



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 05:25 AM UTC

PDB ID : 8OXU / pdb_00008oxu
Title : Crystal Structure of the Hsp90-LA1011 Complex
Authors : Roe, S.M.; Prodromou, C.
Deposited on : 2023-05-02
Resolution : 2.94 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

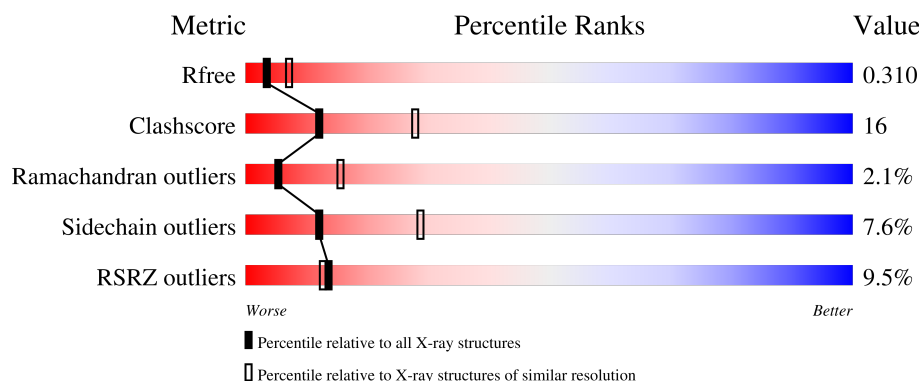
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

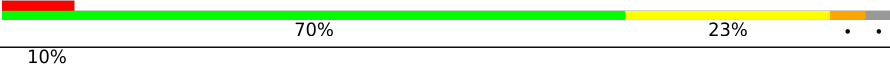
The reported resolution of this entry is 2.94 Å.

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	1159 (2.96-2.92)
Clashscore	190562	1184 (2.96-2.92)
Ramachandran outliers	187476	1131 (2.96-2.92)
Sidechain outliers	187428	1131 (2.96-2.92)
RSRZ outliers	180081	1159 (2.96-2.92)

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	241	
1	B	241	
1	C	241	
1	D	241	

2 Entry composition [i](#)

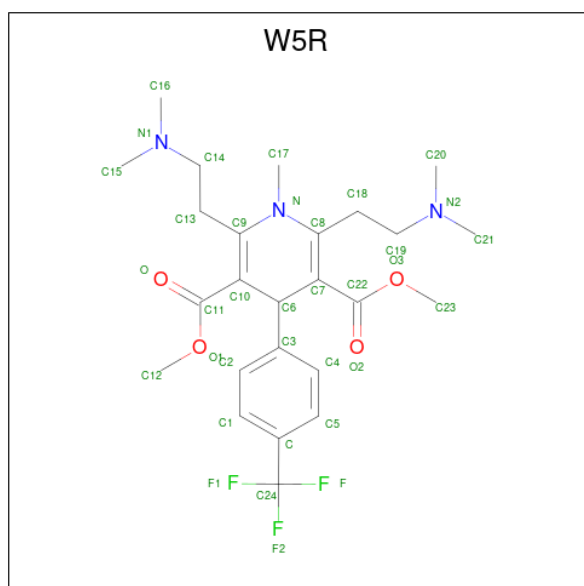
There are 3 unique types of molecules in this entry. The entry contains 6816 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 ATP-dependent molecular chaperone HSP82.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	231	Total	C	N	O	S	0	0	0
			1712	1102	270	336	4			
1	B	233	Total	C	N	O	S	0	0	0
			1692	1086	273	330	3			
1	C	234	Total	C	N	O	S	0	0	0
			1709	1090	272	343	4			
1	D	233	Total	C	N	O	S	0	0	0
			1667	1071	268	325	3			

- Molecule 2 is dimethyl 2,6-bis[2-(dimethylamino)ethyl]-1-methyl-4-[4-(trifluoromethyl)phenyl]-4 {H}-pyridine-3,5-dicarboxylate (CCD ID: W5R) (formula: C₂₅H₃₄F₃N₃O₄) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	C	1	Total	C	F	N	O	0	0
			35	25	3	3	4		

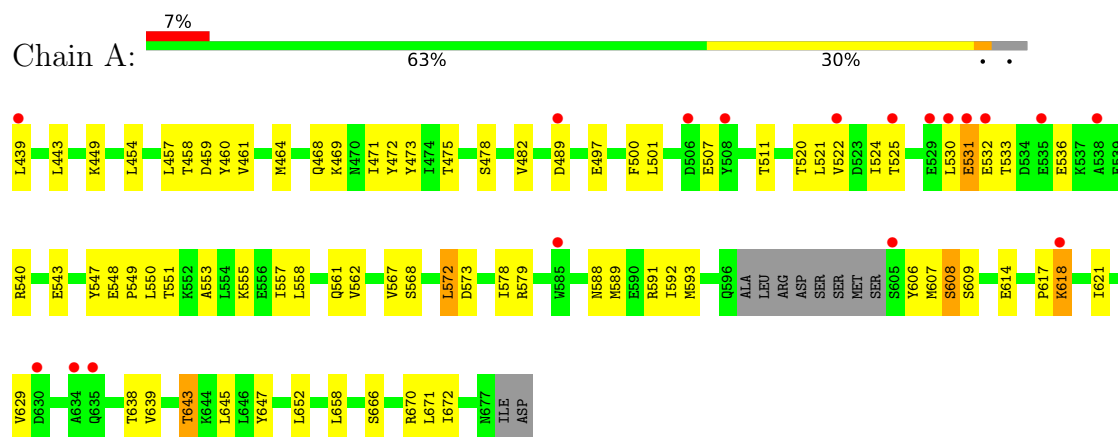
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	O	0	0
			1	1		

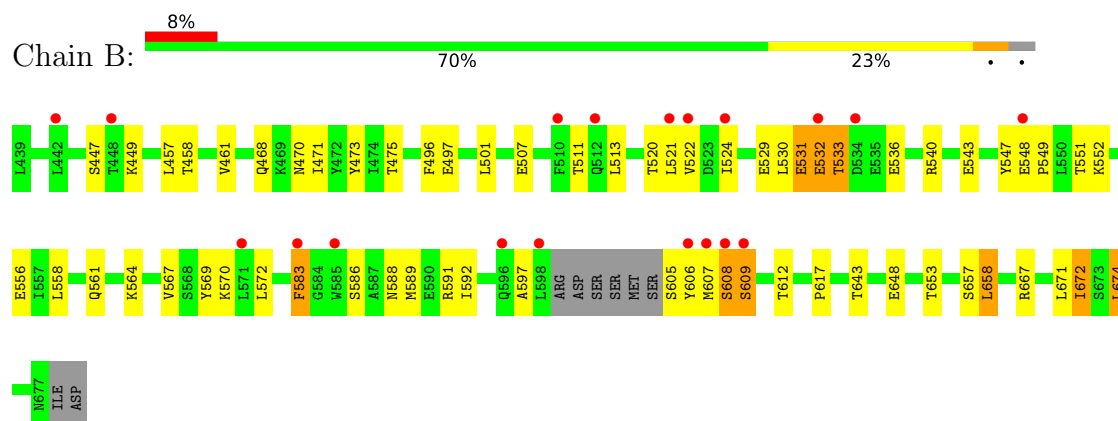
3 Residue-property plots [i](#)

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

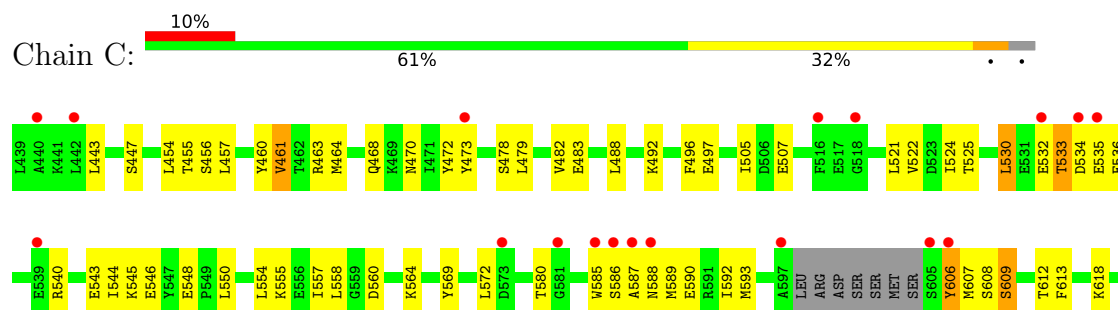
• Molecule 1: ATP-dependent molecular chaperone HSP82

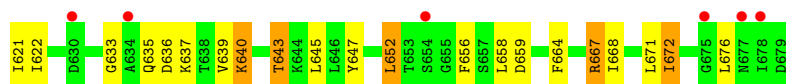


• Molecule 1: ATP-dependent molecular chaperone HSP82

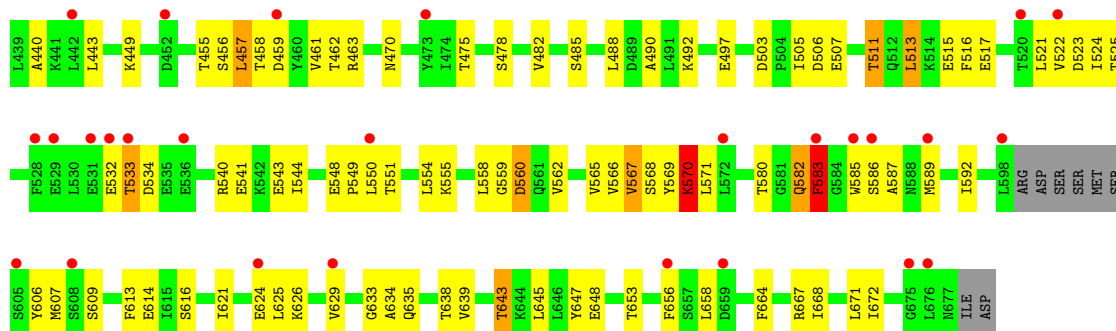


• Molecule 1: ATP-dependent molecular chaperone HSP82





- Molecule 1: ATP-dependent molecular chaperone HSP82



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	97.71 Å 97.71 Å 275.57 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	79.71 – 2.94 79.71 – 2.94	Depositor EDS
% Data completeness (in resolution range)	78.1 (79.71-2.94) 78.4 (79.71-2.94)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.68 (at 2.96 Å)	Xtriage
Refinement program	REFMAC 5.8.0257, PHENIX 1.19_4092	Depositor
R, R_{free}	0.247 , 0.306 0.261 , 0.310	Depositor DCC
R_{free} test set	1125 reflections (3.82%)	wwPDB-VP
Wilson B-factor (Å ²)	70.1	Xtriage
Anisotropy	0.045	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 80.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	6816	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.44% 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: W5R

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	0.21	0/1740	0.54	2/2368 (0.1%)
1	B	0.18	0/1719	0.52	0/2341
1	C	0.28	0/1734	0.59	2/2361 (0.1%)
1	D	0.29	0/1694	0.70	7/2315 (0.3%)
All	All	0.24	0/6887	0.59	11/9385 (0.1%)

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	D	0	1

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	570	LYS	CB-CA-C	-6.05	98.39	110.42
1	D	624	GLU	CB-CG-CD	-5.93	102.52	112.60
1	D	569	TYR	CA-C-N	5.91	132.83	121.54
1	D	569	TYR	C-N-CA	5.91	132.83	121.54
1	D	624	GLU	CB-CA-C	5.76	119.31	109.24
1	C	640	LYS	CA-CB-CG	5.74	125.59	114.10
1	C	640	LYS	CB-CG-CD	-5.42	98.84	111.30
1	D	582	GLN	CA-C-N	5.24	131.54	121.54
1	D	582	GLN	C-N-CA	5.24	131.54	121.54
1	A	617	PRO	CA-C-N	5.16	131.71	122.38
1	A	617	PRO	C-N-CA	5.16	131.71	122.38

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	570	LYS	Peptide

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	1712	0	1602	50	0
1	B	1692	0	1571	42	0
1	C	1709	0	1579	60	0
1	D	1667	0	1520	62	0
2	C	35	0	0	0	0
3	A	1	0	0	0	0
All	All	6816	0	6272	204	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 16.

All (204) 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:475:THR:H	1:A:524:ILE:HD11	1.35	0.89
1:A:469:LYS:H	1:A:530:LEU:HD21	1.37	0.89
1:D:550:LEU:HD11	1:D:625:LEU:HD23	1.63	0.81
1:A:449:LYS:HD2	1:A:497:GLU:HB2	1.63	0.81
1:A:522:VAL:HG12	1:A:525:THR:H	1.44	0.79
1:C:564:LYS:HG3	1:C:612:THR:HG23	1.64	0.79
1:D:645:LEU:HD13	1:D:667:ARG:HG2	1.65	0.77
1:A:522:VAL:HG13	1:A:524:ILE:HD12	1.65	0.77
1:B:530:LEU:O	1:B:532:GLU:N	2.20	0.74
1:A:672:ILE:HG22	1:D:671:LEU:HD21	1.71	0.72
1:D:475:THR:H	1:D:524:ILE:HD11	1.55	0.71
1:D:540:ARG:O	1:D:543:GLU:N	2.25	0.69
1:B:522:VAL:HG12	1:B:524:ILE:H	1.58	0.68
1:C:637:LYS:HA	1:C:640:LYS:HD2	1.74	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:459:ASP:HB3	1:D:463:ARG:HH12	1.60	0.66
1:C:554:LEU:HA	1:C:557:ILE:HD12	1.77	0.65
1:D:571:LEU:HD23	1:D:616:SER:HB2	1.79	0.65
1:B:533:THR:HG23	1:B:536:GLU:H	1.62	0.64
1:C:533:THR:O	1:C:535:GLU:N	2.29	0.64
1:B:608:SER:OG	1:B:609:SER:N	2.27	0.63
1:C:633:GLY:HA2	1:C:636:ASP:HB2	1.80	0.62
1:B:449:LYS:HB2	1:B:497:GLU:HG3	1.82	0.62
1:D:560:ASP:OD1	1:D:560:ASP:N	2.28	0.62
1:D:550:LEU:HD13	1:D:626:LYS:HA	1.81	0.61
1:B:564:LYS:HB2	1:B:612:THR:HG23	1.82	0.61
1:D:507:GLU:O	1:D:511:THR:OG1	2.18	0.60
1:D:589:MET:HA	1:D:589:MET:HE2	1.83	0.60
1:D:555:LYS:NZ	1:D:559:GLY:O	2.35	0.59
1:A:555:LYS:NZ	1:A:562:VAL:O	2.35	0.59
1:C:443:LEU:HA	1:C:456:SER:HA	1.84	0.58
1:C:645:LEU:HG	1:C:667:ARG:HG2	1.85	0.58
1:A:536:GLU:HB3	1:A:540:ARG:HD2	1.85	0.58
1:B:532:GLU:O	1:B:533:THR:HG22	2.03	0.58
1:D:645:LEU:HD11	1:D:668:ILE:HD13	1.86	0.58
1:A:553:ALA:O	1:A:557:ILE:HD12	2.04	0.58
1:A:606:TYR:CE2	1:A:608:SER:HB3	2.39	0.57
1:C:608:SER:OG	1:C:609:SER:N	2.28	0.57
1:D:648:GLU:HG2	1:D:658:LEU:HD21	1.85	0.57
1:B:588:ASN:O	1:B:592:ILE:HG13	2.04	0.57
1:D:525:THR:HA	1:D:580:THR:O	2.05	0.57
1:D:606:TYR:O	1:D:607:MET:HG3	2.05	0.56
1:C:550:LEU:HD22	1:C:554:LEU:HD11	1.87	0.56
1:A:579:ARG:HE	1:A:614:GLU:CD	2.14	0.56
1:A:672:ILE:HD11	1:D:621:ILE:HD13	1.86	0.56
1:A:572:LEU:HA	1:A:618:LYS:HZ2	1.71	0.56
1:C:536:GLU:O	1:C:540:ARG:HG3	2.05	0.56
1:C:470:ASN:HB3	1:C:521:LEU:HD21	1.88	0.55
1:A:652:LEU:HD21	1:A:658:LEU:HG	1.88	0.55
1:C:522:VAL:HG12	1:C:525:THR:H	1.72	0.55
1:D:633:GLY:O	1:D:635:GLN:N	2.40	0.55
1:D:443:LEU:HA	1:D:456:SER:HA	1.88	0.55
1:A:639:VAL:O	1:A:643:THR:HG23	2.07	0.54
1:B:540:ARG:HD2	1:B:569:TYR:CE1	2.42	0.54
1:D:458:THR:O	1:D:461:VAL:HG22	2.08	0.54
1:A:530:LEU:O	1:A:531:GLU:C	2.51	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:648:GLU:HG2	1:D:658:LEU:CD2	2.38	0.54
1:C:588:ASN:O	1:C:592:ILE:HD12	2.07	0.53
1:D:490:ALA:HB2	1:D:570:LYS:O	2.09	0.53
1:D:664:PHE:O	1:D:668:ILE:HG12	2.07	0.53
1:B:470:ASN:HB3	1:B:521:LEU:HD21	1.90	0.53
1:C:618:LYS:C	1:C:618:LYS:HZ3	2.16	0.53
1:D:470:ASN:HB3	1:D:521:LEU:HD21	1.91	0.53
1:D:522:VAL:HG12	1:D:524:ILE:H	1.74	0.53
1:A:645:LEU:HD21	1:A:671:LEU:HD22	1.90	0.52
1:B:648:GLU:HB3	1:B:658:LEU:HD22	1.91	0.52
1:B:536:GLU:O	1:B:540:ARG:HG2	2.10	0.52
1:B:648:GLU:OE1	1:B:667:ARG:NH1	2.43	0.52
1:D:587:ALA:HA	1:D:656:PHE:CE2	2.45	0.52
1:B:547:TYR:O	1:B:551:THR:HG23	2.10	0.52
1:A:588:ASN:OD1	1:A:591:ARG:NH2	2.42	0.52
1:D:582:GLN:O	1:D:583:PHE:HB2	2.10	0.52
1:A:478:SER:O	1:A:482:VAL:HG12	2.11	0.51
1:A:573:ASP:H	1:A:618:LYS:NZ	2.09	0.51
1:B:605:SER:O	1:B:607:MET:N	2.40	0.51
1:C:505:ILE:H	1:C:505:ILE:HD12	1.76	0.51
1:D:522:VAL:HG12	1:D:524:ILE:N	2.26	0.51
1:A:588:ASN:O	1:A:592:ILE:HG12	2.11	0.50
1:B:531:GLU:O	1:B:533:THR:N	2.44	0.50
1:D:554:LEU:HB2	1:D:565:VAL:HG11	1.94	0.50
1:B:552:LYS:O	1:B:556:GLU:HG3	2.12	0.50
1:D:668:ILE:O	1:D:672:ILE:HG23	2.12	0.50
1:B:540:ARG:HH11	1:B:540:ARG:HG3	1.75	0.50
1:C:555:LYS:HD2	1:C:555:LYS:O	2.11	0.50
1:B:522:VAL:HG12	1:B:524:ILE:N	2.27	0.50
1:A:489:ASP:HB3	1:A:572:LEU:HD12	1.94	0.49
1:D:587:ALA:HA	1:D:656:PHE:HE2	1.77	0.49
1:A:568:SER:HB3	1:A:614:GLU:HB3	1.93	0.49
1:C:461:VAL:HA	1:C:464:MET:HG3	1.94	0.49
1:A:666:SER:O	1:A:670:ARG:HB2	2.12	0.49
1:C:478:SER:O	1:C:482:VAL:HG12	2.12	0.49
1:C:543:GLU:HA	1:C:546:GLU:OE2	2.12	0.49
1:A:579:ARG:NE	1:A:614:GLU:OE1	2.43	0.49
1:A:652:LEU:HD11	1:A:658:LEU:HD11	1.94	0.49
1:C:652:LEU:HD11	1:C:658:LEU:HG	1.94	0.49
1:A:621:ILE:HG21	1:D:672:ILE:HD11	1.95	0.49
1:D:475:THR:HG21	1:D:507:GLU:HG3	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:664:PHE:CZ	1:D:668:ILE:HD11	2.48	0.49
1:B:672:ILE:HG12	1:C:621:ILE:HD13	1.95	0.49
1:C:522:VAL:CG1	1:C:524:ILE:HB	2.43	0.49
1:C:664:PHE:CZ	1:C:668:ILE:HD11	2.48	0.49
1:D:516:PHE:HD1	1:D:517:GLU:OE1	1.95	0.49
1:A:540:ARG:O	1:A:543:GLU:N	2.46	0.49
1:B:507:GLU:OE2	1:B:586:SER:OG	2.26	0.49
1:C:635:GLN:O	1:C:640:LYS:HE3	2.12	0.48
1:C:507:GLU:OE2	1:C:589:MET:HG2	2.12	0.48
1:A:530:LEU:O	1:A:532:GLU:N	2.47	0.48
1:A:608:SER:OG	1:A:609:SER:N	2.47	0.48
1:D:507:GLU:OE2	1:D:586:SER:OG	2.28	0.48
1:A:589:MET:HE2	1:A:593:MET:HE3	1.96	0.47
1:D:639:VAL:O	1:D:643:THR:HG23	2.14	0.47
1:A:472:TYR:CE1	1:A:521:LEU:HD11	2.49	0.47
1:A:469:LYS:N	1:A:530:LEU:HD21	2.17	0.47
1:B:588:ASN:HD21	1:B:591:ARG:HH21	1.63	0.47
1:A:558:LEU:O	1:A:561:GLN:HG3	2.13	0.47
1:B:457:LEU:O	1:B:461:VAL:HG23	2.15	0.47
1:D:503:ASP:HB3	1:D:506:ASP:OD1	2.15	0.47
1:C:540:ARG:O	1:C:543:GLU:N	2.48	0.46
1:D:457:LEU:O	1:D:461:VAL:HG13	2.15	0.46
1:D:523:ASP:OD2	1:D:523:ASP:N	2.44	0.46
1:C:455:THR:OG1	1:C:463:ARG:NH2	2.41	0.46
1:B:507:GLU:O	1:B:511:THR:HG23	2.16	0.45
1:B:589:MET:HE3	1:B:589:MET:HA	1.97	0.45
1:D:613:PHE:CE1	1:D:647:TYR:HA	2.51	0.45
1:C:505:ILE:HG23	1:D:585:TRP:CZ2	2.52	0.45
1:A:464:MET:HE3	1:A:468:GLN:HB3	1.97	0.45
1:B:468:GLN:HG3	1:B:496:PHE:CE1	2.51	0.45
1:B:473:TYR:CE2	1:B:522:VAL:HG22	2.51	0.45
1:C:457:LEU:O	1:C:461:VAL:HG23	2.16	0.45
1:C:543:GLU:HB3	1:C:569:TYR:OH	2.16	0.45
1:C:545:LYS:O	1:C:548:GLU:HG2	2.17	0.45
1:C:585:TRP:CE3	1:C:590:GLU:HG3	2.52	0.45
1:C:645:LEU:HD23	1:C:664:PHE:HE1	1.82	0.45
1:A:460:TYR:CE2	1:A:471:ILE:HG23	2.51	0.45
1:B:529:GLU:O	1:B:530:LEU:C	2.60	0.44
1:B:671:LEU:HD23	1:C:671:LEU:HD23	1.98	0.44
1:D:478:SER:O	1:D:482:VAL:HG12	2.17	0.44
1:C:473:TYR:CZ	1:C:522:VAL:HG23	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:439:LEU:HD13	1:B:583:PHE:CE2	2.52	0.44
1:C:479:LEU:O	1:C:483:GLU:HB2	2.17	0.44
1:C:488:LEU:HB3	1:C:492:LYS:HD2	1.99	0.44
1:C:639:VAL:O	1:C:643:THR:HG23	2.16	0.44
1:B:540:ARG:O	1:B:543:GLU:N	2.51	0.44
1:B:671:LEU:HA	1:B:674:LEU:HD12	1.98	0.44
1:A:473:TYR:CZ	1:A:522:VAL:HG23	2.53	0.44
1:A:578:ILE:HD11	1:A:647:TYR:CE1	2.52	0.44
1:B:540:ARG:HG3	1:B:540:ARG:NH1	2.32	0.44
1:C:668:ILE:O	1:C:672:ILE:HG23	2.17	0.44
1:C:606:TYR:O	1:C:607:MET:HB3	2.18	0.44
1:D:544:ILE:HD12	1:D:544:ILE:H	1.83	0.43
1:C:470:ASN:HB3	1:C:521:LEU:CD2	2.47	0.43
1:C:460:TYR:O	1:C:464:MET:HG3	2.19	0.43
1:D:488:LEU:HB3	1:D:492:LYS:HD2	2.00	0.43
1:B:471:ILE:HB	1:B:520:THR:HG22	2.00	0.43
1:C:460:TYR:CE2	1:C:497:GLU:HB3	2.53	0.43
1:C:558:LEU:HD23	1:C:558:LEU:HA	1.75	0.43
1:C:587:ALA:HA	1:C:656:PHE:CE1	2.54	0.43
1:B:548:GLU:N	1:B:549:PRO:HD2	2.34	0.43
1:D:533:THR:OG1	1:D:534:ASP:N	2.52	0.43
1:C:532:GLU:O	1:C:533:THR:O	2.37	0.43
1:C:468:GLN:HG3	1:C:496:PHE:CE1	2.54	0.43
1:C:546:GLU:H	1:C:546:GLU:HG3	1.64	0.43
1:C:668:ILE:O	1:C:671:LEU:N	2.52	0.43
1:D:565:VAL:HG22	1:D:613:PHE:HB3	2.01	0.43
1:B:475:THR:HG21	1:B:507:GLU:HG3	2.01	0.43
1:D:513:LEU:C	1:D:515:GLU:H	2.26	0.43
1:A:547:TYR:O	1:A:551:THR:HG23	2.19	0.42
1:B:558:LEU:O	1:B:561:GLN:HG3	2.19	0.42
1:B:567:VAL:HG23	1:B:617:PRO:HD3	2.01	0.42
1:C:530:LEU:HD12	1:C:530:LEU:HA	1.90	0.42
1:A:531:GLU:O	1:A:532:GLU:C	2.62	0.42
1:A:607:MET:O	1:A:608:SER:C	2.62	0.42
1:B:605:SER:C	1:B:607:MET:H	2.23	0.42
1:D:558:LEU:HD23	1:D:562:VAL:HG11	2.01	0.42
1:A:475:THR:HG21	1:A:507:GLU:HG3	2.02	0.42
1:D:449:LYS:HD2	1:D:497:GLU:HB2	2.00	0.42
1:D:544:ILE:HG13	1:D:567:VAL:CG2	2.49	0.42
1:A:458:THR:O	1:A:461:VAL:HG12	2.20	0.42
1:C:606:TYR:CB	1:D:505:ILE:HG21	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:571:LEU:HD23	1:D:616:SER:CB	2.49	0.41
1:A:671:LEU:HD21	1:D:672:ILE:HG22	2.02	0.41
1:C:645:LEU:HD21	1:C:668:ILE:HD13	2.03	0.41
1:D:540:ARG:O	1:D:544:ILE:HD12	2.21	0.41
1:C:652:LEU:CD1	1:C:658:LEU:HG	2.51	0.41
1:A:443:LEU:HD23	1:A:454:LEU:HB3	2.02	0.41
1:A:548:GLU:N	1:A:549:PRO:HD2	2.35	0.41
1:A:550:LEU:HD13	1:A:629:VAL:HG21	2.03	0.41
1:D:548:GLU:HA	1:D:551:THR:OG1	2.21	0.41
1:B:591:ARG:HH11	1:B:591:ARG:HG2	1.83	0.41
1:A:500:PHE:C	1:A:501:LEU:HD23	2.46	0.41
1:A:607:MET:O	1:A:607:MET:HG2	2.21	0.41
1:B:570:LYS:HE2	1:B:570:LYS:HB2	1.82	0.41
1:C:507:GLU:OE2	1:C:586:SER:OG	2.35	0.41
1:C:672:ILE:O	1:C:676:LEU:HD23	2.21	0.41
1:D:548:GLU:N	1:D:549:PRO:HD2	2.36	0.41
1:B:473:TYR:CD2	1:B:501:LEU:HD22	2.56	0.40
1:C:580:THR:HG21	1:C:585:TRP:O	2.21	0.40
1:C:593:MET:HG3	1:D:505:ILE:HG13	2.03	0.40
1:C:613:PHE:CE1	1:C:647:TYR:HA	2.56	0.40
1:D:541:GLU:HA	1:D:544:ILE:HD13	2.03	0.40
1:C:472:TYR:CE1	1:C:521:LEU:HD11	2.56	0.40
1:D:568:SER:HB3	1:D:614:GLU:HB3	2.03	0.40
1:D:470:ASN:HB3	1:D:521:LEU:CD2	2.50	0.40
1:C:658:LEU:HD23	1:C:658:LEU:HA	1.92	0.40
1:D:555:LYS:HD2	1:D:555:LYS:HA	1.76	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	227/241 (94%)	207 (91%)	17 (8%)	3 (1%)	9	25
1	B	229/241 (95%)	205 (90%)	17 (7%)	7 (3%)	3	8
1	C	230/241 (95%)	207 (90%)	20 (9%)	3 (1%)	9	25
1	D	229/241 (95%)	207 (90%)	16 (7%)	6 (3%)	4	12
All	All	915/964 (95%)	826 (90%)	70 (8%)	19 (2%)	5	15

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	531	GLU
1	A	608	SER
1	B	531	GLU
1	B	532	GLU
1	B	583	PHE
1	C	533	THR
1	C	534	ASP
1	C	606	TYR
1	D	533	THR
1	D	583	PHE
1	D	609	SER
1	B	533	THR
1	B	597	ALA
1	B	606	TYR
1	D	634	ALA
1	A	533	THR
1	D	532	GLU
1	B	608	SER
1	D	440	ALA

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	165/216 (76%)	156 (94%)	9 (6%)	19	41
1	B	159/216 (74%)	148 (93%)	11 (7%)	14	33

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	164/216 (76%)	150 (92%)	14 (8%)	10	24
1	D	153/216 (71%)	138 (90%)	15 (10%)	7	20
All	All	641/864 (74%)	592 (92%)	49 (8%)	12	29

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

Mol	Chain	Res	Type
1	A	457	LEU
1	A	459	ASP
1	A	511	THR
1	A	520	THR
1	A	567	VAL
1	A	572	LEU
1	A	618	LYS
1	A	638	THR
1	A	643	THR
1	B	447	SER
1	B	458	THR
1	B	513	LEU
1	B	572	LEU
1	B	609	SER
1	B	643	THR
1	B	653	THR
1	B	657	SER
1	B	658	LEU
1	B	672	ILE
1	B	674	LEU
1	C	447	SER
1	C	454	LEU
1	C	461	VAL
1	C	530	LEU
1	C	544	ILE
1	C	560	ASP
1	C	572	LEU
1	C	609	SER
1	C	622	ILE
1	C	643	THR
1	C	652	LEU
1	C	659	ASP
1	C	667	ARG
1	C	672	ILE

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Mol	Chain	Res	Type
1	D	455	THR
1	D	457	LEU
1	D	462	THR
1	D	485	SER
1	D	511	THR
1	D	513	LEU
1	D	560	ASP
1	D	566	VAL
1	D	567	VAL
1	D	583	PHE
1	D	592	ILE
1	D	629	VAL
1	D	638	THR
1	D	643	THR
1	D	653	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	588	ASN
1	C	561	GLN
1	C	635	GLN
1	D	588	ASN

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](#)

1 ligand is 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	W5R	C	801	-	36,36,36	0.47	0	43,52,52	0.74	1 (2%)

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	W5R	C	801	-	-	7/32/56/56	0/2/2/2

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	801	W5R	C7-C8-N	-2.05	118.14	120.75

There are no chirality outliers.

All (7) torsion outliers are listed below:

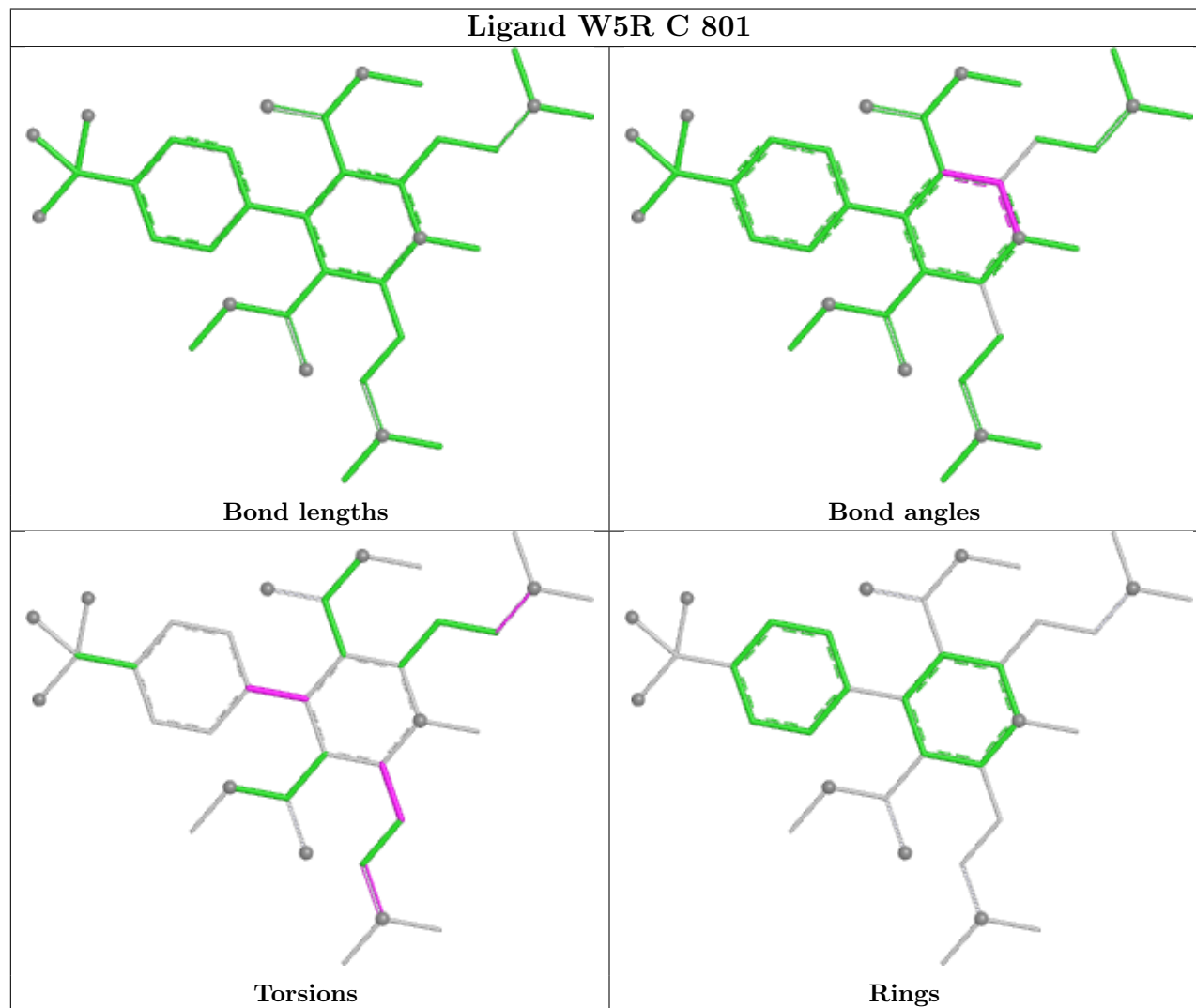
Mol	Chain	Res	Type	Atoms
2	C	801	W5R	C14-C13-C9-N
2	C	801	W5R	C13-C14-N1-C15
2	C	801	W5R	C13-C14-N1-C16
2	C	801	W5R	C18-C19-N2-C21
2	C	801	W5R	C2-C3-C6-C10
2	C	801	W5R	C4-C3-C6-C10
2	C	801	W5R	C14-C13-C9-C10

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 ⓘ

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	231/241 (95%)	0.64	18 (7%) 19 16	26, 48, 83, 111	0
1	B	233/241 (96%)	0.62	19 (8%) 17 15	28, 52, 89, 108	0
1	C	234/241 (97%)	0.85	24 (10%) 12 11	32, 58, 92, 104	0
1	D	233/241 (96%)	0.82	27 (11%) 9 9	30, 62, 100, 116	0
All	All	931/964 (96%)	0.73	88 (9%) 14 13	26, 55, 90, 116	0

All (88) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	532	GLU	6.4
1	C	585	TRP	5.7
1	B	583	PHE	4.9
1	D	585	TRP	4.7
1	A	531	GLU	4.1
1	D	531	GLU	3.9
1	C	442	LEU	3.8
1	A	618	LYS	3.8
1	B	585	TRP	3.7
1	D	624	GLU	3.6
1	A	634	ALA	3.6
1	C	587	ALA	3.5
1	D	520	THR	3.5
1	B	607	MET	3.4
1	C	630	ASP	3.4
1	A	439	LEU	3.3
1	D	583	PHE	3.2
1	A	529	GLU	3.1
1	A	585	TRP	3.0
1	A	535	GLU	3.0
1	D	629	VAL	3.0

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Mol	Chain	Res	Type	RSRZ
1	C	597	ALA	3.0
1	A	530	LEU	3.0
1	B	442	LEU	3.0
1	D	473	TYR	2.9
1	D	550	LEU	2.9
1	D	589	MET	2.9
1	C	606	TYR	2.8
1	C	532	GLU	2.8
1	A	635	GLN	2.8
1	B	532	GLU	2.8
1	C	534	ASP	2.8
1	A	506	ASP	2.8
1	C	535	GLU	2.7
1	B	598	LEU	2.7
1	D	675	GLY	2.7
1	A	605	SER	2.7
1	D	586	SER	2.7
1	B	606	TYR	2.6
1	C	581	GLY	2.6
1	B	521	LEU	2.6
1	B	448	THR	2.6
1	A	489	ASP	2.6
1	D	452	ASP	2.6
1	D	572	LEU	2.5
1	C	586	SER	2.5
1	C	516	PHE	2.5
1	C	678	ILE	2.5
1	D	536	GLU	2.5
1	D	459	ASP	2.5
1	D	659	ASP	2.5
1	B	534	ASP	2.4
1	D	533	THR	2.4
1	C	675	GLY	2.4
1	D	442	LEU	2.4
1	B	512	GLN	2.4
1	B	524	ILE	2.4
1	D	598	LEU	2.3
1	C	654	SER	2.3
1	D	608	SER	2.3
1	C	573	ASP	2.3
1	B	548	GLU	2.3
1	A	522	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
1	D	522	VAL	2.3
1	D	676	LEU	2.2
1	B	596	GLN	2.2
1	A	508	TYR	2.2
1	B	609	SER	2.2
1	B	571	LEU	2.2
1	C	539	GLU	2.2
1	A	525	THR	2.2
1	A	538	ALA	2.2
1	B	522	VAL	2.2
1	D	605	SER	2.2
1	D	528	PHE	2.2
1	C	634	ALA	2.2
1	D	532	GLU	2.2
1	C	588	ASN	2.2
1	C	677	ASN	2.2
1	B	608	SER	2.2
1	D	529	GLU	2.2
1	C	518	GLY	2.1
1	B	510	PHE	2.1
1	C	473	TYR	2.1
1	C	605	SER	2.0
1	D	656	PHE	2.0
1	C	440	ALA	2.0
1	A	630	ASP	2.0

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

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

6.3 Carbohydrates ⓘ

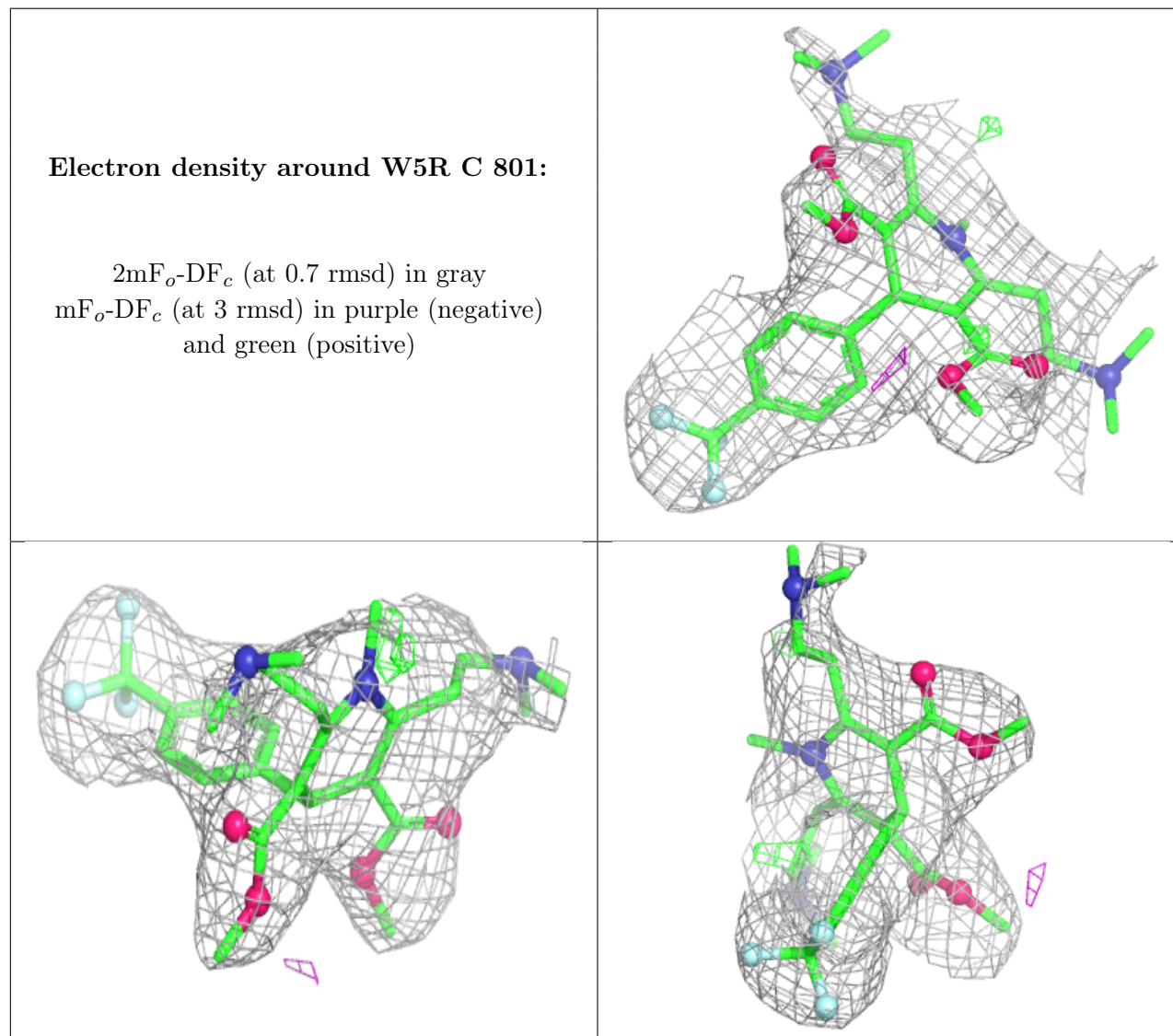
There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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	W5R	C	801	35/35	0.85	0.17	52,65,80,90	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.



6.5 Other polymers ⓘ

There are no such residues in this entry.