



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 08:47 PM UTC

PDB ID : 4LKO / pdb_00004lko
Title : Crystal structure of human DPP-IV in complex with BMS-744891
Authors : Klei, H.E.
Deposited on : 2013-07-08
Resolution : 2.43 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

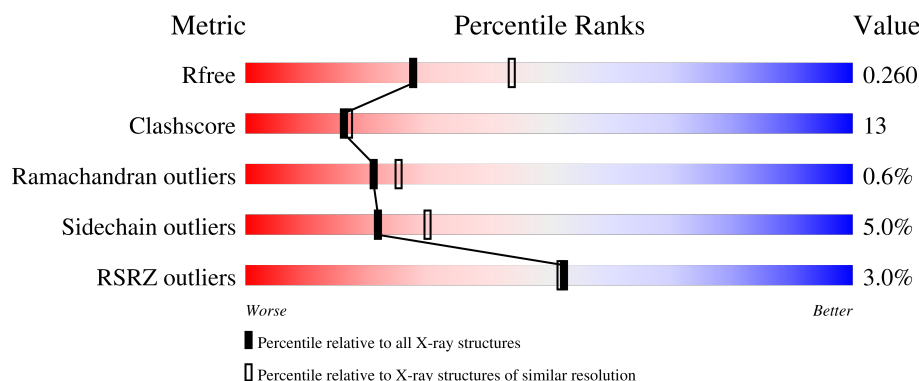
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

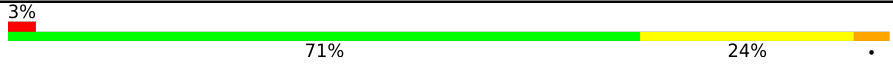
The reported resolution of this entry is 2.43 Å.

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	2340 (2.46-2.42)
Clashscore	190562	2400 (2.46-2.42)
Ramachandran outliers	187476	2379 (2.46-2.42)
Sidechain outliers	187428	2379 (2.46-2.42)
RSRZ outliers	180081	2340 (2.46-2.42)

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	728	
1	B	728	

2 Entry composition [i](#)

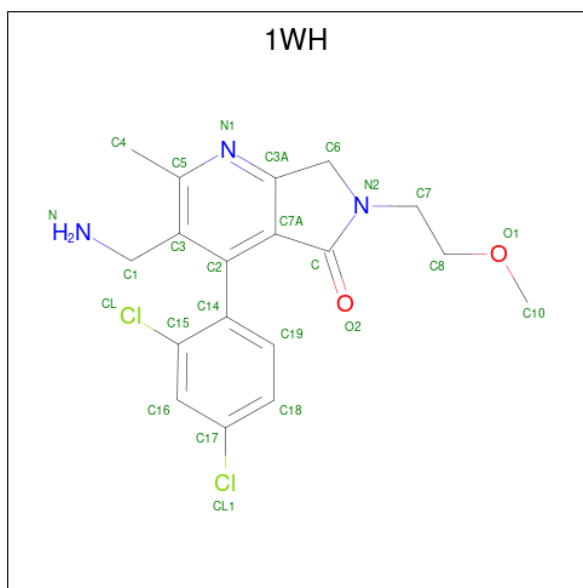
There are 3 unique types of molecules in this entry. The entry contains 12074 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Dipeptidyl peptidase 4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	728	Total	C	N	O	S	0	0	0
			5963	3827	982	1128	26			
1	B	728	Total	C	N	O	S	0	0	0
			5963	3827	982	1128	26			

- Molecule 2 is 3-(aminomethyl)-4-(2,4-dichlorophenyl)-6-(2-methoxyethyl)-2-methyl-6,7-dihydro-5H-pyrrolo[3,4-b]pyridin-5-one (CCD ID: 1WH) (formula: C₁₈H₁₉Cl₂N₃O₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Cl	N	O	0	0
			25	18	2	3	2		
2	B	1	Total	C	Cl	N	O	0	0
			25	18	2	3	2		

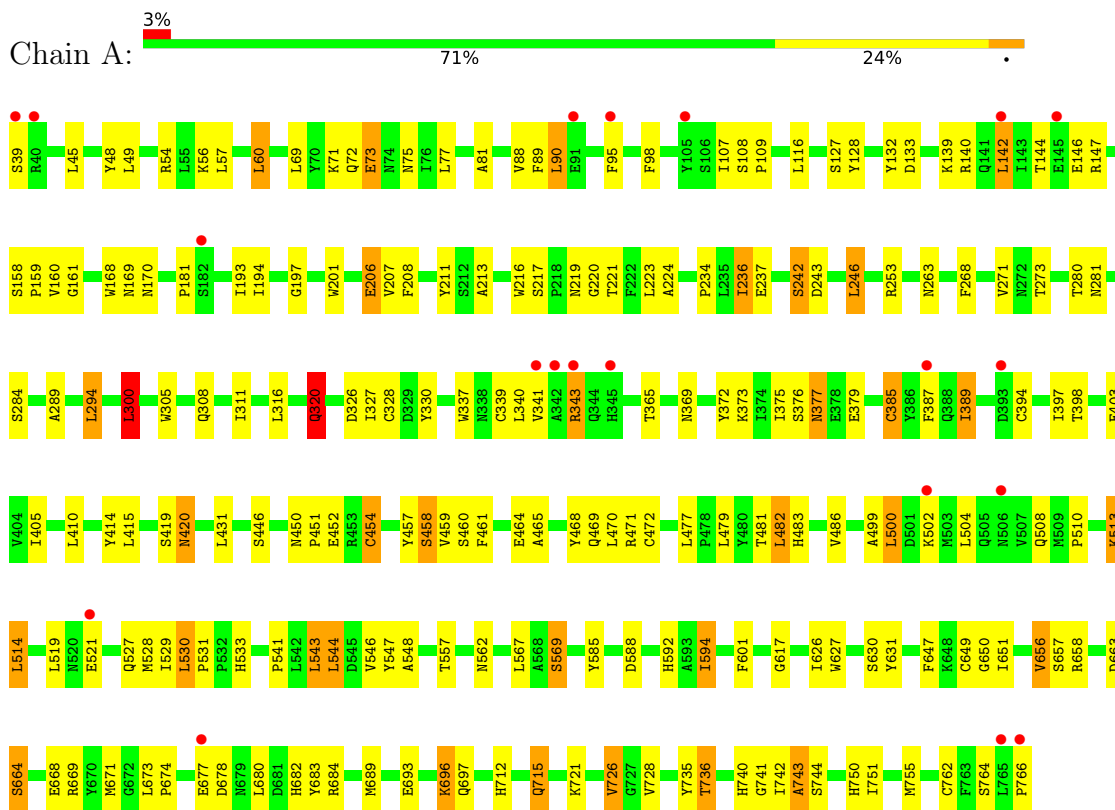
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	63	Total 63	O 63	0	0
3	B	35	Total 35	O 35	0	0

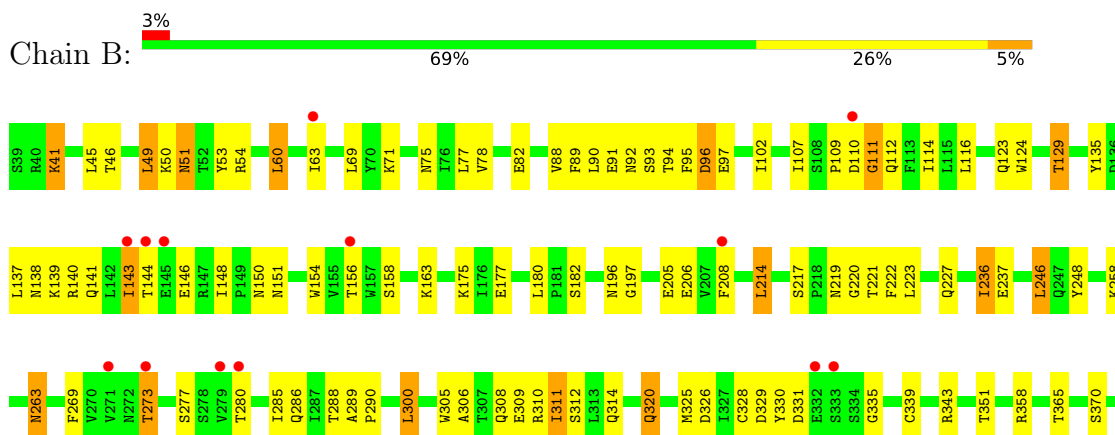
3 Residue-property plots

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: Dipeptidyl peptidase 4



• Molecule 1: Dipeptidyl peptidase 4





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	65.90Å 68.30Å 421.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.43 – 2.43 47.43 – 2.44	Depositor EDS
% Data completeness (in resolution range)	96.4 (47.43-2.43) 96.5 (47.43-2.44)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.63 (at 2.42Å)	Xtriage
Refinement program	CNS, CNX 2002	Depositor
R, R_{free}	0.226 , 0.263 0.226 , 0.260	Depositor DCC
R_{free} test set	4996 reflections (7.12%)	wwPDB-VP
Wilson B-factor (Å ²)	50.5	Xtriage
Anisotropy	0.543	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 23.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.064 for k,h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	12074	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.07% 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: 1WH

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	0.46	0/6135	0.99	35/8344 (0.4%)
1	B	0.43	0/6135	0.97	20/8344 (0.2%)
All	All	0.45	0/12270	0.98	55/16688 (0.3%)

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

There are no bond length outliers.

All (55) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	389	ILE	N-CA-C	12.52	123.39	110.62
1	A	546	VAL	N-CA-C	9.41	122.04	108.48
1	A	220	GLY	N-CA-C	-8.92	102.34	115.63
1	B	148	ILE	N-CA-C	-7.73	101.37	108.95
1	A	458	SER	N-CA-C	-7.63	97.72	109.24
1	A	450	ASN	CA-C-N	7.59	129.33	119.84
1	A	450	ASN	C-N-CA	7.59	129.33	119.84
1	B	546	VAL	N-CA-C	7.35	120.06	108.89
1	B	300	LEU	N-CA-C	-7.21	97.49	108.96
1	A	530	LEU	CA-C-N	6.98	124.75	119.66
1	A	530	LEU	C-N-CA	6.98	124.75	119.66
1	A	300	LEU	N-CA-C	-6.87	97.53	108.73
1	B	458	SER	N-CA-C	-6.79	98.99	109.24
1	B	584	GLY	N-CA-C	6.76	121.83	112.57

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	158	SER	N-CA-C	-6.65	99.09	109.87
1	B	656	VAL	N-CA-C	-6.58	98.72	109.12
1	B	220	GLY	N-CA-C	-6.52	105.91	115.63
1	B	683	TYR	N-CA-C	-6.52	104.10	111.14
1	B	743	ALA	N-CA-C	6.51	121.38	113.38
1	A	656	VAL	N-CA-C	-6.33	99.12	109.12
1	A	521	GLU	N-CA-C	-6.27	103.67	113.02
1	A	465	ALA	N-CA-C	6.17	119.93	111.54
1	A	548	ALA	N-CA-C	6.17	120.39	112.86
1	B	206	GLU	N-CA-C	6.11	120.89	113.38
1	A	454	CYS	N-CA-C	6.01	118.55	108.02
1	A	683	TYR	N-CA-C	-5.97	104.85	111.36
1	B	414	TYR	N-CA-C	5.86	118.08	109.24
1	B	143	ILE	N-CA-C	-5.81	101.53	109.37
1	A	207	VAL	N-CA-C	5.80	116.94	111.48
1	A	201	TRP	N-CA-C	5.77	117.24	111.07
1	A	477	LEU	CA-C-N	5.77	126.10	119.93
1	A	477	LEU	C-N-CA	5.77	126.10	119.93
1	A	664	SER	N-CA-C	5.75	117.22	111.07
1	A	206	GLU	N-CA-C	5.67	120.47	113.50
1	A	320	GLN	N-CA-C	5.64	122.82	110.80
1	A	529	ILE	N-CA-C	-5.62	98.97	107.73
1	A	541	PRO	N-CA-C	-5.53	103.11	111.41
1	A	365	THR	N-CA-C	-5.48	103.15	110.55
1	B	205	GLU	N-CA-C	5.46	117.04	111.14
1	A	414	TYR	N-CA-C	5.46	117.48	109.24
1	B	595	ASN	N-CA-C	5.39	118.39	110.30
1	A	242	SER	CB-CA-C	-5.39	108.77	115.89
1	A	54	ARG	N-CA-C	5.39	118.41	109.46
1	A	211	TYR	N-CA-C	-5.35	105.89	112.90
1	A	569	SER	N-CA-C	5.34	117.19	111.36
1	A	81	ALA	N-CA-C	5.31	117.07	111.28
1	A	217	SER	N-CA-C	-5.28	103.29	110.36
1	B	525	TRP	N-CA-C	5.27	118.28	110.48
1	B	365	THR	N-CA-C	-5.18	103.65	110.33
1	B	454	CYS	N-CA-C	5.09	116.81	108.52
1	A	108	SER	CA-C-N	5.08	124.87	119.28
1	A	108	SER	C-N-CA	5.08	124.87	119.28
1	A	647	PHE	N-CA-C	5.06	117.22	109.07
1	B	547	TYR	N-CA-C	-5.06	100.02	110.80
1	A	743	ALA	N-CA-C	5.01	119.40	113.28

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	547	TYR	Sidechain

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	5963	0	5695	142	0
1	B	5963	0	5695	165	0
2	A	25	0	19	0	0
2	B	25	0	19	0	0
3	A	63	0	0	1	0
3	B	35	0	0	0	0
All	All	12074	0	11428	300	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 13.

All (300) 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:508:GLN:HE21	1:A:533:HIS:CE1	1.76	1.04
1:B:600:THR:HG23	1:B:601:PHE:H	1.22	0.99
1:B:51:ASN:HD22	1:B:54:ARG:HG2	1.32	0.95
1:B:177:GLU:HB2	1:B:180:LEU:HD23	1.49	0.95
1:A:508:GLN:HE21	1:A:533:HIS:HE1	0.99	0.93
1:B:649:CYS:HG	1:B:762:CYS:HG	1.19	0.89
1:A:385:CYS:HG	1:A:394:CYS:HG	1.21	0.89
1:A:528:MET:HE3	1:A:530:LEU:HD21	1.56	0.87
1:A:403:GLU:H	1:A:420:ASN:HD21	1.24	0.85
1:B:600:THR:HG23	1:B:601:PHE:N	1.91	0.83
1:B:77:LEU:HD23	1:B:88:VAL:HA	1.61	0.82
1:A:39:SER:HB2	1:A:508:GLN:HG3	1.59	0.82
1:A:89:PHE:HE2	1:A:107:ILE:HD13	1.45	0.81
1:A:377:ASN:C	1:A:377:ASN:HD22	1.88	0.81
1:A:693:GLU:O	1:A:696:LYS:HG3	1.81	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:328:CYS:HG	1:A:339:CYS:HG	1.16	0.80
1:B:403:GLU:H	1:B:420:ASN:HD21	1.29	0.80
1:A:508:GLN:NE2	1:A:533:HIS:HE1	1.80	0.79
1:B:600:THR:CG2	1:B:601:PHE:H	1.96	0.79
1:B:89:PHE:HE2	1:B:107:ILE:HD13	1.47	0.78
1:A:649:CYS:HG	1:A:762:CYS:HG	0.77	0.77
1:A:340:LEU:HB3	1:A:343:ARG:HD3	1.66	0.77
1:B:471:ARG:HG3	1:B:480:TYR:CE2	2.20	0.76
1:B:221:THR:O	1:B:273:THR:HG22	1.85	0.76
1:B:536:LYS:H	1:B:536:LYS:HD3	1.49	0.76
1:B:320:GLN:OE1	1:B:669:ARG:HD3	1.86	0.76
1:A:657:SER:H	1:A:715:GLN:NE2	1.84	0.75
1:A:142:LEU:HD12	1:A:142:LEU:H	1.52	0.75
1:B:726:VAL:HG13	1:B:728:VAL:HG23	1.69	0.74
1:A:726:VAL:HG12	1:A:728:VAL:HG23	1.69	0.74
1:B:735:TYR:OH	1:B:750:HIS:HD2	1.71	0.73
1:B:597:ARG:O	1:B:600:THR:HG22	1.87	0.73
1:A:499:ALA:HA	1:A:502:LYS:HE3	1.70	0.72
1:A:751:ILE:HG12	1:A:755:MET:HE2	1.71	0.72
1:B:236:ILE:HD13	1:B:237:GLU:N	2.04	0.72
1:B:139:LYS:C	1:B:140:ARG:HD2	2.15	0.71
1:A:236:ILE:HG13	1:A:712:HIS:CE1	2.26	0.71
1:B:377:ASN:C	1:B:377:ASN:HD22	1.97	0.70
1:B:358:ARG:HB3	1:B:358:ARG:HH21	1.55	0.70
1:B:219:ASN:HB2	1:B:308:GLN:OE1	1.91	0.70
1:B:358:ARG:HB3	1:B:358:ARG:NH2	2.07	0.70
1:B:657:SER:H	1:B:715:GLN:NE2	1.89	0.70
1:A:237:GLU:OE2	1:A:253:ARG:HD3	1.92	0.69
1:B:217:SER:OG	1:B:222:PHE:HB2	1.92	0.69
1:B:236:ILE:HG13	1:B:712:HIS:CE1	2.27	0.69
1:B:89:PHE:CE2	1:B:107:ILE:HD13	2.27	0.68
1:B:175:LYS:HG3	1:B:182:SER:HB3	1.75	0.68
1:B:677:GLU:CD	1:B:677:GLU:H	2.01	0.67
1:A:500:LEU:HD22	1:A:504:LEU:HG	1.76	0.67
1:B:481:THR:OG1	1:B:483:HIS:HE1	1.77	0.67
1:A:89:PHE:CE2	1:A:107:ILE:HD13	2.29	0.67
1:A:508:GLN:NE2	1:A:533:HIS:CE1	2.58	0.66
1:B:640:LEU:HD11	1:B:650:GLY:HA3	1.78	0.66
1:B:312:SER:OG	1:B:325:MET:HE3	1.97	0.65
1:A:237:GLU:HG2	1:A:253:ARG:HG2	1.80	0.64
1:A:696:LYS:HD2	1:A:697:GLN:HG3	1.79	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:196:ASN:OD1	1:B:227:GLN:HG3	1.98	0.64
1:A:343:ARG:HD2	1:A:343:ARG:N	2.13	0.64
1:A:481:THR:OG1	1:A:483:HIS:HE1	1.81	0.63
1:B:78:VAL:HG23	1:B:89:PHE:HB2	1.80	0.63
1:B:558:VAL:HG22	1:B:560:ARG:NH1	2.13	0.63
1:A:735:TYR:OH	1:A:750:HIS:HD2	1.81	0.62
1:B:51:ASN:ND2	1:B:54:ARG:HG2	2.12	0.62
1:B:91:GLU:CD	1:B:91:GLU:H	2.07	0.62
1:B:60:LEU:H	1:B:60:LEU:HD22	1.64	0.62
1:A:343:ARG:HD2	1:A:343:ARG:H	1.63	0.62
1:B:102:ILE:HD13	1:B:116:LEU:HD22	1.81	0.62
1:B:680:LEU:HD13	1:B:680:LEU:C	2.26	0.61
1:B:111:GLY:O	1:B:137:LEU:HD12	2.00	0.61
1:B:306:ALA:HB3	1:B:310:ARG:HG2	1.81	0.61
1:A:142:LEU:HD12	1:A:142:LEU:N	2.15	0.60
1:A:377:ASN:ND2	1:A:379:GLU:H	1.99	0.60
1:A:377:ASN:C	1:A:377:ASN:ND2	2.54	0.60
1:A:454:CYS:CB	1:A:472:CYS:HG	2.14	0.60
1:B:143:ILE:HG13	1:B:143:ILE:O	2.01	0.59
1:A:139:LYS:O	1:A:140:ARG:HB2	2.01	0.59
1:B:177:GLU:CB	1:B:180:LEU:HD23	2.30	0.59
1:A:510:PRO:HD3	1:A:569:SER:HB2	1.85	0.58
1:B:514:LEU:HD12	1:B:557:THR:HG22	1.85	0.58
1:A:160:VAL:HG12	1:A:161:GLY:N	2.18	0.58
1:A:193:ILE:HG22	1:A:194:ILE:HG12	1.85	0.57
1:B:269:PHE:CE1	1:B:286:GLN:HB2	2.39	0.57
1:A:405:ILE:HD13	1:A:419:SER:HA	1.84	0.57
1:B:46:THR:CG2	1:B:50:LYS:HD3	2.35	0.57
1:A:289:ALA:HB3	1:A:294:LEU:CD1	2.35	0.57
1:B:415:LEU:C	1:B:415:LEU:HD23	2.29	0.57
1:B:403:GLU:H	1:B:420:ASN:ND2	2.00	0.57
1:B:614:SER:HA	1:B:619:VAL:CG2	2.34	0.57
1:B:454:CYS:CB	1:B:472:CYS:SG	2.92	0.56
1:B:614:SER:HA	1:B:619:VAL:HG21	1.87	0.56
1:A:454:CYS:CB	1:A:472:CYS:SG	2.93	0.56
1:A:242:SER:OG	1:A:243:ASP:N	2.38	0.56
1:B:377:ASN:C	1:B:377:ASN:ND2	2.61	0.56
1:B:459:VAL:HG22	1:B:460:SER:N	2.20	0.56
1:B:454:CYS:SG	1:B:472:CYS:SG	3.04	0.56
1:A:341:VAL:HG12	1:A:341:VAL:O	2.06	0.56
1:B:63:ILE:HD11	1:B:109:PRO:O	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:403:GLU:H	1:A:420:ASN:ND2	2.00	0.56
1:A:236:ILE:HD13	1:A:237:GLU:N	2.20	0.55
1:B:486:VAL:HG12	1:B:487:ASN:OD1	2.05	0.55
1:B:704:HIS:ND1	1:B:716:SER:HB2	2.21	0.55
1:A:72:GLN:O	1:A:73:GLU:HB2	2.07	0.55
1:B:454:CYS:HG	1:B:472:CYS:HG	1.54	0.55
1:A:736:THR:HG21	1:B:717:ALA:O	2.07	0.55
1:B:377:ASN:HD21	1:B:381:TYR:H	1.53	0.55
1:B:373:LYS:HD3	1:B:375:ILE:HD11	1.88	0.54
1:B:562:ASN:C	1:B:562:ASN:HD22	2.15	0.54
1:A:224:ALA:HB1	1:A:268:PHE:CZ	2.43	0.54
1:B:657:SER:H	1:B:715:GLN:HE21	1.55	0.54
1:A:385:CYS:HB3	1:A:387:PHE:CE1	2.41	0.54
1:A:420:ASN:C	1:A:420:ASN:HD22	2.15	0.54
1:B:454:CYS:CB	1:B:472:CYS:HG	2.20	0.54
1:B:45:LEU:HG	1:B:49:LEU:HD22	1.89	0.54
1:A:415:LEU:HD23	1:A:415:LEU:C	2.33	0.54
1:A:514:LEU:HD12	1:A:557:THR:HG22	1.91	0.54
1:B:154:TRP:NE1	1:B:156:THR:OG1	2.41	0.54
1:B:454:CYS:HB3	1:B:472:CYS:SG	2.48	0.54
1:A:736:THR:CG2	1:B:721:LYS:HB2	2.39	0.53
1:A:657:SER:H	1:A:715:GLN:HE21	1.55	0.53
1:B:693:GLU:O	1:B:696:LYS:HG2	2.08	0.53
1:A:242:SER:HB3	1:A:246:LEU:HD12	1.90	0.53
1:B:236:ILE:HD13	1:B:237:GLU:H	1.74	0.53
1:A:271:VAL:HG22	1:A:284:SER:HA	1.90	0.53
1:B:377:ASN:ND2	1:B:381:TYR:H	2.06	0.53
1:B:741:GLY:O	1:B:742:ILE:C	2.52	0.53
1:A:77:LEU:HD23	1:A:88:VAL:HA	1.92	0.52
1:A:48:TYR:CD2	1:A:49:LEU:HD12	2.45	0.52
1:A:451:PRO:HG2	1:A:452:GLU:OE2	2.09	0.52
1:A:454:CYS:HB3	1:A:472:CYS:SG	2.49	0.52
1:A:674:PRO:O	1:A:680:LEU:HD13	2.10	0.51
1:A:671:MET:HE1	1:A:682:HIS:CD2	2.46	0.51
1:B:82:GLU:HB2	1:B:467:TYR:OH	2.11	0.51
1:B:459:VAL:HG22	1:B:460:SER:H	1.76	0.51
1:A:197:GLY:C	1:A:213:ALA:HB3	2.36	0.51
1:A:369:ASN:O	1:A:389:ILE:HG12	2.10	0.51
1:B:403:GLU:OE1	1:B:585:TYR:HA	2.10	0.50
1:B:726:VAL:CG1	1:B:728:VAL:HG23	2.39	0.50
1:A:684:ARG:HG3	1:A:684:ARG:HH11	1.76	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:197:GLY:HA2	1:B:214:LEU:HD13	1.94	0.50
1:A:543:LEU:HD12	1:A:567:LEU:HD13	1.94	0.50
1:B:144:THR:HG22	1:B:144:THR:O	2.11	0.50
1:A:45:LEU:O	1:A:49:LEU:HD13	2.12	0.49
1:B:150:ASN:O	1:B:151:ASN:HB2	2.11	0.49
1:A:446:SER:HB2	1:A:457:TYR:CE2	2.47	0.49
1:B:139:LYS:HD3	1:B:141:GLN:NE2	2.27	0.49
1:A:289:ALA:HB3	1:A:294:LEU:HD11	1.94	0.49
1:B:290:PRO:HG3	1:B:326:ASP:OD2	2.13	0.49
1:B:314:GLN:HG2	1:B:325:MET:HG3	1.95	0.48
1:B:422:TYR:CD1	1:B:447:CYS:SG	3.06	0.48
1:B:69:LEU:CD2	1:B:78:VAL:HG22	2.44	0.48
1:B:546:VAL:HG22	1:B:547:TYR:N	2.28	0.48
1:B:648:LYS:HE3	1:B:762:CYS:SG	2.53	0.48
1:A:454:CYS:SG	1:A:472:CYS:SG	3.03	0.48
1:A:109:PRO:HG2	1:A:158:SER:O	2.14	0.48
1:B:129:THR:HG23	1:B:151:ASN:HA	1.96	0.48
1:A:750:HIS:HE1	1:B:728:VAL:O	1.97	0.48
1:A:219:ASN:N	1:A:308:GLN:OE1	2.47	0.47
1:B:146:GLU:O	1:B:175:LYS:NZ	2.41	0.47
1:A:630:SER:OG	1:A:631:TYR:N	2.47	0.47
1:B:461:PHE:CD1	1:B:468:TYR:HB3	2.49	0.47
1:A:60:LEU:C	1:A:60:LEU:HD22	2.39	0.47
1:A:169:ASN:O	1:A:170:ASN:HB2	2.15	0.47
1:A:458:SER:OG	1:A:471:ARG:HD3	2.15	0.47
1:A:658:ARG:HB2	1:A:689:MET:HE1	1.97	0.47
1:A:236:ILE:HD13	1:A:236:ILE:C	2.39	0.47
1:A:403:GLU:OE1	1:A:585:TYR:HA	2.15	0.47
1:A:519:LEU:HD23	1:A:519:LEU:HA	1.71	0.47
1:B:536:LYS:H	1:B:536:LYS:CD	2.22	0.47
1:A:530:LEU:HA	1:A:531:PRO:HD3	1.82	0.47
1:B:75:ASN:HD21	1:B:92:ASN:ND2	2.13	0.47
1:B:285:ILE:HD12	1:B:285:ILE:N	2.29	0.47
1:B:521:GLU:O	1:B:521:GLU:HG2	2.15	0.47
1:A:146:GLU:HG3	1:A:181:PRO:N	2.30	0.46
1:A:630:SER:HB2	1:A:740:HIS:NE2	2.29	0.46
1:A:60:LEU:HD13	3:A:960:HOH:O	2.16	0.46
1:A:671:MET:HE1	1:A:682:HIS:HD2	1.79	0.46
1:B:60:LEU:H	1:B:60:LEU:CD2	2.26	0.46
1:B:140:ARG:HD2	1:B:140:ARG:N	2.29	0.46
1:A:459:VAL:HG22	1:A:460:SER:N	2.31	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:377:ASN:ND2	1:A:379:GLU:N	2.64	0.46
1:A:696:LYS:CD	1:A:697:GLN:HG3	2.46	0.46
1:B:305:TRP:CZ3	1:B:311:ILE:HG12	2.50	0.46
1:B:608:GLU:OE1	1:B:611:ARG:HD2	2.15	0.46
1:A:160:VAL:CG1	1:A:161:GLY:N	2.79	0.46
1:A:372:TYR:HA	1:A:385:CYS:O	2.16	0.46
1:A:48:TYR:CE1	1:A:562:ASN:HA	2.50	0.46
1:B:110:ASP:O	1:B:112:GLN:N	2.49	0.46
1:B:498:SER:O	1:B:502:LYS:HG3	2.15	0.46
1:B:627:TRP:HB2	1:B:651:ILE:HB	1.99	0.45
1:B:677:GLU:CD	1:B:677:GLU:N	2.72	0.45
1:A:90:LEU:HD23	1:A:90:LEU:HA	1.70	0.45
1:A:236:ILE:HG13	1:A:712:HIS:ND1	2.30	0.45
1:B:597:ARG:O	1:B:600:THR:CG2	2.63	0.45
1:B:263:ASN:HD22	1:B:263:ASN:HA	1.60	0.45
1:B:671:MET:HE1	1:B:682:HIS:ND1	2.32	0.45
1:B:94:THR:O	1:B:95:PHE:HB2	2.16	0.45
1:A:375:ILE:HG22	1:A:376:SER:O	2.17	0.45
1:A:736:THR:HG23	1:B:721:LYS:HB2	1.98	0.45
1:A:159:PRO:HD3	1:A:216:TRP:CB	2.47	0.45
1:A:461:PHE:CD1	1:A:468:TYR:HB3	2.52	0.44
1:B:351:THR:HG22	1:B:592:HIS:HB3	1.99	0.44
1:B:658:ARG:HB3	1:B:661:TYR:CD2	2.53	0.44
1:A:60:LEU:HD13	1:A:60:LEU:H	1.83	0.44
1:A:95:PHE:HB3	1:A:98:PHE:HB2	1.99	0.44
1:B:102:ILE:HD11	1:B:116:LEU:HD13	1.99	0.44
1:B:288:THR:HG22	1:B:289:ALA:O	2.18	0.44
1:B:558:VAL:CG2	1:B:560:ARG:CZ	2.94	0.44
1:B:123:GLN:HG2	1:B:124:TRP:N	2.32	0.44
1:B:434:ILE:HG23	1:B:442:VAL:HG22	1.98	0.44
1:A:280:THR:HG22	1:A:281:ASN:N	2.33	0.44
1:A:482:LEU:HD23	1:A:483:HIS:H	1.81	0.44
1:A:513:LYS:O	1:A:527:GLN:HA	2.18	0.44
1:B:208:PHE:HZ	1:B:300:LEU:CD1	2.31	0.44
1:B:402:TRP:CD2	1:B:421:GLU:HB2	2.53	0.44
1:A:56:LYS:O	1:A:57:LEU:HD23	2.18	0.44
1:A:116:LEU:O	1:A:132:TYR:HA	2.18	0.44
1:B:137:LEU:O	1:B:139:LYS:O	2.36	0.44
1:A:741:GLY:O	1:A:742:ILE:C	2.59	0.44
1:B:46:THR:HG23	1:B:50:LYS:HD3	1.99	0.44
1:B:331:ASP:O	1:B:335:GLY:N	2.50	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:594:ILE:HG23	1:B:594:ILE:O	2.17	0.44
1:B:664:SER:HB2	1:B:668:GLU:OE2	2.18	0.43
1:A:127:SER:O	1:A:128:TYR:HB3	2.18	0.43
1:B:258:LYS:NZ	1:B:712:HIS:ND1	2.63	0.43
1:B:657:SER:HB3	1:B:719:ILE:HD11	2.01	0.43
1:A:305:TRP:CH2	1:A:311:ILE:HD11	2.53	0.43
1:A:594:ILE:O	1:A:594:ILE:HG23	2.17	0.43
1:A:594:ILE:CD1	1:A:601:PHE:HB2	2.48	0.43
1:A:743:ALA:O	1:A:744:SER:C	2.60	0.43
1:B:328:CYS:SG	1:B:339:CYS:SG	3.04	0.43
1:A:343:ARG:H	1:A:343:ARG:CD	2.24	0.43
1:B:435:GLN:HE22	1:B:441:LYS:HD2	1.83	0.43
1:B:596:ARG:HA	1:B:670:TYR:O	2.19	0.43
1:B:598:LEU:O	1:B:682:HIS:HE1	2.01	0.43
1:B:616:MET:HB3	1:B:618:PHE:CE2	2.54	0.43
1:B:680:LEU:CD2	1:B:684:ARG:HD3	2.48	0.43
1:B:309:GLU:HB3	1:B:330:TYR:HB3	2.01	0.43
1:B:660:GLU:HG3	1:B:683:TYR:CD2	2.54	0.43
1:A:627:TRP:HB2	1:A:651:ILE:HB	2.01	0.43
1:A:208:PHE:HZ	1:A:300:LEU:HD13	1.83	0.42
1:A:668:GLU:HG2	1:A:673:LEU:HD23	2.01	0.42
1:A:60:LEU:HD11	1:A:469:GLN:CD	2.43	0.42
1:B:60:LEU:HD23	1:B:60:LEU:O	2.19	0.42
1:B:163:LYS:HZ3	1:B:273:THR:HG23	1.84	0.42
1:B:742:ILE:O	1:B:742:ILE:HG22	2.18	0.42
1:B:543:LEU:HD12	1:B:567:LEU:HD13	2.01	0.42
1:B:626:ILE:O	1:B:650:GLY:HA2	2.19	0.42
1:A:397:ILE:HG13	1:A:398:THR:HG23	2.00	0.42
1:A:736:THR:HG21	1:B:721:LYS:HB2	2.00	0.42
1:B:217:SER:HG	1:B:222:PHE:HB2	1.82	0.42
1:B:611:ARG:HG2	1:B:611:ARG:HH11	1.83	0.42
1:A:71:LYS:HA	1:A:75:ASN:O	2.19	0.42
1:A:588:ASP:HB3	1:A:592:HIS:CD2	2.54	0.42
1:B:41:LYS:HE3	1:B:53:TYR:OH	2.19	0.42
1:B:46:THR:HG22	1:B:50:LYS:HD3	2.00	0.42
1:A:234:PRO:HB2	1:B:248:TYR:CZ	2.54	0.42
1:A:764:SER:O	1:A:766:PRO:HD3	2.19	0.42
1:B:93:SER:O	1:B:96:ASP:HB2	2.19	0.42
1:B:558:VAL:HG22	1:B:560:ARG:CZ	2.49	0.42
1:A:60:LEU:C	1:A:60:LEU:CD2	2.93	0.42
1:B:91:GLU:C	1:B:93:SER:N	2.78	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:385:CYS:HB3	1:B:387:PHE:CE1	2.54	0.42
1:A:69:LEU:CD1	1:A:107:ILE:HD12	2.50	0.42
1:A:142:LEU:H	1:A:142:LEU:CD1	2.27	0.42
1:A:320:GLN:OE1	1:A:669:ARG:HD3	2.20	0.42
1:B:114:ILE:HG23	1:B:135:TYR:HB3	2.02	0.42
1:B:656:VAL:HG13	1:B:715:GLN:HE22	1.85	0.41
1:A:464:GLU:O	1:A:464:GLU:HG2	2.20	0.41
1:A:664:SER:HB2	1:A:668:GLU:OE2	2.20	0.41
1:B:654:ALA:N	1:B:655:PRO:CD	2.83	0.41
1:A:168:TRP:O	1:A:169:ASN:HB2	2.21	0.41
1:A:728:VAL:O	1:B:750:HIS:HE1	2.03	0.41
1:A:60:LEU:HD22	1:A:60:LEU:O	2.20	0.41
1:A:452:GLU:N	1:A:452:GLU:CD	2.79	0.41
1:B:734:TRP:CD1	1:B:734:TRP:C	2.98	0.41
1:A:327:ILE:HG13	1:A:343:ARG:HB2	2.03	0.41
1:A:677:GLU:HG3	1:A:678:ASP:N	2.36	0.41
1:B:95:PHE:C	1:B:97:GLU:N	2.77	0.41
1:B:329:ASP:OD2	1:B:343:ARG:NH1	2.53	0.41
1:B:433:LYS:HG2	1:B:443:THR:HG23	2.03	0.41
1:B:500:LEU:HD13	1:B:500:LEU:O	2.20	0.41
1:B:536:LYS:HD3	1:B:536:LYS:N	2.27	0.41
1:A:369:ASN:C	1:A:389:ILE:HG12	2.46	0.40
1:A:330:TYR:HB2	1:A:337:TRP:CH2	2.56	0.40
1:A:470:LEU:HD23	1:A:470:LEU:HA	1.99	0.40
1:A:626:ILE:O	1:A:650:GLY:HA2	2.21	0.40
1:B:137:LEU:HD23	1:B:137:LEU:HA	1.86	0.40
1:B:620:ASP:OD1	1:B:622:LYS:HE2	2.21	0.40
1:A:144:THR:HG22	1:A:147:ARG:HH21	1.86	0.40
1:A:206:GLU:OE2	1:A:663:ASP:OD2	2.39	0.40
1:A:544:LEU:HA	1:A:544:LEU:HD23	1.73	0.40
1:B:246:LEU:HD23	1:B:246:LEU:HA	1.92	0.40
1:B:277:SER:HB3	1:B:280:THR:HB	2.02	0.40
1:B:594:ILE:CD1	1:B:601:PHE:HB2	2.52	0.40
1:B:137:LEU:O	1:B:138:ASN:C	2.64	0.40
1:A:221:THR:O	1:A:273:THR:HB	2.21	0.40
1:A:481:THR:OG1	1:A:483:HIS:CE1	2.69	0.40
1:B:397:ILE:HG13	1:B:439:TYR:CE1	2.56	0.40
1:B:546:VAL:CG2	1:B:547:TYR:N	2.84	0.40
1:B:596:ARG:N	1:B:670:TYR:O	2.51	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	726/728 (100%)	675 (93%)	47 (6%)	4 (1%)	21	25
1	B	726/728 (100%)	682 (94%)	39 (5%)	5 (1%)	18	21
All	All	1452/1456 (100%)	1357 (94%)	86 (6%)	9 (1%)	21	25

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	389	ILE
1	B	111	GLY
1	B	320	GLN
1	A	320	GLN
1	A	617	GLY
1	B	617	GLY
1	B	393	ASP
1	A	486	VAL
1	B	486	VAL

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	653/653 (100%)	619 (95%)	34 (5%)	21	28
1	B	653/653 (100%)	622 (95%)	31 (5%)	23	33
All	All	1306/1306 (100%)	1241 (95%)	65 (5%)	22	30

All (65) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	60	LEU
1	A	73	GLU
1	A	90	LEU
1	A	133	ASP
1	A	142	LEU
1	A	223	LEU
1	A	236	ILE
1	A	246	LEU
1	A	263	ASN
1	A	294	LEU
1	A	300	LEU
1	A	316	LEU
1	A	326	ASP
1	A	343	ARG
1	A	373	LYS
1	A	377	ASN
1	A	385	CYS
1	A	410	LEU
1	A	420	ASN
1	A	431	LEU
1	A	479	LEU
1	A	482	LEU
1	A	500	LEU
1	A	513	LYS
1	A	514	LEU
1	A	543	LEU
1	A	544	LEU
1	A	594	ILE
1	A	656	VAL
1	A	696	LYS
1	A	715	GLN
1	A	721	LYS
1	A	726	VAL
1	A	736	THR
1	B	41	LYS
1	B	49	LEU
1	B	51	ASN
1	B	60	LEU
1	B	71	LYS
1	B	90	LEU
1	B	96	ASP
1	B	129	THR

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Mol	Chain	Res	Type
1	B	214	LEU
1	B	223	LEU
1	B	236	ILE
1	B	246	LEU
1	B	263	ASN
1	B	273	THR
1	B	311	ILE
1	B	370	SER
1	B	373	LYS
1	B	377	ASN
1	B	385	CYS
1	B	420	ASN
1	B	479	LEU
1	B	507	VAL
1	B	514	LEU
1	B	536	LYS
1	B	543	LEU
1	B	558	VAL
1	B	562	ASN
1	B	594	ILE
1	B	619	VAL
1	B	679	ASN
1	B	715	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (42) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	72	GLN
1	A	123	GLN
1	A	141	GLN
1	A	169	ASN
1	A	247	GLN
1	A	263	ASN
1	A	298	HIS
1	A	363	HIS
1	A	377	ASN
1	A	388	GLN
1	A	420	ASN
1	A	483	HIS
1	A	506	ASN
1	A	508	GLN
1	A	533	HIS

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Mol	Chain	Res	Type
1	A	592	HIS
1	A	679	ASN
1	A	697	GLN
1	A	715	GLN
1	A	750	HIS
1	B	51	ASN
1	B	75	ASN
1	B	100	HIS
1	B	103	ASN
1	B	141	GLN
1	B	169	ASN
1	B	263	ASN
1	B	363	HIS
1	B	377	ASN
1	B	383	HIS
1	B	420	ASN
1	B	435	GLN
1	B	483	HIS
1	B	533	HIS
1	B	562	ASN
1	B	586	GLN
1	B	612	GLN
1	B	679	ASN
1	B	715	GLN
1	B	731	GLN
1	B	748	HIS
1	B	750	HIS

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

2 ligands are modelled in this entry.

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	1WH	A	801	-	25,27,27	1.95	6 (24%)	33,39,39	3.20	12 (36%)
2	1WH	B	801	-	25,27,27	1.77	9 (36%)	33,39,39	2.24	10 (30%)

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	1WH	A	801	-	-	1/9/22/22	0/3/3/3
2	1WH	B	801	-	-	1/9/22/22	0/3/3/3

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	801	1WH	C15-CL	-4.33	1.63	1.73
2	A	801	1WH	C17-CL1	-4.30	1.64	1.74
2	B	801	1WH	C18-C17	4.15	1.45	1.38
2	A	801	1WH	C18-C17	3.68	1.44	1.38
2	A	801	1WH	C19-C18	3.16	1.43	1.38
2	B	801	1WH	C14-C15	2.77	1.44	1.39
2	B	801	1WH	C16-C17	2.64	1.42	1.38
2	B	801	1WH	C16-C15	2.52	1.42	1.38
2	A	801	1WH	C14-C15	2.44	1.44	1.39
2	B	801	1WH	C3-C5	2.39	1.43	1.39
2	B	801	1WH	C15-CL	-2.37	1.68	1.73
2	B	801	1WH	C19-C18	2.25	1.42	1.38
2	B	801	1WH	C5-N1	2.18	1.37	1.34
2	B	801	1WH	C17-CL1	-2.14	1.69	1.74
2	A	801	1WH	C3-C5	2.06	1.42	1.39

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	801	1WH	C16-C17-CL1	-7.50	109.77	119.17
2	A	801	1WH	C16-C15-C14	7.09	127.48	121.97
2	A	801	1WH	C16-C15-CL	-6.77	107.37	118.45
2	A	801	1WH	C8-C7-N2	-5.43	101.90	113.41
2	A	801	1WH	C15-C16-C17	-5.41	112.61	118.72
2	B	801	1WH	C16-C15-C14	5.39	126.16	121.97
2	B	801	1WH	C15-C16-C17	-5.19	112.86	118.72
2	B	801	1WH	C8-C7-N2	-4.78	103.28	113.41
2	A	801	1WH	C18-C17-CL1	4.62	126.17	119.36
2	A	801	1WH	C4-C5-C3	-4.36	116.52	122.31
2	B	801	1WH	C16-C17-CL1	-4.08	114.05	119.17
2	A	801	1WH	C7A-C-N2	3.65	109.05	106.40
2	A	801	1WH	C2-C7A-C	3.36	133.56	129.65
2	B	801	1WH	C4-C5-C3	-3.28	117.95	122.31
2	A	801	1WH	C14-C15-CL	3.26	125.14	120.67
2	A	801	1WH	C1-C3-C5	-3.17	115.88	119.69
2	A	801	1WH	O2-C-N2	-3.00	119.08	125.26
2	B	801	1WH	C16-C15-CL	-2.70	114.02	118.45
2	B	801	1WH	C18-C17-CL1	2.32	122.78	119.36
2	B	801	1WH	C10-O1-C8	2.20	126.16	112.90
2	B	801	1WH	C5-N1-C3A	-2.11	115.98	118.66
2	B	801	1WH	C1-C3-C5	-2.02	117.27	119.69

There are no chirality outliers.

All (2) torsion outliers are listed below:

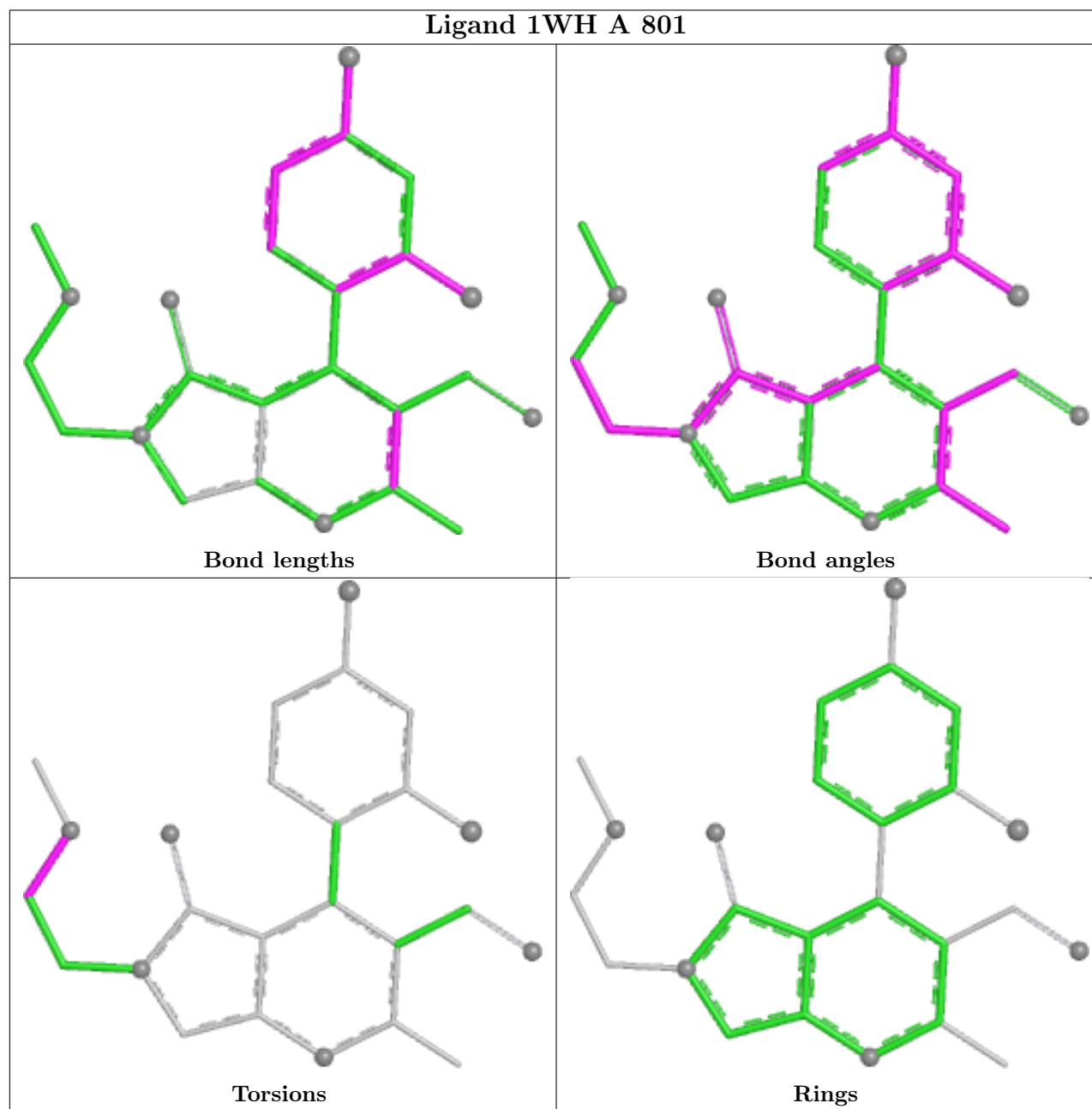
Mol	Chain	Res	Type	Atoms
2	A	801	1WH	C7-C8-O1-C10
2	B	801	1WH	C7-C8-O1-C10

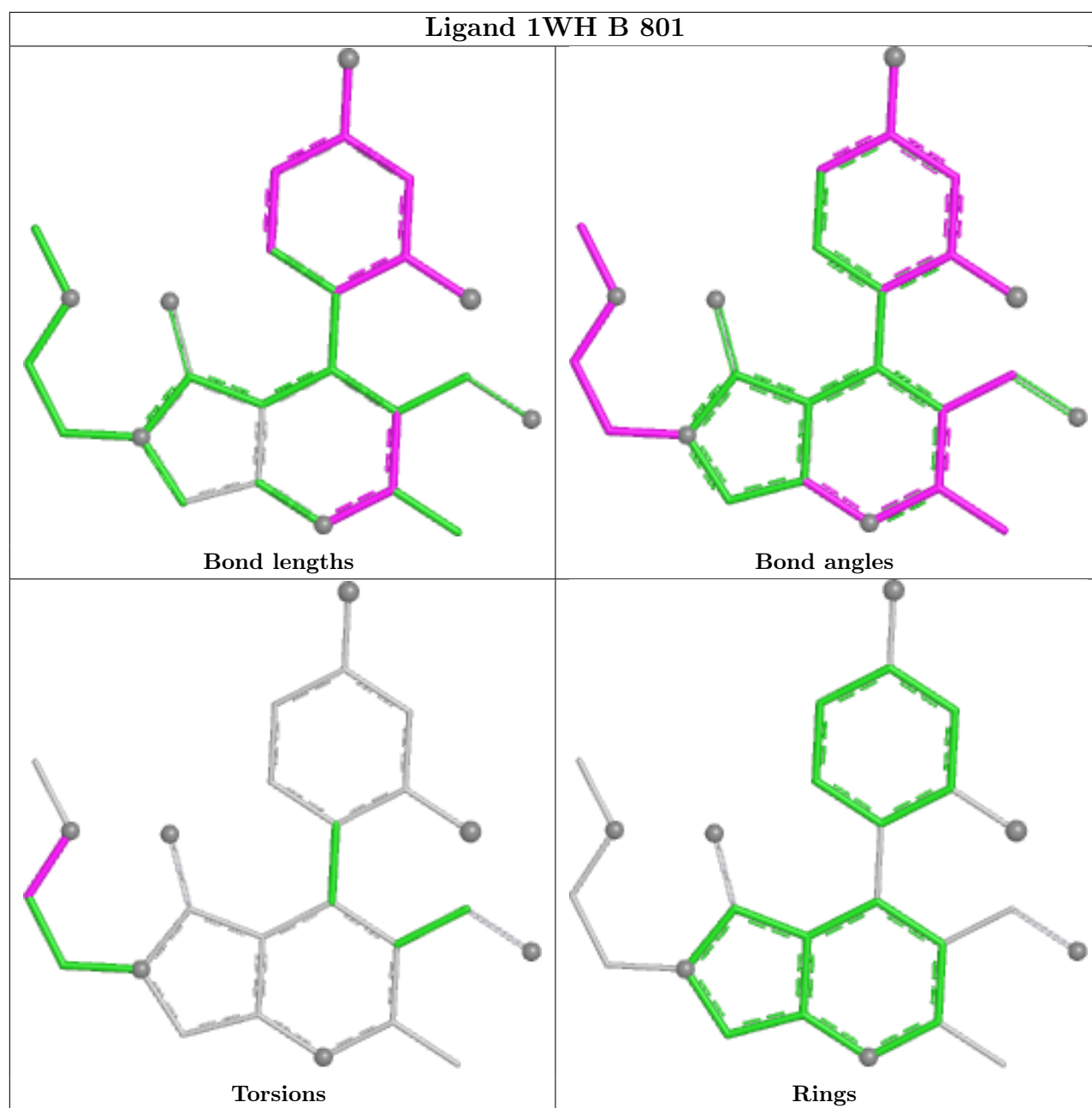
There are no ring outliers.

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	728/728 (100%)	0.15	20 (2%) 56 55	34, 47, 69, 84	0
1	B	728/728 (100%)	0.36	23 (3%) 50 49	33, 54, 75, 84	0
All	All	1456/1456 (100%)	0.26	43 (2%) 52 52	33, 50, 73, 84	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	39	SER	6.6
1	A	342	ALA	6.0
1	A	766	PRO	5.7
1	B	766	PRO	4.2
1	A	521	GLU	3.8
1	A	341	VAL	3.7
1	B	143	ILE	3.7
1	A	765	LEU	3.5
1	B	279	VAL	3.4
1	A	105	TYR	3.0
1	B	145	GLU	2.9
1	B	547	TYR	2.8
1	A	502	LYS	2.8
1	B	271	VAL	2.8
1	B	521	GLU	2.8
1	A	343	ARG	2.7
1	A	40	ARG	2.6
1	B	389	ILE	2.6
1	A	91	GLU	2.6
1	B	397	ILE	2.5
1	B	156	THR	2.5
1	B	332	GLU	2.5
1	A	142	LEU	2.4
1	B	144	THR	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	677	GLU	2.4
1	B	63	ILE	2.4
1	A	387	PHE	2.3
1	B	471	ARG	2.3
1	A	345	HIS	2.3
1	B	280	THR	2.3
1	B	273	THR	2.2
1	A	393	ASP	2.2
1	A	95	PHE	2.2
1	A	145	GLU	2.2
1	A	506	ASN	2.2
1	B	208	PHE	2.1
1	B	463	LYS	2.1
1	B	333	SER	2.1
1	B	506	ASN	2.1
1	B	110	ASP	2.0
1	B	502	LYS	2.0
1	B	390	ASP	2.0
1	A	182	SER	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.

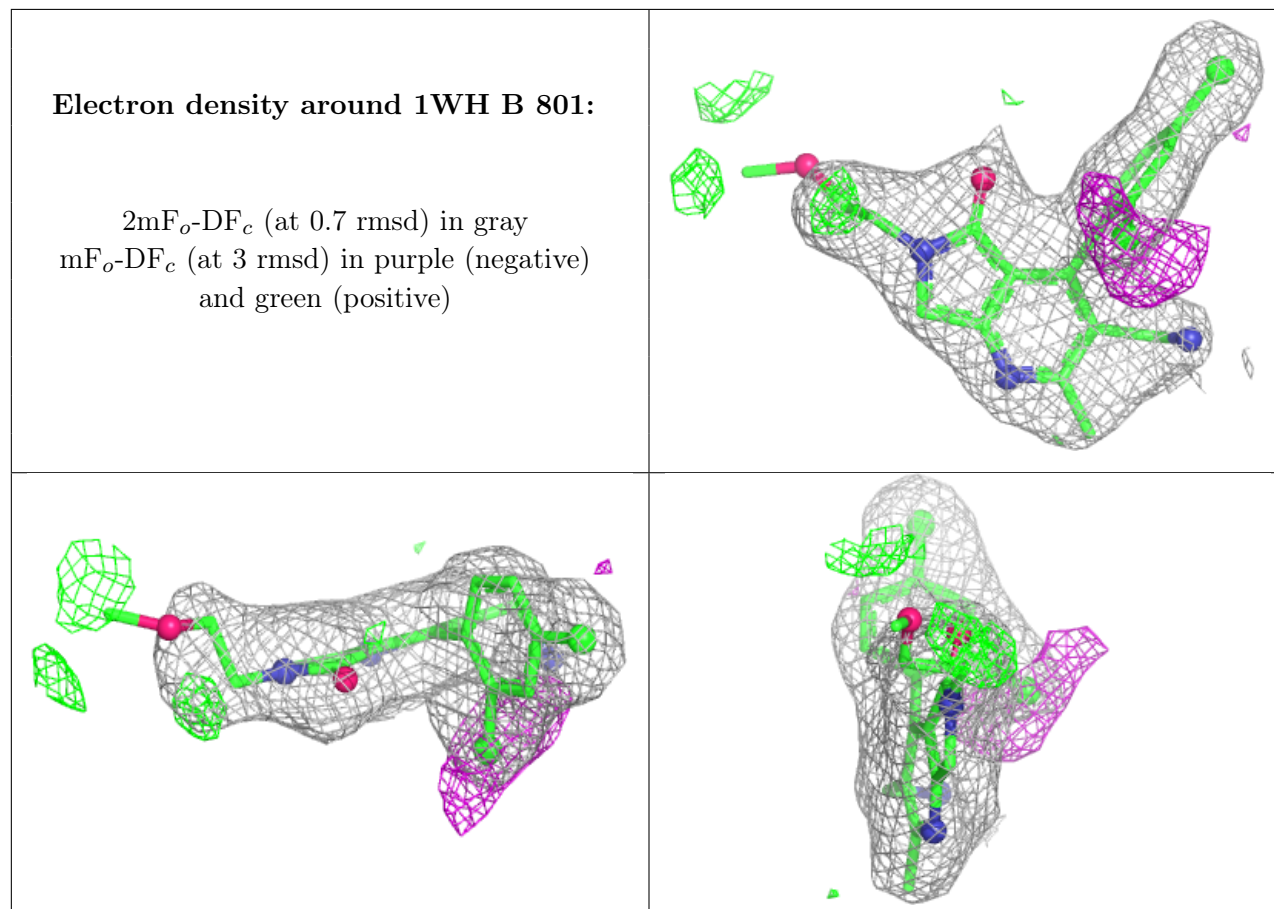
6.4 Ligands [i](#)

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	1WH	B	801	25/25	0.84	0.16	55,58,67,74	0
2	1WH	A	801	25/25	0.87	0.14	52,56,63,68	0

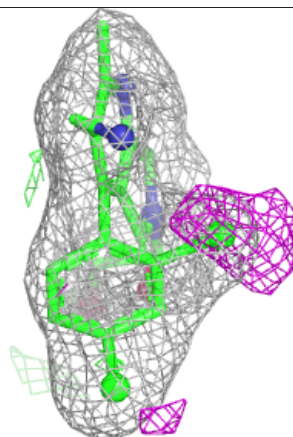
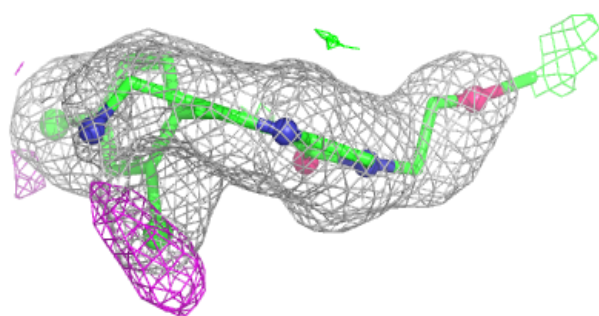
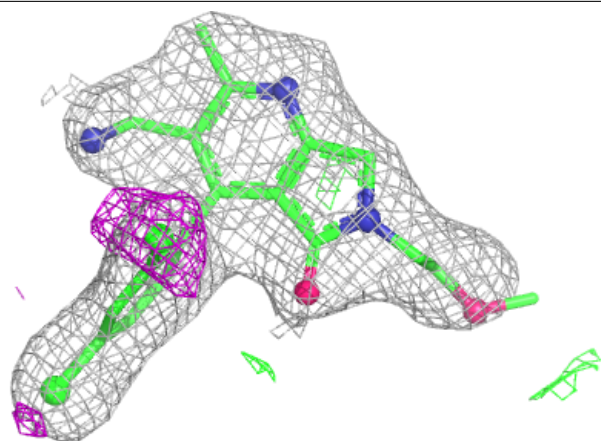
The following is a graphical depiction of the model fit to experimental electron density of all

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 1WH A 801:

$2mF_o - DF_c$ (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.