



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:55 PM UTC

PDB ID : 3Q4C / pdb_00003q4c
Title : Crystal Structure of Wild Type BRAF kinase domain in complex with organometallic inhibitor CNS292
Authors : Xie, P.; Streu, C.; Qin, J.; Pregman, H.; Pagano, N.; Meggers, E.; Marmorstein, R.
Deposited on : 2010-12-23
Resolution : 3.20 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

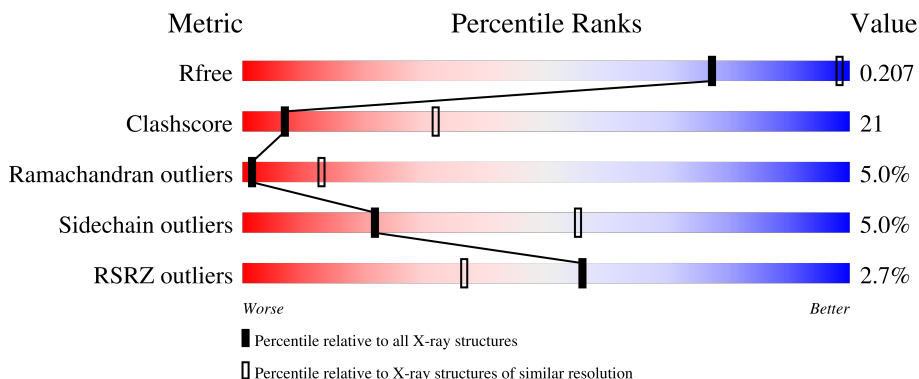
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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	307	
1	B	307	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	RSW	A	1	-	-	-	X

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4250 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.

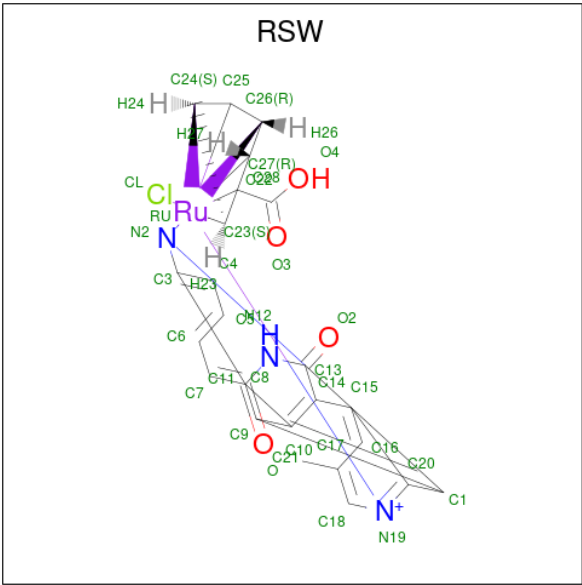
- Molecule 1 is a protein called Serine/threonine-protein kinase B-raf.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	264	Total	C	N	O	S	0	0	0
			2108	1352	367	376	13			
1	B	264	Total	C	N	O	S	0	0	0
			2108	1352	367	376	13			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	419	MET	-	expression tag	UNP P15056
A	420	ASP	-	expression tag	UNP P15056
A	421	ARG	-	expression tag	UNP P15056
A	422	GLY	-	expression tag	UNP P15056
A	423	SER	-	expression tag	UNP P15056
A	424	HIS	-	expression tag	UNP P15056
A	425	HIS	-	expression tag	UNP P15056
A	426	HIS	-	expression tag	UNP P15056
A	427	HIS	-	expression tag	UNP P15056
A	428	HIS	-	expression tag	UNP P15056
A	429	HIS	-	expression tag	UNP P15056
A	430	GLY	-	expression tag	UNP P15056
B	419	MET	-	expression tag	UNP P15056
B	420	ASP	-	expression tag	UNP P15056
B	421	ARG	-	expression tag	UNP P15056
B	422	GLY	-	expression tag	UNP P15056
B	423	SER	-	expression tag	UNP P15056
B	424	HIS	-	expression tag	UNP P15056
B	425	HIS	-	expression tag	UNP P15056
B	426	HIS	-	expression tag	UNP P15056
B	427	HIS	-	expression tag	UNP P15056
B	428	HIS	-	expression tag	UNP P15056
B	429	HIS	-	expression tag	UNP P15056
B	430	GLY	-	expression tag	UNP P15056

- Molecule 2 is [(1,2,3,4,5,6-eta)-(1S,2R,3R,4R,5S,6S)-1-carboxycyclohexane-1,2,3,4,5,6-hexayl](chloro)(3-methyl-5,7-dioxo-6,7-dihydro-5H-pyrido[2,3-a]pyrrolo[3,4-c]carbazol-12-ide-kappa a 2 N 1 ,N 12)ruthenium(1+) (CCD ID: RSW) (formula: C₂₅H₁₆ClN₃O₄Ru).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	C	Cl	N	O	Ru		
2	A	1	34	25	1	3	4	1	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	94.91Å 94.91Å 163.35Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 3.20 50.00 – 3.20	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-3.20) 99.1 (50.00-3.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.59 (at 3.07Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.232 , 0.280 0.208 , 0.207	Depositor DCC
R_{free} test set	712 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	51.0	Xtriage
Anisotropy	0.126	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 43.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	4250	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.06% 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: RSW

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	0.48	0/2153	0.98	11/2905 (0.4%)
1	B	0.51	0/2153	1.00	8/2905 (0.3%)
All	All	0.50	0/4306	0.99	19/5810 (0.3%)

There are no bond length outliers.

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	455	GLN	N-CA-C	-9.72	102.22	114.56
1	B	540	LEU	N-CA-C	6.48	118.34	111.28
1	A	641	PHE	N-CA-C	-6.10	104.74	111.82
1	A	540	LEU	N-CA-C	6.03	117.65	111.14
1	A	550	ILE	N-CA-C	-5.95	106.69	111.81
1	B	448	ASP	N-CA-C	5.77	123.08	110.80
1	A	541	HIS	N-CA-C	5.65	121.07	113.72
1	A	645	LEU	N-CA-C	-5.65	105.12	111.28
1	A	453	ASP	N-CA-C	5.63	118.39	110.23
1	B	521	LYS	CA-C-N	-5.44	113.04	119.84
1	B	521	LYS	C-N-CA	-5.44	113.04	119.84
1	A	521	LYS	CA-C-N	-5.40	113.08	119.84
1	A	521	LYS	C-N-CA	-5.40	113.08	119.84
1	A	483	MET	N-CA-C	5.37	117.42	109.69
1	A	485	ASN	N-CA-C	5.18	117.83	111.82
1	B	470	VAL	N-CA-C	5.16	116.39	108.71
1	B	568	ALA	N-CA-C	-5.09	106.33	112.54
1	B	543	ILE	N-CA-C	-5.02	107.57	112.29
1	A	496	ALA	N-CA-C	-5.02	105.26	111.33

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	2108	0	2153	94	0
1	B	2108	0	2153	91	0
2	A	34	0	10	0	0
All	All	4250	0	4316	181	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 21.

All (181) 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:667:MET:HE2	1:A:673:LEU:HB2	1.45	0.98
1:A:550:ILE:HD12	1:A:720:LEU:HD21	1.52	0.89
1:A:720:LEU:HB3	1:A:721:PRO:HA	1.54	0.88
1:B:455:GLN:HE21	1:B:475:TRP:HE1	1.14	0.87
1:A:635:GLN:NE2	1:A:705:LEU:HD21	1.93	0.84
1:B:453:ASP:OD1	1:B:521:LYS:HG3	1.80	0.82
1:B:720:LEU:HB3	1:B:721:PRO:O	1.80	0.81
1:A:716:LEU:HD12	1:A:716:LEU:O	1.81	0.79
1:B:483:MET:HE2	1:B:523:GLN:HB2	1.65	0.77
1:A:455:GLN:HE21	1:A:475:TRP:HE1	1.31	0.75
1:B:566:LEU:HD22	1:B:571:ILE:HG21	1.70	0.74
1:B:455:GLN:NE2	1:B:475:TRP:HE1	1.85	0.73
1:B:484:LEU:CD1	1:B:486:VAL:HG22	2.19	0.72
1:A:630:ASN:N	1:A:631:PRO:HD3	2.05	0.72
1:A:550:ILE:CD1	1:A:720:LEU:HD21	2.19	0.72
1:B:484:LEU:HD12	1:B:486:VAL:HG22	1.74	0.70
1:B:453:ASP:CG	1:B:521:LYS:HG3	2.16	0.70
1:A:585:GLU:HG2	1:B:587:LEU:HD21	1.73	0.69
1:B:635:GLN:NE2	1:B:705:LEU:HD21	2.07	0.69
1:A:484:LEU:HD12	1:A:524:LEU:HD12	1.74	0.68
1:A:522:PRO:O	1:A:523:GLN:HG3	1.93	0.68
1:A:451:ILE:HD12	1:A:456:ILE:HD11	1.74	0.68
1:A:720:LEU:HB3	1:A:721:PRO:CA	2.25	0.67
1:A:551:LYS:O	1:A:555:ILE:HG13	1.94	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:503:VAL:HG22	1:B:599:VAL:HG23	1.77	0.66
1:B:663:GLN:O	1:B:667:MET:HG3	1.98	0.64
1:B:673:LEU:HD12	1:B:674:SER:H	1.63	0.63
1:B:550:ILE:H	1:B:550:ILE:HD13	1.63	0.63
1:B:521:LYS:HB2	1:B:522:PRO:HD3	1.80	0.62
1:B:550:ILE:HD13	1:B:550:ILE:N	2.14	0.62
1:B:619:MET:HE2	1:B:623:VAL:HG12	1.81	0.62
1:A:585:GLU:O	1:A:587:LEU:HD22	2.00	0.62
1:A:455:GLN:NE2	1:A:475:TRP:HE1	1.97	0.60
1:A:521:LYS:HB2	1:A:522:PRO:HD3	1.82	0.60
1:B:450:GLU:HB2	1:B:518:TYR:CZ	2.37	0.60
1:B:630:ASN:N	1:B:631:PRO:HD3	2.16	0.60
1:A:532:GLU:H	1:A:584:HIS:CD2	2.19	0.59
1:B:514:LEU:H	1:B:529:GLN:HE21	1.47	0.59
1:B:453:ASP:CB	1:B:521:LYS:HG3	2.33	0.59
1:A:514:LEU:H	1:A:529:GLN:HE21	1.49	0.59
1:B:616:ILE:HG22	1:B:619:MET:SD	2.43	0.58
1:A:490:THR:OG1	1:A:493:GLN:HG3	2.04	0.58
1:A:629:LYS:HA	1:A:629:LYS:HE2	1.85	0.58
1:B:503:VAL:HG22	1:B:599:VAL:CG2	2.34	0.58
1:B:566:LEU:HD22	1:B:571:ILE:CG2	2.34	0.57
1:A:486:VAL:O	1:A:486:VAL:HG23	2.04	0.57
1:B:572:ILE:HG22	1:B:574:ARG:CG	2.35	0.57
1:A:557:ARG:HH11	1:A:557:ARG:HB3	1.70	0.57
1:A:557:ARG:HB3	1:A:557:ARG:NH1	2.21	0.56
1:A:467:PHE:HD2	1:A:467:PHE:H	1.53	0.56
1:A:547:PHE:HB3	1:A:551:LYS:HB3	1.87	0.56
1:A:690:ARG:O	1:A:694:GLU:HG3	2.05	0.56
1:B:550:ILE:HD12	1:B:720:LEU:HD21	1.88	0.55
1:A:500:GLU:OE1	1:A:595:GLY:HA2	2.07	0.55
1:A:568:ALA:C	1:A:570:SER:H	2.13	0.55
1:A:648:LEU:HB3	1:A:649:MET:HE2	1.88	0.55
1:A:532:GLU:H	1:A:584:HIS:HD2	1.55	0.55
1:B:520:THR:HG22	1:B:524:LEU:CD2	2.38	0.54
1:A:531:CYS:SG	1:A:582:PHE:HB3	2.48	0.54
1:B:453:ASP:HB3	1:B:521:LYS:HG3	1.88	0.54
1:B:577:LYS:HD2	1:B:615:SER:HB2	1.90	0.53
1:A:508:ARG:HD3	1:B:515:PHE:O	2.08	0.53
1:B:621:PRO:O	1:B:624:ILE:HG22	2.09	0.53
1:B:720:LEU:HB3	1:B:721:PRO:C	2.33	0.53
1:B:572:ILE:HG22	1:B:574:ARG:HG3	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:627:GLN:HG3	1:B:628:ASP:OD1	2.09	0.52
1:A:718:ARG:O	1:A:720:LEU:N	2.43	0.52
1:B:490:THR:HG23	1:B:493:GLN:OE1	2.10	0.52
1:A:680:VAL:HG21	1:A:689:LYS:HD2	1.91	0.52
1:A:585:GLU:HB3	1:A:587:LEU:CD2	2.40	0.52
1:A:643:ILE:O	1:A:647:GLU:HG2	2.10	0.51
1:B:619:MET:HE2	1:B:623:VAL:CG1	2.40	0.51
1:A:492:GLN:O	1:A:495:GLN:HB2	2.11	0.51
1:B:550:ILE:H	1:B:550:ILE:CD1	2.22	0.51
1:B:673:LEU:HD12	1:B:674:SER:N	2.23	0.51
1:A:599:VAL:O	1:A:599:VAL:HG12	2.11	0.51
1:A:660:ASN:O	1:A:664:ILE:HG13	2.11	0.51
1:B:501:VAL:O	1:B:505:ARG:HG2	2.11	0.51
1:B:475:TRP:O	1:B:476:HIS:HB2	2.11	0.50
1:A:484:LEU:CD2	1:A:486:VAL:HG22	2.41	0.50
1:A:625:ARG:O	1:A:626:MET:C	2.54	0.50
1:B:484:LEU:HD11	1:B:486:VAL:HG22	1.92	0.50
1:A:621:PRO:HD3	1:A:639:TYR:CE1	2.46	0.50
1:B:549:MET:O	1:B:553:ILE:HG13	2.11	0.50
1:B:616:ILE:HA	1:B:619:MET:HG3	1.93	0.49
1:A:638:VAL:HG13	1:A:709:ILE:HD11	1.95	0.49
1:A:464:SER:HA	1:A:469:THR:HA	1.95	0.49
1:B:453:ASP:HB3	1:B:521:LYS:CB	2.43	0.49
1:A:521:LYS:O	1:A:522:PRO:C	2.54	0.49
1:B:667:MET:HE2	1:B:672:TYR:HB2	1.94	0.48
1:B:722:LYS:HG2	1:B:722:LYS:O	2.14	0.48
1:A:572:ILE:HG22	1:A:574:ARG:HG2	1.96	0.48
1:B:483:MET:CE	1:B:523:GLN:HB2	2.39	0.48
1:B:548:GLU:H	1:B:548:GLU:CD	2.21	0.48
1:A:509:HIS:HB2	1:A:565:TYR:CE2	2.49	0.47
1:A:616:ILE:HG22	1:A:619:MET:SD	2.55	0.47
1:B:594:PHE:HB3	1:B:597:ALA:HB3	1.97	0.47
1:B:464:SER:HA	1:B:469:THR:HA	1.97	0.47
1:B:487:THR:C	1:B:489:PRO:HD3	2.40	0.47
1:B:635:GLN:HE21	1:B:705:LEU:HD21	1.76	0.47
1:A:585:GLU:CG	1:B:587:LEU:HD11	2.45	0.47
1:B:573:HIS:O	1:B:574:ARG:HB2	2.14	0.47
1:B:577:LYS:C	1:B:579:ASN:H	2.22	0.46
1:B:594:PHE:HD1	1:B:597:ALA:HB2	1.79	0.46
1:A:630:ASN:N	1:A:631:PRO:CD	2.76	0.46
1:A:474:LYS:HD2	1:A:474:LYS:C	2.40	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:577:LYS:C	1:A:579:ASN:H	2.23	0.46
1:B:572:ILE:HG22	1:B:574:ARG:HG2	1.97	0.46
1:A:722:LYS:O	1:A:722:LYS:HG2	2.14	0.46
1:B:554:ASP:OD1	1:B:557:ARG:NH1	2.49	0.46
1:B:494:LEU:HD12	1:B:494:LEU:O	2.15	0.46
1:A:460:GLN:HG3	1:A:461:ARG:N	2.31	0.46
1:B:573:HIS:HE1	1:B:593:ASP:O	1.99	0.45
1:B:661:ARG:O	1:B:665:ILE:HG13	2.15	0.45
1:A:699:LYS:HE3	1:A:702:GLU:OE1	2.15	0.45
1:A:464:SER:CB	1:A:469:THR:HG22	2.47	0.45
1:B:705:LEU:O	1:B:709:ILE:HG13	2.15	0.45
1:A:522:PRO:C	1:A:523:GLN:HG3	2.41	0.45
1:B:532:GLU:H	1:B:584:HIS:CD2	2.34	0.45
1:A:549:MET:HE3	1:A:684:CYS:HA	1.97	0.45
1:A:568:ALA:C	1:A:570:SER:N	2.74	0.45
1:A:643:ILE:O	1:A:646:TYR:HB3	2.17	0.45
1:B:591:ILE:HG22	1:B:592:GLY:N	2.31	0.45
1:A:583:LEU:HD23	1:A:583:LEU:HA	1.71	0.45
1:B:573:HIS:CE1	1:B:593:ASP:O	2.70	0.45
1:A:511:ASN:ND2	1:A:558:GLN:HB3	2.32	0.45
1:B:599:VAL:HG12	1:B:599:VAL:O	2.17	0.45
1:A:696:LEU:O	1:A:697:LYS:C	2.61	0.44
1:B:537:TYR:N	1:B:578:SER:O	2.50	0.44
1:A:521:LYS:HA	1:A:521:LYS:HD3	1.84	0.44
1:A:515:PHE:O	1:B:508:ARG:HD3	2.17	0.44
1:A:599:VAL:O	1:A:600:LYS:C	2.61	0.44
1:B:635:GLN:CD	1:B:635:GLN:H	2.25	0.44
1:B:471:TYR:CD2	1:B:471:TYR:N	2.85	0.44
1:B:493:GLN:C	1:B:495:GLN:N	2.74	0.44
1:A:626:MET:O	1:A:627:GLN:C	2.61	0.44
1:A:562:GLY:O	1:A:565:TYR:HB3	2.17	0.43
1:B:513:LEU:HD12	1:B:513:LEU:HA	1.89	0.43
1:A:577:LYS:HD2	1:A:618:TRP:CE3	2.54	0.43
1:A:706:PHE:O	1:A:707:PRO:C	2.59	0.43
1:A:493:GLN:C	1:A:495:GLN:N	2.76	0.43
1:B:646:TYR:O	1:B:650:THR:HG23	2.18	0.43
1:B:449:TRP:HD1	1:B:505:ARG:NH2	2.16	0.43
1:B:667:MET:HE2	1:B:672:TYR:CB	2.48	0.43
1:A:584:HIS:HB3	1:A:588:THR:HB	2.01	0.43
1:B:578:SER:HB3	1:B:647:GLU:OE1	2.19	0.43
1:A:616:ILE:HD12	1:A:617:LEU:N	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:550:ILE:N	1:B:550:ILE:CD1	2.80	0.42
1:A:451:ILE:O	1:A:452:PRO:C	2.62	0.42
1:A:537:TYR:CD1	1:A:541:HIS:HD2	2.38	0.42
1:A:577:LYS:CD	1:A:615:SER:HB2	2.49	0.42
1:B:629:LYS:HD3	1:B:629:LYS:HA	1.75	0.42
1:B:554:ASP:O	1:B:558:GLN:HG3	2.19	0.42
1:B:660:ASN:OD1	1:B:662:ASP:HB2	2.19	0.42
1:B:720:LEU:N	1:B:721:PRO:HA	2.34	0.42
1:A:475:TRP:O	1:A:476:HIS:HB2	2.20	0.42
1:A:559:THR:HG23	1:A:591:ILE:HD11	2.01	0.42
1:A:629:LYS:C	1:A:631:PRO:HD3	2.45	0.42
1:A:620:ALA:HB1	1:A:622:GLU:OE1	2.19	0.41
1:A:481:VAL:HG22	1:A:527:VAL:HG22	2.03	0.41
1:A:661:ARG:O	1:A:665:ILE:HG13	2.20	0.41
1:A:712:SER:C	1:A:714:GLU:N	2.78	0.41
1:B:464:SER:O	1:B:467:PHE:HE2	2.03	0.41
1:B:532:GLU:H	1:B:584:HIS:HD2	1.68	0.41
1:B:547:PHE:HB2	1:B:552:LEU:CD2	2.50	0.41
1:B:715:LEU:HD23	1:B:715:LEU:C	2.45	0.41
1:A:633:SER:O	1:A:634:PHE:C	2.63	0.41
1:A:646:TYR:CD2	1:A:654:PRO:HG3	2.55	0.41
1:A:557:ARG:HH12	1:A:558:GLN:HG2	1.86	0.41
1:A:511:ASN:HD21	1:A:558:GLN:HB3	1.86	0.41
1:A:536:LEU:HD21	1:A:648:LEU:HD21	2.02	0.41
1:A:639:TYR:CD2	1:A:639:TYR:C	2.98	0.41
1:B:621:PRO:O	1:B:625:ARG:HB2	2.20	0.41
1:B:646:TYR:CD2	1:B:654:PRO:HG3	2.57	0.40
1:A:497:PHE:O	1:A:498:LYS:C	2.64	0.40
1:B:455:GLN:HE21	1:B:475:TRP:NE1	1.96	0.40
1:A:577:LYS:C	1:A:579:ASN:N	2.79	0.40
1:B:463:GLY:O	1:B:464:SER:HB3	2.22	0.40
1:A:619:MET:HE2	1:A:623:VAL:HG12	2.03	0.40
1:A:709:ILE:O	1:A:713:ILE:HG13	2.22	0.40
1:B:495:GLN:OE1	1:B:495:GLN:HA	2.21	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	260/307 (85%)	218 (84%)	29 (11%)	13 (5%)	1	13
1	B	260/307 (85%)	227 (87%)	20 (8%)	13 (5%)	1	13
All	All	520/614 (85%)	445 (86%)	49 (9%)	26 (5%)	1	13

All (26) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	627	GLN
1	A	721	PRO
1	B	448	ASP
1	B	464	SER
1	B	721	PRO
1	A	681	ARG
1	B	586	ASP
1	B	593	ASP
1	B	627	GLN
1	B	718	ARG
1	A	448	ASP
1	A	464	SER
1	A	700	ARG
1	A	719	SER
1	B	575	ASP
1	B	701	ASP
1	A	487	THR
1	A	626	MET
1	A	701	ASP
1	A	497	PHE
1	B	454	GLY
1	A	486	VAL
1	A	720	LEU
1	B	630	ASN
1	B	522	PRO

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Mol	Chain	Res	Type
1	B	489	PRO

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	232/271 (86%)	223 (96%)	9 (4%)	28	62
1	B	232/271 (86%)	218 (94%)	14 (6%)	17	50
All	All	464/542 (86%)	441 (95%)	23 (5%)	22	55

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	447	ASP
1	A	467	PHE
1	A	484	LEU
1	A	499	ASN
1	A	540	LEU
1	A	548	GLU
1	A	587	LEU
1	A	624	ILE
1	A	716	LEU
1	B	453	ASP
1	B	457	THR
1	B	467	PHE
1	B	471	TYR
1	B	474	LYS
1	B	484	LEU
1	B	486	VAL
1	B	495	GLN
1	B	540	LEU
1	B	548	GLU
1	B	550	ILE
1	B	587	LEU
1	B	628	ASP

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Mol	Chain	Res	Type
1	B	716	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	455	GLN
1	A	511	ASN
1	A	529	GLN
1	A	538	HIS
1	A	558	GLN
1	A	579	ASN
1	A	580	ASN
1	A	584	HIS
1	B	455	GLN
1	B	485	ASN
1	B	529	GLN
1	B	538	HIS
1	B	561	GLN
1	B	579	ASN
1	B	584	HIS

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	RSW	A	1	-	30,45,45	3.36	20 (66%)	35,106,106	3.09	13 (37%)

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	RSW	A	1	-	-	0/0/152/152	0/6/12/12

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1	RSW	RU-N2	-7.58	1.91	2.08
2	A	1	RSW	C18-N19	5.89	1.43	1.35
2	A	1	RSW	C9-C10	5.33	1.50	1.40
2	A	1	RSW	C14-C13	-5.09	1.41	1.49
2	A	1	RSW	C16-C17	4.55	1.45	1.37
2	A	1	RSW	C4-C3	4.17	1.46	1.39
2	A	1	RSW	C6-C7	4.15	1.46	1.38
2	A	1	RSW	C8-C3	3.89	1.47	1.41
2	A	1	RSW	C5-C4	3.64	1.45	1.38
2	A	1	RSW	C7-C8	3.42	1.45	1.40
2	A	1	RSW	C3-N2	3.40	1.45	1.39
2	A	1	RSW	C14-C10	3.34	1.44	1.38
2	A	1	RSW	C18-C17	3.09	1.45	1.39
2	A	1	RSW	C15-C20	3.09	1.47	1.42
2	A	1	RSW	C6-C5	3.00	1.44	1.38
2	A	1	RSW	C14-C15	2.57	1.48	1.43
2	A	1	RSW	C20-N19	2.42	1.40	1.37
2	A	1	RSW	C10-C11	-2.39	1.41	1.49
2	A	1	RSW	O2-C13	2.37	1.28	1.23
2	A	1	RSW	O-C11	2.31	1.28	1.23

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1	RSW	C14-C13-N12	9.82	114.41	105.83
2	A	1	RSW	C13-N12-C11	-9.05	104.72	112.59
2	A	1	RSW	C15-C14-C13	6.97	135.43	129.01
2	A	1	RSW	C15-C20-N19	4.10	125.28	122.26
2	A	1	RSW	O2-C13-C14	-3.93	123.50	129.10
2	A	1	RSW	C16-C15-C14	3.19	128.87	123.21
2	A	1	RSW	O4-C28-C22	3.11	122.85	113.69
2	A	1	RSW	C16-C15-C20	-3.05	115.21	119.49
2	A	1	RSW	O4-C28-O3	-2.48	115.92	123.86
2	A	1	RSW	C9-C1-C20	2.38	126.52	124.32
2	A	1	RSW	C9-C8-C3	2.32	108.83	106.49
2	A	1	RSW	C8-C9-C10	2.32	142.22	134.87
2	A	1	RSW	O-C11-C10	-2.18	125.31	128.47

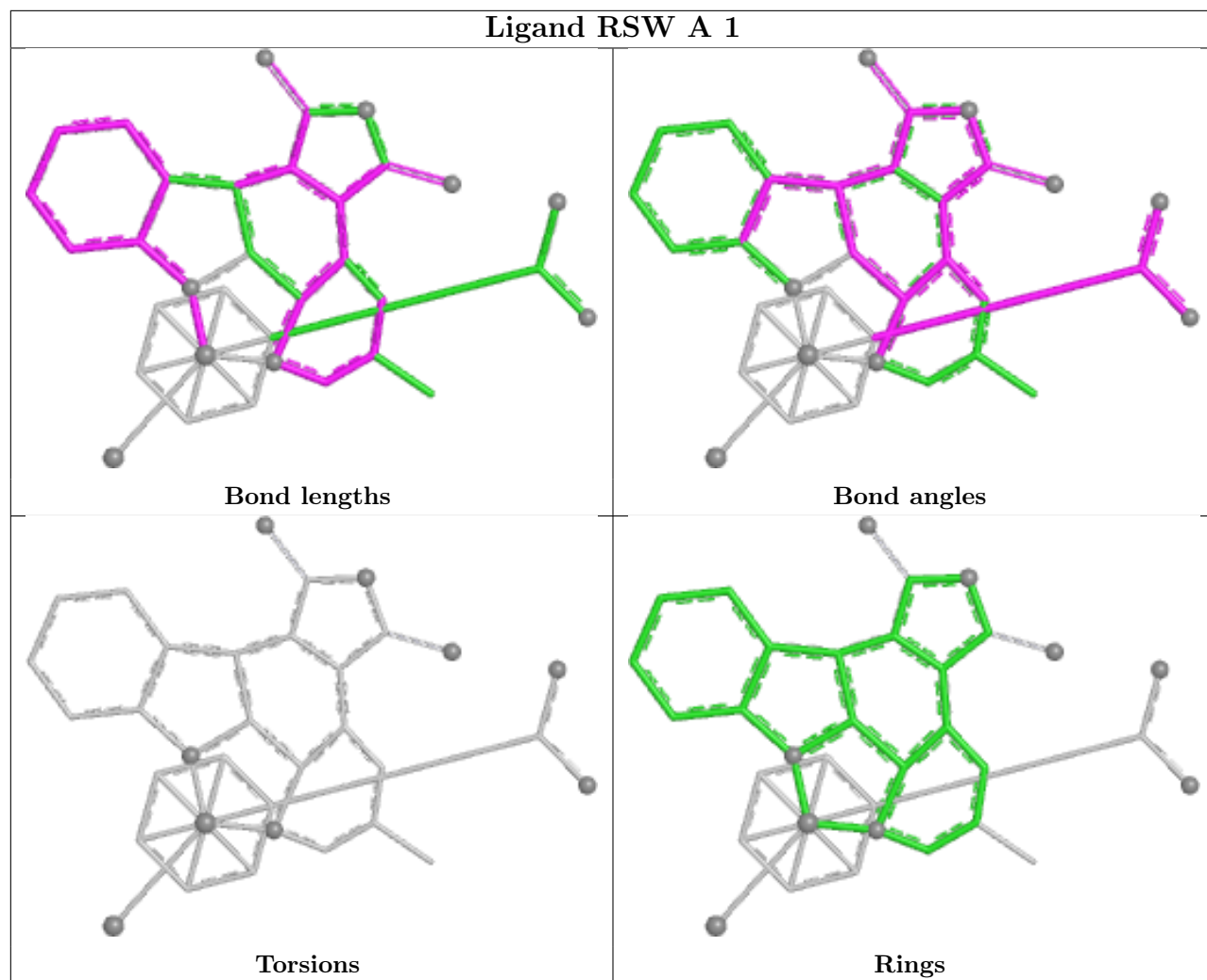
There are no chirality outliers.

There are no torsion outliers.

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 [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	264/307 (85%)	0.05	6 (2%) 61 41	16, 41, 81, 104	0
1	B	264/307 (85%)	0.02	8 (3%) 52 33	18, 40, 77, 133	0
All	All	528/614 (85%)	0.04	14 (2%) 56 36	16, 41, 78, 133	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	465	GLY	4.1
1	B	466	SER	3.9
1	B	720	LEU	3.7
1	A	720	LEU	3.3
1	A	721	PRO	2.9
1	A	626	MET	2.8
1	B	721	PRO	2.8
1	B	467	PHE	2.7
1	A	467	PHE	2.5
1	A	629	LYS	2.5
1	B	626	MET	2.4
1	B	624	ILE	2.2
1	A	466	SER	2.2
1	B	722	LYS	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 ⓘ

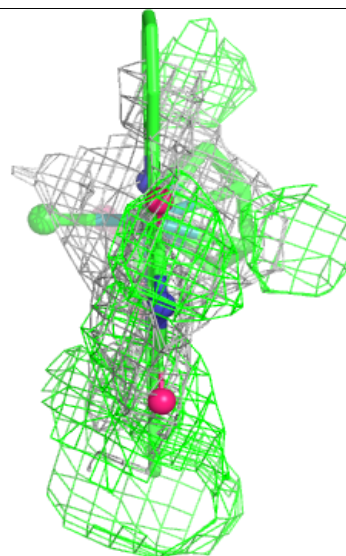
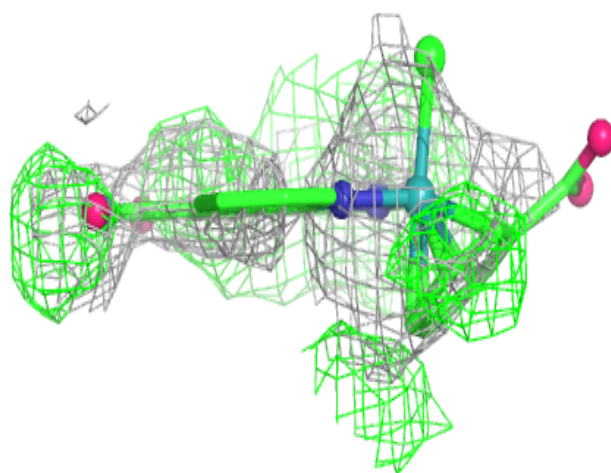
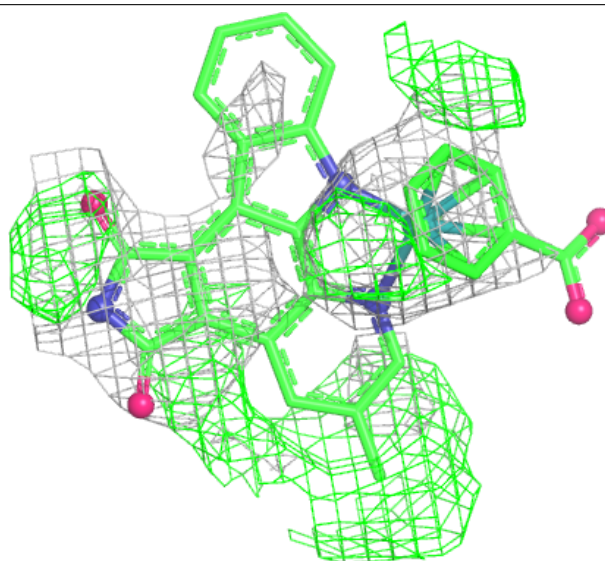
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists 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	RSW	A	1	34/34	0.76	0.42	84,89,92,93	34

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 RSW A 1:

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.