



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:08 AM UTC

PDB ID : 4MLP / pdb_00004mlp
Title : Mammalian cryptochrome in complex with a small molecule competitor of its ubiquitin ligase
Authors : Nangle, S.; Xing, W.; Zheng, N.
Deposited on : 2013-09-06
Resolution : 1.94 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

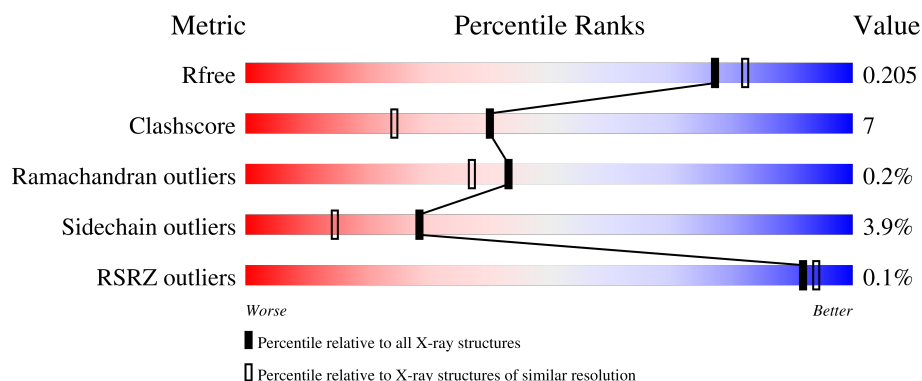
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.94 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	1452 (1.94-1.94)
Clashscore	190562	1494 (1.94-1.94)
Ramachandran outliers	187476	1479 (1.94-1.94)
Sidechain outliers	187428	1479 (1.94-1.94)
RSRZ outliers	180081	1453 (1.94-1.94)

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	512	
1	B	512	
1	C	512	
1	D	512	

2 Entry composition [i](#)

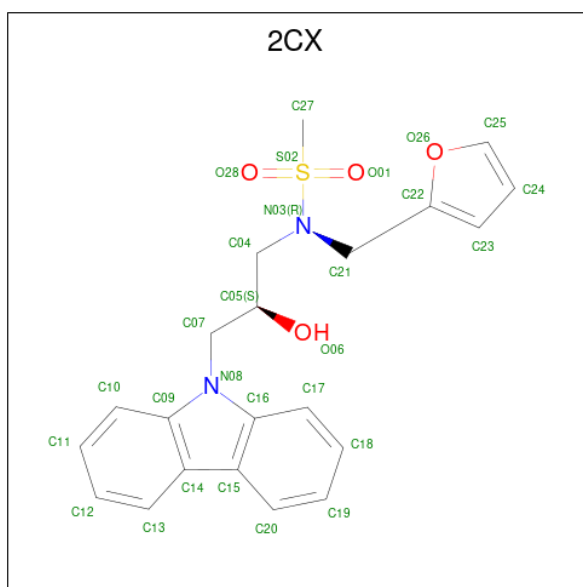
There are 3 unique types of molecules in this entry. The entry contains 17726 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Cryptochrome-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	492	Total	C	N	O	S	0	0	0
			4015	2577	713	705	20			
1	A	492	Total	C	N	O	S	0	0	0
			4015	2577	713	705	20			
1	B	492	Total	C	N	O	S	0	0	0
			4015	2577	713	705	20			
1	D	492	Total	C	N	O	S	0	0	0
			4015	2577	713	705	20			

- Molecule 2 is N-[(2S)-3-(9H-carbazol-9-yl)-2-hydroxypropyl]-N-(furan-2-ylmethyl)methanesulfonamide (CCD ID: 2CX) (formula: C₂₁H₂₂N₂O₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	C	1	Total	C	N	O	S	0	0
			28	21	2	4	1		
2	A	1	Total	C	N	O	S	0	0
			28	21	2	4	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	S	0	0
			28	21	2	4	1		
2	D	1	Total	C	N	O	S	0	0
			28	21	2	4	1		

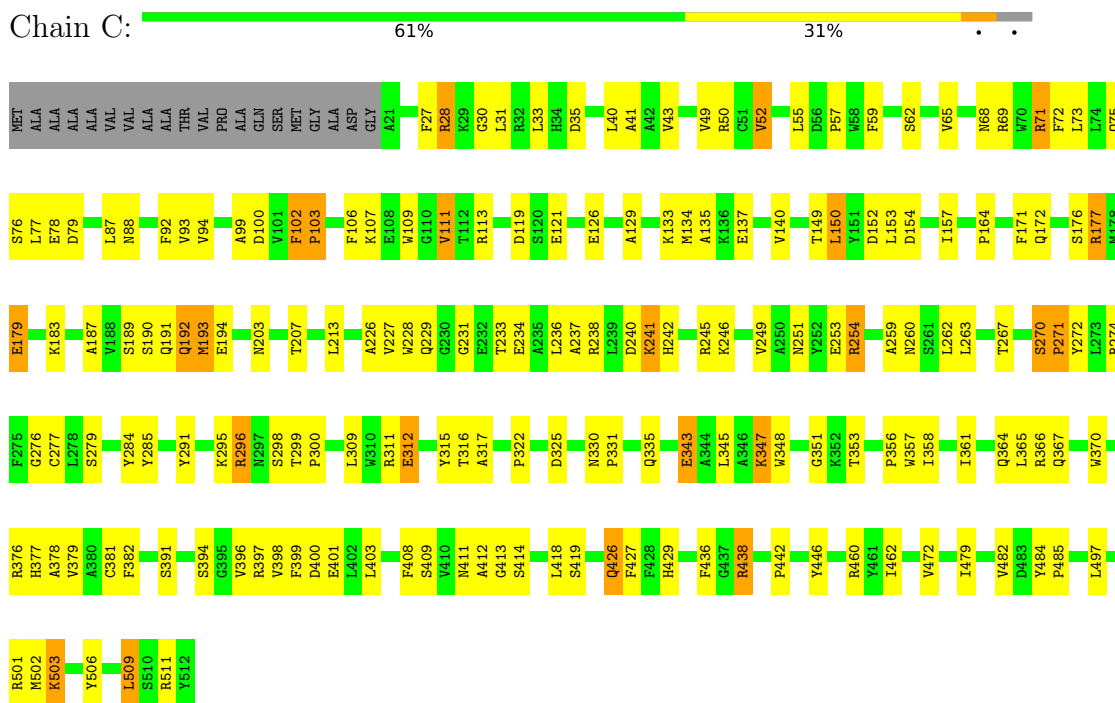
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	434	Total	O	0	0
			434	434		
3	A	461	Total	O	0	0
			461	461		
3	B	322	Total	O	0	0
			322	322		
3	D	337	Total	O	0	0
			337	337		

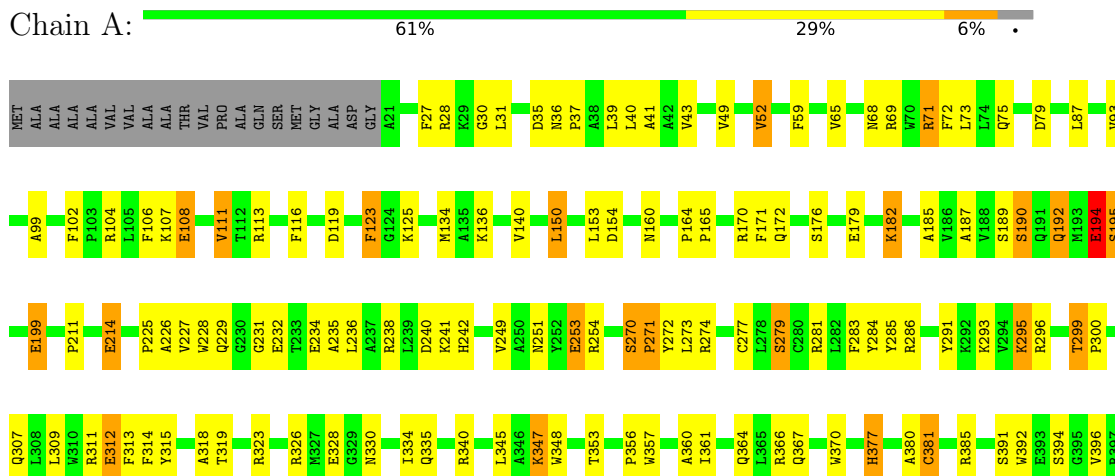
3 Residue-property plots

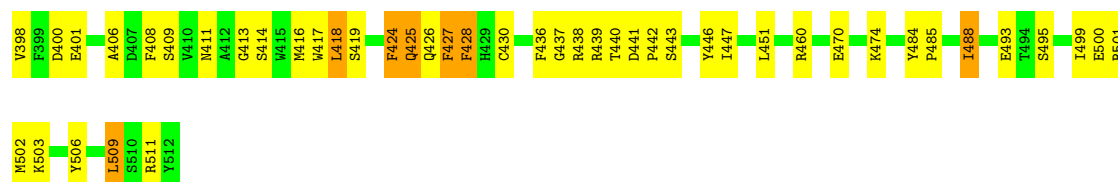
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Cryptochrome-2



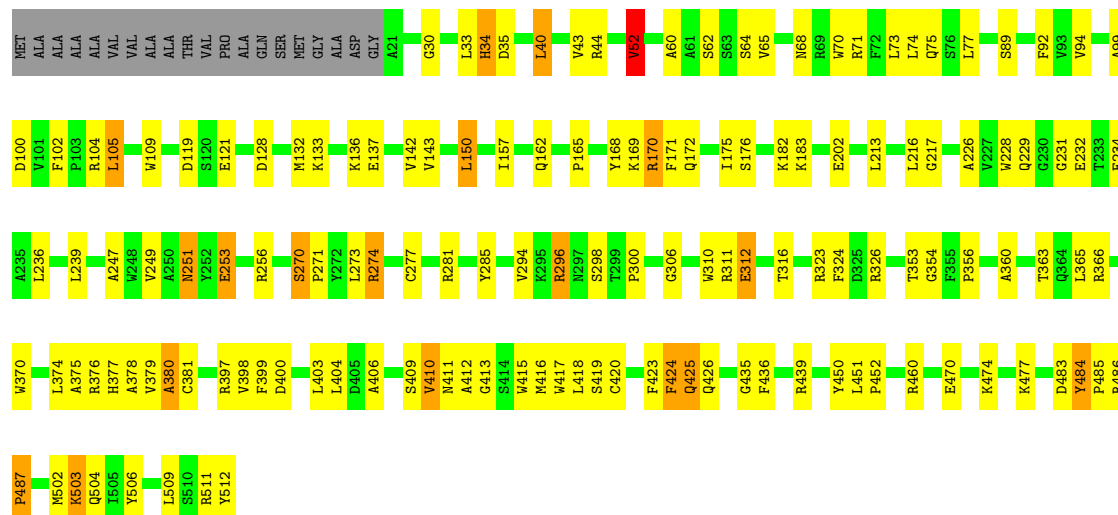
• Molecule 1: Cryptochrome-2





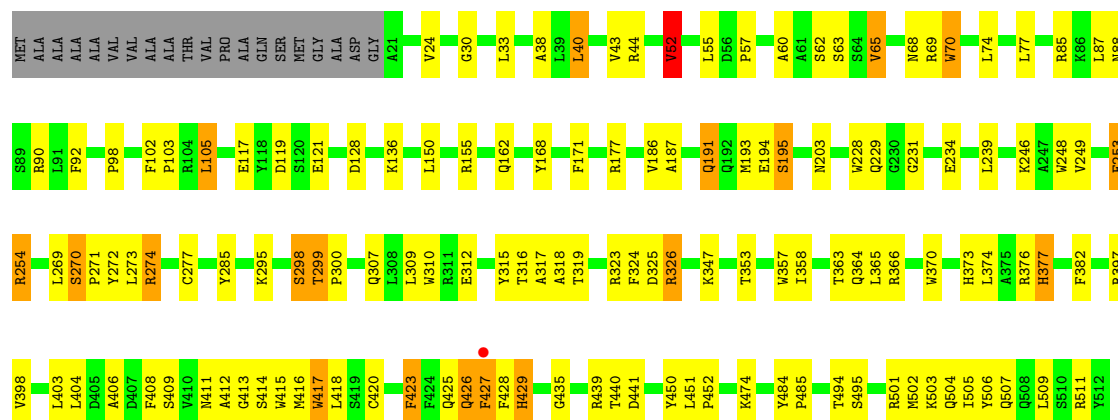
• Molecule 1: Cryptochrome-2

Chain B: 68% 25%



• Molecule 1: Cryptochrome-2

Chain D: 69% 23%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	48.21Å 93.65Å 141.18Å 89.94° 90.09° 90.25°	Depositor
Resolution (Å)	44.46 – 1.94 44.46 – 1.94	Depositor EDS
% Data completeness (in resolution range)	(Not available) (44.46-1.94) 93.8 (44.46-1.94)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.04 (at 1.94Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1168)	Depositor
R, R_{free}	0.179 , 0.217 0.186 , 0.205	Depositor DCC
R_{free} test set	1530 reflections (0.65%)	wwPDB-VP
Wilson B-factor (Å ²)	26.0	Xtriage
Anisotropy	0.337	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 37.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.109 for h,-k,-l 0.110 for -h,k,-l 0.457 for -h,-k,l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	17726	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.09% 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: 2CX

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	1.96	122/4133 (3.0%)	1.05	12/5608 (0.2%)
1	B	1.83	78/4133 (1.9%)	1.04	12/5608 (0.2%)
1	C	1.97	128/4133 (3.1%)	1.06	7/5608 (0.1%)
1	D	1.82	73/4133 (1.8%)	1.04	9/5608 (0.2%)
All	All	1.90	401/16532 (2.4%)	1.05	40/22432 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	C	0	1
All	All	0	2

All (401) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	300	PRO	C-O	-9.01	1.17	1.24
1	D	300	PRO	C-O	-8.06	1.18	1.24
1	A	300	PRO	C-O	-7.90	1.18	1.24
1	C	309	LEU	C-O	-7.83	1.14	1.24
1	A	234	GLU	C-O	-7.75	1.15	1.24
1	B	300	PRO	C-O	-7.71	1.18	1.24
1	B	239	LEU	C-O	-7.50	1.15	1.24
1	A	102	PHE	C-O	-7.06	1.17	1.24
1	B	409	SER	C-O	-6.89	1.16	1.24
1	C	113	ARG	C-O	-6.88	1.16	1.24
1	A	164	PRO	C-O	-6.87	1.19	1.24
1	A	284	TYR	C-O	-6.86	1.16	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	150	LEU	C-O	-6.85	1.16	1.24
1	D	312	GLU	C-O	-6.84	1.16	1.24
1	C	400	ASP	C-O	-6.81	1.16	1.24
1	C	382	PHE	C-O	-6.75	1.16	1.24
1	A	271	PRO	C-O	-6.75	1.15	1.24
1	C	176	SER	C-O	-6.75	1.15	1.24
1	A	40	LEU	C-O	-6.74	1.16	1.24
1	C	171	PHE	C-O	-6.73	1.16	1.24
1	A	398	VAL	C-O	-6.71	1.16	1.24
1	C	150	LEU	C-O	-6.70	1.16	1.24
1	C	436	PHE	C-O	-6.70	1.16	1.24
1	B	366	ARG	C-O	-6.62	1.16	1.24
1	A	366	ARG	C-O	-6.61	1.16	1.24
1	C	73	LEU	C-O	-6.60	1.16	1.24
1	A	172	GLN	C-O	-6.59	1.16	1.24
1	C	79	ASP	C-O	-6.58	1.16	1.24
1	C	72	PHE	C-O	-6.57	1.16	1.24
1	A	370	TRP	C-O	-6.57	1.15	1.23
1	C	227	VAL	C-O	-6.57	1.16	1.24
1	D	239	LEU	C-O	-6.56	1.16	1.24
1	C	40	LEU	C-O	-6.54	1.16	1.24
1	A	72	PHE	C-O	-6.45	1.16	1.24
1	C	242	HIS	C-O	-6.45	1.16	1.24
1	A	377	HIS	C-O	-6.44	1.16	1.24
1	C	164	PRO	C-O	-6.43	1.19	1.24
1	C	240	ASP	C-O	-6.43	1.16	1.24
1	C	394	SER	C-O	-6.42	1.16	1.24
1	D	65	VAL	C-O	-6.38	1.17	1.24
1	D	40	LEU	C-O	-6.38	1.16	1.24
1	C	93	VAL	C-O	-6.37	1.16	1.24
1	D	171	PHE	C-O	-6.37	1.16	1.24
1	C	381	CYS	C-O	-6.35	1.16	1.24
1	A	394	SER	C-O	-6.35	1.16	1.24
1	C	397	ARG	C-O	-6.35	1.16	1.24
1	D	325	ASP	C-N	-6.35	1.24	1.33
1	D	119	ASP	C-O	-6.35	1.16	1.24
1	C	367	GLN	C-O	-6.32	1.16	1.24
1	A	367	GLN	C-O	-6.30	1.16	1.24
1	A	314	PHE	C-O	-6.30	1.16	1.24
1	C	345	LEU	C-O	-6.29	1.16	1.24
1	A	408	PHE	C-O	-6.29	1.16	1.24
1	D	102	PHE	C-O	-6.29	1.18	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	274	ARG	C-O	-6.29	1.16	1.24
1	C	462	ILE	C-O	-6.27	1.16	1.24
1	B	353	THR	C-O	-6.27	1.16	1.24
1	B	417	TRP	C-O	-6.26	1.17	1.24
1	A	281	ARG	C-O	-6.25	1.16	1.24
1	D	274	ARG	C-O	-6.25	1.17	1.24
1	B	171	PHE	C-O	-6.24	1.17	1.24
1	A	73	LEU	C-O	-6.23	1.17	1.24
1	A	238	ARG	C-O	-6.23	1.16	1.24
1	A	307	GLN	C-O	-6.20	1.16	1.24
1	B	416	MET	C-O	-6.20	1.16	1.24
1	C	353	THR	C-O	-6.19	1.16	1.24
1	A	171	PHE	C-O	-6.18	1.16	1.24
1	B	451	LEU	C-O	-6.18	1.17	1.24
1	B	30	GLY	C-O	-6.17	1.17	1.24
1	B	65	VAL	C-O	-6.16	1.17	1.24
1	D	74	LEU	C-O	-6.16	1.17	1.24
1	B	274	ARG	C-O	-6.16	1.17	1.24
1	A	356	PRO	C-O	-6.14	1.16	1.24
1	C	438	ARG	C-O	-6.14	1.17	1.24
1	A	345	LEU	C-O	-6.13	1.17	1.24
1	D	398	VAL	C-O	-6.13	1.16	1.24
1	A	240	ASP	C-O	-6.12	1.17	1.24
1	C	315	TYR	C-O	-6.12	1.17	1.24
1	A	272	TYR	C-O	-6.12	1.17	1.24
1	C	234	GLU	C-O	-6.12	1.17	1.24
1	B	168	TYR	C-O	-6.11	1.16	1.24
1	D	121	GLU	C-O	-6.11	1.16	1.24
1	C	325	ASP	C-O	-6.09	1.16	1.24
1	C	68	ASN	C-O	-6.09	1.17	1.24
1	C	370	TRP	C-O	-6.07	1.16	1.23
1	B	277	CYS	C-O	-6.07	1.16	1.24
1	C	30	GLY	C-O	-6.06	1.17	1.24
1	C	102	PHE	C-O	-6.06	1.18	1.24
1	A	59	PHE	C-O	-6.05	1.17	1.24
1	A	313	PHE	C-O	-6.04	1.17	1.24
1	A	242	HIS	C-O	-6.03	1.16	1.24
1	C	41	ALA	C-O	-6.03	1.17	1.24
1	A	30	GLY	C-O	-6.03	1.17	1.24
1	A	309	LEU	C-O	-6.02	1.17	1.24
1	A	401	GLU	C-O	-6.00	1.17	1.24
1	A	315	TYR	C-O	-5.99	1.17	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	436	PHE	C-O	-5.99	1.17	1.24
1	B	232	GLU	C-O	-5.98	1.17	1.24
1	D	270	SER	C-O	-5.98	1.19	1.24
1	C	52	VAL	C-O	-5.97	1.18	1.24
1	C	271	PRO	C-O	-5.96	1.16	1.24
1	C	365	LEU	C-O	-5.96	1.17	1.24
1	D	77	LEU	C-O	-5.95	1.17	1.24
1	B	381	CYS	C-O	-5.95	1.17	1.24
1	B	150	LEU	C-O	-5.94	1.17	1.24
1	C	311	ARG	C-O	-5.93	1.17	1.24
1	B	374	LEU	C-O	-5.91	1.17	1.24
1	A	79	ASP	C-O	-5.90	1.17	1.24
1	B	74	LEU	C-O	-5.88	1.17	1.24
1	C	262	LEU	C-O	-5.88	1.16	1.24
1	B	102	PHE	C-O	-5.87	1.19	1.24
1	B	273	LEU	C-O	-5.87	1.17	1.24
1	C	409	SER	C-O	-5.87	1.17	1.24
1	A	49	VAL	C-O	-5.86	1.17	1.24
1	C	413	GLY	C-O	-5.85	1.16	1.23
1	A	31	LEU	C-O	-5.85	1.16	1.23
1	D	409	SER	C-O	-5.85	1.17	1.24
1	D	319	THR	C-O	-5.84	1.16	1.24
1	B	316	THR	C-O	-5.84	1.17	1.24
1	B	77	LEU	C-O	-5.83	1.17	1.24
1	A	273	LEU	C-O	-5.83	1.17	1.24
1	B	234	GLU	C-O	-5.83	1.17	1.24
1	B	370	TRP	C-O	-5.83	1.17	1.23
1	A	361	ILE	C-O	-5.81	1.17	1.24
1	A	438	ARG	C-O	-5.81	1.17	1.24
1	D	417	TRP	C-O	-5.81	1.17	1.24
1	B	398	VAL	C-O	-5.80	1.17	1.24
1	D	404	LEU	C-O	-5.80	1.16	1.24
1	D	413	GLY	C-O	-5.79	1.16	1.23
1	A	414	SER	C-O	-5.79	1.17	1.24
1	C	316	THR	C-O	-5.78	1.17	1.24
1	A	123	PHE	C-O	-5.78	1.17	1.24
1	D	370	TRP	C-O	-5.78	1.17	1.23
1	A	381	CYS	C-O	-5.77	1.17	1.24
1	A	391	SER	C-O	-5.77	1.17	1.23
1	A	227	VAL	C-O	-5.77	1.17	1.24
1	A	330	ASN	C-O	-5.77	1.17	1.24
1	C	366	ARG	C-O	-5.76	1.17	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	107	LYS	C-O	-5.75	1.17	1.24
1	A	65	VAL	C-O	-5.73	1.18	1.24
1	C	446	TYR	C-O	-5.72	1.17	1.24
1	A	413	GLY	C-O	-5.72	1.16	1.23
1	B	363	THR	C-O	-5.71	1.17	1.24
1	A	153	LEU	C-O	-5.71	1.17	1.24
1	C	270	SER	C-O	-5.70	1.19	1.24
1	A	107	LYS	C-O	-5.70	1.17	1.24
1	D	68	ASN	C-O	-5.70	1.17	1.24
1	B	128	ASP	C-O	-5.70	1.17	1.24
1	A	113	ARG	C-O	-5.69	1.17	1.24
1	A	409	SER	C-O	-5.69	1.17	1.24
1	A	312	GLU	C-O	-5.68	1.17	1.24
1	D	128	ASP	C-O	-5.68	1.17	1.24
1	D	168	TYR	C-O	-5.68	1.17	1.24
1	A	235	ALA	C-O	-5.65	1.17	1.24
1	A	460	ARG	C-O	-5.64	1.16	1.24
1	C	272	TYR	C-O	-5.64	1.17	1.24
1	A	229	GLN	C-O	-5.64	1.17	1.23
1	D	310	TRP	C-O	-5.64	1.17	1.24
1	C	119	ASP	C-O	-5.63	1.17	1.24
1	C	411	ASN	C-O	-5.63	1.17	1.24
1	C	414	SER	C-O	-5.63	1.17	1.24
1	D	366	ARG	C-O	-5.62	1.17	1.24
1	A	400	ASP	C-O	-5.62	1.17	1.24
1	C	291	TYR	C-O	-5.61	1.17	1.24
1	D	411	ASN	C-O	-5.61	1.17	1.24
1	A	154	ASP	C-O	-5.61	1.17	1.24
1	A	364	GLN	C-O	-5.61	1.17	1.24
1	C	153	LEU	C-O	-5.60	1.17	1.24
1	D	451	LEU	C-O	-5.60	1.18	1.24
1	C	376	ARG	C-O	-5.60	1.17	1.24
1	B	310	TRP	C-O	-5.60	1.17	1.24
1	C	408	PHE	C-O	-5.60	1.17	1.24
1	D	38	ALA	C-O	-5.59	1.17	1.24
1	A	335	GLN	C-O	-5.59	1.17	1.24
1	D	382	PHE	C-O	-5.59	1.17	1.24
1	C	228	TRP	C-O	-5.58	1.17	1.23
1	A	41	ALA	C-O	-5.58	1.17	1.24
1	C	241	LYS	C-O	-5.58	1.17	1.24
1	B	365	LEU	C-O	-5.57	1.17	1.24
1	D	353	THR	C-O	-5.57	1.16	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	70	TRP	C-O	-5.57	1.17	1.24
1	C	312	GLU	C-O	-5.57	1.17	1.24
1	B	172	GLN	C-O	-5.57	1.17	1.24
1	A	228	TRP	C-O	-5.57	1.17	1.23
1	B	412	ALA	C-O	-5.57	1.17	1.24
1	C	238	ARG	C-O	-5.56	1.17	1.24
1	A	396	VAL	C-O	-5.55	1.17	1.24
1	B	415	TRP	C-O	-5.54	1.17	1.24
1	B	399	PHE	C-O	-5.54	1.17	1.24
1	A	283	PHE	C-O	-5.53	1.17	1.24
1	D	412	ALA	C-O	-5.53	1.17	1.24
1	A	440	THR	C-O	-5.53	1.17	1.24
1	D	373	HIS	C-O	-5.53	1.17	1.24
1	A	411	ASN	C-O	-5.53	1.17	1.24
1	D	358	ILE	C-O	-5.52	1.17	1.24
1	A	69	ARG	C-O	-5.51	1.17	1.24
1	B	75	GLN	C-O	-5.51	1.17	1.24
1	A	286	ARG	C-O	-5.51	1.17	1.24
1	C	379	VAL	C-O	-5.50	1.17	1.24
1	C	129	ALA	C-O	-5.50	1.17	1.24
1	B	121	GLU	C-O	-5.50	1.16	1.24
1	D	364	GLN	C-O	-5.49	1.17	1.24
1	A	279	SER	C-O	-5.48	1.17	1.23
1	D	228	TRP	C-O	-5.48	1.17	1.24
1	C	401	GLU	C-O	-5.47	1.17	1.24
1	C	364	GLN	C-O	-5.47	1.17	1.24
1	D	103	PRO	C-O	-5.47	1.17	1.24
1	B	379	VAL	C-O	-5.47	1.17	1.24
1	B	410	VAL	C-O	-5.46	1.17	1.24
1	A	441	ASP	C-O	-5.45	1.18	1.24
1	D	376	ARG	C-O	-5.45	1.17	1.24
1	C	229	GLN	C-O	-5.45	1.17	1.23
1	C	317	ALA	C-O	-5.45	1.17	1.24
1	D	63	SER	C-O	-5.45	1.17	1.23
1	A	277	CYS	C-O	-5.45	1.17	1.24
1	A	488	ILE	C-O	-5.44	1.16	1.24
1	B	411	ASN	C-O	-5.44	1.17	1.24
1	C	347	LYS	C-O	-5.44	1.17	1.24
1	D	317	ALA	C-O	-5.44	1.17	1.24
1	D	150	LEU	C-O	-5.43	1.17	1.24
1	A	392	TRP	C-O	-5.43	1.17	1.24
1	D	309	LEU	C-O	-5.43	1.17	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	211	PRO	C-O	-5.43	1.17	1.23
1	B	484	TYR	C-O	-5.43	1.18	1.24
1	B	143	VAL	C-O	-5.43	1.18	1.24
1	A	319	THR	C-O	-5.42	1.17	1.24
1	B	270	SER	C-O	-5.42	1.19	1.24
1	A	417	TRP	C-O	-5.41	1.17	1.24
1	C	31	LEU	C-O	-5.41	1.16	1.23
1	D	117	GLU	C-O	-5.41	1.17	1.23
1	A	71	ARG	C-O	-5.41	1.17	1.24
1	A	495	SER	C-O	-5.41	1.17	1.24
1	C	103	PRO	C-O	-5.41	1.17	1.24
1	C	59	PHE	C-O	-5.41	1.17	1.24
1	D	273	LEU	C-O	-5.40	1.17	1.24
1	C	398	VAL	C-O	-5.39	1.18	1.24
1	C	94	VAL	C-O	-5.39	1.18	1.24
1	A	179	GLU	C-O	-5.39	1.16	1.23
1	A	499	ILE	C-O	-5.39	1.17	1.24
1	B	40	LEU	C-O	-5.39	1.17	1.24
1	C	172	GLN	C-O	-5.38	1.17	1.24
1	B	312	GLU	C-O	-5.38	1.17	1.24
1	A	37	PRO	C-O	-5.38	1.17	1.24
1	B	68	ASN	C-O	-5.38	1.17	1.24
1	A	357	TRP	C-O	-5.37	1.18	1.24
1	A	419	SER	C-O	-5.37	1.16	1.24
1	A	451	LEU	C-O	-5.37	1.17	1.24
1	A	274	ARG	C-O	-5.37	1.17	1.24
1	D	377	HIS	C-O	-5.36	1.17	1.24
1	C	399	PHE	C-O	-5.36	1.17	1.24
1	A	119	ASP	C-O	-5.36	1.17	1.24
1	A	348	TRP	C-O	-5.36	1.18	1.24
1	C	403	LEU	C-O	-5.36	1.17	1.24
1	D	98	PRO	C-O	-5.36	1.17	1.24
1	D	365	LEU	C-O	-5.36	1.18	1.24
1	C	331	PRO	C-O	-5.35	1.17	1.24
1	C	246	LYS	C-O	-5.35	1.18	1.24
1	C	140	VAL	C-O	-5.35	1.18	1.24
1	A	503	LYS	C-O	-5.35	1.17	1.24
1	C	135	ALA	C-O	-5.34	1.17	1.24
1	C	71	ARG	C-O	-5.34	1.17	1.24
1	A	39	LEU	C-O	-5.34	1.17	1.24
1	A	437	GLY	C-O	-5.34	1.17	1.23
1	B	413	GLY	C-O	-5.33	1.17	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	226	ALA	C-O	-5.33	1.17	1.23
1	C	236	LEU	C-O	-5.32	1.17	1.24
1	B	378	ALA	C-O	-5.32	1.18	1.24
1	C	472	VAL	C-O	-5.32	1.18	1.24
1	D	70	TRP	C-O	-5.32	1.17	1.24
1	A	226	ALA	C-O	-5.32	1.17	1.23
1	C	76	SER	C-O	-5.31	1.17	1.24
1	B	375	ALA	C-O	-5.31	1.18	1.24
1	C	192	GLN	C-O	-5.31	1.18	1.24
1	B	360	ALA	C-O	-5.31	1.18	1.24
1	A	360	ALA	C-O	-5.30	1.18	1.24
1	D	277	CYS	C-O	-5.30	1.17	1.24
1	C	260	ASN	C-O	-5.30	1.17	1.24
1	C	377	HIS	C-O	-5.29	1.18	1.24
1	C	391	SER	C-O	-5.29	1.17	1.23
1	C	442	PRO	C-O	-5.29	1.17	1.24
1	C	152	ASP	C-O	-5.29	1.17	1.24
1	D	441	ASP	C-O	-5.28	1.18	1.24
1	C	277	CYS	C-O	-5.28	1.17	1.24
1	C	356	PRO	C-O	-5.28	1.17	1.24
1	C	482	VAL	C-O	-5.28	1.17	1.24
1	A	442	PRO	C-O	-5.28	1.17	1.24
1	A	27	PHE	C-O	-5.28	1.17	1.23
1	A	347	LYS	C-O	-5.28	1.18	1.24
1	A	443	SER	C-O	-5.27	1.17	1.24
1	A	185	ALA	C-O	-5.27	1.16	1.23
1	D	414	SER	C-O	-5.27	1.18	1.24
1	B	94	VAL	C-O	-5.27	1.18	1.24
1	C	35	ASP	C-O	-5.27	1.17	1.23
1	B	100	ASP	C-O	-5.26	1.17	1.24
1	B	236	LEU	C-O	-5.26	1.18	1.24
1	B	229	GLN	C-O	-5.26	1.17	1.23
1	C	28	ARG	C-O	-5.25	1.17	1.24
1	D	69	ARG	C-O	-5.25	1.18	1.24
1	A	116	PHE	C-O	-5.24	1.17	1.23
1	B	311	ARG	C-O	-5.24	1.18	1.24
1	C	121	GLU	C-O	-5.24	1.16	1.24
1	B	170	ARG	C-O	-5.24	1.18	1.24
1	D	105	LEU	C-O	-5.24	1.18	1.24
1	D	357	TRP	C-O	-5.24	1.18	1.24
1	A	270	SER	C-O	-5.23	1.19	1.24
1	A	311	ARG	C-O	-5.23	1.18	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	73	LEU	C-O	-5.23	1.18	1.24
1	A	406	ALA	C-O	-5.23	1.17	1.23
1	C	263	LEU	C-O	-5.23	1.17	1.23
1	C	412	ALA	C-O	-5.23	1.18	1.24
1	D	24	VAL	C-O	-5.23	1.18	1.24
1	A	241	LYS	C-O	-5.23	1.18	1.24
1	D	248	TRP	C-O	-5.23	1.18	1.24
1	D	408	PHE	C-O	-5.22	1.18	1.24
1	A	182	LYS	C-O	-5.22	1.17	1.23
1	B	157	ILE	C-O	-5.22	1.18	1.24
1	B	216	LEU	C-O	-5.21	1.17	1.24
1	C	361	ILE	C-O	-5.21	1.18	1.24
1	D	307	GLN	C-O	-5.20	1.18	1.24
1	C	460	ARG	C-O	-5.20	1.17	1.24
1	D	234	GLU	C-O	-5.20	1.18	1.24
1	C	157	ILE	C-O	-5.19	1.18	1.24
1	D	315	TYR	C-O	-5.18	1.18	1.24
1	A	75	GLN	C-O	-5.18	1.18	1.24
1	C	335	GLN	C-O	-5.18	1.17	1.24
1	C	358	ILE	C-O	-5.18	1.18	1.24
1	C	351	GLY	C-O	-5.18	1.18	1.24
1	C	65	VAL	C-O	-5.17	1.19	1.24
1	C	27	PHE	C-O	-5.17	1.17	1.23
1	A	68	ASN	C-O	-5.17	1.18	1.24
1	C	100	ASP	C-O	-5.17	1.18	1.24
1	C	322	PRO	C-O	-5.16	1.17	1.24
1	C	357	TRP	C-O	-5.16	1.18	1.24
1	A	225	PRO	C-O	-5.16	1.17	1.23
1	A	353	THR	C-O	-5.16	1.17	1.24
1	C	259	ALA	C-O	-5.16	1.18	1.24
1	A	291	TYR	C-O	-5.16	1.18	1.24
1	D	416	MET	C-O	-5.16	1.18	1.24
1	A	93	VAL	C-O	-5.16	1.18	1.24
1	C	419	SER	C-O	-5.15	1.17	1.24
1	B	119	ASP	C-O	-5.15	1.18	1.24
1	C	284	TYR	C-O	-5.15	1.18	1.24
1	C	330	ASN	C-O	-5.15	1.17	1.24
1	C	378	ALA	C-O	-5.15	1.18	1.24
1	D	374	LEU	C-O	-5.14	1.18	1.24
1	B	377	HIS	C-O	-5.14	1.18	1.24
1	D	363	THR	C-O	-5.14	1.18	1.24
1	B	175	ILE	C-O	-5.14	1.17	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	415	TRP	C-O	-5.13	1.18	1.24
1	A	176	SER	C-O	-5.13	1.17	1.24
1	B	306	GLY	C-O	-5.13	1.17	1.23
1	A	214	GLU	C-O	-5.12	1.18	1.24
1	B	487	PRO	C-O	-5.12	1.17	1.23
1	B	404	LEU	C-O	-5.12	1.17	1.24
1	B	400	ASP	C-O	-5.11	1.18	1.24
1	B	60	ALA	C-O	-5.11	1.17	1.24
1	D	229	GLN	C-O	-5.11	1.17	1.23
1	C	75	GLN	C-O	-5.09	1.18	1.24
1	C	154	ASP	C-O	-5.09	1.18	1.24
1	C	503	LYS	C-O	-5.09	1.18	1.24
1	B	376	ARG	C-O	-5.09	1.18	1.24
1	C	279	SER	C-O	-5.09	1.17	1.23
1	B	34	HIS	C-O	-5.09	1.17	1.23
1	B	176	SER	C-O	-5.09	1.17	1.24
1	C	479	ILE	C-O	-5.09	1.18	1.24
1	A	318	ALA	C-O	-5.09	1.18	1.24
1	D	30	GLY	C-O	-5.08	1.18	1.24
1	C	276	GLY	C-O	-5.08	1.17	1.24
1	A	446	TYR	C-O	-5.08	1.18	1.24
1	C	77	LEU	C-O	-5.07	1.18	1.24
1	C	233	THR	C-O	-5.07	1.18	1.24
1	B	183	LYS	C-O	-5.07	1.17	1.23
1	C	237	ALA	C-O	-5.07	1.18	1.24
1	D	60	ALA	C-O	-5.06	1.17	1.24
1	C	69	ARG	C-O	-5.06	1.18	1.24
1	B	436	PHE	C-O	-5.06	1.18	1.24
1	B	99	ALA	C-O	-5.05	1.18	1.24
1	B	228	TRP	C-O	-5.05	1.18	1.23
1	C	343	GLU	C-O	-5.05	1.18	1.24
1	B	226	ALA	C-O	-5.05	1.17	1.23
1	D	62	SER	C-O	-5.05	1.17	1.23
1	B	281	ARG	C-O	-5.04	1.18	1.24
1	A	236	LEU	C-O	-5.04	1.18	1.24
1	C	177	ARG	C-O	-5.04	1.17	1.24
1	A	447	ILE	C-O	-5.04	1.18	1.24
1	B	380	ALA	C-O	-5.04	1.18	1.24
1	D	347	LYS	C-O	-5.04	1.18	1.24
1	B	324	PHE	C-O	-5.03	1.17	1.24
1	C	396	VAL	C-O	-5.03	1.18	1.24
1	A	232	GLU	C-O	-5.03	1.18	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	316	THR	C-O	-5.03	1.18	1.24
1	A	140	VAL	C-O	-5.02	1.18	1.24
1	C	49	VAL	C-O	-5.02	1.18	1.24
1	D	272	TYR	C-O	-5.01	1.18	1.24
1	D	324	PHE	C-O	-5.01	1.17	1.24
1	C	348	TRP	C-O	-5.01	1.18	1.24
1	A	328	GLU	C-O	-5.01	1.17	1.23
1	A	36	ASN	C-O	-5.00	1.18	1.24
1	A	416	MET	C-O	-5.00	1.18	1.24
1	D	269	LEU	C-O	-5.00	1.17	1.24
1	A	192	GLN	C-O	-5.00	1.18	1.24

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	231	GLY	N-CA-C	10.85	124.44	112.29
1	B	231	GLY	N-CA-C	10.59	122.01	111.95
1	C	231	GLY	N-CA-C	9.99	123.48	112.29
1	D	231	GLY	N-CA-C	9.31	122.72	112.29
1	D	326	ARG	O-C-N	-8.24	113.10	123.24
1	B	251	ASN	N-CA-C	7.55	119.59	111.36
1	A	52	VAL	N-CA-C	7.23	118.90	108.48
1	C	251	ASN	N-CA-C	6.98	118.97	111.36
1	C	52	VAL	N-CA-C	6.88	118.39	108.48
1	D	253	GLU	N-CA-C	6.84	118.53	111.14
1	C	43	VAL	N-CA-C	6.67	117.46	110.72
1	C	33	LEU	N-CA-C	6.61	118.57	111.36
1	B	294	VAL	N-CA-C	6.59	117.38	110.72
1	A	43	VAL	N-CA-C	6.50	117.29	110.72
1	D	33	LEU	N-CA-C	6.30	118.22	111.36
1	A	251	ASN	N-CA-C	6.05	117.96	111.36
1	A	108	GLU	N-CA-C	6.04	118.66	111.71
1	C	426	GLN	N-CA-C	-6.04	100.72	109.59
1	D	440	THR	N-CA-C	6.00	117.49	111.07
1	D	52	VAL	N-CA-C	5.95	117.33	108.46
1	A	367	GLN	N-CA-C	5.90	117.79	111.36
1	B	33	LEU	N-CA-C	5.88	117.77	111.36
1	A	285	TYR	N-CA-C	5.79	117.67	111.36
1	A	409	SER	N-CA-C	5.77	117.38	111.14
1	B	424	PHE	CB-CA-C	-5.74	109.97	116.63
1	A	425	GLN	CB-CA-C	-5.68	110.04	116.63
1	B	425	GLN	N-CA-C	5.67	116.40	108.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	52	VAL	N-CA-C	5.61	116.81	108.46
1	B	285	TYR	N-CA-C	5.57	117.43	111.36
1	B	483	ASP	N-CA-C	5.54	118.25	111.82
1	C	285	TYR	N-CA-C	5.53	117.39	111.36
1	D	285	TYR	N-CA-C	5.44	117.29	111.36
1	D	318	ALA	N-CA-C	5.34	117.19	111.36
1	B	426	GLN	N-CA-C	5.22	117.41	108.90
1	B	253	GLU	N-CA-C	5.20	116.81	111.03
1	D	423	PHE	N-CA-C	5.17	118.84	111.92
1	A	424	PHE	N-CA-C	5.13	116.74	108.08
1	B	217	GLY	N-CA-C	5.11	123.10	115.64
1	A	427	PHE	N-CA-C	5.11	121.67	110.80
1	A	194	GLU	N-CA-C	-5.08	100.22	108.41

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	28	ARG	Peptide
1	C	28	ARG	Peptide

5.2 Too-close contacts [i](#)

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	4015	0	3944	54	0
1	B	4015	0	3944	47	0
1	C	4015	0	3944	48	1
1	D	4015	0	3944	65	0
2	A	28	0	22	1	0
2	B	28	0	22	1	0
2	C	28	0	22	0	0
2	D	28	0	22	0	0
3	A	461	0	0	11	1
3	B	322	0	0	8	0
3	C	434	0	0	11	2
3	D	337	0	0	5	2

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	17726	0	15864	212	3

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 7.

All (212) 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:189:SER:OG	1:A:192:GLN:HG3	1.18	1.27
1:C:254:ARG:HG3	1:C:254:ARG:HH11	1.24	1.02
1:A:199:GLU:O	3:A:1201:HOH:O	1.83	0.96
1:A:296:ARG:HH11	1:A:296:ARG:HG2	1.31	0.94
1:D:506:TYR:O	1:D:511:ARG:NH2	2.02	0.93
1:A:189:SER:OG	1:A:192:GLN:CG	2.15	0.92
1:D:420:CYS:SG	3:D:1239:HOH:O	2.28	0.91
1:C:502:MET:HE1	1:C:506:TYR:HE1	1.36	0.90
1:A:439:ARG:NH2	3:A:1339:HOH:O	2.01	0.90
1:C:343:GLU:OE2	3:C:1360:HOH:O	1.88	0.90
1:A:506:TYR:O	1:A:511:ARG:NH2	2.04	0.89
1:A:189:SER:HG	1:A:192:GLN:HG3	1.36	0.88
1:B:296:ARG:HH11	1:B:296:ARG:HG2	1.40	0.86
1:B:503:LYS:HA	1:B:503:LYS:HE2	1.61	0.83
1:C:347:LYS:NZ	3:C:1279:HOH:O	2.12	0.82
1:D:435:GLY:O	1:D:439:ARG:HG3	1.79	0.82
1:D:423:PHE:O	3:D:1104:HOH:O	1.97	0.81
1:B:506:TYR:O	1:B:511:ARG:NH2	2.14	0.81
1:D:417:TRP:CZ3	1:D:428:PHE:HE2	2.01	0.79
1:D:503:LYS:HE2	1:D:507:GLN:NE2	1.99	0.78
1:D:507:GLN:C	1:D:511:ARG:HH22	1.91	0.78
1:A:334:ILE:HD11	1:A:424:PHE:HE1	1.49	0.78
1:B:150:LEU:N	1:B:312:GLU:OE2	2.15	0.77
1:B:133:LYS:O	1:B:137:GLU:HG3	1.85	0.77
1:C:506:TYR:O	1:C:511:ARG:NH2	2.18	0.76
1:B:182:LYS:NZ	3:B:1205:HOH:O	1.75	0.76
1:C:502:MET:HE1	1:C:506:TYR:CE1	2.21	0.75
1:A:506:TYR:HA	1:A:509:LEU:HD22	1.70	0.74
1:A:296:ARG:HG2	1:A:296:ARG:NH1	2.00	0.73
1:A:427:PHE:O	1:A:428:PHE:CD2	2.43	0.72
1:B:435:GLY:O	1:B:439:ARG:HG3	1.89	0.72
1:D:417:TRP:CH2	1:D:428:PHE:HE2	2.07	0.72
1:D:426:GLN:HG2	1:D:427:PHE:H	1.56	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:104:ARG:O	1:A:108:GLU:HG2	1.92	0.69
1:D:502:MET:HE1	1:D:506:TYR:HE1	1.56	0.69
1:C:347:LYS:NZ	3:C:1321:HOH:O	2.26	0.69
1:A:334:ILE:CD1	1:A:424:PHE:HE1	2.05	0.69
1:B:44:ARG:NH1	3:B:1249:HOH:O	2.10	0.66
1:B:502:MET:HE1	1:B:506:TYR:CE1	2.31	0.65
1:D:191:GLN:O	1:D:195:SER:OG	2.14	0.65
1:B:169:LYS:NZ	3:B:1288:HOH:O	2.29	0.65
1:D:155:ARG:NH1	3:D:1226:HOH:O	2.22	0.65
1:A:511:ARG:NH1	3:A:1388:HOH:O	2.30	0.64
1:B:486:ARG:HB2	1:B:487:PRO:HD2	1.78	0.64
1:D:502:MET:HE1	1:D:506:TYR:CE1	2.32	0.64
1:C:254:ARG:HG3	1:C:254:ARG:NH1	1.97	0.64
1:B:274:ARG:NH1	1:B:406:ALA:O	2.32	0.63
1:A:323:ARG:HH12	1:A:326:ARG:NH1	1.97	0.62
1:A:150:LEU:N	1:A:312:GLU:OE2	2.25	0.62
1:D:274:ARG:NH1	1:D:406:ALA:O	2.33	0.62
1:B:296:ARG:HH11	1:B:296:ARG:CG	2.10	0.62
1:D:52:VAL:HG12	1:D:92:PHE:HB2	1.81	0.61
1:B:419:SER:O	1:B:420:CYS:HB2	1.99	0.61
1:C:52:VAL:HG12	1:C:92:PHE:HB2	1.83	0.60
1:D:426:GLN:OE1	1:D:426:GLN:N	2.25	0.60
1:A:165:PRO:HB3	1:A:170:ARG:HG3	1.83	0.60
1:B:274:ARG:CZ	1:B:406:ALA:O	2.50	0.60
1:A:334:ILE:HD11	1:A:424:PHE:CE1	2.33	0.60
1:A:123:PHE:HB2	3:A:1346:HOH:O	2.02	0.60
1:D:203:ASN:OD1	1:D:203:ASN:N	2.31	0.59
1:B:323:ARG:CZ	1:B:326:ARG:HD2	2.32	0.59
1:A:500:GLU:OE1	3:A:1397:HOH:O	2.16	0.59
1:C:253:GLU:O	1:C:254:ARG:HD3	2.03	0.59
1:D:377:HIS:HA	1:D:418:LEU:CD1	2.33	0.59
1:C:426:GLN:H	1:C:426:GLN:CD	2.11	0.59
1:B:504:GLN:NE2	3:B:1198:HOH:O	2.36	0.58
1:A:249:VAL:HA	1:A:253:GLU:HG3	1.86	0.58
1:B:64:SER:OG	1:B:397:ARG:NH2	2.36	0.58
1:A:340:ARG:NH2	3:A:1166:HOH:O	2.36	0.57
1:C:106:PHE:HA	1:C:111:VAL:HG13	1.86	0.57
1:D:274:ARG:CZ	1:D:406:ALA:O	2.52	0.57
1:C:177:ARG:NH2	3:C:1209:HOH:O	2.29	0.57
1:A:106:PHE:HA	1:A:111:VAL:HG13	1.87	0.56
1:D:85:ARG:NH2	1:D:90:ARG:NH1	2.53	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:506:TYR:HA	1:C:509:LEU:HD22	1.88	0.56
1:D:323:ARG:NH1	1:D:326:ARG:CD	2.69	0.56
1:D:426:GLN:H	1:D:426:GLN:CD	2.09	0.56
1:B:512:TYR:O	3:B:1229:HOH:O	2.18	0.56
1:D:427:PHE:CD2	1:D:428:PHE:N	2.69	0.56
1:D:85:ARG:NH2	1:D:90:ARG:CZ	2.69	0.55
1:C:502:MET:CE	1:C:506:TYR:CE1	2.89	0.55
1:B:40:LEU:HB3	1:B:44:ARG:HH12	1.72	0.55
1:C:189:SER:HB3	1:C:192:GLN:HG3	1.88	0.55
1:B:71:ARG:HD3	1:B:213:LEU:HD11	1.89	0.55
1:C:191:GLN:NE2	1:C:194:GLU:OE1	2.40	0.54
1:B:423:PHE:O	1:B:424:PHE:HB2	2.08	0.53
1:D:323:ARG:CZ	1:D:326:ARG:CD	2.86	0.53
1:D:323:ARG:NH1	1:D:326:ARG:NE	2.57	0.53
1:C:179:GLU:H	1:C:179:GLU:CD	2.17	0.53
1:A:182:LYS:NZ	3:A:1376:HOH:O	2.31	0.53
1:B:502:MET:HE1	1:B:506:TYR:HE1	1.72	0.53
1:C:426:GLN:O	1:C:426:GLN:NE2	2.42	0.53
1:C:187:ALA:HB1	3:C:1266:HOH:O	2.09	0.52
1:C:177:ARG:NE	3:C:1209:HOH:O	2.31	0.52
1:D:417:TRP:CH2	1:D:428:PHE:CE2	2.93	0.52
1:C:50:ARG:HH21	1:C:109:TRP:HB3	1.75	0.52
1:D:507:GLN:CA	1:D:511:ARG:HH22	2.23	0.52
1:A:347:LYS:HE2	3:A:1113:HOH:O	2.09	0.52
1:B:484:TYR:CD2	1:B:485:PRO:HD2	2.44	0.51
1:C:427:PHE:HE1	1:C:429:HIS:CG	2.29	0.51
1:D:504:GLN:NE2	1:D:507:GLN:OE1	2.44	0.51
1:D:428:PHE:C	1:D:428:PHE:CD1	2.88	0.51
1:A:190:SER:O	1:A:194:GLU:HG3	2.11	0.51
1:D:187:ALA:HA	3:D:1183:HOH:O	2.11	0.51
1:A:502:MET:HE1	1:A:506:TYR:HE1	1.76	0.50
1:B:132:MET:HG2	1:B:142:VAL:HG11	1.93	0.50
1:B:450:TYR:C	1:B:452:PRO:HD3	2.37	0.50
1:A:71:ARG:HD2	3:A:1434:HOH:O	2.12	0.50
1:D:43:VAL:HG12	1:D:43:VAL:O	2.11	0.50
1:D:428:PHE:C	1:D:428:PHE:HD1	2.19	0.50
1:D:377:HIS:ND1	1:D:418:LEU:HD12	2.26	0.50
1:C:438:ARG:HD3	3:C:1174:HOH:O	2.11	0.50
1:A:484:TYR:CD1	1:A:485:PRO:HD2	2.47	0.50
1:C:99:ALA:O	1:C:134:MET:HE1	2.11	0.50
1:A:424:PHE:O	1:A:425:GLN:HB2	2.10	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:427:PHE:H	1:D:427:PHE:HD2	1.59	0.49
1:C:189:SER:CB	1:C:192:GLN:HG3	2.42	0.49
1:A:194:GLU:O	1:A:195:SER:CB	2.60	0.49
1:D:249:VAL:HA	1:D:253:GLU:HB2	1.95	0.49
1:D:323:ARG:CZ	1:D:326:ARG:HD3	2.43	0.49
1:B:296:ARG:HG2	1:B:296:ARG:NH1	2.19	0.49
1:D:507:GLN:C	1:D:511:ARG:NH2	2.68	0.49
1:B:165:PRO:HB3	1:B:170:ARG:HG3	1.95	0.49
1:C:87:LEU:HD13	1:C:193:MET:HG2	1.94	0.48
1:D:377:HIS:HA	1:D:418:LEU:HD11	1.94	0.48
1:D:270:SER:OG	1:D:271:PRO:HD3	2.14	0.48
1:D:418:LEU:C	1:D:418:LEU:HD23	2.38	0.48
1:A:187:ALA:HB2	3:A:1300:HOH:O	2.13	0.48
1:C:426:GLN:CD	1:C:426:GLN:N	2.72	0.48
1:C:484:TYR:CD2	1:C:485:PRO:HD2	2.49	0.48
1:A:99:ALA:O	1:A:134:MET:HE1	2.14	0.48
1:B:354:GLY:O	1:B:356:PRO:HD3	2.14	0.48
1:C:133:LYS:HE2	1:C:137:GLU:OE2	2.14	0.47
1:C:183:LYS:NZ	3:C:1260:HOH:O	2.27	0.47
1:A:377:HIS:CE1	2:A:900:2CX:H1	2.49	0.47
1:A:270:SER:OG	1:A:271:PRO:HD3	2.14	0.47
1:B:296:ARG:CG	1:B:296:ARG:NH1	2.73	0.47
1:D:246:LYS:HB2	1:D:246:LYS:HE2	1.63	0.47
1:B:503:LYS:HA	1:B:503:LYS:CE	2.36	0.47
1:D:295:LYS:O	1:D:298:SER:OG	2.33	0.47
1:B:270:SER:OG	1:B:271:PRO:HD3	2.15	0.47
1:D:427:PHE:CG	1:D:428:PHE:N	2.83	0.47
1:D:450:TYR:C	1:D:452:PRO:HD3	2.39	0.47
1:D:503:LYS:HE2	1:D:507:GLN:HE22	1.80	0.46
1:C:150:LEU:N	1:C:312:GLU:OE2	2.35	0.46
1:A:295:LYS:O	1:A:296:ARG:C	2.57	0.46
1:D:40:LEU:O	1:D:44:ARG:HG3	2.15	0.46
1:A:87:LEU:HD22	1:A:190:SER:HA	1.97	0.46
1:A:502:MET:HE1	1:A:506:TYR:CE1	2.50	0.46
1:D:417:TRP:CZ3	1:D:428:PHE:CE2	2.92	0.45
1:C:241:LYS:HB3	1:C:267:THR:HG22	1.98	0.45
1:C:312:GLU:HG2	3:C:1036:HOH:O	2.17	0.45
1:B:136:LYS:HG3	1:B:137:GLU:N	2.31	0.45
1:C:179:GLU:OE1	3:C:1249:HOH:O	2.21	0.45
1:D:254:ARG:HH11	1:D:254:ARG:CG	2.30	0.45
1:C:189:SER:HB3	1:C:192:GLN:H	1.82	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:381:CYS:HG	1:A:430:CYS:HG	1.63	0.44
1:B:105:LEU:HD13	1:B:109:TRP:CZ3	2.53	0.44
1:B:425:GLN:NE2	3:B:1227:HOH:O	2.51	0.44
1:B:380:ALA:HB1	1:B:418:LEU:HD22	2.00	0.44
1:C:55:LEU:O	1:C:57:PRO:HD3	2.18	0.44
1:C:270:SER:OG	1:C:271:PRO:HD3	2.17	0.44
1:A:347:LYS:HD3	1:A:488:ILE:HG21	1.99	0.44
1:D:484:TYR:CG	1:D:485:PRO:HD2	2.52	0.44
1:C:71:ARG:HB2	1:C:213:LEU:HD11	2.00	0.44
1:B:247:ALA:O	1:B:251:ASN:HB2	2.17	0.44
1:D:417:TRP:CE2	1:D:425:GLN:NE2	2.87	0.43
1:D:162:GLN:HA	1:D:162:GLN:OE1	2.19	0.43
1:B:43:VAL:O	1:B:43:VAL:HG12	2.19	0.43
1:D:87:LEU:HD13	1:D:193:MET:CB	2.49	0.43
1:B:410:VAL:CG1	2:B:900:2CX:H19	2.48	0.43
1:C:78:GLU:CD	3:C:1339:HOH:O	2.61	0.43
1:A:299:THR:HG21	1:B:62:SER:OG	2.18	0.43
1:D:55:LEU:C	1:D:55:LEU:HD23	2.44	0.43
1:D:155:ARG:HG2	1:D:155:ARG:HH11	1.84	0.42
1:A:199:GLU:H	1:A:199:GLU:HG2	1.70	0.42
1:C:55:LEU:C	1:C:55:LEU:HD23	2.44	0.42
1:D:270:SER:N	1:D:271:PRO:CD	2.82	0.42
1:C:427:PHE:CE1	1:C:429:HIS:HA	2.55	0.42
1:D:484:TYR:CD1	1:D:485:PRO:HD2	2.54	0.42
1:C:62:SER:OG	1:D:299:THR:HG21	2.19	0.42
1:C:295:LYS:O	1:C:296:ARG:C	2.63	0.42
1:A:35:ASP:HB3	1:A:279:SER:HB2	2.01	0.42
1:B:484:TYR:CG	1:B:485:PRO:HD2	2.55	0.42
1:D:507:GLN:HA	1:D:511:ARG:HH22	1.83	0.42
1:D:426:GLN:N	1:D:426:GLN:CD	2.74	0.42
1:A:509:LEU:HB2	1:A:511:ARG:HH21	1.85	0.41
1:A:380:ALA:HB1	1:A:418:LEU:HD22	2.02	0.41
1:B:477:LYS:NZ	3:B:1295:HOH:O	2.53	0.41
1:A:160:ASN:OD1	1:A:170:ARG:NH1	2.52	0.41
1:B:249:VAL:HA	1:B:253:GLU:HB2	2.02	0.41
1:D:494:THR:HG23	1:D:495:SER:N	2.34	0.41
1:A:385:ARG:NH2	1:A:428:PHE:CZ	2.89	0.41
1:A:509:LEU:O	1:A:511:ARG:NH2	2.54	0.41
1:C:149:THR:HB	1:C:312:GLU:OE2	2.21	0.41
1:D:503:LYS:HE2	1:D:507:GLN:HE21	1.78	0.41
1:C:503:LYS:HD2	1:C:503:LYS:HA	1.89	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:323:ARG:CZ	1:A:326:ARG:HD2	2.50	0.41
1:A:296:ARG:NH1	1:A:296:ARG:CG	2.74	0.41
1:A:470:GLU:CD	1:A:474:LYS:HZ1	2.27	0.41
1:B:298:SER:N	3:B:1244:HOH:O	2.54	0.41
1:B:323:ARG:NH1	1:B:326:ARG:HD2	2.36	0.41
1:D:177:ARG:NE	3:D:1261:HOH:O	2.44	0.41
1:C:245:ARG:O	1:C:249:VAL:HG23	2.21	0.41
1:A:380:ALA:CB	1:A:418:LEU:HD22	2.50	0.41
1:D:55:LEU:O	1:D:57:PRO:HD3	2.21	0.40
1:A:293:LYS:NZ	3:A:1394:HOH:O	2.54	0.40
1:B:34:HIS:O	1:B:35:ASP:C	2.64	0.40
1:B:52:VAL:HG12	1:B:92:PHE:HB2	2.03	0.40
1:C:102:PHE:N	1:C:103:PRO:CD	2.84	0.40
1:A:87:LEU:O	1:A:194:GLU:HG2	2.22	0.40
1:D:65:VAL:HB	1:D:70:TRP:HE1	1.86	0.40

All (3) 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)
3:C:1335:HOH:O	3:A:1352:HOH:O[1_554]	1.50	0.70
3:C:1414:HOH:O	3:D:1289:HOH:O[1_565]	2.18	0.02
1:C:177:ARG:NH1	3:D:1287:HOH:O[1_565]	2.19	0.01

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	490/512 (96%)	474 (97%)	14 (3%)	2 (0%)	30	22
1	B	490/512 (96%)	480 (98%)	10 (2%)	0	100	100
1	C	490/512 (96%)	480 (98%)	10 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	490/512 (96%)	478 (98%)	10 (2%)	2 (0%)	30	22
All	All	1960/2048 (96%)	1912 (98%)	44 (2%)	4 (0%)	43	37

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	426	GLN
1	A	195	SER
1	A	428	PHE
1	D	429	HIS

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	424/434 (98%)	407 (96%)	17 (4%)	28	14
1	B	424/434 (98%)	410 (97%)	14 (3%)	33	20
1	C	424/434 (98%)	408 (96%)	16 (4%)	29	15
1	D	424/434 (98%)	405 (96%)	19 (4%)	24	11
All	All	1696/1736 (98%)	1630 (96%)	66 (4%)	28	15

All (66) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	88	ASN
1	C	111	VAL
1	C	126	GLU
1	C	179	GLU
1	C	190	SER
1	C	193	MET
1	C	203	ASN
1	C	207	THR
1	C	254	ARG
1	C	296	ARG

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Mol	Chain	Res	Type
1	C	298	SER
1	C	299	THR
1	C	418	LEU
1	C	497	LEU
1	C	501	ARG
1	C	509	LEU
1	A	52	VAL
1	A	111	VAL
1	A	125	LYS
1	A	136	LYS
1	A	190	SER
1	A	194	GLU
1	A	199	GLU
1	A	214	GLU
1	A	253	GLU
1	A	254	ARG
1	A	295	LYS
1	A	299	THR
1	A	418	LEU
1	A	426	GLN
1	A	493	GLU
1	A	501	ARG
1	A	509	LEU
1	B	52	VAL
1	B	89	SER
1	B	104	ARG
1	B	105	LEU
1	B	162	GLN
1	B	202	GLU
1	B	256	ARG
1	B	296	ARG
1	B	403	LEU
1	B	460	ARG
1	B	470	GLU
1	B	474	LYS
1	B	503	LYS
1	B	509	LEU
1	D	52	VAL
1	D	88	ASN
1	D	105	LEU
1	D	136	LYS
1	D	186	VAL

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Mol	Chain	Res	Type
1	D	191	GLN
1	D	194	GLU
1	D	195	SER
1	D	254	ARG
1	D	298	SER
1	D	299	THR
1	D	397	ARG
1	D	403	LEU
1	D	427	PHE
1	D	429	HIS
1	D	474	LYS
1	D	501	ARG
1	D	505	ILE
1	D	509	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	297	ASN
1	A	507	GLN
1	B	201	GLN
1	B	297	ASN
1	B	307	GLN
1	B	425	GLN
1	B	504	GLN
1	D	88	ASN
1	D	297	ASN
1	D	307	GLN
1	D	373	HIS
1	D	504	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 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	2CX	A	900	-	31,31,31	2.37	13 (41%)	41,45,45	2.47	8 (19%)
2	2CX	C	900	-	31,31,31	2.33	13 (41%)	41,45,45	2.55	7 (17%)
2	2CX	D	900	-	31,31,31	2.32	14 (45%)	41,45,45	2.64	8 (19%)
2	2CX	B	900	-	31,31,31	2.29	14 (45%)	41,45,45	2.75	9 (21%)

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	2CX	A	900	-	-	5/18/18/18	0/4/4/4
2	2CX	C	900	-	-	5/18/18/18	0/4/4/4
2	2CX	D	900	-	-	4/18/18/18	0/4/4/4
2	2CX	B	900	-	-	4/18/18/18	0/4/4/4

All (54) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	900	2CX	C13-C14	5.86	1.48	1.40
2	B	900	2CX	C13-C14	5.73	1.48	1.40
2	A	900	2CX	O28-S02	5.66	1.52	1.43
2	C	900	2CX	C13-C14	5.64	1.48	1.40
2	A	900	2CX	C13-C14	5.61	1.48	1.40
2	C	900	2CX	O28-S02	5.56	1.51	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	900	2CX	O28-S02	4.85	1.50	1.43
2	B	900	2CX	O28-S02	4.80	1.50	1.43
2	A	900	2CX	S02-N03	3.95	1.70	1.63
2	C	900	2CX	S02-N03	3.91	1.70	1.63
2	A	900	2CX	C19-C20	3.81	1.45	1.38
2	D	900	2CX	S02-N03	3.57	1.69	1.63
2	A	900	2CX	C11-C10	3.47	1.44	1.38
2	D	900	2CX	C11-C10	3.44	1.44	1.38
2	B	900	2CX	C19-C20	3.41	1.44	1.38
2	B	900	2CX	C18-C17	3.39	1.44	1.38
2	B	900	2CX	S02-N03	3.38	1.69	1.63
2	D	900	2CX	C19-C20	3.34	1.44	1.38
2	C	900	2CX	C19-C20	3.26	1.44	1.38
2	D	900	2CX	C18-C17	3.24	1.44	1.38
2	B	900	2CX	C11-C10	3.15	1.44	1.38
2	D	900	2CX	C04-C05	3.11	1.60	1.52
2	C	900	2CX	C11-C10	3.08	1.44	1.38
2	C	900	2CX	C18-C17	3.03	1.44	1.38
2	C	900	2CX	C04-C05	2.96	1.59	1.52
2	A	900	2CX	C18-C17	2.94	1.43	1.38
2	A	900	2CX	C04-C05	2.84	1.59	1.52
2	B	900	2CX	C04-C05	2.72	1.59	1.52
2	D	900	2CX	C23-C22	2.53	1.41	1.34
2	C	900	2CX	C21-C22	2.53	1.53	1.48
2	B	900	2CX	C23-C22	2.50	1.41	1.34
2	D	900	2CX	C27-S02	2.49	1.81	1.75
2	C	900	2CX	C23-C22	2.48	1.41	1.34
2	B	900	2CX	C04-N03	-2.46	1.43	1.47
2	A	900	2CX	C17-C16	2.46	1.43	1.39
2	A	900	2CX	C23-C22	2.43	1.41	1.34
2	A	900	2CX	C21-C22	2.42	1.53	1.48
2	D	900	2CX	C21-C22	2.36	1.53	1.48
2	B	900	2CX	O01-S02	-2.36	1.40	1.43
2	B	900	2CX	C27-S02	2.32	1.80	1.75
2	C	900	2CX	C27-S02	2.30	1.80	1.75
2	D	900	2CX	C04-N03	-2.29	1.44	1.47
2	B	900	2CX	C21-C22	2.27	1.53	1.48
2	D	900	2CX	O01-S02	-2.25	1.40	1.43
2	A	900	2CX	C24-C25	2.23	1.40	1.32
2	A	900	2CX	C27-S02	2.23	1.80	1.75
2	A	900	2CX	C04-N03	-2.21	1.44	1.47
2	B	900	2CX	C21-N03	-2.21	1.44	1.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	900	2CX	C21-N03	-2.21	1.44	1.47
2	C	900	2CX	C24-C25	2.19	1.40	1.32
2	C	900	2CX	C04-N03	-2.19	1.44	1.47
2	D	900	2CX	C24-C25	2.15	1.40	1.32
2	C	900	2CX	C14-C15	2.12	1.51	1.46
2	B	900	2CX	C24-C25	2.08	1.40	1.32

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	900	2CX	O28-S02-O01	-12.16	100.28	118.46
2	D	900	2CX	O28-S02-O01	-11.48	101.29	118.46
2	A	900	2CX	O28-S02-O01	-11.32	101.53	118.46
2	B	900	2CX	O28-S02-O01	-11.19	101.72	118.46
2	B	900	2CX	C05-C04-N03	-7.58	102.91	113.24
2	A	900	2CX	O01-S02-N03	6.94	112.82	107.02
2	C	900	2CX	O01-S02-N03	6.13	112.13	107.02
2	B	900	2CX	O01-S02-N03	6.09	112.10	107.02
2	D	900	2CX	C05-C04-N03	-6.03	105.02	113.24
2	D	900	2CX	O28-S02-N03	5.86	111.91	107.02
2	D	900	2CX	O01-S02-N03	5.36	111.50	107.02
2	B	900	2CX	O28-S02-N03	5.04	111.22	107.02
2	A	900	2CX	C27-S02-N03	4.45	112.51	107.27
2	C	900	2CX	C27-S02-N03	4.32	112.35	107.27
2	C	900	2CX	O28-S02-N03	4.13	110.47	107.02
2	B	900	2CX	O28-S02-C27	3.05	112.48	108.30
2	D	900	2CX	O28-S02-C27	2.92	112.31	108.30
2	A	900	2CX	O28-S02-N03	2.72	109.29	107.02
2	C	900	2CX	C25-O26-C22	2.64	111.48	106.22
2	C	900	2CX	C16-N08-C09	2.48	111.03	108.43
2	A	900	2CX	C25-O26-C22	2.47	111.15	106.22
2	A	900	2CX	C16-N08-C09	2.45	110.99	108.43
2	D	900	2CX	C25-O26-C22	2.30	110.81	106.22
2	B	900	2CX	C27-S02-N03	2.30	109.98	107.27
2	B	900	2CX	O26-C22-C21	2.30	125.94	118.01
2	B	900	2CX	C17-C16-N08	2.29	132.26	129.08
2	D	900	2CX	O26-C22-C21	2.23	125.70	118.01
2	A	900	2CX	O28-S02-C27	2.18	111.29	108.30
2	D	900	2CX	C27-S02-N03	2.18	109.84	107.27
2	B	900	2CX	C25-O26-C22	2.17	110.55	106.22
2	C	900	2CX	O26-C22-C21	2.08	125.18	118.01
2	A	900	2CX	O26-C22-C21	2.01	124.93	118.01

There are no chirality outliers.

All (18) torsion outliers are listed below:

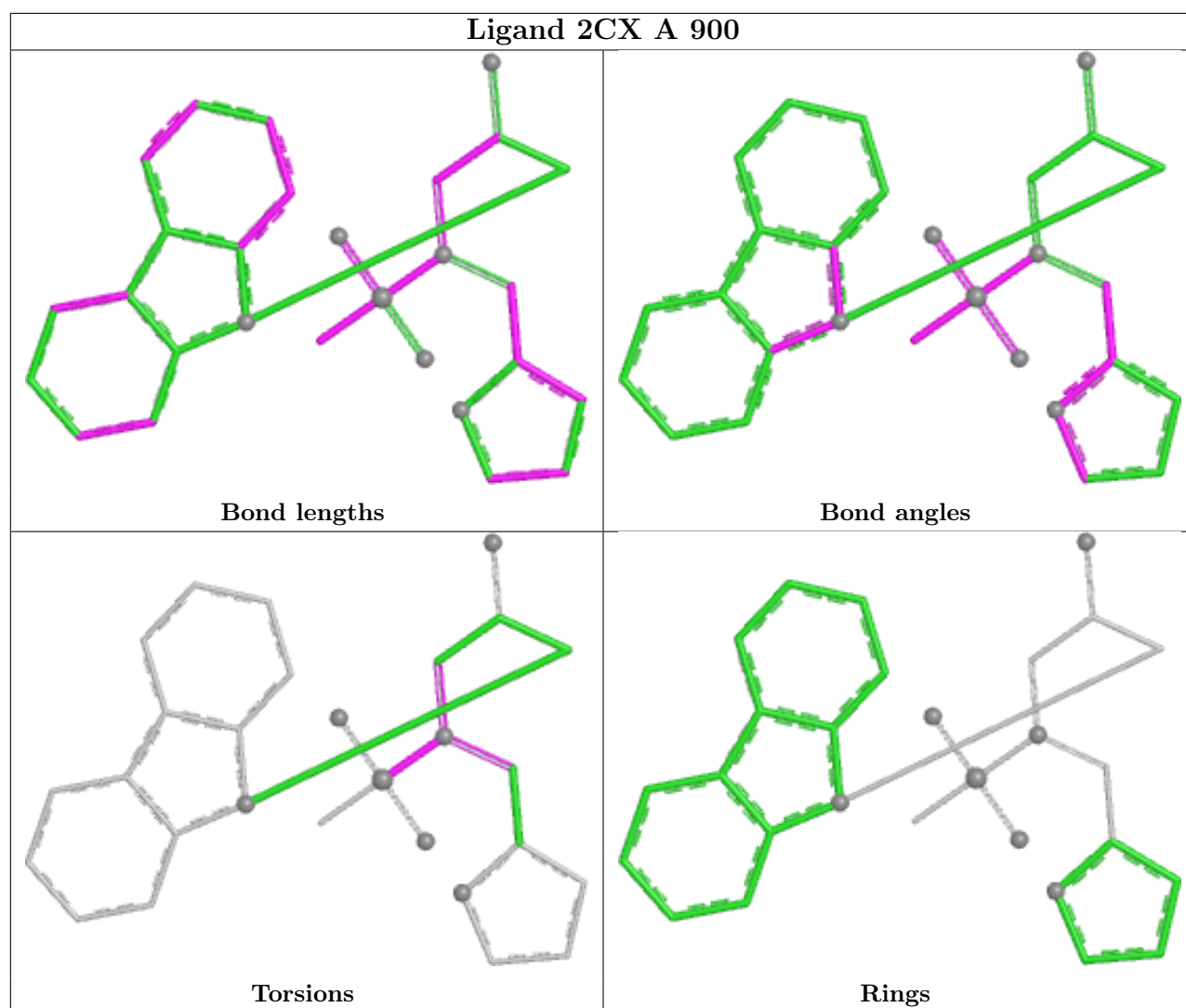
Mol	Chain	Res	Type	Atoms
2	C	900	2CX	C22-C21-N03-S02
2	C	900	2CX	C21-N03-S02-C27
2	C	900	2CX	C21-N03-S02-O28
2	C	900	2CX	C05-C04-N03-C21
2	A	900	2CX	C22-C21-N03-S02
2	A	900	2CX	C21-N03-S02-C27
2	A	900	2CX	C21-N03-S02-O28
2	A	900	2CX	C05-C04-N03-C21
2	B	900	2CX	C04-N03-S02-C27
2	B	900	2CX	C04-N03-S02-O28
2	D	900	2CX	C04-N03-S02-O28
2	B	900	2CX	N03-C21-C22-O26
2	D	900	2CX	N03-C21-C22-O26
2	A	900	2CX	C04-N03-S02-O28
2	C	900	2CX	C04-N03-S02-O28
2	B	900	2CX	C04-N03-S02-O01
2	D	900	2CX	C04-N03-S02-O01
2	D	900	2CX	C04-N03-S02-C27

There are no ring outliers.

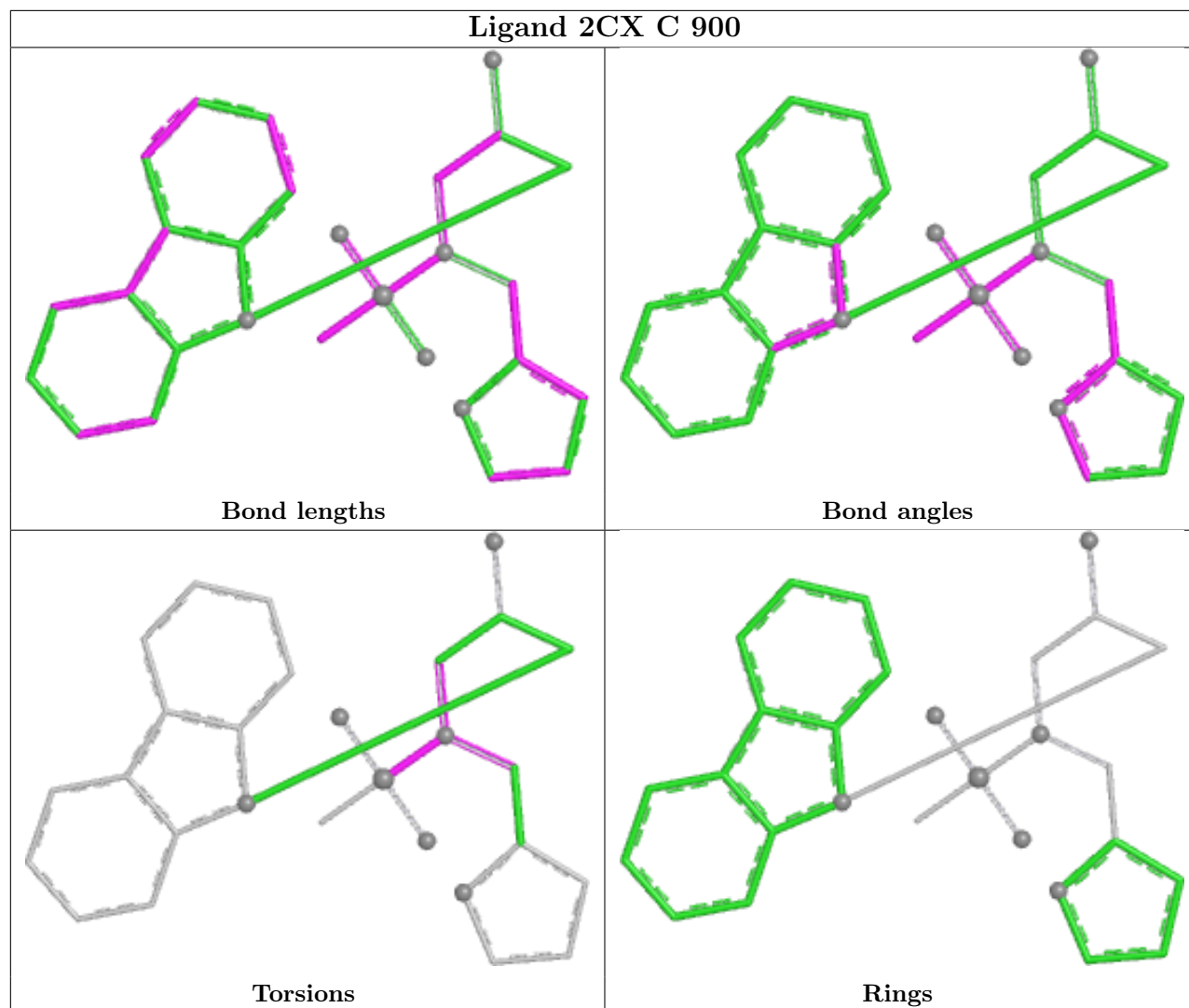
2 monomers are involved in 2 short contacts:

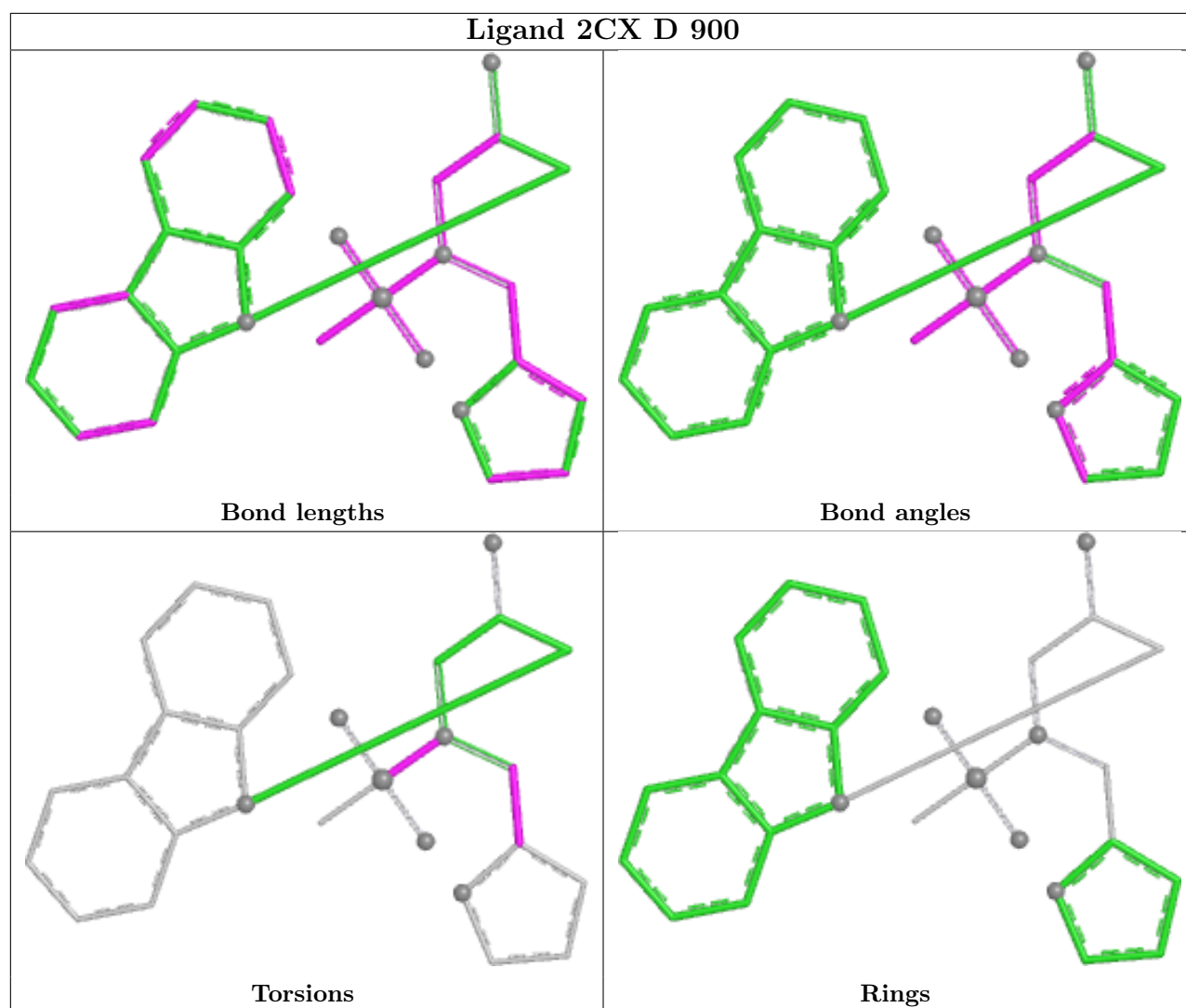
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	900	2CX	1	0
2	B	900	2CX	1	0

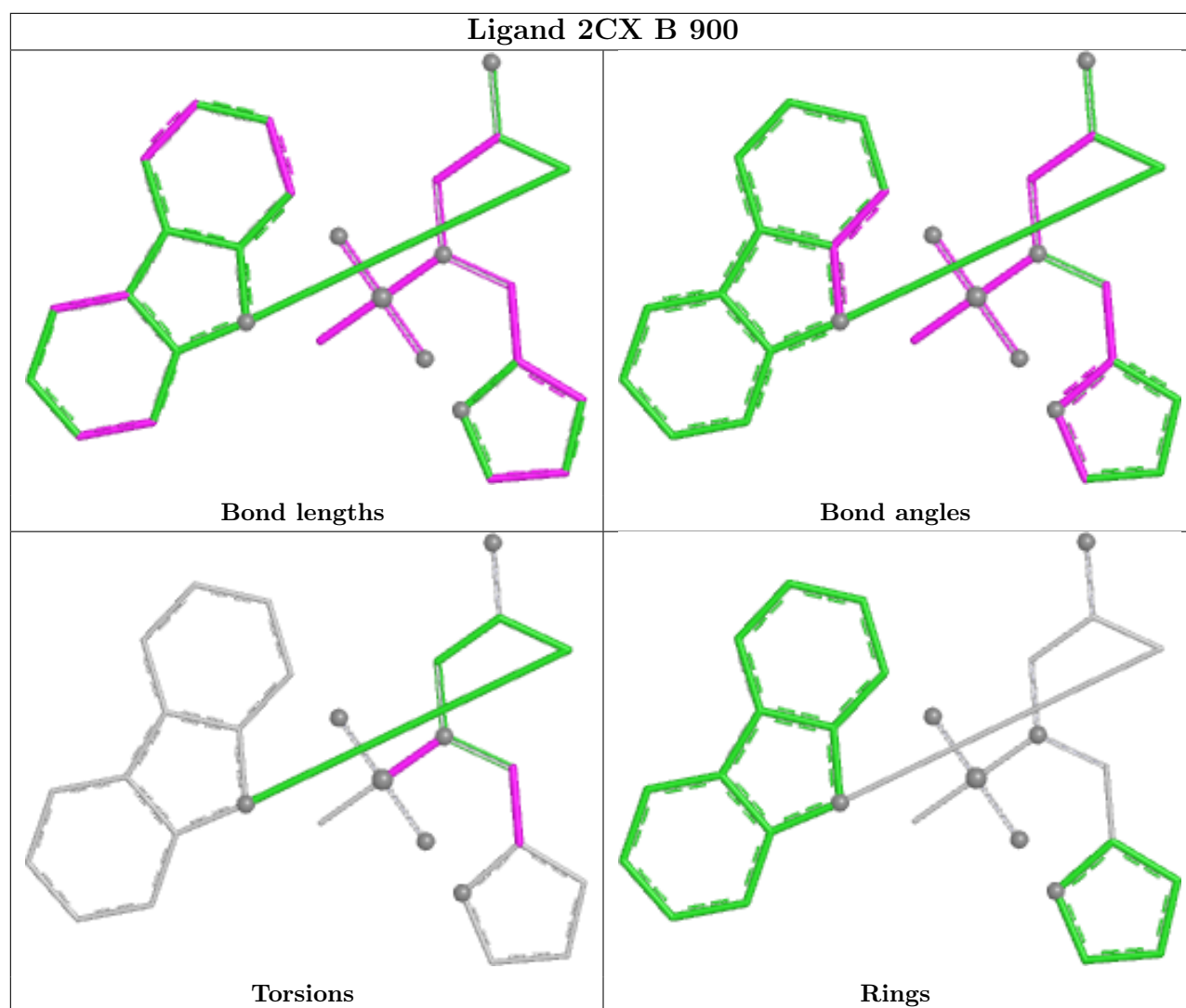
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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



Ligand 2CX C 900







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(Å²)	Q<0.9
1	A	492/512 (96%)	-1.03	0	100100	15, 26, 50, 74	0
1	B	492/512 (96%)	-0.89	0	100100	19, 33, 54, 75	0
1	C	492/512 (96%)	-1.03	0	100100	15, 26, 48, 71	0
1	D	492/512 (96%)	-0.89	1 (0%)	9194	19, 33, 54, 81	0
All	All	1968/2048 (96%)	-0.96	1 (0%)	9294	15, 29, 53, 81	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	427	PHE	2.4

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(Å ²)	Q<0.9
2	2CX	C	900	28/28	0.98	0.05	22,28,41,43	0

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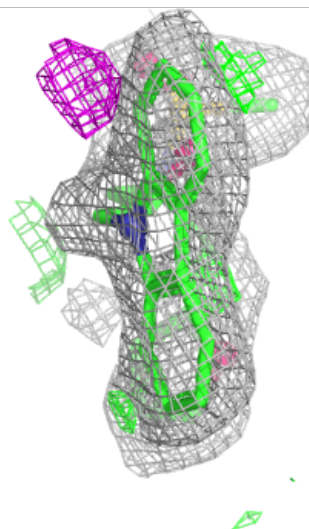
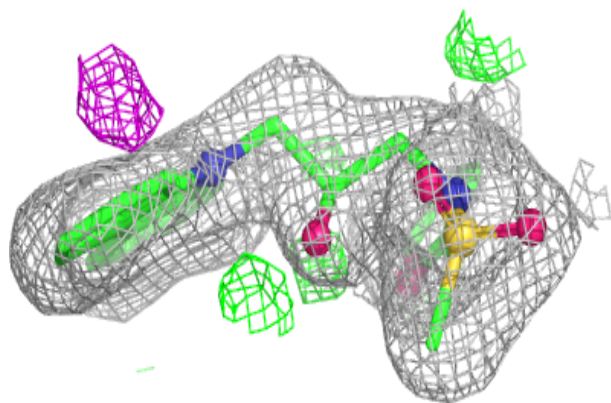
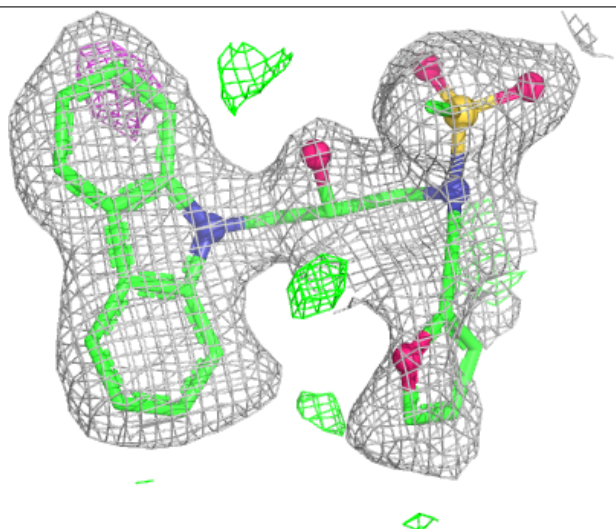
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	2CX	D	900	28/28	0.98	0.05	25,32,44,47	0
2	2CX	B	900	28/28	0.99	0.05	24,31,42,43	0
2	2CX	A	900	28/28	0.99	0.04	20,27,40,41	0

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.

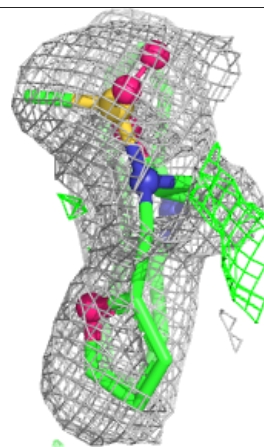
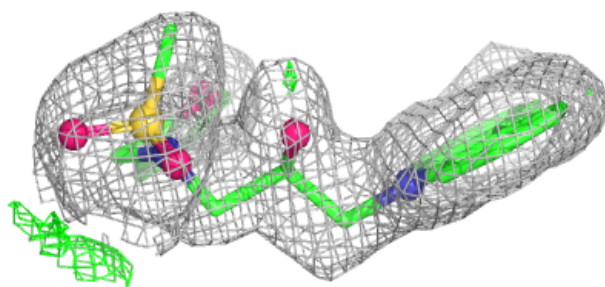
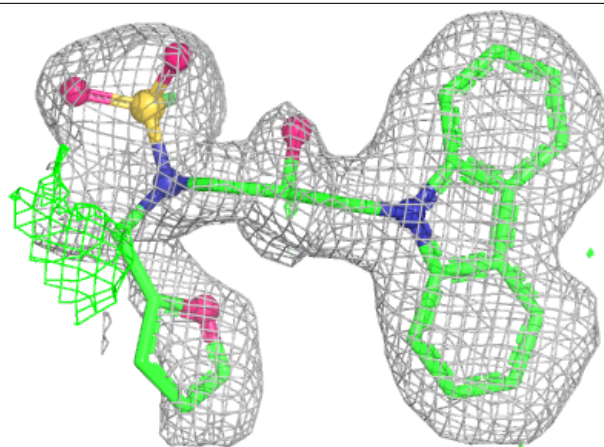
Electron density around 2CX C 900:

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



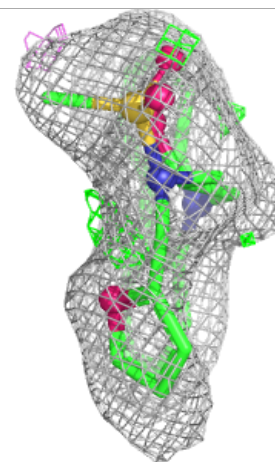
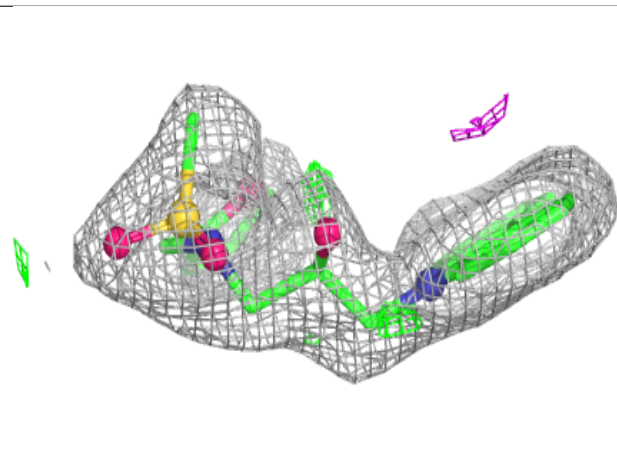
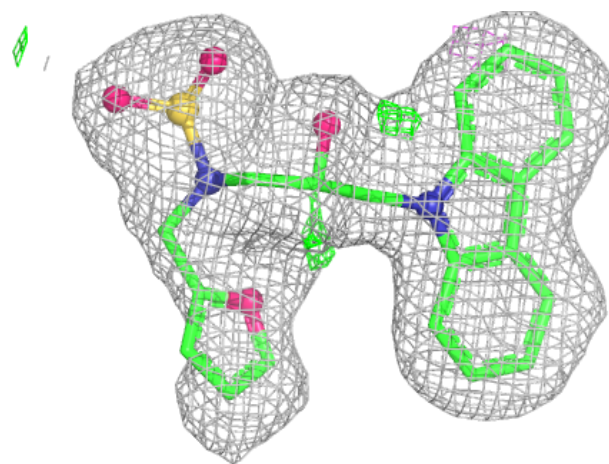
Electron density around 2CX D 900:

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



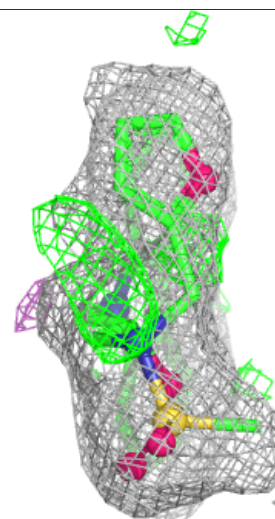
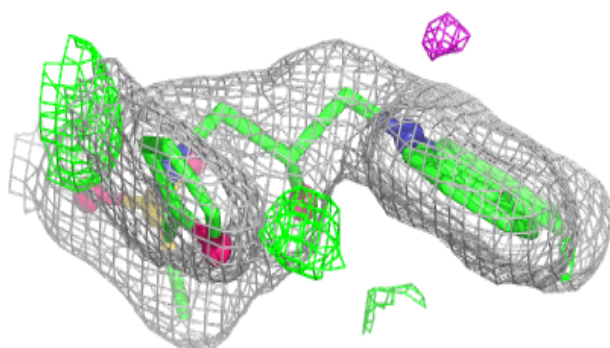
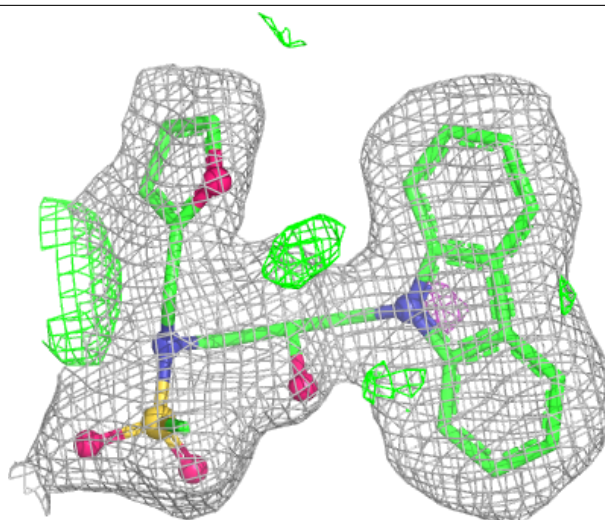
Electron density around 2CX B 900:

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



Electron density around 2CX A 900:

$2mF_o-DF_c$ (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.