



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 11:24 PM UTC

PDB ID : 9Q5Y / pdb_00009q5y
Title : Human prolyl endopeptidase (PREP) - complex with YX45
Authors : Fucci, I.J.; Xie, Y.; Yoo, E.; Monteiro, D.C.F.
Deposited on : 2025-08-21
Resolution : 1.49 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

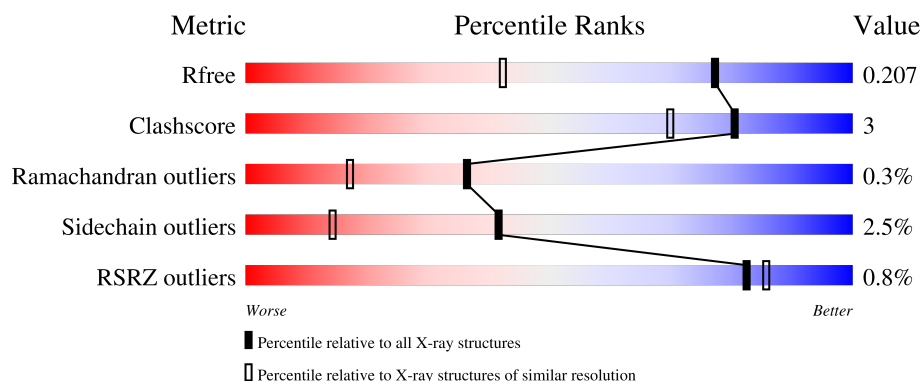
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.49 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	4037 (1.50-1.50)
Clashscore	190562	4235 (1.50-1.50)
Ramachandran outliers	187476	4153 (1.50-1.50)
Sidechain outliers	187428	4150 (1.50-1.50)
RSRZ outliers	180081	4039 (1.50-1.50)

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	711	90% 8% ..
1	B	711	90% 8% .
1	C	711	86% 11% ..

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
4	GOL	A	806	-	X	-	-
4	GOL	B	810	-	X	-	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 36295 atoms, of which 16962 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.

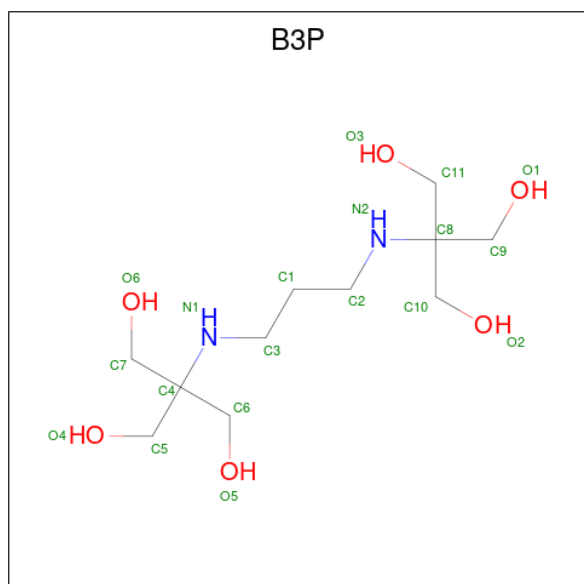
- Molecule 1 is a protein called Prolyl endopeptidase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	706	Total	C	H	N	O	S	178	6	0
			11267	3659	5564	949	1067	28			
1	B	708	Total	C	H	N	O	S	178	5	0
			11282	3664	5570	950	1071	27			
1	C	707	Total	C	H	N	O	S	177	3	0
			11230	3652	5539	942	1070	27			

There are 3 discrepancies between the modelled and reference sequences:

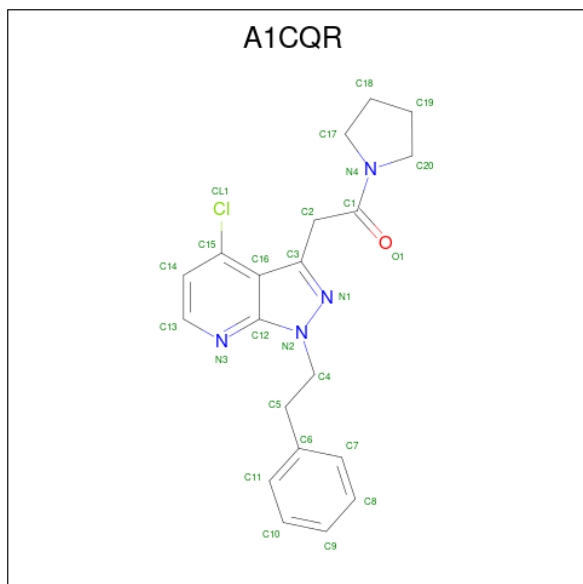
Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	expression tag	UNP P48147
B	0	GLY	-	expression tag	UNP P48147
C	0	GLY	-	expression tag	UNP P48147

- Molecule 2 is 2-[3-(2-HYDROXY-1,1-DIHYDROXYMETHYL-ETHYLAMINO)-PROPYLAMINO]-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (CCD ID: B3P) (formula: C₁₁H₂₆N₂O₆).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	H	N	O	8	0
			45	11	26	2	6		

- Molecule 3 is 2-[4-chloro-1-(2-phenylethyl)-1H-pyrazolo[3,4-b]pyridin-3-yl]-1-(pyrrolidin-1-yl)ethan-1-one (CCD ID: A1CQR) (formula: C₂₀H₂₁ClN₄O).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total 47	C 20	Cl 1	H 21	N 4	O 1	0	0
3	B	1	Total 47	C 20	Cl 1	H 21	N 4	O 1	0	0
3	C	1	Total 47	C 20	Cl 1	H 21	N 4	O 1	0	0

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



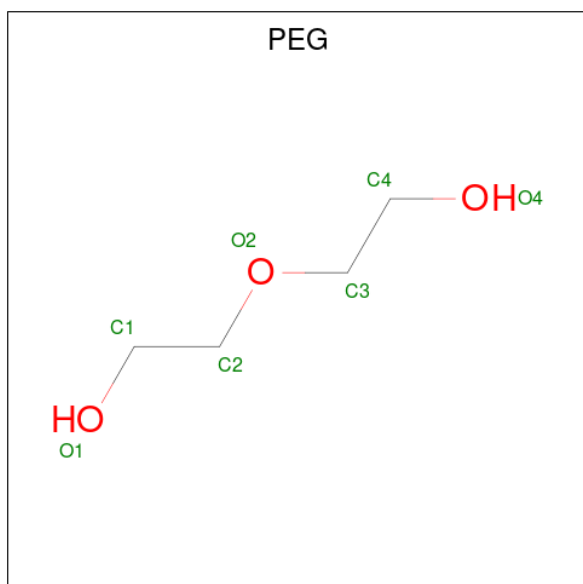
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	H	O	3	0
			14	3	8	3		
4	A	1	Total	C	H	O	3	0
			14	3	8	3		
4	A	1	Total	C	H	O	3	0
			14	3	8	3		
4	A	1	Total	C	H	O	3	0
			14	3	8	3		
4	A	1	Total	C	H	O	3	0
			14	3	8	3		
4	B	1	Total	C	H	O	3	0
			14	3	8	3		
4	B	1	Total	C	H	O	3	0
			14	3	8	3		
4	B	1	Total	C	H	O	3	0
			14	3	8	3		
4	B	1	Total	C	H	O	3	0
			14	3	8	3		
4	B	1	Total	C	H	O	3	0
			14	3	8	3		
4	B	1	Total	C	H	O	3	0
			14	3	8	3		
4	B	1	Total	C	H	O	3	0
			14	3	8	3		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	C	1	Total	C	H	O	3	0
			14	3	8	3		
4	C	1	Total	C	H	O	3	0
			14	3	8	3		
4	C	1	Total	C	H	O	3	0
			14	3	8	3		
4	C	1	Total	C	H	O	3	0
			14	3	8	3		
4	C	1	Total	C	H	O	3	0
			14	3	8	3		

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	H	O	2	0
			17	4	10	3		
5	B	1	Total	C	H	O	2	0
			17	4	10	3		
5	C	1	Total	C	H	O	2	0
			17	4	10	3		
5	C	1	Total	C	H	O	2	0
			17	4	10	3		

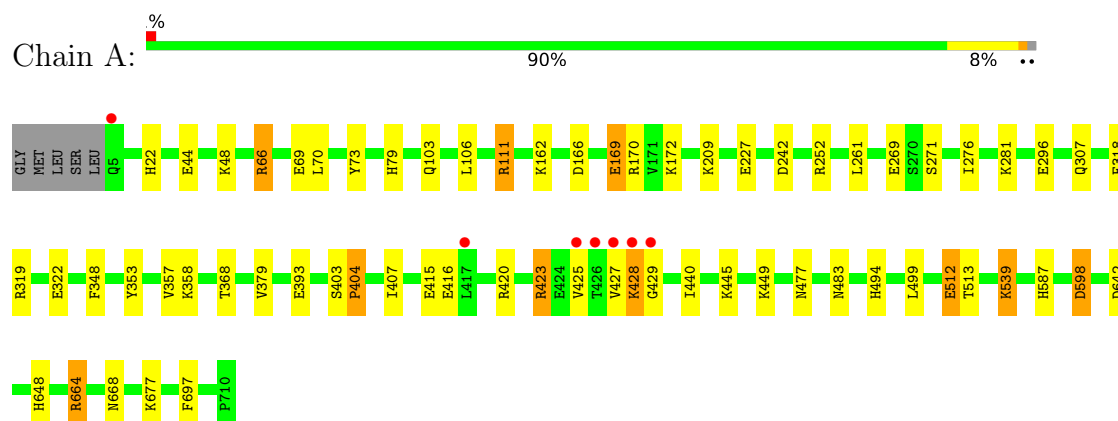
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	764	Total 768	O 768	0	4
6	B	658	Total 660	O 660	0	2
6	C	553	Total 554	O 554	0	1

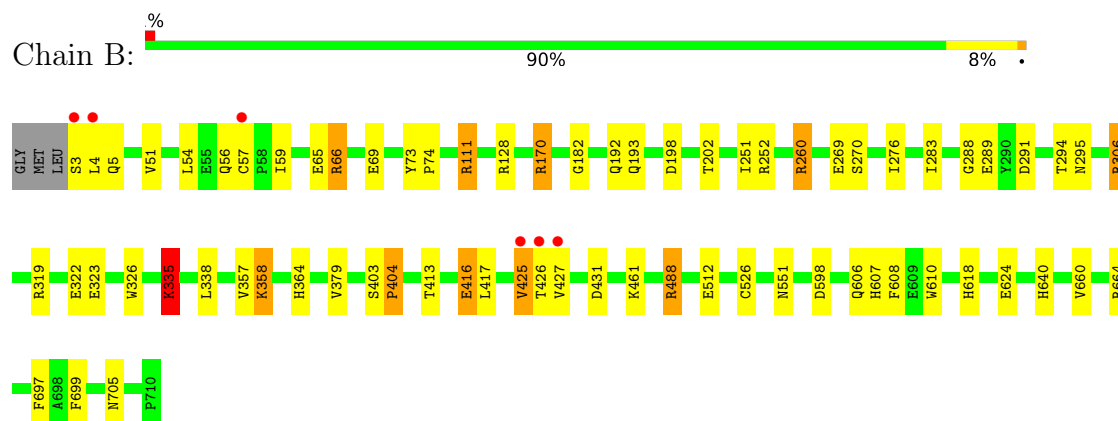
3 Residue-property plots

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.

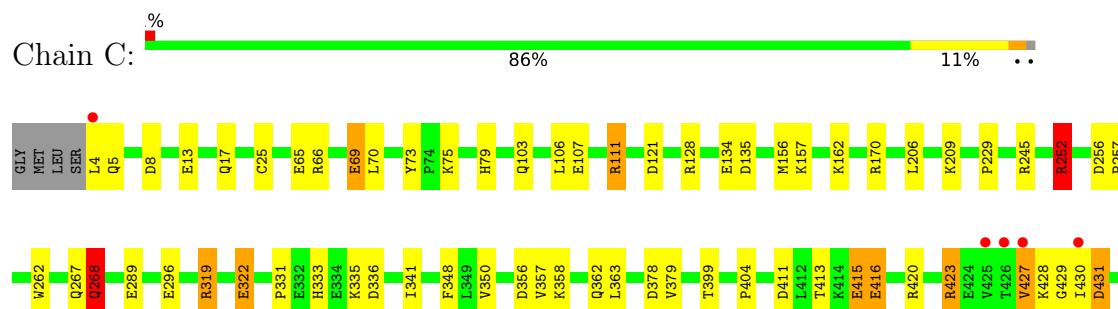
• Molecule 1: Prolyl endopeptidase



• Molecule 1: Prolyl endopeptidase



• Molecule 1: Prolyl endopeptidase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	106.22Å 66.88Å 156.91Å 90.00° 99.29° 90.00°	Depositor
Resolution (Å)	34.94 – 1.49 34.94 – 1.49	Depositor EDS
% Data completeness (in resolution range)	73.4 (34.94-1.49) 73.6 (34.94-1.49)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.85 (at 1.49Å)	Xtriage
Refinement program	REFMAC 5.8.0430 (refmacat 0.4.100)	Depositor
R, R_{free}	0.163 , 0.206 0.164 , 0.207	Depositor DCC
R_{free} test set	17662 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	17.0	Xtriage
Anisotropy	0.102	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 39.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	36295	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.17% 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: GOL, PEG, A1CQR, B3P

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	0.98	3/5873 (0.1%)	1.36	24/7959 (0.3%)
1	B	0.96	6/5879 (0.1%)	1.36	33/7967 (0.4%)
1	C	0.93	5/5852 (0.1%)	1.41	46/7932 (0.6%)
All	All	0.96	14/17604 (0.1%)	1.38	103/23858 (0.4%)

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	7
1	B	0	11
1	C	0	7
All	All	0	25

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	170	ARG	NE-CZ	-8.95	1.23	1.33
1	A	648	HIS	CE1-NE2	8.06	1.40	1.32
1	B	640	HIS	ND1-CE1	6.00	1.38	1.32
1	A	587	HIS	CD2-NE2	5.70	1.44	1.37
1	B	512	GLU	CD-OE1	5.68	1.36	1.25
1	C	229	PRO	CA-CB	-5.36	1.46	1.53
1	A	172	LYS	CD-CE	-5.34	1.36	1.52
1	C	289	GLU	CD-OE1	5.26	1.35	1.25
1	C	544	SER	CA-CB	-5.25	1.47	1.54
1	B	417	LEU	C-N	5.22	1.37	1.33
1	B	338	LEU	C-O	-5.20	1.17	1.24
1	C	615	SER	CA-CB	-5.17	1.45	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	289	GLU	CD-OE1	5.05	1.34	1.25
1	B	618	HIS	CG-CD2	-5.02	1.30	1.35

All (103) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	322	GLU	CB-CG-CD	10.64	130.69	112.60
1	A	512	GLU	CB-CG-CD	10.63	130.68	112.60
1	C	13	GLU	CB-CG-CD	9.13	128.12	112.60
1	C	268	GLN	CB-CA-C	-9.04	93.64	110.36
1	C	268	GLN	N-CA-CB	8.97	124.28	110.46
1	B	697	PHE	CA-CB-CG	-8.22	105.58	113.80
1	A	252	ARG	CD-NE-CZ	8.16	135.83	124.40
1	B	306	ARG	CB-CA-C	7.80	121.39	110.16
1	C	512	GLU	CB-CG-CD	7.52	125.39	112.60
1	C	670	LEU	N-CA-CB	-7.49	99.45	110.84
1	B	260	ARG	NE-CZ-NH2	7.45	125.91	119.20
1	C	296	GLU	CB-CG-CD	7.45	125.26	112.60
1	C	378	ASP	CA-CB-CG	7.38	119.97	112.60
1	B	461	LYS	CB-CA-C	-7.29	97.50	109.53
1	B	51	VAL	O-C-N	-7.12	115.86	120.42
1	A	404	PRO	CA-C-N	-7.08	107.54	121.41
1	A	404	PRO	C-N-CA	-7.08	107.54	121.41
1	C	623	PRO	CB-CA-C	6.98	120.42	111.56
1	C	423	ARG	N-CA-CB	6.92	123.22	111.66
1	C	662	ARG	CD-NE-CZ	6.88	134.03	124.40
1	C	603	ASP	CA-CB-CG	6.86	119.46	112.60
1	C	135	ASP	CA-CB-CG	6.85	119.45	112.60
1	A	348	PHE	CA-CB-CG	6.75	120.55	113.80
1	C	128	ARG	NE-CZ-NH1	-6.71	114.79	121.50
1	C	705	ASN	CB-CA-C	-6.65	101.85	112.09
1	C	8	ASP	CB-CA-C	-6.64	99.26	109.89
1	C	618	HIS	CA-CB-CG	-6.58	107.22	113.80
1	B	170	ARG	CD-NE-CZ	6.58	133.61	124.40
1	C	442	TYR	CB-CA-C	6.39	118.47	108.63
1	A	513	THR	N-CA-CB	6.38	119.60	110.16
1	A	697	PHE	CA-CB-CG	-6.35	107.45	113.80
1	C	642	ASP	CA-CB-CG	6.25	118.85	112.60
1	C	423	ARG	CB-CA-C	-6.25	97.02	110.45
1	B	404	PRO	CA-C-N	-6.15	109.36	121.41
1	B	404	PRO	C-N-CA	-6.15	109.36	121.41
1	C	598	ASP	CA-CB-CG	6.09	118.69	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	128	ARG	NE-CZ-NH2	6.06	124.66	119.20
1	C	252	ARG	CG-CD-NE	-6.06	98.67	112.00
1	A	281	LYS	N-CA-CB	-6.06	103.14	110.53
1	B	335	LYS	CB-CG-CD	5.96	125.00	111.30
1	A	477	ASN	CA-CB-CG	5.95	118.55	112.60
1	B	699	PHE	CA-CB-CG	5.92	119.72	113.80
1	B	294	THR	CA-CB-OG1	-5.92	100.72	109.60
1	C	5	GLN	N-CA-CB	5.91	119.87	110.65
1	B	198	ASP	CA-CB-CG	5.86	118.46	112.60
1	B	608	PHE	CA-CB-CG	-5.86	107.94	113.80
1	A	642	ASP	CA-CB-CG	5.85	118.45	112.60
1	B	526	CYS	N-CA-CB	5.83	118.63	109.94
1	B	65	GLU	CB-CG-CD	-5.82	102.70	112.60
1	A	494	HIS	CB-CG-CD2	-5.80	123.67	131.20
1	C	468	ALA	O-C-N	5.74	130.30	123.24
1	C	695	ASP	CA-CB-CG	5.72	118.32	112.60
1	C	111	ARG	CG-CD-NE	-5.71	99.43	112.00
1	C	411	ASP	O-C-N	-5.71	115.40	122.82
1	C	466	HIS	CA-CB-CG	-5.65	108.15	113.80
1	B	413	THR	CA-CB-OG1	-5.65	101.13	109.60
1	B	260	ARG	NE-CZ-NH1	-5.64	115.86	121.50
1	C	547	ARG	CB-CA-C	5.62	120.07	110.85
1	A	393	GLU	CB-CG-CD	5.57	122.08	112.60
1	C	156	MET	CG-SD-CE	-5.57	88.64	100.90
1	C	639	ASP	CA-CB-CG	5.57	118.17	112.60
1	B	291	ASP	CA-CB-CG	5.57	118.17	112.60
1	A	483	ASN	CA-CB-CG	-5.51	107.09	112.60
1	B	193	GLN	OE1-CD-NE2	-5.50	117.10	122.60
1	C	121	ASP	CB-CA-C	-5.47	99.22	109.72
1	A	539	LYS	CG-CD-CE	5.46	123.86	111.30
1	C	331	PRO	CB-CA-C	5.46	118.14	110.98
1	B	202	THR	CA-CB-OG1	-5.44	101.44	109.60
1	A	420	ARG	CD-NE-CZ	5.44	132.01	124.40
1	A	169	GLU	CA-CB-CG	-5.40	103.31	114.10
1	A	166	ASP	CA-CB-CG	5.38	117.98	112.60
1	B	705	ASN	CB-CA-C	-5.37	103.81	112.09
1	A	281	LYS	CB-CA-C	5.36	120.16	112.12
1	C	431	ASP	CA-CB-CG	5.36	117.96	112.60
1	B	322	GLU	CG-CD-OE1	5.35	130.71	118.40
1	B	364	HIS	CB-CG-ND1	-5.34	114.69	122.70
1	B	295	ASN	O-C-N	5.34	129.46	123.27
1	A	307	GLN	N-CA-C	-5.28	106.78	113.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	51	VAL	N-CA-CB	-5.26	106.99	110.52
1	B	364	HIS	CB-CG-CD2	5.25	138.03	131.20
1	A	598	ASP	CA-CB-CG	5.24	117.84	112.60
1	C	209	LYS	CB-CG-CD	5.24	123.35	111.30
1	C	436	GLN	N-CA-CB	-5.20	102.64	111.69
1	C	170	ARG	NE-CZ-NH1	-5.19	116.31	121.50
1	B	660	VAL	CA-C-N	5.18	126.66	120.13
1	B	660	VAL	C-N-CA	5.18	126.66	120.13
1	A	668	ASN	OD1-CG-ND2	5.17	127.77	122.60
1	C	399	THR	OG1-CB-CG2	-5.16	98.97	109.30
1	C	157	LYS	CB-CA-C	5.16	118.26	109.80
1	B	56	GLN	CB-CA-C	-5.12	99.28	109.99
1	B	289	GLU	CG-CD-OE2	-5.12	106.62	118.40
1	B	288	GLY	O-C-N	-5.11	118.84	123.54
1	A	318	PHE	CA-CB-CG	5.10	118.90	113.80
1	A	209	LYS	CA-CB-CG	-5.10	103.89	114.10
1	C	134	GLU	CB-CG-CD	5.09	121.26	112.60
1	C	413	THR	CA-CB-OG1	-5.08	101.97	109.60
1	C	415	GLU	N-CA-CB	-5.08	102.32	110.19
1	C	660	VAL	CA-C-N	5.08	125.72	120.03
1	C	660	VAL	C-N-CA	5.08	125.72	120.03
1	C	319	ARG	CD-NE-CZ	5.07	131.50	124.40
1	A	269	GLU	CB-CG-CD	5.07	121.22	112.60
1	B	322	GLU	CB-CG-CD	5.04	121.17	112.60
1	B	322	GLU	CG-CD-OE2	-5.03	106.83	118.40

There are no chirality outliers.

All (25) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	111	ARG	Sidechain
1	A	319	ARG	Sidechain
1	A	423	ARG	Sidechain
1	A	66[A]	ARG	Sidechain
1	A	66[B]	ARG	Sidechain
1	A	664[A]	ARG	Sidechain
1	A	664[B]	ARG	Sidechain
1	B	111	ARG	Sidechain
1	B	128	ARG	Sidechain
1	B	170	ARG	Sidechain
1	B	252[A]	ARG	Sidechain
1	B	252[B]	ARG	Sidechain

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Mol	Chain	Res	Type	Group
1	B	3	SER	Peptide
1	B	306	ARG	Sidechain
1	B	488	ARG	Sidechain
1	B	66[A]	ARG	Sidechain
1	B	66[B]	ARG	Sidechain
1	B	664	ARG	Sidechain
1	C	111	ARG	Sidechain
1	C	245	ARG	Sidechain
1	C	252	ARG	Sidechain
1	C	319	ARG	Sidechain
1	C	423	ARG	Sidechain
1	C	428	LYS	Peptide
1	C	493	ARG	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	5703	5564	5541	27	1
1	B	5712	5570	5547	26	0
1	C	5691	5539	5516	31	0
2	A	19	26	26	1	0
3	A	26	21	0	0	0
3	B	26	21	0	0	0
3	C	26	21	0	0	0
4	A	30	40	39	3	0
4	B	54	72	71	9	1
4	C	36	48	48	5	0
5	A	7	10	10	0	0
5	B	7	10	10	0	0
5	C	14	20	20	3	0
6	A	768	0	0	7	2
6	B	660	0	0	6	1
6	C	554	0	0	7	2
All	All	19333	16962	16828	88	4

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 3.

All (88) 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:C:25:CYS:HB3	6:C:1318:HOH:O	1.53	1.07
1:A:296:GLU:HG3	6:A:1497:HOH:O	1.58	1.00
1:A:162:LYS:NZ	6:A:901:HOH:O	1.88	0.95
5:C:808:PEG:H32	6:C:1236:HOH:O	1.78	0.83
1:C:336:ASP:OD2	4:C:809:GOL:O3	1.97	0.83
1:A:170:ARG:NH2	6:A:902:HOH:O	2.11	0.79
1:C:466:HIS:HB2	1:C:498:ILE:HD12	1.68	0.75
1:B:606:GLN:HB2	6:B:1385:HOH:O	1.89	0.72
1:B:269:GLU:O	4:B:804:GOL:H12	1.90	0.71
1:B:319:ARG:NH2	4:B:811:GOL:H11	2.06	0.71
1:C:333:HIS:CD2	4:C:809:GOL:O2	2.44	0.71
5:C:808:PEG:C3	6:C:1236:HOH:O	2.36	0.71
1:A:66[B]:ARG:HH12	1:A:69:GLU:HG2	1.59	0.67
1:A:404:PRO:HG2	1:A:425:VAL:O	1.94	0.67
1:B:379:VAL:HG12	6:B:1475:HOH:O	1.94	0.67
1:A:664[B]:ARG:HG2	1:A:664[B]:ARG:HH21	1.59	0.66
1:C:416:GLU:O	1:C:416:GLU:CD	2.38	0.66
1:C:335:LYS:HE3	6:C:1249:HOH:O	1.95	0.66
1:C:73:TYR:O	1:C:75:LYS:HE2	1.96	0.65
1:A:73:TYR:CD2	1:A:425:VAL:HG11	2.34	0.63
1:C:461:LYS:HD2	1:C:461:LYS:N	2.15	0.61
1:A:664[B]:ARG:NH2	1:A:664[B]:ARG:CG	2.61	0.61
4:B:807:GOL:O3	6:B:901:HOH:O	2.09	0.60
1:B:319:ARG:HH22	4:B:811:GOL:H11	1.65	0.60
1:A:379:VAL:HG12	6:A:1565:HOH:O	2.02	0.60
1:C:17:GLN:NE2	6:C:908:HOH:O	2.35	0.59
1:A:664[B]:ARG:HH21	1:A:664[B]:ARG:CG	2.16	0.58
1:C:335:LYS:NZ	1:C:356:ASP:OD1	2.32	0.57
1:C:427:VAL:HG23	1:C:484:TYR:OH	2.03	0.57
1:A:664[B]:ARG:HG2	1:A:664[B]:ARG:NH2	2.19	0.57
1:C:358:LYS:HD2	1:C:379:VAL:HG13	1.87	0.56
1:A:276:ILE:HD12	4:A:806:GOL:H12	1.87	0.56
1:C:362:GLN:HE22	4:C:809:GOL:C1	2.18	0.56
1:B:54:LEU:HA	1:B:57:CYS:SG	2.46	0.56
1:B:276:ILE:HD12	4:B:810:GOL:O1	2.06	0.55
1:A:353:TYR:CG	4:A:807:GOL:H32	2.42	0.54
1:B:57:CYS:HB3	6:B:1352:HOH:O	2.07	0.54
1:B:488:ARG:HE	4:B:808:GOL:H11	1.73	0.54
1:B:66[B]:ARG:NE	1:B:66[B]:ARG:HA	2.23	0.53
1:B:488:ARG:HE	4:B:808:GOL:C1	2.20	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:22:HIS:HE1	2:A:801:B3P:O5	1.91	0.53
1:A:66[B]:ARG:NH1	1:A:69:GLU:HG2	2.24	0.53
1:B:404:PRO:HG3	1:B:427:VAL:HG23	1.91	0.53
1:C:404:PRO:HG3	1:C:484:TYR:CZ	2.44	0.52
1:B:551:ASN:ND2	4:B:808:GOL:O2	2.43	0.52
1:B:73:TYR:CD1	1:B:425:VAL:HG11	2.45	0.51
1:B:251:ILE:HB	1:B:260:ARG:HB2	1.91	0.51
1:C:431:ASP:HB3	1:C:434:ASP:OD2	2.11	0.51
4:C:803:GOL:H12	6:C:1043:HOH:O	2.12	0.50
1:A:48:LYS:HG3	6:A:983:HOH:O	2.11	0.50
1:C:79:HIS:NE2	1:C:103:GLN:NE2	2.60	0.49
1:B:358:LYS:HD2	1:B:379:VAL:HG13	1.95	0.48
1:B:182:GLY:O	6:B:902:HOH:O	2.20	0.47
1:C:65:GLU:O	1:C:69[B]:GLU:OE1	2.31	0.47
1:C:539:LYS:HB3	1:C:539:LYS:HE3	1.67	0.47
1:C:206:LEU:O	4:C:803:GOL:H31	2.15	0.47
1:A:70:LEU:CD2	1:A:429:GLY:HA3	2.45	0.46
1:A:440:ILE:C	1:A:440:ILE:HD12	2.40	0.46
1:B:607:HIS:CD2	1:B:610:TRP:CH2	3.03	0.46
1:C:416:GLU:O	1:C:416:GLU:OE2	2.34	0.45
1:C:341:ILE:HA	1:C:350:VAL:O	2.16	0.45
1:C:348:PHE:CD2	1:C:363:LEU:HD21	2.52	0.45
1:B:270:SER:HA	4:B:804:GOL:H12	1.97	0.45
1:B:416:GLU:H	1:B:416:GLU:HG2	1.57	0.45
1:C:70:LEU:CD2	1:C:429:GLY:HA3	2.48	0.44
1:C:538:ILE:HD13	1:C:545:PRO:HG3	2.00	0.44
1:A:227:GLU:O	4:A:806:GOL:H11	2.18	0.43
1:A:415:GLU:HG3	1:A:416:GLU:N	2.33	0.43
1:C:268:GLN:OE1	1:C:268:GLN:CA	2.66	0.43
1:A:242:ASP:OD2	6:A:904:HOH:O	2.21	0.43
1:A:322:GLU:HG3	6:A:1077:HOH:O	2.18	0.42
1:C:70:LEU:HD22	1:C:429:GLY:HA3	2.01	0.42
1:C:162:LYS:HE2	1:C:162:LYS:HB3	1.59	0.42
1:B:323:GLU:HA	1:B:326:TRP:CE2	2.54	0.42
1:C:268:GLN:OE1	1:C:268:GLN:HA	2.19	0.42
1:A:79:HIS:NE2	1:A:103:GLN:NE2	2.67	0.42
1:C:252:ARG:NH2	1:C:256:ASP:O	2.52	0.42
1:B:73:TYR:HA	1:B:74:PRO:HD3	1.92	0.42
1:B:319:ARG:HD3	1:B:319:ARG:HA	1.81	0.42
1:C:483:ASN:HB2	6:C:1399:HOH:O	2.20	0.41
1:A:407:ILE:HD12	1:A:423:ARG:HD2	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:358:LYS:HE2	1:B:358:LYS:HB2	1.90	0.41
1:A:44:GLU:CG	1:A:48:LYS:HE2	2.51	0.41
1:B:4:LEU:HD23	1:B:4:LEU:HA	1.97	0.41
1:B:335:LYS:HE2	6:B:1132:HOH:O	2.20	0.40
1:C:262:TRP:HH2	5:C:808:PEG:H11	1.86	0.40
1:A:427:VAL:C	1:A:428:LYS:O	2.64	0.40
1:A:677:LYS:HB3	1:A:677:LYS:HE2	1.88	0.40

All (4) 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)
6:A:1584:HOH:O	6:A:1608:HOH:O[2_546]	2.10	0.10
6:B:1498:HOH:O	6:C:1337:HOH:O[2_455]	2.11	0.09
6:A:1023:HOH:O	6:C:1011:HOH:O[2_445]	2.12	0.08
1:A:368:THR:O	4:B:804:GOL:H11[2_545]	1.57	0.03

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	710/711 (100%)	688 (97%)	20 (3%)	2 (0%)	36	17
1	B	711/711 (100%)	686 (96%)	23 (3%)	2 (0%)	36	17
1	C	708/711 (100%)	685 (97%)	21 (3%)	2 (0%)	36	17
All	All	2129/2133 (100%)	2059 (97%)	64 (3%)	6 (0%)	36	17

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	428	LYS
1	B	5	GLN

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Mol	Chain	Res	Type
1	A	357	VAL
1	C	554	SER
1	B	357	VAL
1	C	357	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	619/617 (100%)	605 (98%)	14 (2%)	44	16
1	B	620/617 (100%)	605 (98%)	15 (2%)	43	15
1	C	617/617 (100%)	597 (97%)	20 (3%)	34	8
All	All	1856/1851 (100%)	1807 (97%)	49 (3%)	42	13

All (49) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	106	LEU
1	A	111	ARG
1	A	169	GLU
1	A	261	LEU
1	A	271	SER
1	A	358	LYS
1	A	403	SER
1	A	445	LYS
1	A	449	LYS
1	A	499[A]	LEU
1	A	499[B]	LEU
1	A	512	GLU
1	A	539	LYS
1	A	598	ASP
1	B	59	ILE
1	B	69	GLU
1	B	111	ARG
1	B	192[A]	GLN

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Mol	Chain	Res	Type
1	B	192[B]	GLN
1	B	283	ILE
1	B	335	LYS
1	B	358	LYS
1	B	403	SER
1	B	416	GLU
1	B	425	VAL
1	B	426	THR
1	B	431	ASP
1	B	598	ASP
1	B	624	GLU
1	C	4	LEU
1	C	66	ARG
1	C	69[A]	GLU
1	C	69[B]	GLU
1	C	106	LEU
1	C	107	GLU
1	C	257	PRO
1	C	267	GLN
1	C	268	GLN
1	C	322	GLU
1	C	415	GLU
1	C	416	GLU
1	C	420	ARG
1	C	427	VAL
1	C	430	ILE
1	C	461	LYS
1	C	512	GLU
1	C	539	LYS
1	C	598	ASP
1	C	621	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	GLN
1	A	22	HIS
1	A	103	GLN
1	A	388	GLN
1	A	577	GLN
1	B	17	GLN
1	B	103	GLN

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Mol	Chain	Res	Type
1	B	477	ASN
1	B	483	ASN
1	B	531	GLN
1	B	551	ASN
1	B	577	GLN
1	C	103	GLN
1	C	362	GLN
1	C	388	GLN
1	C	483	ASN
1	C	524	GLN
1	C	531	GLN
1	C	566	GLN
1	C	606	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](#)

28 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	$\# Z > 2$
4	GOL	B	801	-	5,5,5	0.24	0	5,5,5	0.55	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	GOL	B	803	-	5,5,5	0.16	0	5,5,5	0.46	0
4	GOL	C	805	-	5,5,5	0.38	0	5,5,5	0.51	0
4	GOL	B	811	-	5,5,5	0.17	0	5,5,5	0.79	0
4	GOL	A	803	-	5,5,5	0.25	0	5,5,5	1.02	0
4	GOL	B	809	-	5,5,5	0.25	0	5,5,5	0.15	0
4	GOL	A	806	-	5,5,5	1.10	1 (20%)	5,5,5	2.64	2 (40%)
5	PEG	A	805	-	6,6,6	0.32	0	5,5,5	0.13	0
4	GOL	A	808	-	5,5,5	0.45	0	5,5,5	1.07	0
4	GOL	C	802	-	5,5,5	0.45	0	5,5,5	0.61	0
3	A1CQR	A	802	-	29,29,29	1.18	2 (6%)	34,40,40	1.38	6 (17%)
5	PEG	B	806	-	6,6,6	0.53	0	5,5,5	0.30	0
4	GOL	C	809	-	5,5,5	0.45	0	5,5,5	0.82	0
5	PEG	C	806	-	6,6,6	0.35	0	5,5,5	0.21	0
4	GOL	B	808	-	5,5,5	0.30	0	5,5,5	0.53	0
2	B3P	A	801	-	18,18,18	0.82	0	23,23,23	1.02	1 (4%)
4	GOL	B	810	-	5,5,5	0.45	0	5,5,5	1.90	3 (60%)
4	GOL	B	805	-	5,5,5	0.36	0	5,5,5	0.99	0
4	GOL	C	803	-	5,5,5	0.52	0	5,5,5	0.76	0
4	GOL	A	807	-	5,5,5	0.16	0	5,5,5	0.80	0
4	GOL	B	807	-	5,5,5	0.10	0	5,5,5	0.49	0
3	A1CQR	B	802	-	29,29,29	1.30	3 (10%)	34,40,40	1.52	7 (20%)
3	A1CQR	C	801	-	29,29,29	1.21	2 (6%)	34,40,40	1.24	4 (11%)
4	GOL	A	804	-	5,5,5	0.50	0	5,5,5	0.73	0
4	GOL	B	804	-	5,5,5	1.43	1 (20%)	5,5,5	1.06	1 (20%)
5	PEG	C	808	-	6,6,6	0.48	0	5,5,5	0.53	0
4	GOL	C	807	-	5,5,5	0.30	0	5,5,5	0.59	0
4	GOL	C	804	-	5,5,5	0.78	0	5,5,5	1.06	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
4	GOL	B	801	-	-	2/4/4/4	-
4	GOL	B	803	-	-	0/4/4/4	-
4	GOL	C	805	-	-	4/4/4/4	-
4	GOL	B	811	-	-	4/4/4/4	-
4	GOL	A	803	-	-	2/4/4/4	-
4	GOL	B	809	-	-	0/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GOL	A	806	-	-	4/4/4/4	-
5	PEG	A	805	-	-	3/4/4/4	-
4	GOL	A	808	-	-	1/4/4/4	-
4	GOL	C	802	-	-	0/4/4/4	-
3	A1CQR	A	802	-	-	0/13/20/20	0/4/4/4
5	PEG	B	806	-	-	2/4/4/4	-
4	GOL	C	809	-	-	4/4/4/4	-
5	PEG	C	806	-	-	3/4/4/4	-
4	GOL	B	808	-	-	1/4/4/4	-
2	B3P	A	801	-	-	0/28/28/28	-
4	GOL	B	810	-	-	4/4/4/4	-
4	GOL	B	805	-	-	2/4/4/4	-
4	GOL	C	803	-	-	4/4/4/4	-
4	GOL	A	807	-	-	3/4/4/4	-
4	GOL	B	807	-	-	2/4/4/4	-
3	A1CQR	B	802	-	-	1/13/20/20	0/4/4/4
3	A1CQR	C	801	-	-	0/13/20/20	0/4/4/4
4	GOL	A	804	-	-	0/4/4/4	-
4	GOL	B	804	-	-	0/4/4/4	-
5	PEG	C	808	-	-	3/4/4/4	-
4	GOL	C	807	-	-	2/4/4/4	-
4	GOL	C	804	-	-	3/4/4/4	-

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	801	A1CQR	C3-N1	4.67	1.36	1.29
3	A	802	A1CQR	C3-N1	3.79	1.34	1.29
3	B	802	A1CQR	C3-N1	3.49	1.34	1.29
3	B	802	A1CQR	C16-C12	3.21	1.45	1.40
4	B	804	GOL	O2-C2	-3.02	1.34	1.43
3	A	802	A1CQR	C12-N2	2.49	1.37	1.35
4	A	806	GOL	O2-C2	-2.37	1.36	1.43
3	B	802	A1CQR	N2-N1	2.14	1.41	1.36
3	C	801	A1CQR	C2-C3	-2.01	1.47	1.50

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	806	GOL	O2-C2-C1	-3.88	93.12	109.18
3	B	802	A1CQR	C16-C12-N3	-3.80	122.94	126.97
4	A	806	GOL	C3-C2-C1	3.69	125.34	111.80
3	A	802	A1CQR	O1-C1-N4	-3.61	115.69	122.12
3	A	802	A1CQR	C16-C12-N3	-3.59	123.16	126.97
3	B	802	A1CQR	O1-C1-N4	-3.58	115.75	122.12
3	C	801	A1CQR	C16-C12-N3	-3.41	123.35	126.97
3	B	802	A1CQR	C14-C15-C16	-3.11	117.69	121.86
3	C	801	A1CQR	O1-C1-N4	-2.98	116.81	122.12
4	B	810	GOL	O2-C2-C1	-2.90	97.15	109.18
2	A	801	B3P	C11-C8-C9	-2.64	104.22	110.02
3	B	802	A1CQR	N3-C12-N2	2.60	129.40	125.34
3	A	802	A1CQR	O1-C1-C2	2.57	127.21	121.28
3	B	802	A1CQR	C5-C4-N2	2.43	114.49	111.93
3	A	802	A1CQR	C14-C15-C16	-2.41	118.64	121.86
3	A	802	A1CQR	N3-C12-N2	2.36	129.03	125.34
3	C	801	A1CQR	N3-C12-N2	2.36	129.03	125.34
3	B	802	A1CQR	C12-N2-N1	-2.26	109.78	111.05
3	A	802	A1CQR	C4-N2-N1	2.25	121.69	119.75
4	B	810	GOL	C3-C2-C1	2.21	119.92	111.80
4	B	804	GOL	C3-C2-C1	2.08	119.42	111.80
3	B	802	A1CQR	O1-C1-C2	2.08	126.06	121.28
3	C	801	A1CQR	C2-C3-N1	2.07	123.09	118.75
4	B	810	GOL	O1-C1-C2	-2.04	101.21	110.38

There are no chirality outliers.

All (54) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	802	A1CQR	O1-C1-N4-C20
4	A	807	GOL	O1-C1-C2-C3
4	B	807	GOL	C1-C2-C3-O3
4	B	810	GOL	O1-C1-C2-C3
4	B	810	GOL	C1-C2-C3-O3
4	B	811	GOL	O1-C1-C2-C3
4	C	803	GOL	C1-C2-C3-O3
4	C	804	GOL	C1-C2-C3-O3
4	C	804	GOL	O2-C2-C3-O3
4	C	805	GOL	O1-C1-C2-O2
4	C	805	GOL	O1-C1-C2-C3
4	C	805	GOL	C1-C2-C3-O3
4	C	807	GOL	C1-C2-C3-O3
4	C	809	GOL	O1-C1-C2-C3

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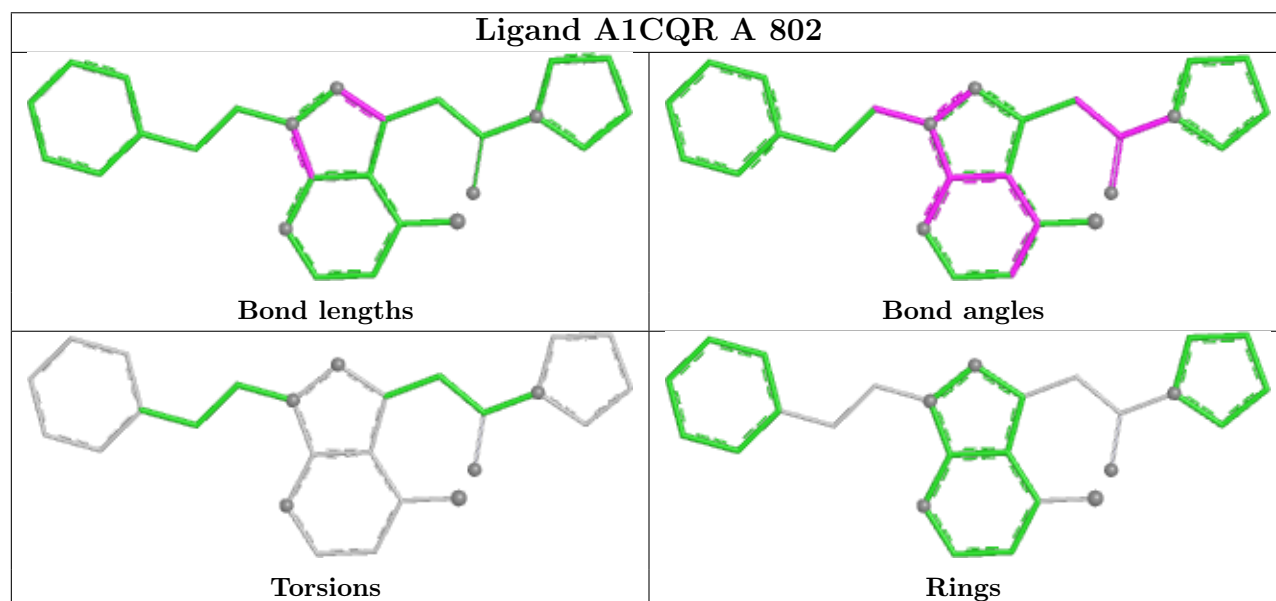
Mol	Chain	Res	Type	Atoms
4	C	809	GOL	C1-C2-C3-O3
4	C	809	GOL	O2-C2-C3-O3
5	A	805	PEG	O2-C3-C4-O4
4	B	810	GOL	O2-C2-C3-O3
4	C	803	GOL	O2-C2-C3-O3
4	C	805	GOL	O2-C2-C3-O3
5	C	806	PEG	O1-C1-C2-O2
4	A	806	GOL	O1-C1-C2-C3
4	A	806	GOL	C1-C2-C3-O3
4	A	808	GOL	O1-C1-C2-C3
4	B	801	GOL	C1-C2-C3-O3
4	B	805	GOL	C1-C2-C3-O3
4	B	811	GOL	C1-C2-C3-O3
4	C	803	GOL	O1-C1-C2-C3
5	A	805	PEG	O1-C1-C2-O2
4	A	806	GOL	O2-C2-C3-O3
4	A	807	GOL	O1-C1-C2-O2
4	C	809	GOL	O1-C1-C2-O2
5	C	808	PEG	O1-C1-C2-O2
5	B	806	PEG	O2-C3-C4-O4
5	C	806	PEG	O2-C3-C4-O4
4	B	801	GOL	O2-C2-C3-O3
4	B	810	GOL	O1-C1-C2-O2
4	B	805	GOL	O2-C2-C3-O3
4	B	807	GOL	O2-C2-C3-O3
4	B	811	GOL	O2-C2-C3-O3
4	B	808	GOL	O2-C2-C3-O3
4	B	811	GOL	O1-C1-C2-O2
4	C	807	GOL	O2-C2-C3-O3
5	C	806	PEG	C4-C3-O2-C2
5	C	808	PEG	C1-C2-O2-C3
4	A	807	GOL	O2-C2-C3-O3
5	A	805	PEG	C1-C2-O2-C3
5	B	806	PEG	C1-C2-O2-C3
4	C	803	GOL	O1-C1-C2-O2
4	A	803	GOL	O1-C1-C2-C3
5	C	808	PEG	C4-C3-O2-C2
4	C	804	GOL	O1-C1-C2-O2
4	A	803	GOL	O1-C1-C2-O2
4	A	806	GOL	O1-C1-C2-O2

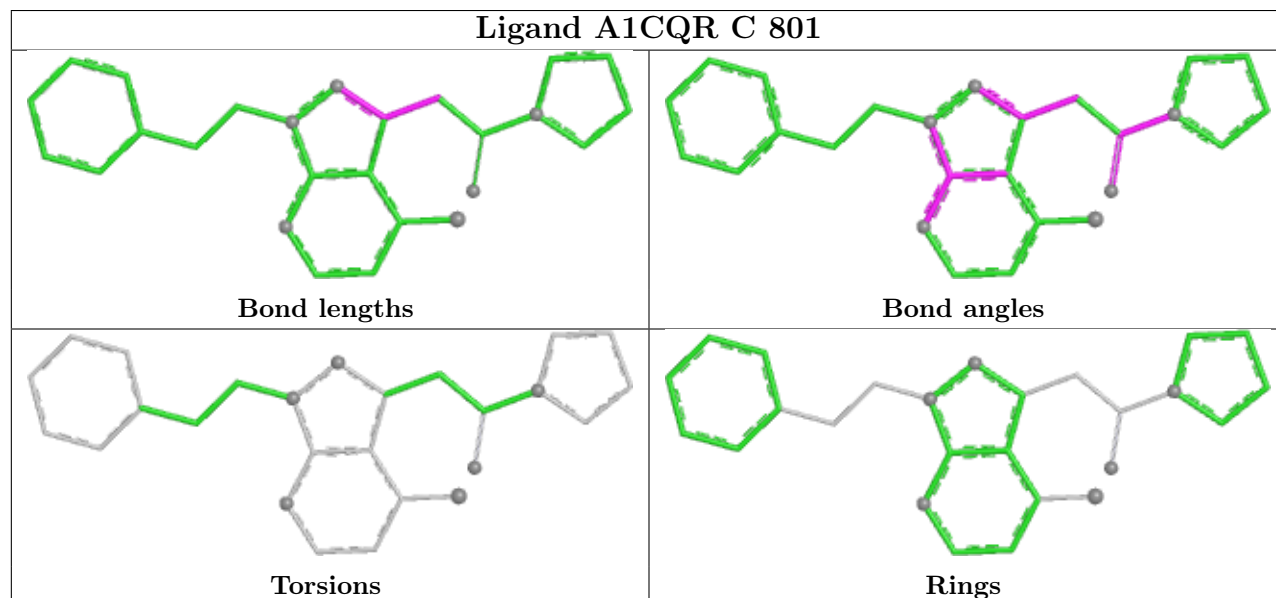
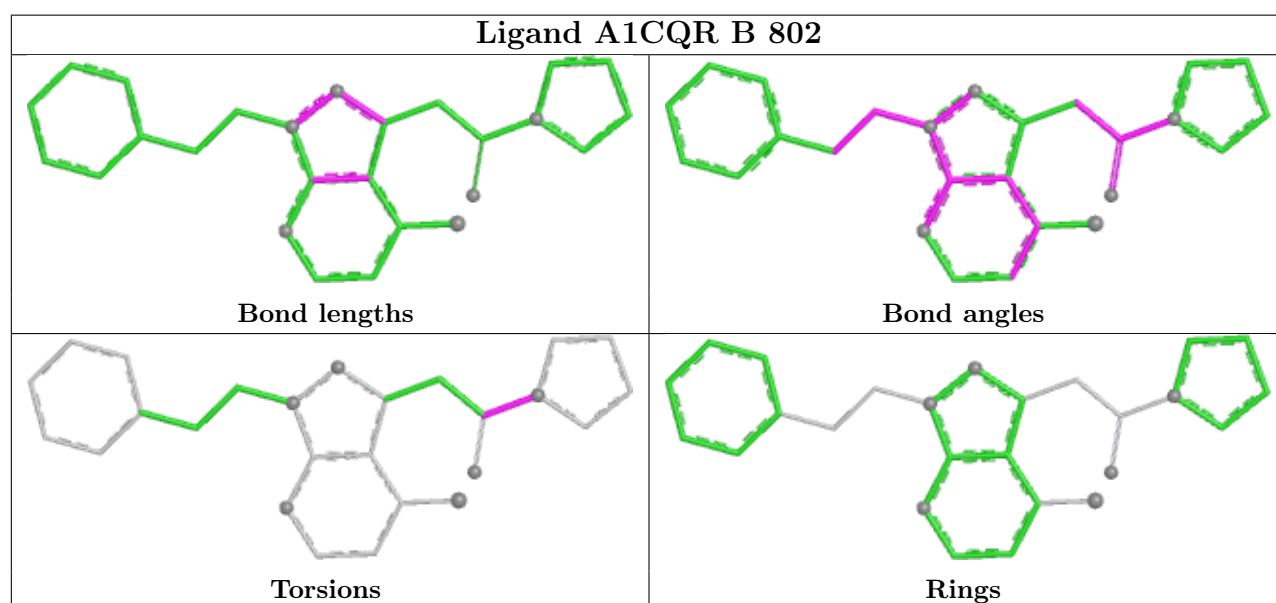
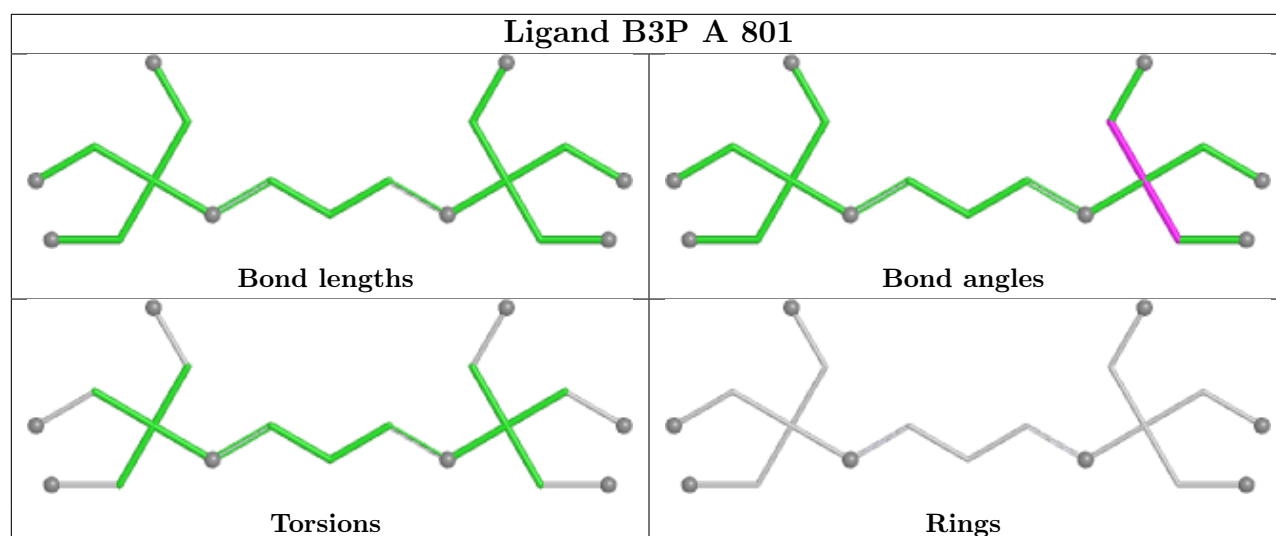
There are no ring outliers.

11 monomers are involved in 22 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	811	GOL	2	0
4	A	806	GOL	2	0
4	C	809	GOL	3	0
4	B	808	GOL	3	0
2	A	801	B3P	1	0
4	B	810	GOL	1	0
4	C	803	GOL	2	0
4	A	807	GOL	1	0
4	B	807	GOL	1	0
4	B	804	GOL	2	1
5	C	808	PEG	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	706/711 (99%)	-0.66	7 (0%) 79 82	7, 17, 33, 73	6 (0%)
1	B	708/711 (99%)	-0.42	6 (0%) 82 86	11, 20, 39, 74	5 (0%)
1	C	707/711 (99%)	-0.17	5 (0%) 84 87	9, 24, 47, 81	3 (0%)
All	All	2121/2133 (99%)	-0.41	18 (0%) 82 86	7, 20, 42, 81	14 (0%)

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	425	VAL	4.9
1	B	426	THR	4.9
1	A	429	GLY	4.3
1	C	4	LEU	4.2
1	C	426	THR	4.0
1	A	427	VAL	4.0
1	B	4	LEU	3.8
1	C	427	VAL	3.4
1	C	430	ILE	3.2
1	A	426	THR	3.0
1	B	427	VAL	2.9
1	A	428	LYS	2.8
1	A	425	VAL	2.7
1	B	3	SER	2.4
1	B	57	CYS	2.2
1	C	425	VAL	2.2
1	A	417	LEU	2.1
1	A	5	GLN	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.

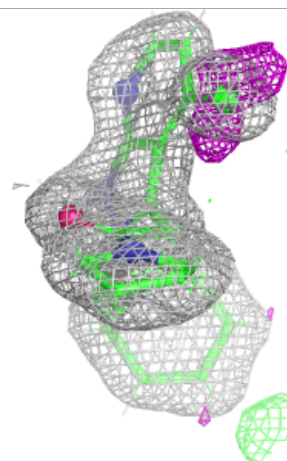
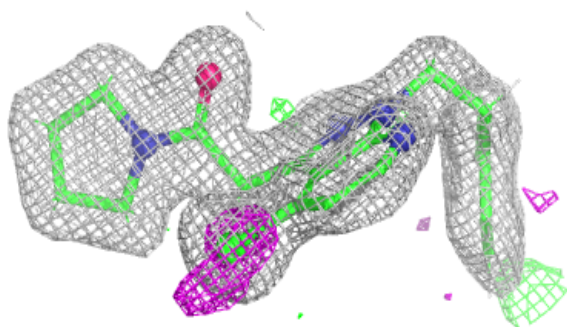
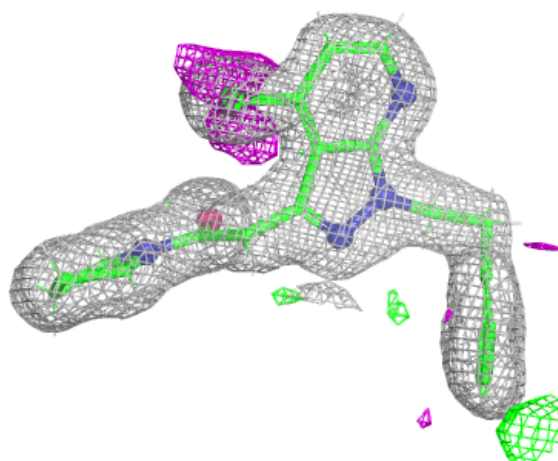
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	GOL	C	805	6/6	0.71	0.25	25,26,39,39	14
5	PEG	C	806	7/7	0.79	0.19	21,25,39,39	17
4	GOL	C	809	6/6	0.81	0.23	20,24,39,39	14
4	GOL	C	804	6/6	0.82	0.15	32,39,46,50	3
4	GOL	B	807	6/6	0.83	0.14	39,47,52,54	3
4	GOL	B	808	6/6	0.83	0.21	19,23,39,39	14
4	GOL	B	810	6/6	0.85	0.13	31,41,45,48	3
5	PEG	A	805	7/7	0.87	0.11	38,47,50,55	2
4	GOL	A	806	6/6	0.87	0.12	20,27,39,39	14
4	GOL	B	811	6/6	0.88	0.12	28,39,47,51	3
5	PEG	B	806	7/7	0.88	0.12	36,50,55,59	2
4	GOL	B	809	6/6	0.88	0.17	18,26,39,39	14
4	GOL	C	803	6/6	0.89	0.12	26,36,39,54	3
4	GOL	C	807	6/6	0.89	0.11	39,41,44,46	3
4	GOL	B	801	6/6	0.89	0.10	28,37,39,41	3
4	GOL	A	808	6/6	0.90	0.10	26,35,39,44	3
4	GOL	B	805	6/6	0.91	0.10	32,38,39,43	3
5	PEG	C	808	7/7	0.91	0.10	19,32,39,52	2
4	GOL	C	802	6/6	0.93	0.08	21,31,39,39	3
4	GOL	A	803	6/6	0.94	0.07	29,33,39,39	3
4	GOL	B	804	6/6	0.94	0.12	12,23,39,39	14
3	A1CQR	C	801	26/26	0.95	0.07	18,25,32,36	0
4	GOL	A	807	6/6	0.95	0.07	24,31,39,39	3
4	GOL	A	804	6/6	0.95	0.06	20,25,39,39	3
3	A1CQR	B	802	26/26	0.96	0.07	16,22,31,31	0
4	GOL	B	803	6/6	0.97	0.05	23,25,39,39	3
3	A1CQR	A	802	26/26	0.97	0.06	13,21,26,27	0
2	B3P	A	801	19/19	0.98	0.04	11,14,39,39	8

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.

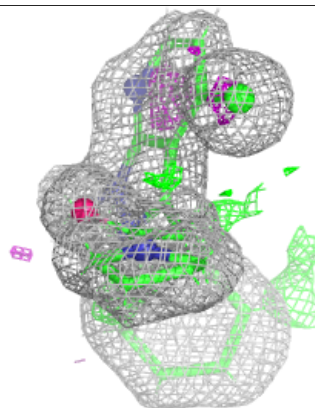
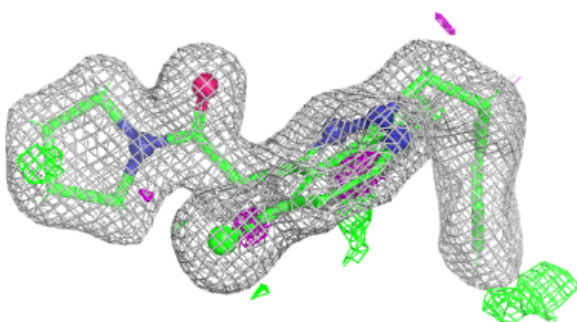
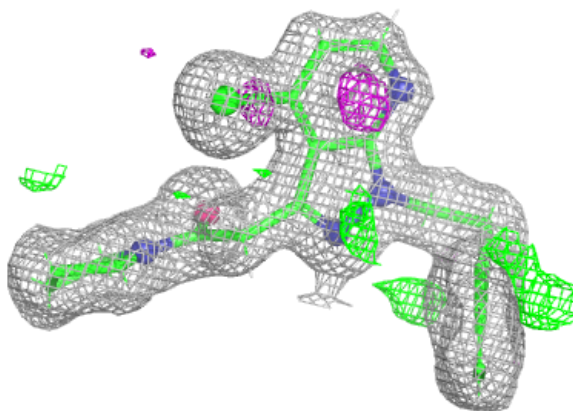
Electron density around A1CQR C 801:

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

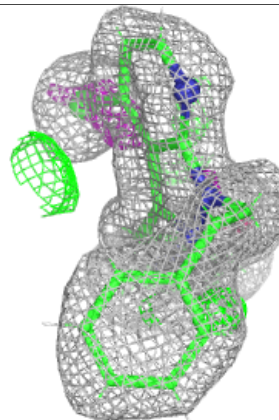
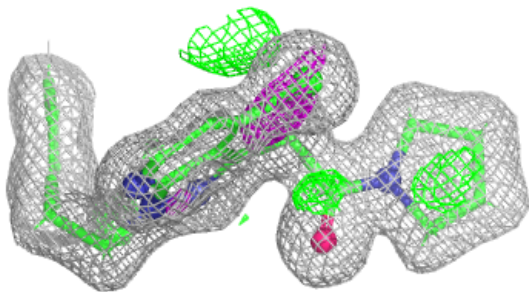
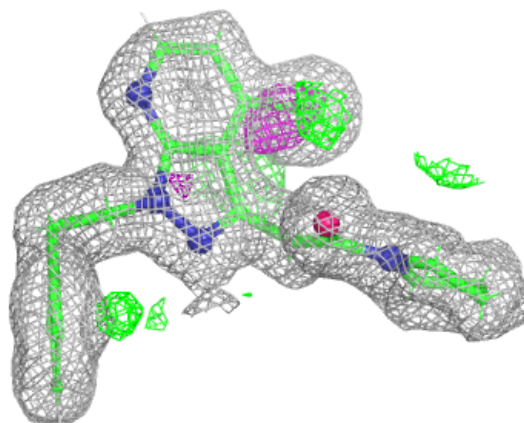


Electron density around A1CQR B 802:

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

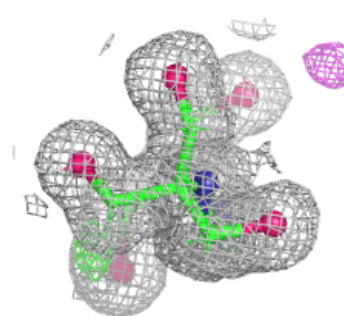
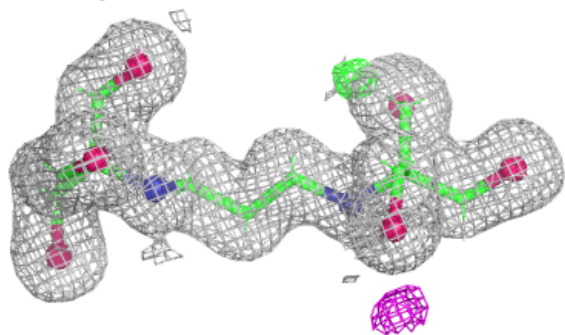
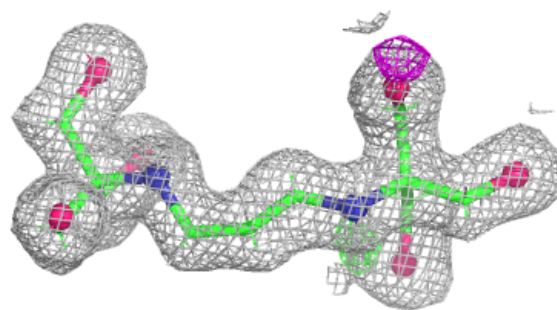
**Electron density around A1CQR A 802:**

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



Electron density around B3P 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.