159:HOH:O	1.69	0.76
1:A:988:GLN:CB	4:A:2174:HOH:O	2.33	0.76
1:B:441:LEU:HD23	1:B:449:VAL:HG11	1.69	0.75
1:A:635:ASN:HD21	1:A:732:ASN:HD22	1.35	0.74
3:B:1999:I41:C18	3:B:1999:I41:O14	2.37	0.73
1:B:671:ASN:OD1	1:B:701:LYS:HD3	1.87	0.73
1:B:93:HIS:CE1	1:B:368:ARG:HH21	2.07	0.72
1:B:635:ASN:ND2	1:B:732:ASN:HD22	1.88	0.72
1:A:300:GLN:HE21	1:A:502:GLN:HE21	1.40	0.69
1:B:711:ARG:HG3	4:B:2098:HOH:O	1.92	0.68
1:A:204:LEU:CD2	1:A:204:LEU:HG	2.12	0.68
1:B:184:ASN:ND2	1:B:223:LYS:NZ	2.39	0.68
1:A:196:ASN:HD21	1:A:199:TRP:HD1	1.42	0.68
1:A:196:ASN:ND2	1:A:199:TRP:HD1	1.91	0.67
1:B:635:ASN:HD21	1:B:732:ASN:HD22	1.42	0.66
3:A:1999:I41:C08	4:A:2177:HOH:O	2.43	0.65
1:A:291:HIS:CD2	1:A:370:PHE:HB2	2.32	0.65
1:A:309:ASP:H	1:A:672:ASN:ND2	1.91	0.65
1:B:776:TRP:CD1	1:B:953:LYS:HD3	2.33	0.64
1:A:600:LEU:HD11	1:A:648:LYS:HB3	1.81	0.63
1:B:285:LEU:HD12	1:B:286:PRO:HD2	1.81	0.62
1:B:727:ALA:CB	1:B:742:MET:HE1	2.30	0.62
1:A:204:LEU:CD2	1:A:204:LEU:CD1	2.75	0.61
1:A:93:HIS:HE1	1:A:368:ARG:HH21	1.49	0.61
1:A:583:ALA:CB	1:A:626:GLY:HA2	2.31	0.60
1:A:308:LYS:HE3	1:A:672:ASN:HB3	1.82	0.60
1:A:300:GLN:NE2	1:A:502:GLN:HE21	1.99	0.60
1:B:559:LEU:HD13	1:B:742:MET:HE2	1.84	0.60
1:A:557:SER:CB	1:A:742:MET:HE3	2.29	0.59
3:B:1999:I41:O14	3:B:1999:I41:O20	2.20	0.59
1:B:623:THR:HG22	1:B:624:ILE:N	2.18	0.59
1:A:776:TRP:NE1	1:A:953:LYS:HE2	2.18	0.58
1:B:689:LEU:CD2	1:B:995:MET:HG2	2.34	0.58
1:A:1007:LEU:HD12	1:B:1000:ARG:HG2	1.86	0.58
1:B:266:ASP:HB2	4:B:2040:HOH:O	2.02	0.58
1:A:359:LEU:O	3:A:1999:I41:H10	2.04	0.58
1:B:52:ASN:O	1:B:53:HIS:C	2.46	0.58
1:A:709:LEU:HB3	1:A:710:PRO:HD3	1.86	0.57
1:B:341:GLU:OE2	3:B:1999:I41:H10	2.04	0.57
1:A:843:ILE:HG22	1:A:844:GLN:H	1.70	0.57
1:A:441:LEU:HD23	1:A:449:VAL:HG11	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:795:TYR:CE1	1:A:953:LYS:HD2	2.40	0.56
1:A:688:LEU:HD13	1:A:995:MET:HE1	1.86	0.56
1:B:689:LEU:HD23	1:B:995:MET:HG2	1.86	0.56
1:A:782:ARG:HH12	1:A:963:MET:H	1.54	0.56
1:A:223:LYS:HE2	1:A:227:GLU:OE2	2.06	0.55
1:B:300:GLN:HE21	1:B:502:GLN:HE21	1.55	0.55
1:B:184:ASN:ND2	1:B:223:LYS:HZ3	2.02	0.55
1:A:660:GLU:HG3	4:A:2126:HOH:O	2.06	0.55
1:A:557:SER:HB2	1:A:742:MET:CE	2.32	0.55
1:A:671:ASN:O	1:A:674:ARG:HG2	2.07	0.55
1:A:801:SER:O	1:A:802:THR:C	2.50	0.54
1:B:559:LEU:HD22	1:B:742:MET:HB2	1.88	0.54
1:A:933:LYS:O	1:A:937:ILE:HG12	2.07	0.54
1:B:709:LEU:HB3	1:B:710:PRO:HD3	1.89	0.54
1:B:294:GLN:H	1:B:297:HIS:HD2	1.55	0.54
1:A:632:LYS:HD2	4:A:2118:HOH:O	2.07	0.54
1:A:636:ASP:OD1	1:A:636:ASP:O	2.27	0.53
1:A:691:THR:HG23	1:A:841:ASN:HD21	1.73	0.53
1:A:940:TYR:CE2	1:A:945:ALA:HB2	2.43	0.53
1:A:523:LYS:HD3	1:A:524:LEU:N	2.24	0.53
1:A:321:ASP:OD1	1:A:371:MET:HE3	2.08	0.53
1:A:196:ASN:HD21	1:A:199:TRP:CD1	2.23	0.52
1:A:94:ILE:HG13	1:A:248:TYR:HB3	1.92	0.52
1:A:988:GLN:HB3	4:A:2174:HOH:O	2.06	0.52
1:A:303:LYS:HD3	1:A:485:PHE:CE1	2.44	0.52
1:B:559:LEU:HB2	1:B:742:MET:HE3	1.91	0.52
1:B:285:LEU:HD12	1:B:286:PRO:CD	2.39	0.52
1:A:937:ILE:HG22	1:A:941:LYS:HD2	1.93	0.51
1:B:236:ASP:OD1	1:B:236:ASP:C	2.53	0.51
1:A:294:GLN:H	1:A:297:HIS:HD2	1.57	0.51
1:B:388:GLU:OE2	1:B:509:VAL:HG22	2.11	0.51
1:B:308:LYS:HD3	1:B:672:ASN:HB3	1.92	0.51
1:A:586:ASP:HB2	1:A:587:PRO:HD2	1.92	0.50
1:A:722:ARG:HB2	1:A:758:LEU:HD23	1.91	0.50
1:A:722:ARG:HB2	1:A:758:LEU:CD2	2.41	0.50
1:A:815:ILE:HG22	1:A:870:MET:HG3	1.92	0.50
1:B:93:HIS:HE1	1:B:368:ARG:NH2	2.05	0.50
1:B:93:HIS:HD2	1:B:145:GLU:O	1.93	0.50
1:B:184:ASN:ND2	1:B:223:LYS:HZ2	2.02	0.50
1:A:332:HIS:HD2	3:A:1999:I41:H012	1.73	0.50
1:A:151:PHE:CD1	1:A:151:PHE:C	2.89	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:583:ALA:CB	1:B:626:GLY:HA2	2.42	0.49
1:A:533:THR:N	1:A:636:ASP:OD2	2.39	0.49
1:A:552:LYS:HB3	1:A:559:LEU:HB3	1.94	0.49
1:A:285:LEU:HD12	1:A:286:PRO:HD2	1.93	0.49
1:A:311:ARG:NH2	1:A:664:GLU:OE2	2.45	0.49
1:B:110:LEU:HD23	1:B:110:LEU:C	2.38	0.49
1:B:864:GLU:OE2	1:B:953:LYS:HE3	2.12	0.49
1:B:906:LYS:HE3	1:B:921:ASP:OD2	2.13	0.49
1:B:147:THR:HG22	1:B:149:TYR:CE1	2.47	0.49
1:B:272:VAL:O	1:B:276:SER:OG	2.30	0.49
1:B:311:ARG:HB3	1:B:379:LEU:HB2	1.94	0.49
1:B:623:THR:HG22	1:B:625:TYR:H	1.78	0.49
1:A:290:GLU:HB3	4:A:2049:HOH:O	2.12	0.49
1:A:359:LEU:C	1:A:359:LEU:HD23	2.38	0.49
1:A:843:ILE:HG22	1:A:844:GLN:N	2.27	0.49
1:B:328:SER:HB2	1:B:458:GLU:O	2.12	0.49
1:A:114:LEU:HD13	1:A:168:PHE:HB3	1.94	0.49
1:B:776:TRP:NE1	1:B:953:LYS:HD3	2.28	0.48
1:A:136:GLY:HA3	1:A:152:ASP:O	2.14	0.48
1:A:418:ASN:HB3	1:A:454:TYR:O	2.13	0.48
1:A:857:HIS:CD2	4:A:2159:HOH:O	2.54	0.48
1:B:601:LYS:NZ	1:B:620:LEU:O	2.36	0.48
1:A:196:ASN:ND2	1:A:199:TRP:CD1	2.79	0.48
1:B:547:TYR:HB3	1:B:548:PRO:CD	2.43	0.48
1:A:880:GLU:HB2	4:A:2163:HOH:O	2.13	0.48
1:A:245:HIS:O	1:A:249:TYR:HB2	2.14	0.48
1:A:586:ASP:HA	1:A:695:TRP:CZ2	2.48	0.48
1:A:793:ILE:O	1:A:847:ARG:HA	2.14	0.47
1:B:559:LEU:HD13	1:B:742:MET:CE	2.44	0.47
1:B:425:LYS:HD2	1:B:428:GLU:OE2	2.14	0.47
1:A:93:HIS:HD2	1:A:145:GLU:O	1.97	0.47
1:A:583:ALA:HB2	1:A:626:GLY:HA2	1.97	0.47
1:B:676:GLU:OE1	1:B:676:GLU:HA	2.14	0.47
1:B:418:ASN:HB3	1:B:454:TYR:O	2.15	0.46
1:A:93:HIS:CE1	1:A:368:ARG:HH21	2.31	0.46
1:B:311:ARG:HH22	1:B:664:GLU:CD	2.23	0.46
1:A:262:GLU:HB3	1:A:266:ASP:HB2	1.98	0.46
1:A:718:GLN:HE22	1:B:711:ARG:HH12	1.62	0.46
1:B:73:ILE:HG13	1:B:251:SER:HB2	1.98	0.46
1:B:346:LEU:HA	1:B:522:PHE:CE1	2.50	0.46
1:B:441:LEU:CD2	1:B:449:VAL:HG11	2.41	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:860:GLU:OE1	1:B:953:LYS:NZ	2.49	0.46
1:A:134:HIS:O	1:A:154:SER:HB3	2.15	0.46
1:A:643:LYS:HE2	1:A:647:GLU:OE1	2.16	0.46
1:A:906:LYS:NZ	1:A:921:ASP:OD2	2.39	0.46
1:B:131:LEU:CD1	1:B:138:SER:HB2	2.46	0.46
1:B:202:PHE:CZ	1:B:206:LYS:HE3	2.51	0.45
1:B:843:ILE:HG22	1:B:844:GLN:H	1.80	0.45
1:B:359:LEU:C	1:B:359:LEU:HD23	2.42	0.45
1:B:417:LEU:HA	1:B:417:LEU:HD23	1.73	0.45
1:A:771:LEU:HB2	1:A:952:HIS:HB3	1.98	0.45
1:B:259:LEU:HD23	1:B:259:LEU:C	2.42	0.45
1:A:756:LYS:HB2	1:A:756:LYS:HE2	1.41	0.45
1:B:61:LYS:HA	1:B:61:LYS:HD3	1.76	0.45
1:B:194:VAL:HG12	1:B:195:MET:HG2	1.99	0.45
1:B:727:ALA:HB3	1:B:742:MET:CE	2.39	0.45
1:A:581:PRO:HG2	1:A:582:PHE:CD2	2.52	0.45
1:B:684:TYR:OH	1:B:697:LYS:HG2	2.17	0.45
1:B:300:GLN:NE2	1:B:502:GLN:HE21	2.15	0.45
1:A:360:VAL:HA	3:A:1999:I41:H10	1.99	0.45
1:B:552:LYS:HB3	1:B:559:LEU:HB3	1.99	0.45
1:A:857:HIS:O	1:A:860:GLU:HB2	2.17	0.45
1:A:94:ILE:HD12	1:A:94:ILE:HG23	1.53	0.44
1:B:556:MET:HE3	1:B:750:ILE:HD11	1.99	0.44
1:B:573:ASN:OD1	1:B:632:LYS:HG2	2.17	0.44
1:A:559:LEU:HD22	1:A:742:MET:HB2	1.99	0.44
1:B:899:LYS:O	1:B:900:LEU:C	2.60	0.44
1:A:378:ASP:OD1	1:A:378:ASP:C	2.60	0.44
1:B:843:ILE:HG22	1:B:844:GLN:N	2.31	0.44
1:A:80:ASP:O	1:A:83:THR:HG22	2.18	0.44
1:A:341:GLU:HG2	1:A:347:LEU:HD23	2.00	0.44
1:A:449:VAL:HG23	1:A:450:LEU:HD13	2.00	0.44
1:A:730:HIS:HD2	1:A:904:SER:OG	2.01	0.44
1:B:864:GLU:HG3	1:B:986:LEU:HD21	1.99	0.43
1:A:649:MET:HE3	1:A:649:MET:HB2	1.85	0.43
1:A:782:ARG:NH1	1:A:963:MET:H	2.15	0.43
1:A:110:LEU:C	1:A:110:LEU:HD23	2.44	0.43
1:A:807:PHE:HE1	1:A:935:ASP:HB3	1.83	0.43
1:B:683:MET:HA	1:B:792:GLU:OE1	2.19	0.43
1:A:552:LYS:NZ	1:A:746:GLU:OE2	2.50	0.43
1:B:391:ILE:O	1:B:394:MET:HB2	2.19	0.43
1:A:362:GLY:HA2	1:A:373:PHE:CZ	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:948:ALA:HB3	1:A:951:ARG:HB2	2.00	0.43
1:B:583:ALA:HB2	1:B:626:GLY:HA2	2.01	0.43
1:B:623:THR:N	1:B:626:GLY:O	2.44	0.43
1:A:824:ARG:O	1:A:828:GLN:HA	2.19	0.43
1:B:692:GLU:HG2	1:B:693:VAL:HG23	2.01	0.43
1:A:311:ARG:HH22	1:A:664:GLU:CD	2.26	0.42
1:B:708:THR:HB	1:B:710:PRO:HD2	2.01	0.42
1:A:299:LYS:HE3	1:A:299:LYS:HB2	1.89	0.42
1:A:638:GLN:N	1:A:639:PRO:CD	2.82	0.42
1:A:776:TRP:HA	1:A:953:LYS:O	2.18	0.42
1:A:960:ALA:HB3	1:A:963:MET:HE3	2.01	0.42
1:B:931:LEU:HD23	1:B:931:LEU:HA	1.81	0.42
1:A:349:GLU:OE1	1:A:521:LYS:HG2	2.19	0.42
1:A:459:PHE:CE2	1:A:461:PRO:HG3	2.55	0.42
1:A:642:LEU:C	1:A:642:LEU:HD12	2.44	0.42
1:B:62:ARG:HG2	1:B:80:ASP:HB2	2.00	0.42
1:B:294:GLN:H	1:B:297:HIS:CD2	2.35	0.42
1:A:118:THR:HG22	1:A:172:PRO:HA	2.01	0.42
1:B:58:PRO:HG2	1:B:423:ARG:HG3	2.00	0.42
1:B:672:ASN:HD22	1:B:672:ASN:HA	1.60	0.42
1:A:360:VAL:HA	3:A:1999:I41:C10	2.50	0.42
1:A:526:THR:O	1:A:527:LYS:C	2.63	0.42
1:B:570:PRO:C	1:B:635:ASN:HD22	2.28	0.42
1:B:153:VAL:HG22	1:B:154:SER:N	2.35	0.41
1:A:361:GLY:H	3:A:1999:I41:H11	1.68	0.41
1:A:744:MET:O	1:A:744:MET:HG2	2.20	0.41
1:B:90:LEU:HD23	1:B:91:ASP:N	2.35	0.41
1:A:179:LYS:HE3	1:A:179:LYS:HB3	1.91	0.41
1:A:600:LEU:CD1	1:A:648:LYS:HB3	2.49	0.41
1:A:586:ASP:HB2	1:A:587:PRO:CD	2.50	0.41
1:A:788:ASN:ND2	4:A:2027:HOH:O	2.53	0.41
1:A:44:ASN:HA	1:A:45:PRO:HD2	1.82	0.41
1:A:534:ASN:OD1	1:A:534:ASN:C	2.64	0.41
1:B:586:ASP:HB2	1:B:587:PRO:HD2	2.03	0.41
1:B:623:THR:CG2	1:B:624:ILE:N	2.82	0.41
1:B:386:HIS:HD2	1:B:389:ASP:OD2	2.03	0.41
1:B:896:LYS:HA	1:B:897:PRO:HD2	1.90	0.41
1:A:868:ILE:O	1:A:869:THR:C	2.64	0.40
1:B:627:MET:HE3	1:B:627:MET:HB2	1.61	0.40
1:A:281:LYS:HZ3	1:A:281:LYS:HG3	1.63	0.40
1:A:722:ARG:HA	1:A:756:LYS:O	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:559:LEU:HD22	1:B:742:MET:CB	2.50	0.40
1:B:569:LEU:O	1:B:571:LYS:N	2.54	0.40
1:A:770:GLN:HB3	1:A:1003:PRO:HB2	2.03	0.40
1:B:203:GLN:O	1:B:203:GLN:HG3	2.21	0.40
1:B:318:PRO:HD2	1:B:475:ASN:O	2.22	0.40
1:B:948:ALA:HB3	1:B:951:ARG:HB2	2.02	0.40
1:A:341:GLU:HB2	1:A:609:TYR:CD1	2.57	0.40
1:A:425:LYS:HB3	1:A:425:LYS:NZ	2.37	0.40
1:B:309:ASP:N	1:B:672:ASN:HD21	1.95	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:934:GLU:CB	4:B:2001:HOH:O[5_555]	1.50	0.70

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	952/990 (96%)	912 (96%)	39 (4%)	1 (0%)	48	77
1	B	950/990 (96%)	910 (96%)	39 (4%)	1 (0%)	48	77
All	All	1902/1980 (96%)	1822 (96%)	78 (4%)	2 (0%)	48	77

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	1010	PRO
1	A	103	ILE

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	837/879 (95%)	790 (94%)	47 (6%)	19	50
1	B	837/879 (95%)	778 (93%)	59 (7%)	14	40
All	All	1674/1758 (95%)	1568 (94%)	106 (6%)	16	45

All (106) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	50	ILE
1	A	57	SER
1	A	67	LEU
1	A	96	SER
1	A	111	GLN
1	A	154	SER
1	A	196	ASN
1	A	204	LEU
1	A	239	GLN
1	A	277	GLU
1	A	281	LYS
1	A	282	ASN
1	A	324	LYS
1	A	348	SER
1	A	356	VAL
1	A	425	LYS
1	A	440	ILE
1	A	450	LEU
1	A	476	VAL
1	A	486	GLU
1	A	488	LYS
1	A	499	GLN
1	A	521	LYS
1	A	523	LYS
1	A	524	LEU
1	A	527	LYS
1	A	558	LYS

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Mol	Chain	Res	Type
1	A	566	LYS
1	A	590	SER
1	A	603	SER
1	A	612	GLU
1	A	656	GLU
1	A	710	PRO
1	A	711	ARG
1	A	722	ARG
1	A	751	GLU
1	A	756	LYS
1	A	762	GLN
1	A	853	GLU
1	A	871	GLU
1	A	872	LYS
1	A	896	LYS
1	A	906	LYS
1	A	937	ILE
1	A	962	GLU
1	A	988	GLN
1	A	1009	LYS
1	B	57	SER
1	B	61	LYS
1	B	67	LEU
1	B	94	ILE
1	B	96	SER
1	B	111	GLN
1	B	119	LYS
1	B	139	ASN
1	B	156	GLU
1	B	177	SER
1	B	192	LYS
1	B	194	VAL
1	B	229	ARG
1	B	250	SER
1	B	276	SER
1	B	277	GLU
1	B	281	LYS
1	B	282	ASN
1	B	287	GLU
1	B	290	GLU
1	B	299	LYS
1	B	307	ILE

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Mol	Chain	Res	Type
1	B	327	LYS
1	B	353	LYS
1	B	356	VAL
1	B	440	ILE
1	B	450	LEU
1	B	456	LEU
1	B	486	GLU
1	B	512	LYS
1	B	523	LYS
1	B	524	LEU
1	B	612	GLU
1	B	644	LYS
1	B	653	GLU
1	B	668	ARG
1	B	676	GLU
1	B	677	GLN
1	B	702	GLU
1	B	725	ILE
1	B	733	ILE
1	B	736	GLN
1	B	744	MET
1	B	746	GLU
1	B	748	THR
1	B	780	GLN
1	B	802	THR
1	B	832	ILE
1	B	853	GLU
1	B	872	LYS
1	B	875	GLU
1	B	876	ASP
1	B	880	GLU
1	B	896	LYS
1	B	901	SER
1	B	906	LYS
1	B	930	THR
1	B	954	VAL
1	B	993	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (40) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	93	HIS

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Mol	Chain	Res	Type
1	A	111	GLN
1	A	231	ASN
1	A	297	HIS
1	A	329	ASN
1	A	332	HIS
1	A	363	GLN
1	A	399	GLN
1	A	407	GLN
1	A	499	GLN
1	A	502	GLN
1	A	635	ASN
1	A	672	ASN
1	A	677	GLN
1	A	724	HIS
1	A	730	HIS
1	A	787	ASN
1	A	813	GLN
1	A	841	ASN
1	A	857	HIS
1	B	52	ASN
1	B	93	HIS
1	B	139	ASN
1	B	184	ASN
1	B	231	ASN
1	B	297	HIS
1	B	300	GLN
1	B	386	HIS
1	B	407	GLN
1	B	499	GLN
1	B	635	ASN
1	B	672	ASN
1	B	677	GLN
1	B	680	GLN
1	B	730	HIS
1	B	788	ASN
1	B	841	ASN
1	B	914	GLN
1	B	922	ASN
1	B	993	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 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	I41	A	1999	-	28,28,28	1.83	5 (17%)	35,36,36	2.13	8 (22%)
3	I41	B	1999	-	28,28,28	1.81	6 (21%)	35,36,36	2.20	10 (28%)

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	I41	A	1999	-	-	15/26/26/26	0/2/2/2
3	I41	B	1999	-	-	11/26/26/26	0/2/2/2

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1999	I41	O02-C03	5.94	1.47	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1999	I41	O02-C03	5.75	1.47	1.33
3	A	1999	I41	C08-C07	3.14	1.42	1.35
3	B	1999	I41	C17-C18	2.57	1.56	1.51
3	B	1999	I41	C15-N16	2.56	1.52	1.47
3	A	1999	I41	C23-C22	2.37	1.43	1.38
3	B	1999	I41	C21-C22	2.29	1.55	1.51
3	B	1999	I41	C08-C07	2.28	1.40	1.35
3	A	1999	I41	C27-C22	2.18	1.43	1.38
3	A	1999	I41	C26-C27	2.15	1.42	1.38
3	B	1999	I41	C26-C27	2.10	1.42	1.38

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1999	I41	C05-C06-C07	7.15	125.37	113.27
3	A	1999	I41	O02-C03-C05	6.01	126.78	111.49
3	A	1999	I41	O02-C03-O04	-5.00	114.12	123.85
3	A	1999	I41	C10-N11-C07	-4.43	102.15	107.91
3	B	1999	I41	C06-C05-N12	4.43	119.08	110.64
3	A	1999	I41	C08-C07-N11	4.30	110.72	105.65
3	B	1999	I41	O02-C03-C05	4.15	122.04	111.49
3	B	1999	I41	O14-C13-N12	-3.97	116.24	122.95
3	B	1999	I41	O02-C03-O04	-3.85	116.35	123.85
3	B	1999	I41	C01-O02-C03	3.53	123.92	115.92
3	A	1999	I41	C05-C06-C07	-3.30	107.68	113.27
3	A	1999	I41	C06-C05-N12	-3.01	104.90	110.64
3	B	1999	I41	O19-C18-C17	2.93	124.92	113.38
3	B	1999	I41	O19-C18-O20	-2.66	116.50	123.33
3	B	1999	I41	C15-C13-N12	2.62	122.14	115.70
3	A	1999	I41	C21-N16-C17	-2.55	106.64	111.77
3	A	1999	I41	N11-C10-N09	2.27	115.20	111.00
3	B	1999	I41	C22-C21-N16	2.07	117.39	113.15

There are no chirality outliers.

All (26) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1999	I41	C03-C05-C06-C07
3	A	1999	I41	N12-C05-C06-C07
3	A	1999	I41	C05-C06-C07-C08
3	A	1999	I41	C05-C06-C07-N11
3	A	1999	I41	N16-C17-C18-O19

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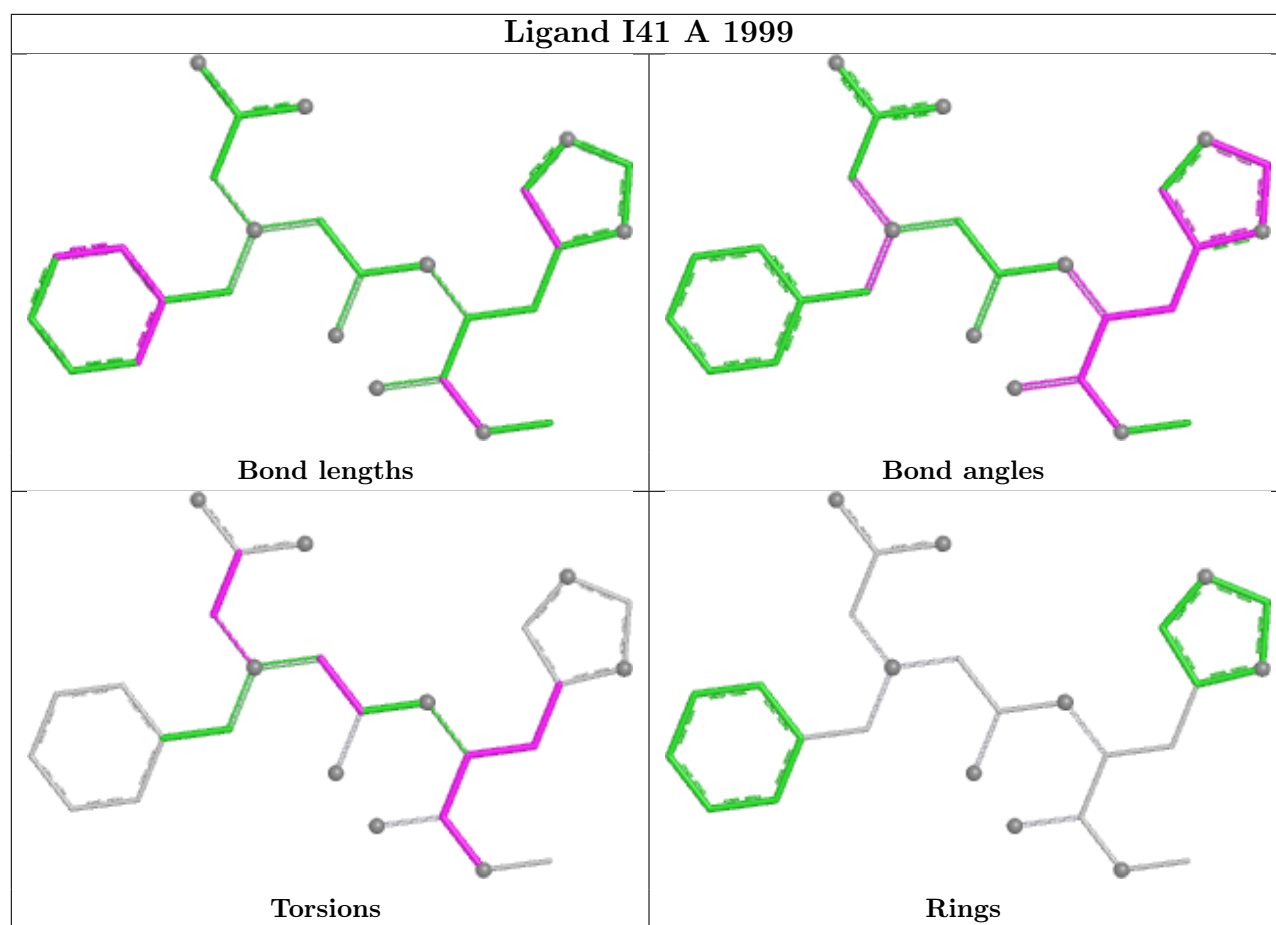
Mol	Chain	Res	Type	Atoms
3	A	1999	I41	N16-C17-C18-O20
3	B	1999	I41	C03-C05-C06-C07
3	B	1999	I41	N12-C05-C06-C07
3	B	1999	I41	C13-C15-N16-C17
3	A	1999	I41	C18-C17-N16-C15
3	A	1999	I41	C05-C03-O02-C01
3	A	1999	I41	O02-C03-C05-N12
3	A	1999	I41	N12-C13-C15-N16
3	A	1999	I41	O04-C03-C05-N12
3	B	1999	I41	N16-C17-C18-O20
3	A	1999	I41	O02-C03-C05-C06
3	A	1999	I41	O14-C13-C15-N16
3	A	1999	I41	O04-C03-O02-C01
3	B	1999	I41	O04-C03-C05-N12
3	A	1999	I41	O04-C03-C05-C06
3	B	1999	I41	O02-C03-C05-N12
3	B	1999	I41	N16-C17-C18-O19
3	B	1999	I41	C05-C06-C07-N11
3	B	1999	I41	C05-C06-C07-C08
3	B	1999	I41	N12-C13-C15-N16
3	B	1999	I41	O14-C13-C15-N16

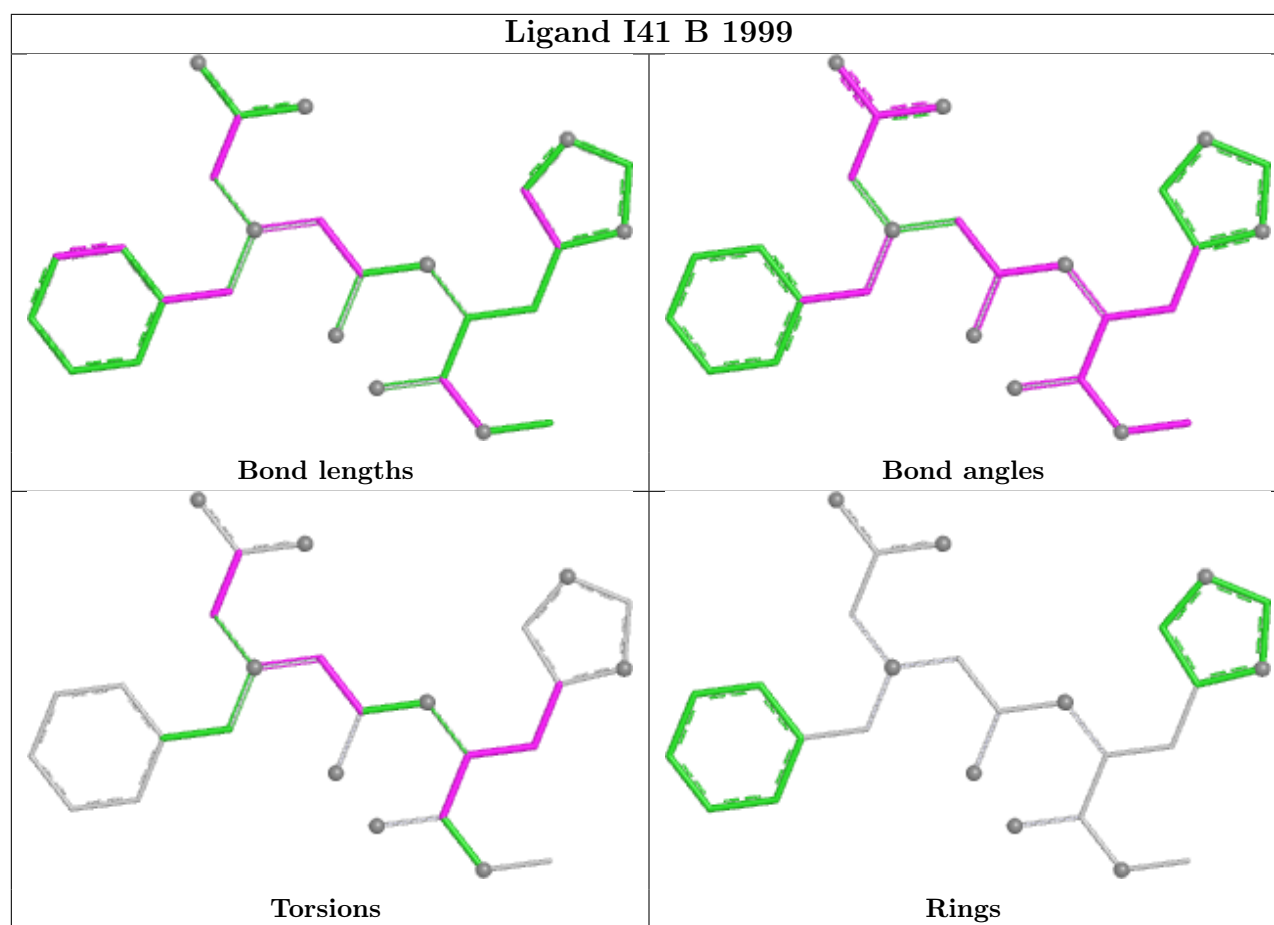
There are no ring outliers.

2 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1999	I41	9	0
3	B	1999	I41	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	956/990 (96%)	-0.74	6 (0%) 85 80	14, 27, 43, 64	0
1	B	954/990 (96%)	-0.63	6 (0%) 85 80	17, 32, 47, 68	0
All	All	1910/1980 (96%)	-0.69	12 (0%) 85 80	14, 30, 45, 68	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	964	ASP	6.8
1	A	964	ASP	6.3
1	A	978	ILE	4.0
1	A	979	ASN	3.8
1	B	53	HIS	3.5
1	B	705	ASP	2.9
1	A	764	VAL	2.7
1	A	1011	HIS	2.6
1	A	43	ASN	2.3
1	B	52	ASN	2.2
1	B	1011	HIS	2.2
1	B	51	GLY	2.1

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](#)

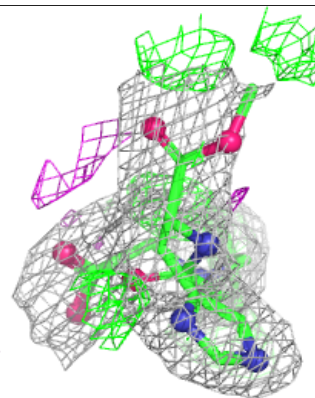
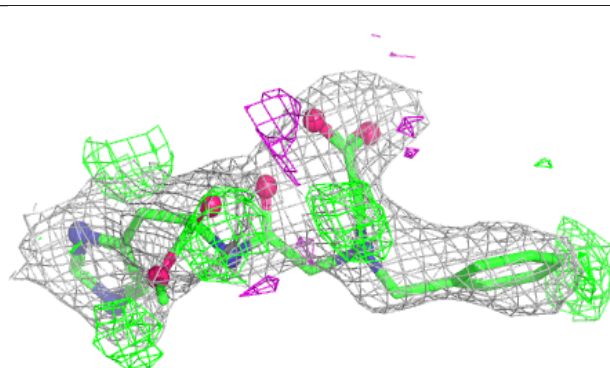
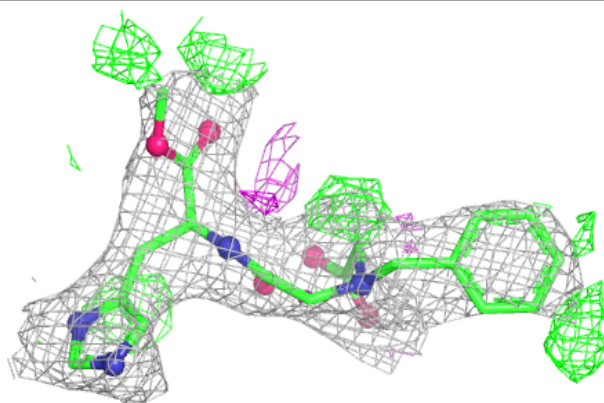
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
3	I41	B	1999	27/27	0.82	0.23	53,63,76,76	0
3	I41	A	1999	27/27	0.83	0.23	35,56,83,84	0
2	ZN	B	1998	1/1	0.99	0.02	33,33,33,33	0
2	ZN	A	1998	1/1	1.00	0.02	31,31,31,31	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.

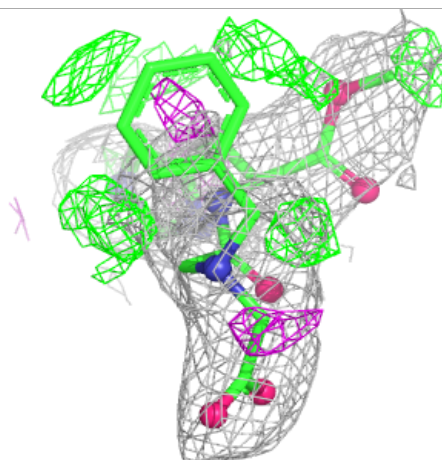
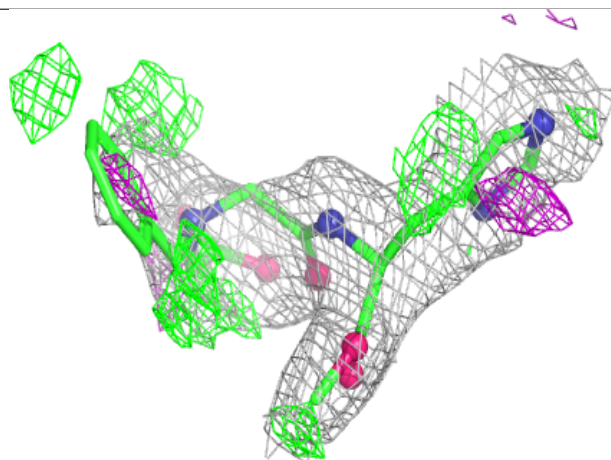
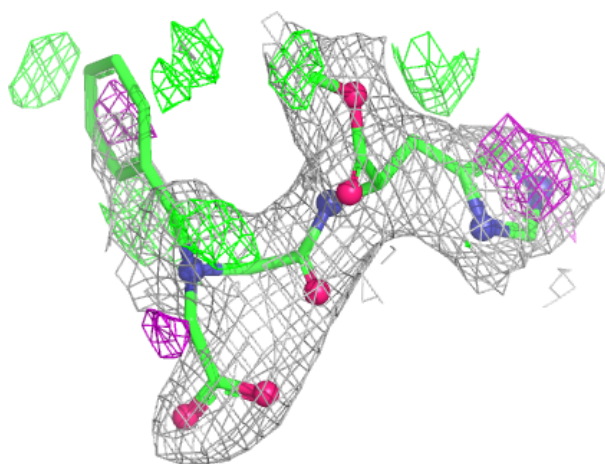
Electron density around I41 B 1999:

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 I41 A 1999:

$2mF_o-DF_c$ (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.