



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 11:42 AM UTC

PDB ID : 7AYR / pdb_00007ayr
Title : Crystal structure of DPP8 in complex with a 4-oxo-b-lactam based inhibitor, B115
Authors : Ross, B.H.; Huber, R.
Deposited on : 2020-11-13
Resolution : 2.69 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

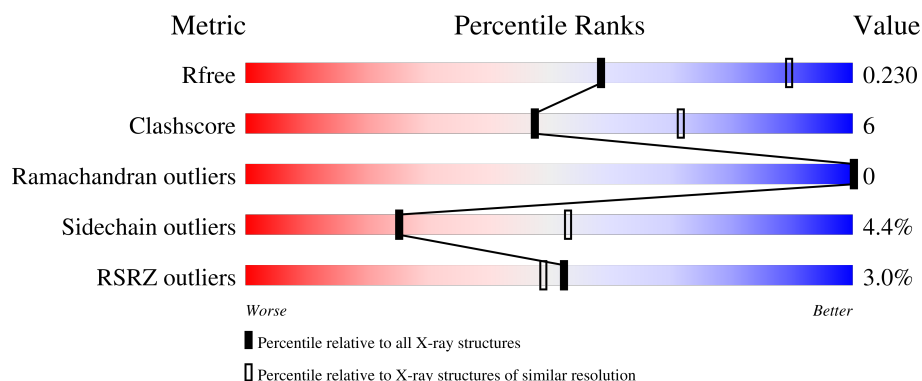
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

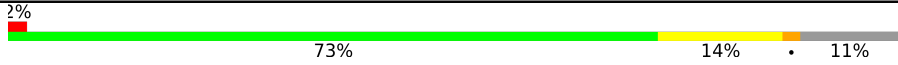
The reported resolution of this entry is 2.69 Å.

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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	898	
1	B	898	

2 Entry composition [i](#)

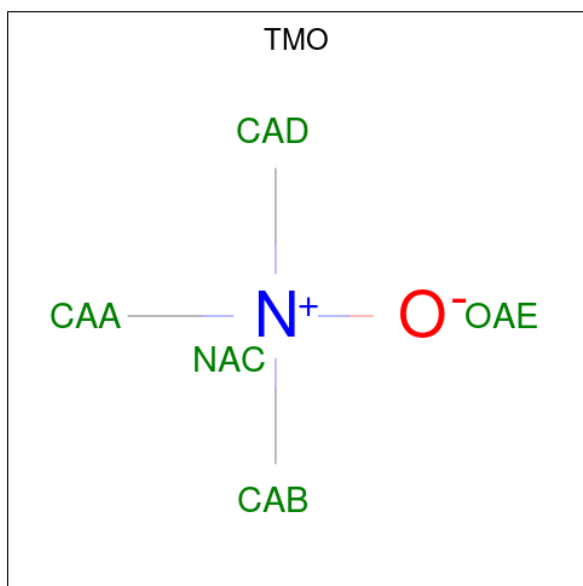
There are 5 unique types of molecules in this entry. The entry contains 13131 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Dipeptidyl peptidase 8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	802	Total	C	N	O	S	0	1	0
			6519	4197	1086	1211	25			
1	B	804	Total	C	N	O	S	0	0	0
			6521	4199	1084	1212	26			

- Molecule 2 is trimethylamine oxide (CCD ID: TMO) (formula: C_3H_9NO).



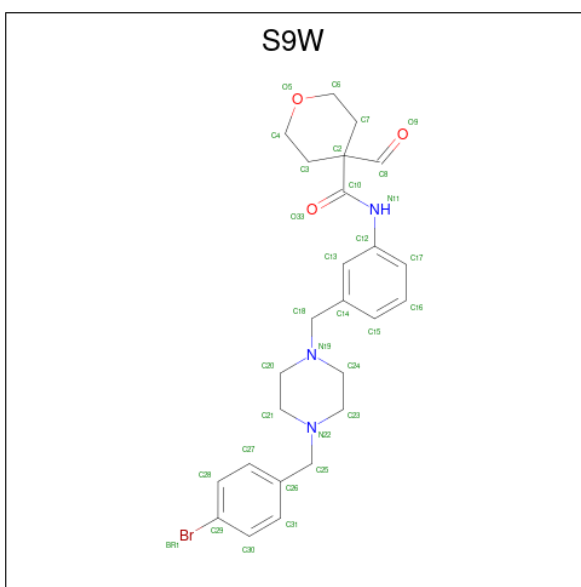
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			5	3	1	1		
2	B	1	Total	C	N	O	0	0
			5	3	1	1		

- Molecule 3 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	P	0	0
			5	4	1		
3	B	1	Total	O	P	0	0
			5	4	1		

- Molecule 4 is {N}-[3-[[4-[(4-bromophenyl)methyl]piperazin-1-yl]methyl]phenyl]-4-methanoyl-oxane-4-carboxamide (CCD ID: S9W) (formula: C₂₅H₃₀BrN₃O₃) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	Br	C	N	O	
			32	1	25	3	3	0

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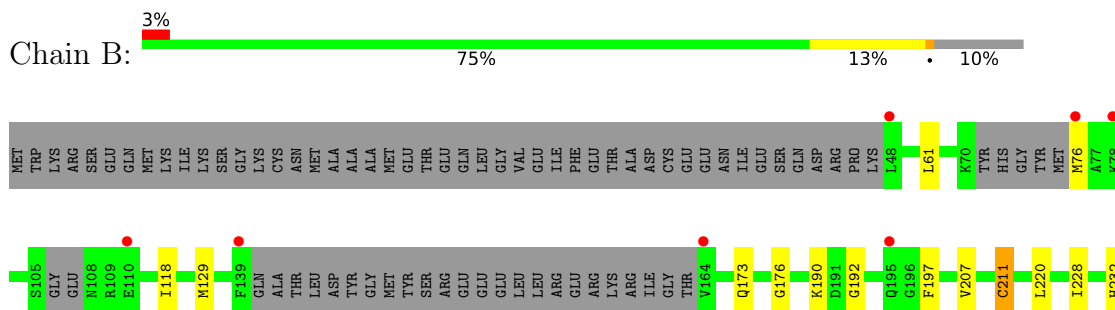
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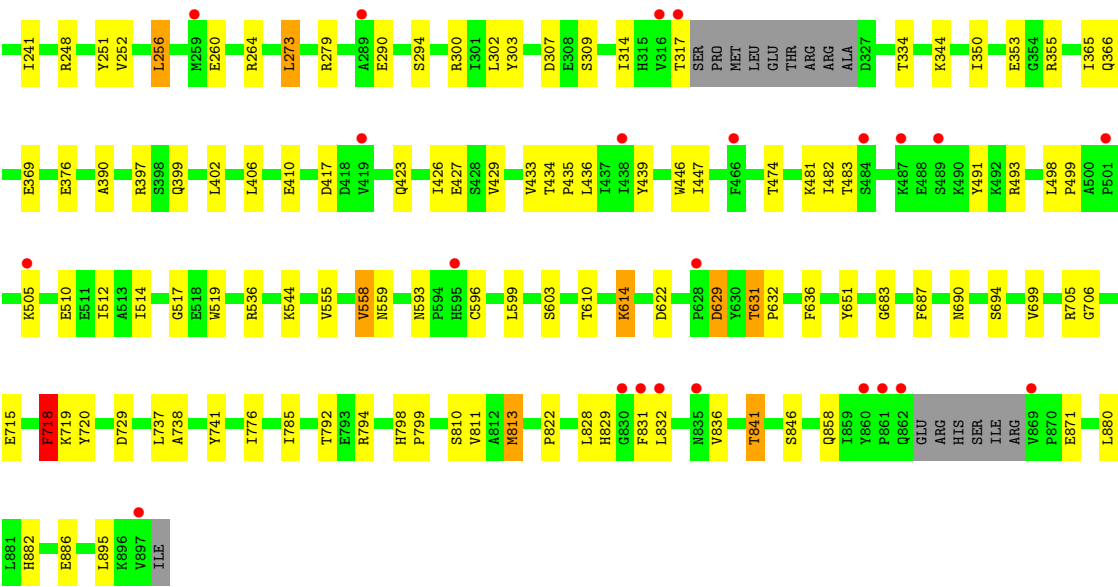
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	B	1	Total	Br	C	N	O	0	0
			32	1	25	3	3		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	4	Total	O	0	0
			4	4		
5	B	3	Total	O	0	0
			3	3		

- Molecule 1: Dipeptidyl peptidase 8





4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	149.56Å 149.56Å 269.63Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.96 – 2.69 48.96 – 2.69	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.96-2.69) 100.0 (48.96-2.69)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.07 (at 2.69Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.187 , 0.230 0.191 , 0.230	Depositor DCC
R_{free} test set	4718 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	75.8	Xtriage
Anisotropy	0.330	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 38.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.029 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	13131	wwPDB-VP
Average B, all atoms (Å ²)	83.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.08% 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: PO4, TMO, S9W

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	1.03	1/6702 (0.0%)	1.31	13/9092 (0.1%)
1	B	1.04	1/6701 (0.0%)	1.34	10/9091 (0.1%)
All	All	1.03	2/13403 (0.0%)	1.33	23/18183 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	517	GLY	C-O	5.35	1.28	1.23
1	A	118	ILE	N-CA	5.21	1.50	1.46

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	631	THR	CA-C-N	13.11	129.23	119.66
1	B	631	THR	C-N-CA	13.11	129.23	119.66
1	A	631	THR	CA-C-N	9.26	129.91	120.38
1	A	631	THR	C-N-CA	9.26	129.91	120.38
1	B	211	CYS	CA-C-N	7.55	127.43	119.05
1	B	211	CYS	C-N-CA	7.55	127.43	119.05
1	A	593	ASN	CA-C-N	7.27	128.92	119.84
1	A	593	ASN	C-N-CA	7.27	128.92	119.84
1	B	593	ASN	CA-C-N	7.23	128.88	119.84
1	B	593	ASN	C-N-CA	7.23	128.88	119.84
1	A	444	ASP	N-CA-C	-6.25	105.29	113.17
1	A	555	VAL	CA-C-O	-5.87	116.03	120.96
1	A	172	HIS	CA-C-N	5.75	129.45	120.82
1	A	172	HIS	C-N-CA	5.75	129.45	120.82
1	B	895	LEU	CA-C-N	5.52	128.24	120.28
1	B	895	LEU	C-N-CA	5.52	128.24	120.28
1	A	242	VAL	CA-C-O	-5.46	116.06	120.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	718	PHE	CA-CB-CG	5.31	119.11	113.80
1	A	242	VAL	O-C-N	5.29	125.87	121.85
1	A	375	VAL	CA-C-O	-5.26	115.05	121.04
1	A	860	TYR	CB-CA-C	5.14	115.99	110.65
1	B	555	VAL	CA-C-O	-5.13	116.65	120.96
1	A	540	PHE	CA-C-O	-5.10	115.77	121.23

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	6519	0	6340	75	0
1	B	6521	0	6340	73	0
2	A	5	0	9	0	0
2	B	5	0	9	0	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
4	A	32	0	0	0	0
4	B	32	0	0	0	0
5	A	4	0	0	1	0
5	B	3	0	0	0	0
All	All	13131	0	12698	145	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 6.

All (145) 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:B:705:ARG:HD3	1:B:729:ASP:OD2	1.77	0.83
1:A:829:HIS:HD2	1:A:841:THR:OG1	1.65	0.78
1:A:882:HIS:CE1	1:B:822:PRO:HG3	2.18	0.78
1:B:365:ILE:HD12	1:B:429:VAL:HG21	1.66	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:434:THR:HG22	1:B:435:PRO:HD2	1.70	0.71
1:A:683:GLY:O	1:A:687:PHE:HB2	1.90	0.70
1:B:252:VAL:HG13	1:B:264:ARG:HB3	1.73	0.69
1:A:705:ARG:HD3	1:A:729:ASP:OD2	1.93	0.69
1:A:334:THR:CG2	1:A:785:ILE:HA	2.24	0.68
1:A:376:GLU:HG3	1:A:397:ARG:HB2	1.74	0.68
1:B:423:GLN:HA	1:B:426:ILE:HD12	1.76	0.68
1:A:290:GLU:OE2	1:A:300:ARG:NH2	2.27	0.66
1:A:828:LEU:HD23	1:A:858:GLN:HB2	1.76	0.65
1:B:211:CYS:HB3	1:B:232:HIS:CD2	2.31	0.65
1:B:399:GLN:OE1	1:B:794:ARG:HD3	1.97	0.65
1:A:273:LEU:HD22	1:A:279:ARG:NH2	2.12	0.64
1:A:822:PRO:HG3	1:B:882:HIS:CE1	2.33	0.62
1:A:399:GLN:OE1	1:A:794:ARG:HD3	1.99	0.61
1:A:139:PHE:HD1	1:A:139:PHE:H	1.48	0.61
1:B:651:TYR:HB2	1:B:699:VAL:HB	1.82	0.61
1:A:104:MET:HG3	1:A:111:ASN:HD22	1.66	0.60
1:A:123:ASN:HD21	1:A:125:ALA:HB3	1.65	0.60
1:B:776:ILE:HD13	1:B:880:LEU:CD1	2.32	0.60
1:A:552:LEU:HB3	1:A:567:THR:HG23	1.83	0.60
1:B:683:GLY:O	1:B:687:PHE:HB2	2.02	0.60
1:B:260:GLU:N	1:B:260:GLU:OE2	2.34	0.58
1:A:719:LYS:HG3	1:A:720:TYR:CD2	2.39	0.58
1:B:302:LEU:HD22	1:B:390:ALA:HB1	1.86	0.57
1:B:446:TRP:HZ2	1:B:715:GLU:HG2	1.68	0.57
1:B:402:LEU:HD12	1:B:447:ILE:HG21	1.87	0.57
1:B:614:LYS:HG3	1:B:614:LYS:O	2.04	0.57
1:A:627:LEU:HD12	1:A:681:PHE:HZ	1.69	0.56
1:B:828:LEU:HD23	1:B:858:GLN:HB2	1.87	0.56
1:A:211:CYS:HB3	1:A:232:HIS:CD2	2.42	0.54
1:B:220:LEU:HD23	1:B:228:ILE:HG22	1.88	0.54
1:A:434:THR:HG21	1:A:489:SER:HB2	1.90	0.54
1:B:434:THR:HG22	1:B:435:PRO:CD	2.37	0.53
1:B:491:TYR:CD2	1:B:499:PRO:HA	2.43	0.53
1:B:481:LYS:HB2	1:B:514:ILE:HD11	1.90	0.53
1:A:139:PHE:CD1	1:A:139:PHE:N	2.78	0.52
1:A:238:ILE:HG23	1:A:249:LEU:HD11	1.90	0.52
1:A:348:ILE:HG22	1:A:350:ILE:CD1	2.39	0.52
1:B:705:ARG:NH1	1:B:729:ASP:OD1	2.42	0.52
1:A:292:THR:HB	1:A:293:PRO:CD	2.40	0.52
1:B:118:ILE:HD12	1:B:599:LEU:CD2	2.40	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:629:ASP:OD1	1:A:629:ASP:N	2.26	0.52
1:B:252:VAL:CG1	1:B:264:ARG:HB3	2.41	0.51
1:B:303:TYR:CZ	1:B:344:LYS:HB2	2.45	0.51
1:B:687:PHE:O	1:B:690:ASN:HB3	2.11	0.51
1:A:465:GLU:OE2	1:A:481:LYS:NZ	2.44	0.51
1:B:290:GLU:OE2	1:B:300:ARG:NH2	2.44	0.50
1:B:629:ASP:OD1	1:B:629:ASP:N	2.22	0.50
1:B:129:MET:HE1	1:B:622:ASP:HB2	1.93	0.49
1:B:248:ARG:HH21	1:B:251:TYR:HA	1.77	0.49
1:A:65:LEU:HD21	1:A:881:LEU:HD12	1.93	0.49
1:B:798:HIS:ND1	1:B:799:PRO:HD2	2.28	0.49
1:A:680:ARG:HD3	5:A:1004:HOH:O	2.12	0.49
1:B:353:GLU:HG3	1:B:355:ARG:NH1	2.28	0.48
1:B:317:THR:HB	1:B:832:LEU:HD21	1.96	0.48
1:B:536:ARG:CZ	1:B:558:VAL:HG21	2.42	0.48
1:A:634:GLU:OE1	1:A:652:LYS:HE2	2.13	0.48
1:A:636:PHE:CZ	1:A:737:LEU:HD11	2.48	0.48
1:B:776:ILE:HD13	1:B:880:LEU:HD11	1.95	0.48
1:A:782:THR:H	1:A:840:HIS:HD2	1.62	0.47
1:B:334:THR:CG2	1:B:785:ILE:HA	2.44	0.47
1:B:176:GLY:O	1:B:190:LYS:HA	2.14	0.47
1:B:307:ASP:OD1	1:B:309:SER:OG	2.33	0.47
1:B:410:GLU:HB3	1:B:433:VAL:HG11	1.96	0.47
1:B:510:GLU:OE2	1:B:512:ILE:HD11	2.13	0.47
1:B:829:HIS:HD2	1:B:841:THR:OG1	1.98	0.47
1:A:314:ILE:HD11	1:A:329:PHE:CZ	2.50	0.47
1:B:406:LEU:HB2	1:B:436:LEU:HB2	1.96	0.47
1:B:314:ILE:C	1:B:314:ILE:HD12	2.39	0.46
1:A:292:THR:HB	1:A:293:PRO:HD2	1.97	0.46
1:A:719:LYS:HE3	1:A:720:TYR:CE2	2.51	0.46
1:A:424:ARG:HB3	1:A:424:ARG:NH1	2.30	0.46
1:B:882:HIS:CE1	1:B:886:GLU:HG3	2.51	0.46
1:A:69:ARG:CZ	1:A:874:GLU:HG2	2.45	0.46
1:A:302:LEU:HD22	1:A:390:ALA:HB1	1.96	0.46
1:A:782:THR:H	1:A:840:HIS:CD2	2.33	0.46
1:B:207:VAL:HG22	1:B:241:ILE:HG22	1.98	0.46
1:B:256:LEU:N	1:B:256:LEU:HD13	2.30	0.46
1:A:65:LEU:HD21	1:A:881:LEU:CD1	2.46	0.45
1:A:402:LEU:C	1:A:402:LEU:HD23	2.41	0.45
1:B:376:GLU:HG2	1:B:397:ARG:HB2	1.97	0.45
1:B:369:GLU:H	1:B:369:GLU:CD	2.25	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:687:PHE:O	1:A:690:ASN:HB3	2.17	0.45
1:B:491:TYR:HE2	1:B:498:LEU:C	2.24	0.45
1:B:810:SER:HB2	1:B:813:MET:HG3	1.99	0.45
1:A:651:TYR:HB2	1:A:699:VAL:HB	1.99	0.44
1:A:273:LEU:HD22	1:A:279:ARG:HH21	1.81	0.44
1:A:69:ARG:NH1	1:A:874:GLU:HG2	2.33	0.44
1:A:463:GLU:OE2	1:A:509:LYS:NZ	2.36	0.44
1:B:376:GLU:CG	1:B:397:ARG:HB2	2.48	0.44
1:A:631:THR:HA	1:A:632:PRO:HD2	1.77	0.44
1:B:536:ARG:NH1	1:B:558:VAL:HG21	2.33	0.44
1:A:706:GLY:HA3	1:A:718:PHE:CE2	2.53	0.43
1:B:719:LYS:HG2	1:B:720:TYR:CD2	2.52	0.43
1:A:217:ASP:N	1:A:218:PRO:CD	2.81	0.43
1:B:273:LEU:HD22	1:B:279:ARG:NH1	2.33	0.43
1:B:76:MET:HG2	1:B:76:MET:O	2.17	0.43
1:A:208:GLU:HG3	1:A:244:ARG:NH2	2.33	0.43
1:A:754:TRP:HA	1:A:778:GLY:O	2.17	0.43
1:B:439:TYR:CD1	1:B:482:ILE:HD13	2.54	0.43
1:B:491:TYR:CE2	1:B:499:PRO:N	2.87	0.43
1:A:252:VAL:HG13	1:A:264:ARG:HB3	2.01	0.43
1:A:84:MET:HE1	1:A:171:TYR:CG	2.54	0.43
1:A:306:ASN:OD1	1:A:341:VAL:HG12	2.18	0.43
1:A:770:ASP:OD1	1:A:770:ASP:N	2.51	0.43
1:B:519:TRP:CG	1:B:544:LYS:HA	2.54	0.43
1:B:831:PHE:HD1	1:B:831:PHE:O	2.01	0.43
1:A:421:GLU:OE1	1:A:424:ARG:NH1	2.52	0.42
1:A:829:HIS:HD2	1:A:841:THR:CB	2.30	0.42
1:A:500:ALA:HB1	1:A:501:PRO:HD2	2.01	0.42
1:A:84:MET:SD	1:A:171:TYR:CD2	3.12	0.42
1:B:129:MET:HE3	1:B:129:MET:HB3	1.88	0.42
1:A:882:HIS:CE1	1:B:822:PRO:CG	2.97	0.42
1:A:353:GLU:HG2	1:A:355:ARG:HG2	2.01	0.42
1:A:524:ARG:HH12	1:A:673:GLN:HG2	1.84	0.42
1:A:123:ASN:C	1:A:123:ASN:ND2	2.77	0.42
1:A:406:LEU:HB2	1:A:436:LEU:HB2	2.01	0.42
1:A:416:GLU:HG3	1:A:421:GLU:HG2	2.02	0.42
1:A:522:LEU:HD13	1:A:550:HIS:CD2	2.55	0.42
1:A:829:HIS:HE1	1:A:836:VAL:O	2.02	0.42
1:B:631:THR:HA	1:B:632:PRO:HD3	1.82	0.42
1:A:519:TRP:CG	1:A:544:LYS:HA	2.54	0.42
1:A:99:ILE:O	1:A:115:TYR:HA	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:276:GLU:OE2	1:A:788:ASP:OD2	2.37	0.41
1:A:627:LEU:HD12	1:A:681:PHE:CZ	2.52	0.41
1:B:192:GLY:HA2	1:B:197:PHE:CE2	2.55	0.41
1:B:706:GLY:HA3	1:B:718:PHE:CE2	2.56	0.41
1:A:79:ALA:O	1:A:80:PRO:C	2.63	0.41
1:B:636:PHE:CZ	1:B:737:LEU:HD11	2.55	0.41
1:A:693:ALA:HA	1:A:697:TYR:O	2.20	0.41
1:B:433:VAL:O	1:B:433:VAL:HG12	2.21	0.41
1:B:61:LEU:HD23	1:B:61:LEU:HA	1.84	0.41
1:A:334:THR:HG21	1:A:784:TRP:O	2.21	0.41
1:B:738:ALA:O	1:B:741:TYR:O	2.39	0.41
1:B:836:VAL:O	1:B:836:VAL:HG12	2.20	0.41
1:B:173:GLN:HE21	1:B:173:GLN:HB3	1.66	0.41
1:B:365:ILE:HG23	1:B:366:GLN:HG3	2.04	0.40
1:A:314:ILE:CD1	1:A:837:HIS:CE1	3.05	0.40
1:A:465:GLU:OE2	1:A:481:LYS:CE	2.69	0.40
1:A:829:HIS:CD2	1:A:841:THR:OG1	2.57	0.40
1:B:433:VAL:O	1:B:493:ARG:NH2	2.54	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	791/898 (88%)	764 (97%)	27 (3%)	0	100	100
1	B	792/898 (88%)	758 (96%)	34 (4%)	0	100	100
All	All	1583/1796 (88%)	1522 (96%)	61 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	713/795 (90%)	674 (94%)	39 (6%)	19	45
1	B	713/795 (90%)	689 (97%)	24 (3%)	32	62
All	All	1426/1590 (90%)	1363 (96%)	63 (4%)	25	53

All (63) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	108	ASN
1	A	123	ASN
1	A	139	PHE
1	A	184	SER
1	A	242	VAL
1	A	252	VAL
1	A	254	ASN
1	A	273	LEU
1	A	291	THR
1	A	294	SER
1	A	312	GLU
1	A	315	HIS
1	A	316	VAL
1	A	328	SER
1	A	340	LYS
1	A	346	SER
1	A	350	ILE
1	A	376	GLU
1	A	424	ARG
1	A	434	THR
1	A	459	SER
1	A	464	ILE
1	A	482	ILE
1	A	510	GLU
1	A	538	VAL
1	A	596	CYS
1	A	603	SER

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Mol	Chain	Res	Type
1	A	610	THR
1	A	614	LYS
1	A	629	ASP
1	A	718	PHE
1	A	725	ILE
1	A	770	ASP
1	A	801	GLN
1	A	811	VAL
1	A	829	HIS
1	A	834	GLU
1	A	892	ILE
1	A	897	VAL
1	B	256	LEU
1	B	273	LEU
1	B	294	SER
1	B	350	ILE
1	B	417	ASP
1	B	427	GLU
1	B	474	THR
1	B	483	THR
1	B	505	LYS
1	B	558	VAL
1	B	559	ASN
1	B	596	CYS
1	B	603	SER
1	B	610	THR
1	B	614	LYS
1	B	629	ASP
1	B	694	SER
1	B	718	PHE
1	B	792	THR
1	B	811	VAL
1	B	813	MET
1	B	841	THR
1	B	846	SER
1	B	871	GLU

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

Mol	Chain	Res	Type
1	A	111	ASN
1	A	123	ASN

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Mol	Chain	Res	Type
1	A	200	GLN
1	A	213	ASN
1	A	403	GLN
1	A	460	HIS
1	A	550	HIS
1	A	580	HIS
1	A	657	GLN
1	A	673	GLN
1	A	829	HIS
1	A	840	HIS
1	A	882	HIS
1	B	60	GLN
1	B	89	ASN
1	B	111	ASN
1	B	200	GLN
1	B	306	ASN
1	B	423	GLN
1	B	550	HIS
1	B	654	HIS
1	B	673	GLN
1	B	829	HIS
1	B	840	HIS
1	B	858	GLN
1	B	862	GLN
1	B	879	HIS
1	B	882	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

6 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$
3	PO4	B	902	-	4,4,4	0.80	0	6,6,6	0.57	0
2	TMO	B	901	-	4,4,4	1.24	0	6,6,6	0.51	0
4	S9W	B	903	1	33,35,35	0.88	1 (3%)	41,48,48	1.92	10 (24%)
3	PO4	A	902	-	4,4,4	1.04	0	6,6,6	0.41	0
4	S9W	A	903	1	33,35,35	1.01	2 (6%)	41,48,48	1.98	11 (26%)
2	TMO	A	901	-	4,4,4	1.12	0	6,6,6	0.49	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	S9W	A	903	1	-	3/20/41/41	0/4/4/4
4	S9W	B	903	1	-	3/20/41/41	0/4/4/4

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	903	S9W	O33-C10	-2.80	1.18	1.22
4	B	903	S9W	C13-C12	2.75	1.43	1.39
4	A	903	S9W	C12-N11	-2.14	1.37	1.41

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	903	S9W	BR1-C29-C30	5.81	127.21	119.30
4	B	903	S9W	C2-C10-N11	4.98	122.25	115.68
4	B	903	S9W	BR1-C29-C28	-4.91	112.61	119.30
4	A	903	S9W	BR1-C29-C28	-4.77	112.79	119.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	903	S9W	C2-C10-N11	4.38	121.45	115.68
4	B	903	S9W	BR1-C29-C30	4.32	125.18	119.30
4	B	903	S9W	O5-C6-C7	-3.67	107.84	111.50
4	A	903	S9W	C25-C26-C27	-3.16	114.93	120.75
4	B	903	S9W	C31-C30-C29	-2.94	115.66	119.18
4	B	903	S9W	C7-C2-C3	2.50	111.32	108.15
4	A	903	S9W	C27-C28-C29	2.44	122.10	119.18
4	B	903	S9W	C25-C26-C27	-2.28	116.55	120.75
4	A	903	S9W	C20-C21-N22	2.27	115.23	110.65
4	A	903	S9W	C7-C2-C3	-2.24	105.33	108.15
4	A	903	S9W	C18-N19-C24	-2.19	106.36	111.07
4	A	903	S9W	C17-C12-N11	-2.13	113.26	120.41
4	B	903	S9W	C18-C14-C15	-2.11	116.86	120.75
4	A	903	S9W	C25-C26-C31	2.11	124.62	120.75
4	B	903	S9W	C13-C12-N11	2.10	126.95	120.13
4	A	903	S9W	C13-C12-N11	2.06	126.84	120.13
4	B	903	S9W	C17-C12-N11	-2.01	113.68	120.41

There are no chirality outliers.

All (6) torsion outliers are listed below:

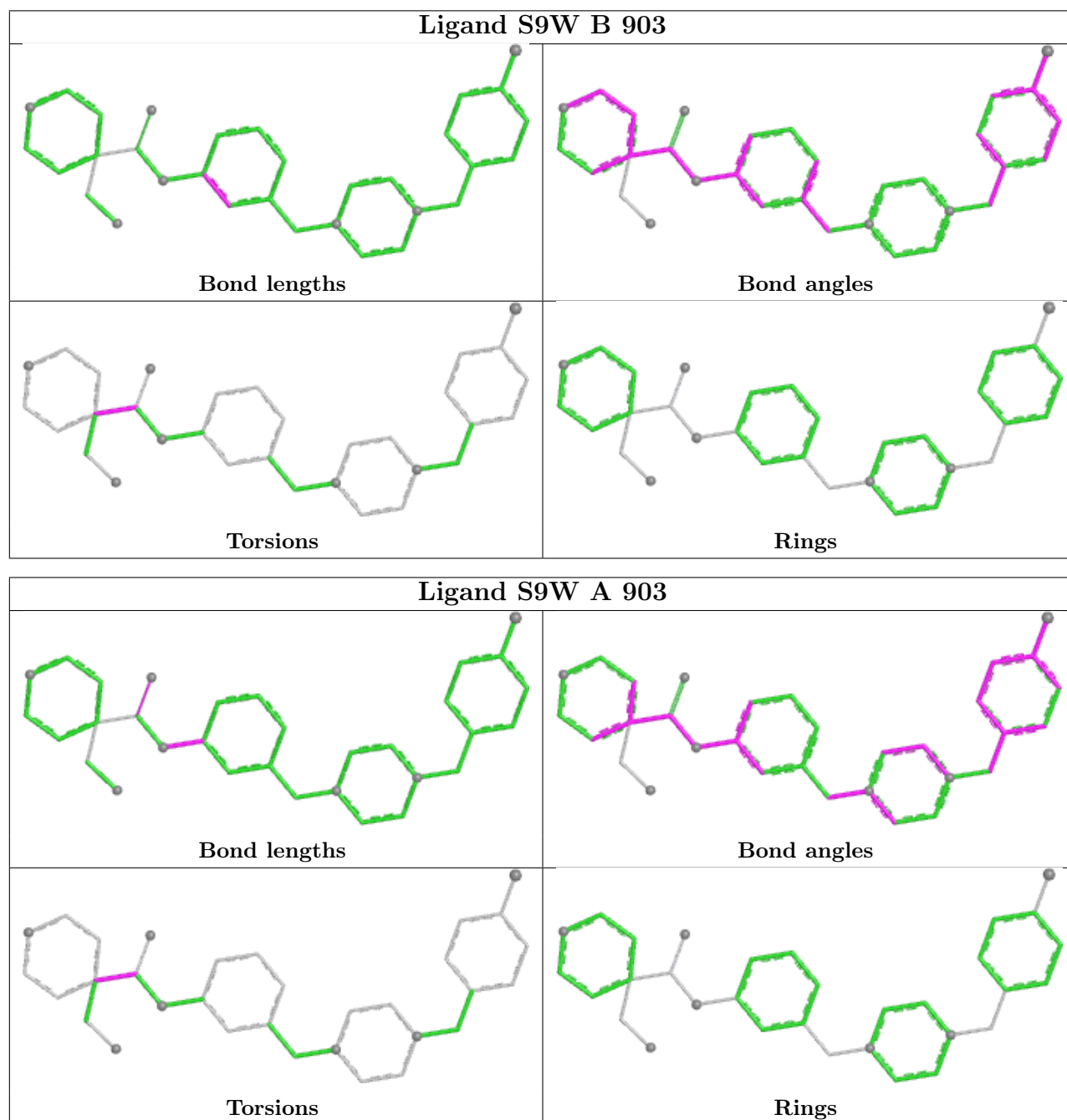
Mol	Chain	Res	Type	Atoms
4	A	903	S9W	N11-C10-C2-C8
4	B	903	S9W	N11-C10-C2-C8
4	A	903	S9W	N11-C10-C2-C7
4	B	903	S9W	N11-C10-C2-C7
4	A	903	S9W	O33-C10-C2-C7
4	B	903	S9W	O33-C10-C2-C7

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier.

The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	802/898 (89%)	-0.10	18 (2%) 62 59	45, 74, 120, 174	1 (0%)
1	B	804/898 (89%)	0.08	30 (3%) 45 41	58, 81, 128, 175	0
All	All	1606/1796 (89%)	-0.01	48 (2%) 52 49	45, 78, 125, 175	1 (0%)

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	897	VAL	5.7
1	A	48	LEU	5.2
1	B	48	LEU	4.4
1	A	869	VAL	4.2
1	A	595	HIS	4.2
1	A	138	LEU	4.1
1	B	869	VAL	4.0
1	A	317	THR	3.8
1	B	861	PRO	3.8
1	B	862	GLN	3.7
1	B	76	MET	3.7
1	A	259	MET	3.2
1	B	139	PHE	3.2
1	A	830	GLY	3.2
1	B	316	VAL	3.2
1	B	419	VAL	3.2
1	B	831	PHE	3.1
1	B	830	GLY	3.0
1	A	862	GLN	3.0
1	B	259	MET	3.0
1	B	317	THR	3.0
1	B	164	VAL	2.9
1	B	438	ILE	2.9
1	B	501	PRO	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	419	VAL	2.8
1	A	487	LYS	2.7
1	A	897	VAL	2.7
1	B	466	PHE	2.7
1	A	77	ALA	2.7
1	A	835	ASN	2.7
1	B	484	SER	2.6
1	B	832	LEU	2.5
1	A	140	GLN	2.4
1	A	834	GLU	2.3
1	B	505	LYS	2.3
1	B	628	PRO	2.3
1	A	70	LYS	2.3
1	B	860	TYR	2.3
1	A	316	VAL	2.3
1	B	595	HIS	2.2
1	B	835	ASN	2.2
1	B	487	LYS	2.2
1	B	195	GLN	2.2
1	B	489	SER	2.1
1	B	78	LYS	2.1
1	B	110	GLU	2.0
1	B	289	ALA	2.0
1	A	195	GLN	2.0

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

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

6.3 Carbohydrates [i](#)

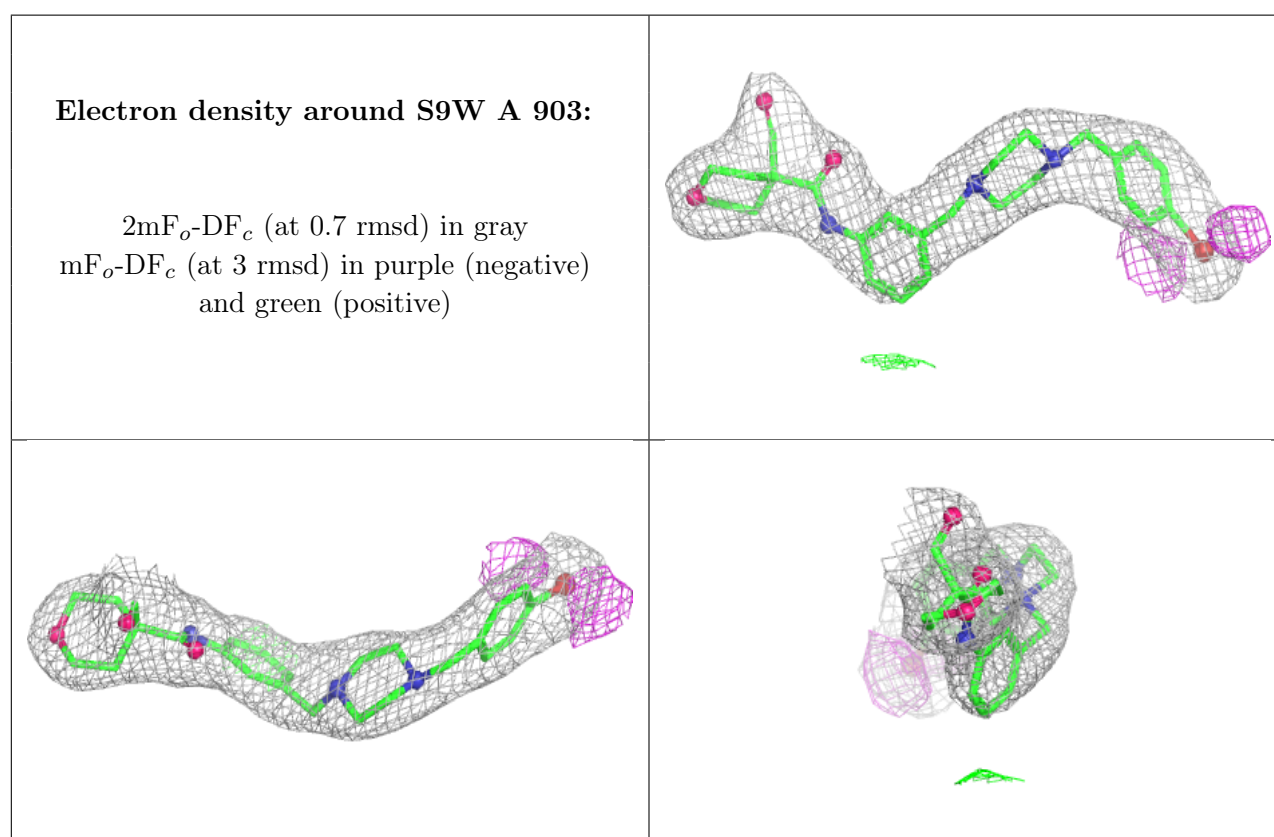
There are no oligosaccharides in this entry.

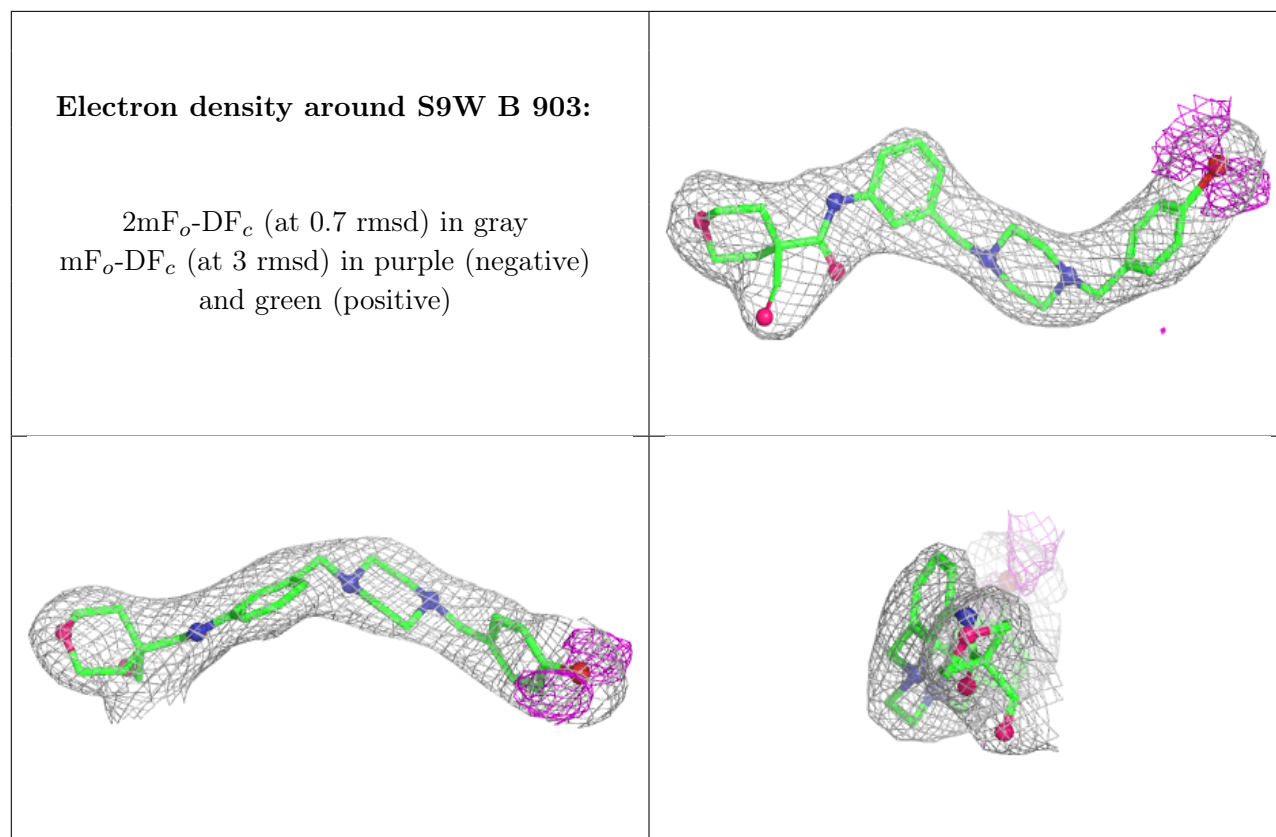
6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	PO4	A	902	5/5	0.84	0.10	103,108,126,129	0
3	PO4	B	902	5/5	0.87	0.08	91,99,111,132	0
4	S9W	A	903	32/32	0.95	0.11	61,81,97,127	0
4	S9W	B	903	32/32	0.96	0.10	62,79,96,113	0
2	TMO	A	901	5/5	0.98	0.12	73,77,79,80	0
2	TMO	B	901	5/5	0.98	0.10	71,76,82,82	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 [i](#)

There are no such residues in this entry.