



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 06:49 AM UTC

PDB ID : 4PF7 / pdb_00004pf7
Title : Crystal structure of insulin degrading enzyme complexed with inhibitor
Authors : Wang, Y.; Guo, S.
Deposited on : 2014-04-28
Resolution : 2.33 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

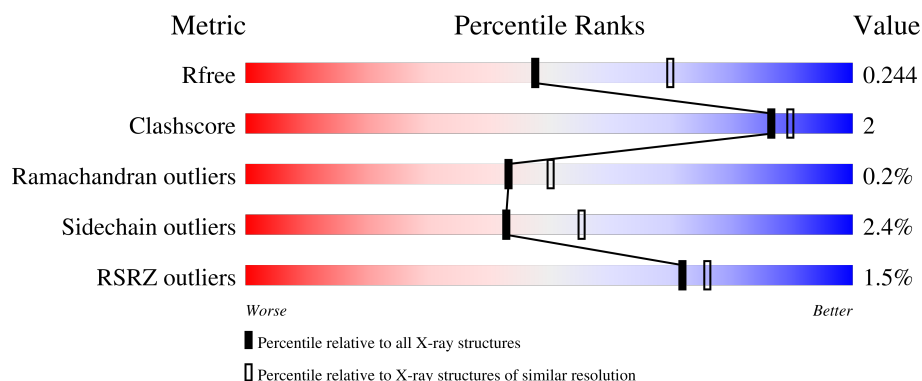
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.33 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	3031 (2.36-2.32)
Clashscore	190562	3127 (2.36-2.32)
Ramachandran outliers	187476	3095 (2.36-2.32)
Sidechain outliers	187428	3095 (2.36-2.32)
RSRZ outliers	180081	3033 (2.36-2.32)

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	989	2% 89% 7% ..
1	B	989	% 88% 8% ..

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 16292 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	955	Total	C	N	O	S	0	1	0
			7812	5035	1310	1445	22			
1	B	959	Total	C	N	O	S	0	1	0
			7848	5054	1317	1455	22			

There are 50 discrepancies between the modelled and reference sequences:

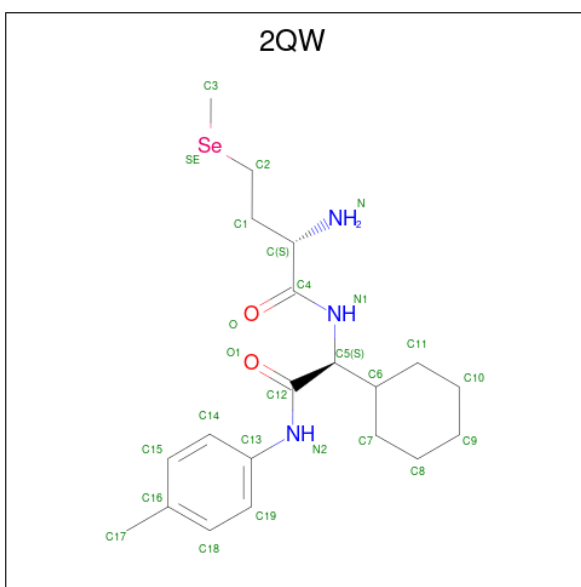
Chain	Residue	Modelled	Actual	Comment	Reference
A	31	MET	-	initiating methionine	UNP P14735
A	32	GLY	-	expression tag	UNP P14735
A	33	HIS	-	expression tag	UNP P14735
A	34	HIS	-	expression tag	UNP P14735
A	35	HIS	-	expression tag	UNP P14735
A	36	HIS	-	expression tag	UNP P14735
A	37	HIS	-	expression tag	UNP P14735
A	38	HIS	-	expression tag	UNP P14735
A	39	GLY	-	expression tag	UNP P14735
A	40	ARG	-	expression tag	UNP P14735
A	41	ALA	-	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
A	974	ALA	CYS	engineered mutation	UNP P14735

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Chain	Residue	Modelled	Actual	Comment	Reference
B	31	MET	-	initiating methionine	UNP P14735
B	32	GLY	-	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	HIS	-	expression tag	UNP P14735
B	38	HIS	-	expression tag	UNP P14735
B	39	GLY	-	expression tag	UNP P14735
B	40	ARG	-	expression tag	UNP P14735
B	41	ALA	-	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 (2S)-2-amino-N-{(1S)-1-cyclohexyl-2-[(4-methylphenyl)amino]-2-oxoethyl}-4-(methylselanyl)butanamide (CCD ID: 2QW) (formula: C₂₀H₃₁N₃O₂Se).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	Se	0	0
			26	20	3	2	1		
2	B	1	Total	C	N	O	Se	0	0
			26	20	3	2	1		

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

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

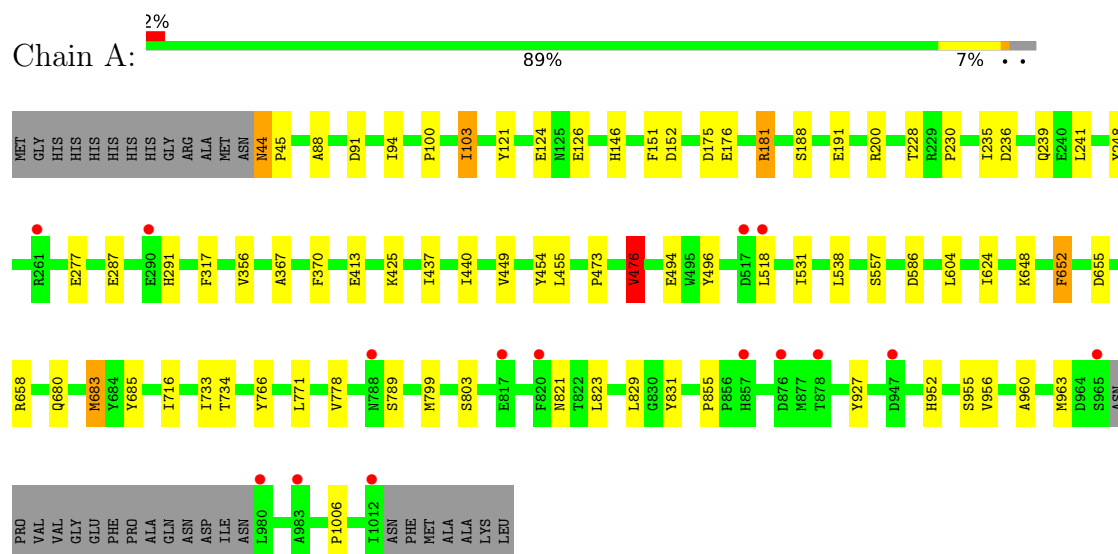
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	289	Total	O	0	0
			289	289		
4	B	289	Total	O	0	0
			289	289		

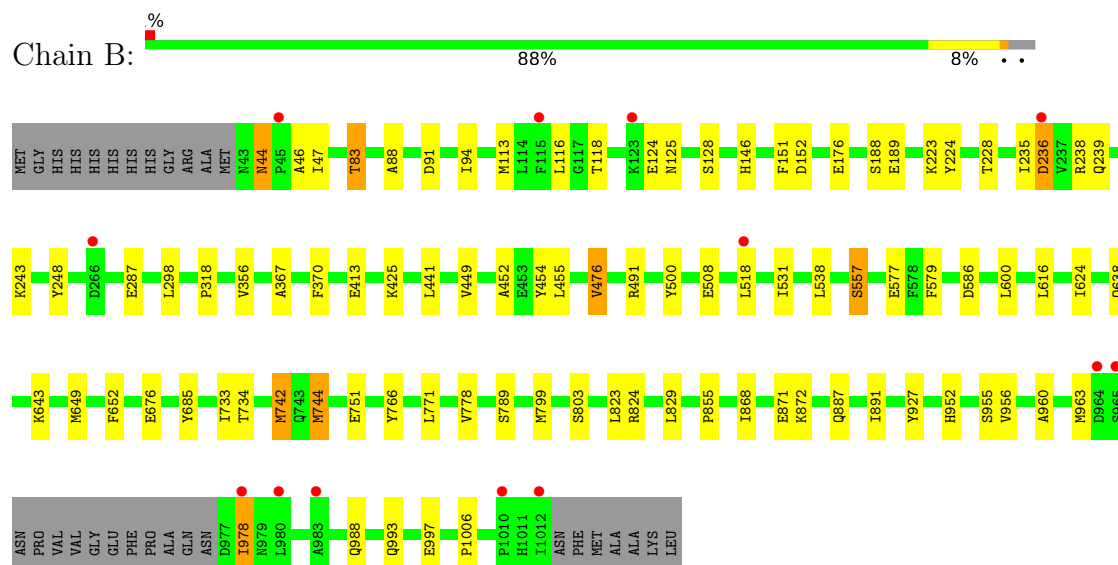
3 Residue-property plots [i](#)

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: Insulin-degrading enzyme



• Molecule 1: Insulin-degrading enzyme



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	78.84Å 115.96Å 124.28Å 90.00° 97.83° 90.00°	Depositor
Resolution (Å)	19.94 – 2.33 19.94 – 2.33	Depositor EDS
% Data completeness (in resolution range)	99.7 (19.94-2.33) 99.5 (19.94-2.33)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.14	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.22 (at 2.33Å)	Xtriage
Refinement program	BUSTER 2.11.2	Depositor
R, R_{free}	0.199 , 0.236 0.205 , 0.244	Depositor DCC
R_{free} test set	1050 reflections (1.12%)	wwPDB-VP
Wilson B-factor (Å ²)	38.2	Xtriage
Anisotropy	0.265	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 35.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	16292	wwPDB-VP
Average B, all atoms (Å ²)	45.0	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: 2QW, 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	0.81	3/8008 (0.0%)	1.30	11/10835 (0.1%)
1	B	0.79	1/8046 (0.0%)	1.30	16/10884 (0.1%)
All	All	0.80	4/16054 (0.0%)	1.30	27/21719 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	683	MET	SD-CE	-7.80	1.60	1.79
1	B	742	MET	SD-CE	-5.52	1.65	1.79
1	A	716	ILE	CA-CB	5.33	1.56	1.54
1	A	44	ASN	C-N	5.25	1.38	1.33

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	152	ASP	CA-CB-CG	7.26	119.86	112.60
1	A	152	ASP	CA-CB-CG	7.10	119.70	112.60
1	A	557	SER	N-CA-C	6.84	120.44	108.75
1	B	557	SER	N-CA-C	6.75	120.29	108.75
1	B	855	PRO	N-CA-C	5.95	116.18	110.47
1	A	855	PRO	N-CA-C	5.89	116.13	110.47
1	A	476	VAL	N-CA-C	5.81	117.56	109.55
1	B	455	LEU	N-CA-C	5.73	118.88	109.76
1	A	455	LEU	N-CA-C	5.70	118.82	109.76
1	B	586	ASP	CA-CB-CG	5.55	118.15	112.60
1	A	586	ASP	CA-CB-CG	5.54	118.14	112.60
1	B	652	PHE	N-CA-C	5.43	118.05	109.96
1	A	494	GLU	N-CA-C	5.41	118.09	111.82
1	B	188	SER	CA-C-N	5.24	127.26	120.44
1	B	188	SER	C-N-CA	5.24	127.26	120.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	871	GLU	CA-C-N	5.18	127.17	120.44
1	B	871	GLU	C-N-CA	5.18	127.17	120.44
1	B	243	LYS	CA-C-N	5.14	127.12	120.44
1	B	243	LYS	C-N-CA	5.14	127.12	120.44
1	A	241	LEU	CA-C-N	5.11	127.13	120.28
1	A	241	LEU	C-N-CA	5.11	127.13	120.28
1	B	476	VAL	N-CA-C	5.11	116.60	109.55
1	B	235	ILE	CA-C-N	5.06	131.22	122.32
1	B	235	ILE	C-N-CA	5.06	131.22	122.32
1	A	191	GLU	CA-C-N	5.02	127.28	120.65
1	A	191	GLU	C-N-CA	5.02	127.28	120.65
1	B	83	THR	N-CA-C	5.01	117.61	110.24

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	7812	0	7741	36	0
1	B	7848	0	7791	31	0
2	A	26	0	26	0	0
2	B	26	0	26	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	289	0	0	1	0
4	B	289	0	0	0	0
All	All	16292	0	15584	67	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 2.

All (67) 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:44:ASN:HB3	1:A:45:PRO:HD3	1.57	0.86
1:B:557:SER:HB2	1:B:742:MET:HE3	1.68	0.75
1:A:648:LYS:O	1:A:652:PHE:HB3	1.88	0.73
1:A:317:PHE:CE2	1:A:476:VAL:HG13	2.31	0.66
1:A:680:GLN:HA	1:A:683:MET:HE3	1.81	0.62
1:A:236:ASP:HB3	1:A:239:GLN:HG2	1.81	0.61
1:A:124:GLU:OE2	1:A:181:ARG:NH1	2.31	0.60
1:A:440:ILE:HG22	1:A:449:VAL:HG23	1.85	0.59
1:A:473:PRO:O	1:A:476:VAL:HG22	2.05	0.56
1:B:298:LEU:HD21	1:B:318:PRO:HG3	1.90	0.54
1:B:425:LYS:HD3	1:B:454:TYR:OH	2.08	0.54
1:B:236:ASP:HB2	1:B:239:GLN:HB3	1.90	0.52
1:A:655:ASP:OD2	1:A:658:ARG:HG2	2.08	0.52
1:A:437:ILE:HG23	1:A:449:VAL:CG2	2.39	0.52
1:A:799:MET:HE1	1:A:1006:PRO:HG2	1.92	0.52
1:A:277:GLU:HB2	4:A:1313:HOH:O	2.10	0.52
1:A:94:ILE:HG13	1:A:248:TYR:HB3	1.92	0.51
1:B:799:MET:HE1	1:B:1006:PRO:HG2	1.92	0.51
1:A:960:ALA:HB3	1:A:963:MET:HB2	1.91	0.51
1:B:643:LYS:HG3	1:B:744:MET:HE3	1.93	0.50
1:B:413:GLU:HG2	1:B:531:ILE:HD11	1.93	0.50
1:B:778:VAL:HG22	1:B:955:SER:HB2	1.92	0.50
1:B:960:ALA:HB3	1:B:963:MET:HB2	1.92	0.50
1:A:437:ILE:HG23	1:A:449:VAL:HG21	1.94	0.50
1:B:224:TYR:HA	1:B:228:THR:HB	1.93	0.49
1:A:648:LYS:HG3	1:A:652:PHE:HB2	1.94	0.49
1:A:771:LEU:HB2	1:A:952:HIS:HB3	1.95	0.49
1:B:367:ALA:HB3	1:B:370:PHE:CE2	2.47	0.49
1:B:771:LEU:HB2	1:B:952:HIS:HB3	1.95	0.48
1:A:44:ASN:HB3	1:A:45:PRO:CD	2.37	0.48
1:A:778:VAL:HG22	1:A:955:SER:HB2	1.95	0.48
1:A:425:LYS:HE2	1:A:454:TYR:OH	2.15	0.47
1:B:643:LYS:HA	1:B:744:MET:HE1	1.96	0.47
1:B:441:LEU:HD23	1:B:449:VAL:HG11	1.96	0.47
1:A:367:ALA:HB3	1:A:370:PHE:CE2	2.50	0.47
1:B:685:TYR:HB2	1:B:956:VAL:HG11	1.96	0.46
1:A:188:SER:HB3	1:A:831:TYR:HD1	1.80	0.46
1:A:413:GLU:HG2	1:A:531:ILE:HD11	1.98	0.46
1:B:491:ARG:HH11	1:B:500:TYR:HE2	1.61	0.46
1:A:648:LYS:HG3	1:A:652:PHE:CB	2.46	0.45
1:B:887:GLN:O	1:B:891:ILE:HG12	2.17	0.45
1:B:91:ASP:OD2	1:B:146:HIS:ND1	2.45	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:100:PRO:HG2	1:A:103:ILE:HD12	1.99	0.44
1:A:685:TYR:HB2	1:A:956:VAL:HG11	1.99	0.44
1:A:538:LEU:HD13	1:A:734:THR:HG23	1.99	0.44
1:B:538:LEU:HD13	1:B:734:THR:HG23	1.99	0.44
1:A:88:ALA:HB3	1:A:151:PHE:CE2	2.53	0.43
1:B:643:LYS:HA	1:B:744:MET:CE	2.49	0.43
1:A:823:LEU:HB3	1:A:829:LEU:HD12	2.01	0.43
1:B:88:ALA:HB3	1:B:151:PHE:CE2	2.53	0.43
1:B:823:LEU:HB3	1:B:829:LEU:HD12	2.00	0.43
1:B:577:GLU:HG2	1:B:579:PHE:CZ	2.54	0.43
1:B:803:SER:HA	1:B:927:TYR:CE2	2.53	0.43
1:B:44:ASN:HD22	1:B:46:ALA:H	1.66	0.43
1:A:291:HIS:CD2	1:A:370:PHE:HB2	2.53	0.42
1:A:624:ILE:HD12	1:A:766:TYR:HD1	1.84	0.42
1:A:200:ARG:HG2	1:A:496:TYR:HB3	2.02	0.42
1:A:803:SER:HA	1:A:927:TYR:CE2	2.54	0.42
1:B:600:LEU:HD21	1:B:649:MET:HG3	2.01	0.42
1:B:94:ILE:HG13	1:B:248:TYR:HB3	2.01	0.42
1:B:113:MET:HA	1:B:116:LEU:HD12	2.01	0.42
1:A:91:ASP:OD2	1:A:146:HIS:ND1	2.46	0.41
1:B:616:LEU:HD11	1:B:638:GLN:HG3	2.02	0.41
1:A:230:PRO:HB3	1:A:235:ILE:HB	2.02	0.41
1:B:624:ILE:HD12	1:B:766:TYR:HD1	1.84	0.41
1:A:121:TYR:HB3	1:A:126:GLU:HG2	2.01	0.41
1:B:44:ASN:HB3	1:B:47:ILE:HD12	2.02	0.40

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	952/989 (96%)	918 (96%)	33 (4%)	1 (0%)	48 57

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	956/989 (97%)	924 (97%)	30 (3%)	2 (0%)	43	50
All	All	1908/1978 (96%)	1842 (96%)	63 (3%)	3 (0%)	43	50

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	175	ASP
1	B	452	ALA
1	B	978	ILE

5.3.2 Protein sidechains [i](#)

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	849/878 (97%)	836 (98%)	13 (2%)	57	69
1	B	856/878 (98%)	828 (97%)	28 (3%)	33	43
All	All	1705/1756 (97%)	1664 (98%)	41 (2%)	43	55

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	103	ILE
1	A	176	GLU
1	A	181	ARG
1	A	228	THR
1	A	287	GLU
1	A	356	VAL
1	A	476	VAL
1	A	518	LEU
1	A	604	LEU
1	A	652	PHE
1	A	733	ILE
1	A	789	SER
1	A	821	ASN

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Mol	Chain	Res	Type
1	B	44	ASN
1	B	83	THR
1	B	118	THR
1	B	124	GLU
1	B	125	ASN
1	B	128	SER
1	B	176	GLU
1	B	189	GLU
1	B	223	LYS
1	B	236	ASP
1	B	238	ARG
1	B	287	GLU
1	B	356	VAL
1	B	476	VAL
1	B	508	GLU
1	B	518	LEU
1	B	676	GLU
1	B	733	ILE
1	B	744	MET
1	B	751	GLU
1	B	789	SER
1	B	824	ARG
1	B	868	ILE
1	B	872	LYS
1	B	978	ILE
1	B	988	GLN
1	B	993	GLN
1	B	997	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	210	ASN
1	A	231	ASN
1	A	232	GLN
1	A	252	ASN
1	A	312	ASN
1	A	376	ASN
1	A	407	GLN
1	A	671	ASN
1	A	788	ASN
1	A	813	GLN

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Mol	Chain	Res	Type
1	A	917	ASN
1	A	988	GLN
1	B	102	ASN
1	B	125	ASN
1	B	129	GLN
1	B	231	ASN
1	B	232	GLN
1	B	239	GLN
1	B	252	ASN
1	B	312	ASN
1	B	376	ASN
1	B	407	GLN
1	B	671	ASN
1	B	770	GLN
1	B	788	ASN
1	B	813	GLN
1	B	917	ASN
1	B	988	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 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	$\# Z > 2$
2	2QW	A	1101	-	25,27,27	1.31	4 (16%)	31,35,35	1.05	2 (6%)
2	2QW	B	1101	-	25,27,27	1.42	3 (12%)	31,35,35	0.95	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
2	2QW	A	1101	-	-	2/23/32/32	0/2/2/2
2	2QW	B	1101	-	-	3/23/32/32	0/2/2/2

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1101	2QW	C4-N1	3.72	1.42	1.34
2	A	1101	2QW	C4-N1	3.70	1.42	1.34
2	B	1101	2QW	C12-N2	2.94	1.42	1.35
2	A	1101	2QW	C14-C13	2.23	1.43	1.39
2	A	1101	2QW	C19-C13	2.21	1.43	1.39
2	A	1101	2QW	C11-C6	2.08	1.58	1.53
2	B	1101	2QW	C14-C15	2.02	1.42	1.38

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1101	2QW	C7-C6-C5	2.16	114.72	111.85
2	A	1101	2QW	C11-C6-C7	2.01	112.87	109.43

There are no chirality outliers.

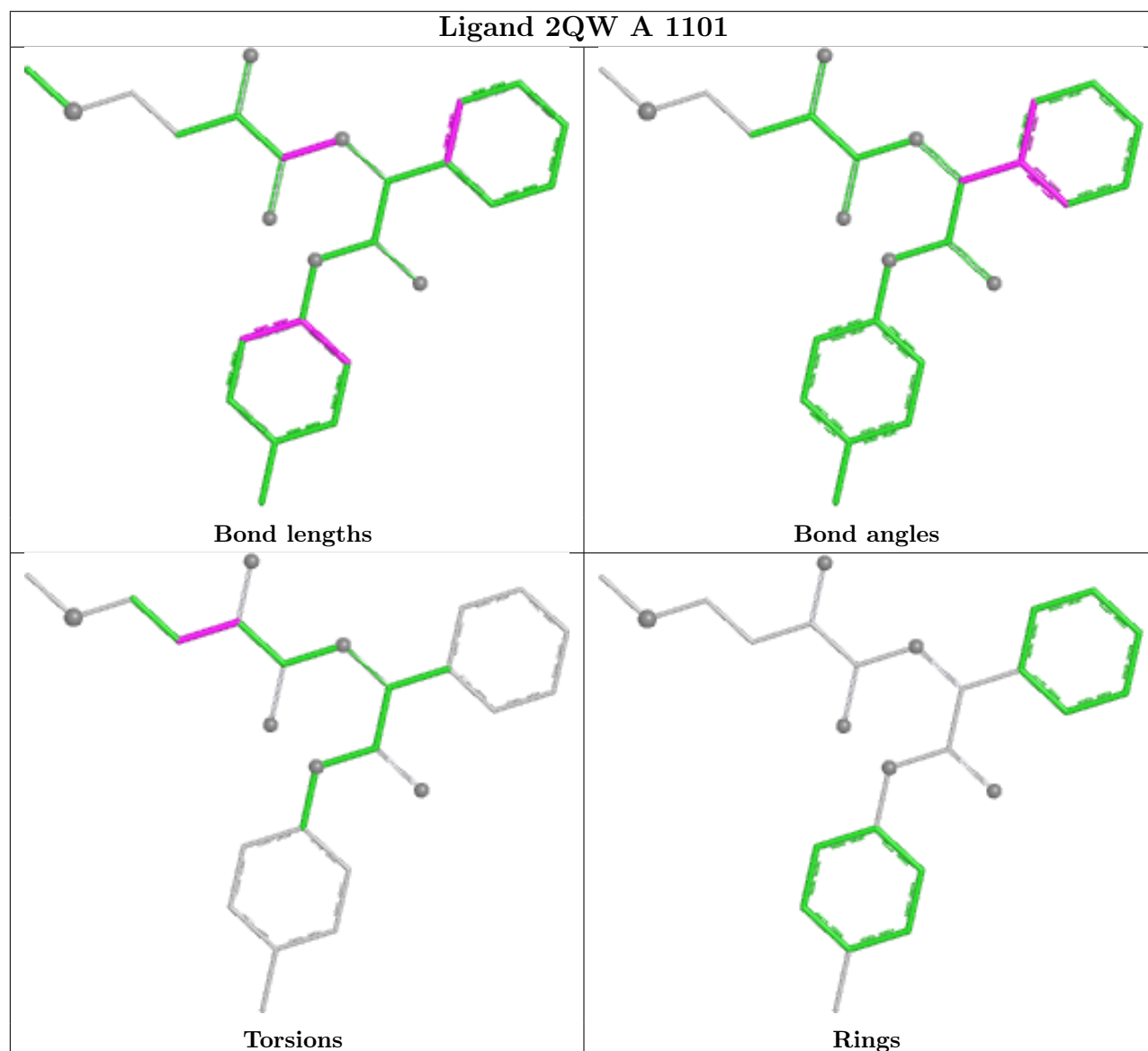
All (5) torsion outliers are listed below:

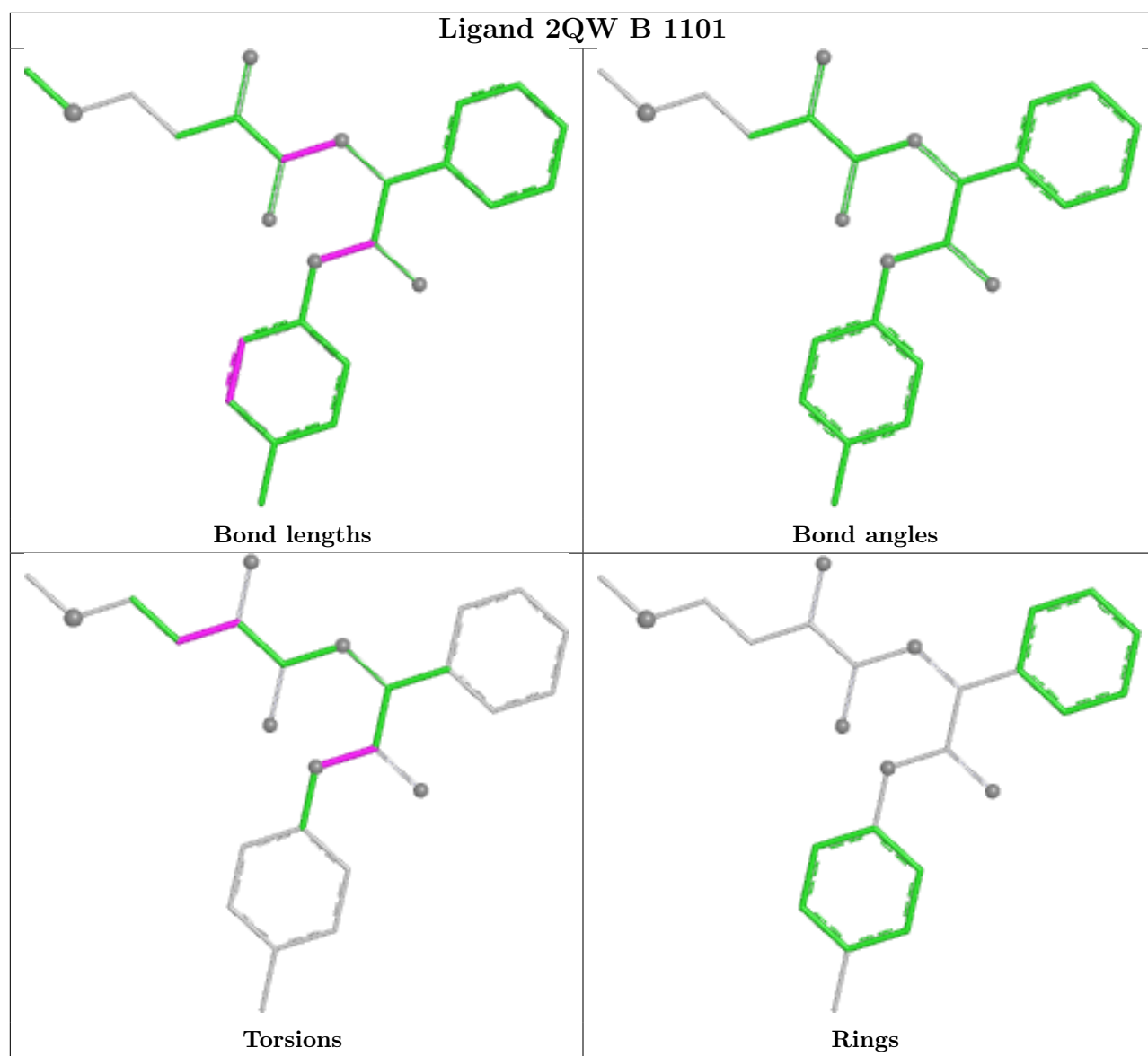
Mol	Chain	Res	Type	Atoms
2	B	1101	2QW	C4-C-C1-C2
2	B	1101	2QW	N-C-C1-C2
2	A	1101	2QW	N-C-C1-C2
2	A	1101	2QW	C4-C-C1-C2
2	B	1101	2QW	O1-C12-N2-C13

There are no ring outliers.

No monomer is involved in short contacts.

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





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	955/989 (96%)	0.17	15 (1%) 70 75	13, 43, 72, 93	1 (0%)
1	B	959/989 (96%)	0.06	13 (1%) 73 77	25, 42, 68, 117	1 (0%)
All	All	1914/1978 (96%)	0.12	28 (1%) 72 76	13, 43, 71, 117	2 (0%)

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1012	ILE	3.6
1	A	980	LEU	3.4
1	A	261	ARG	3.0
1	A	518	LEU	2.9
1	A	817	GLU	2.9
1	A	878	THR	2.9
1	B	965	SER	2.8
1	B	1012	ILE	2.7
1	A	965	SER	2.6
1	A	876	ASP	2.6
1	A	983	ALA	2.6
1	B	978	ILE	2.6
1	A	947	ASP	2.5
1	B	236	ASP	2.4
1	B	45	PRO	2.4
1	A	517	ASP	2.4
1	A	857	HIS	2.4
1	B	518	LEU	2.2
1	B	983	ALA	2.2
1	B	115	PHE	2.2
1	B	980	LEU	2.2
1	A	788	ASN	2.2
1	B	266	ASP	2.2
1	B	123	LYS	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	820	PHE	2.1
1	A	290	GLU	2.0
1	B	1010	PRO	2.0
1	B	964	ASP	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

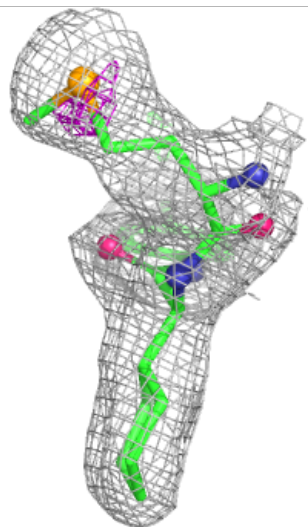
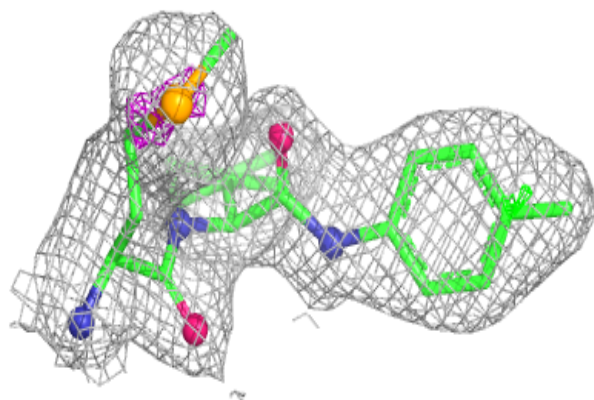
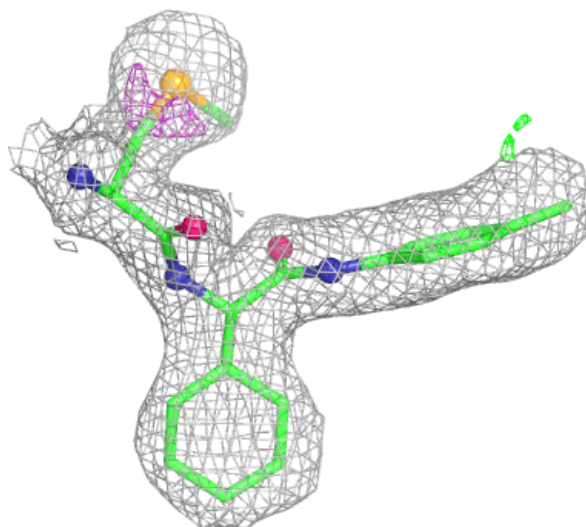
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	2QW	B	1101	26/26	0.93	0.08	30,34,45,53	0
2	2QW	A	1101	26/26	0.95	0.09	32,38,50,56	0
3	ZN	B	1102	1/1	0.98	0.05	86,86,86,86	0
3	ZN	A	1102	1/1	0.99	0.05	72,72,72,72	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.

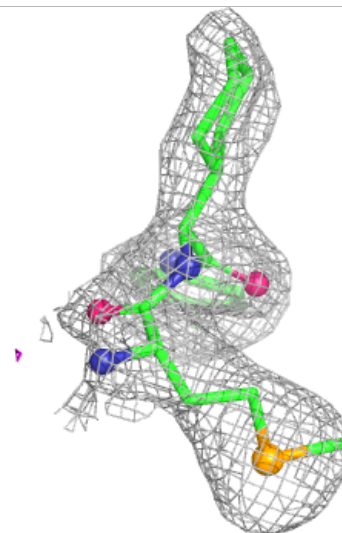
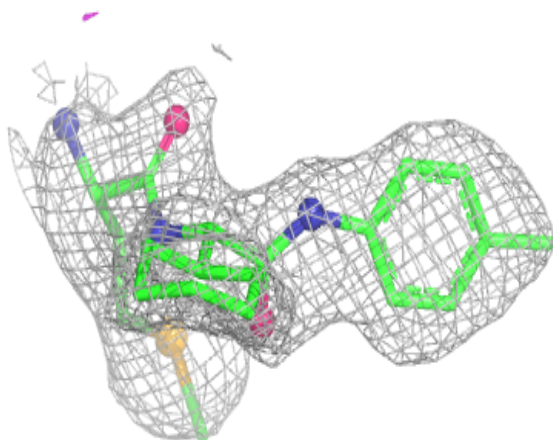
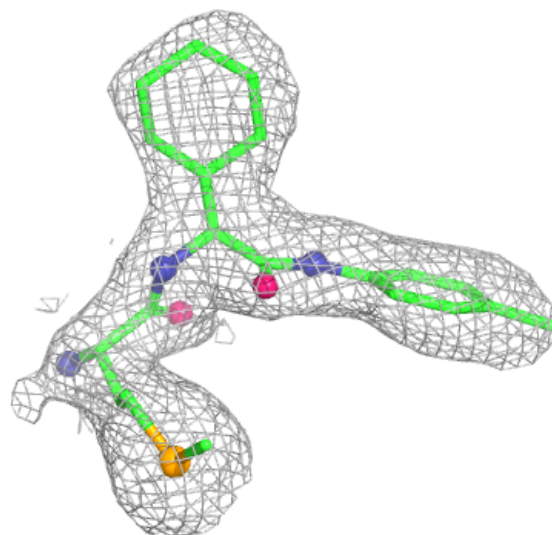
Electron density around 2QW B 1101:

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



Electron density around 2QW A 1101:

$2mF_o-DF_c$ (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.