



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:49 PM UTC

PDB ID : 4UW0 / pdb_00004uw0
Title : Low resolution structure of WbdD with C-terminal bundle ordered to residue 505
Authors : Hagelueken, G.; Huang, H.; Naismith, J.H.
Deposited on : 2014-08-08
Resolution : 3.87 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

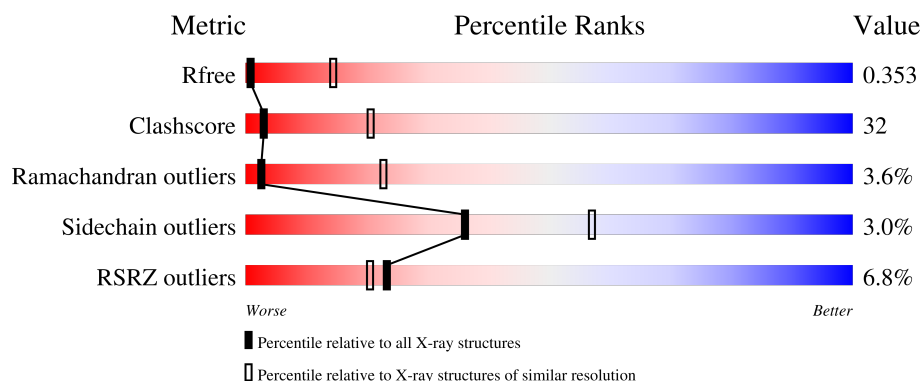
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.87 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	1186 (4.06-3.70)
Clashscore	190562	1012 (4.04-3.72)
Ramachandran outliers	187476	1167 (4.06-3.70)
Sidechain outliers	187428	1160 (4.06-3.70)
RSRZ outliers	180081	1185 (4.06-3.70)

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	505	

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 4074 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.

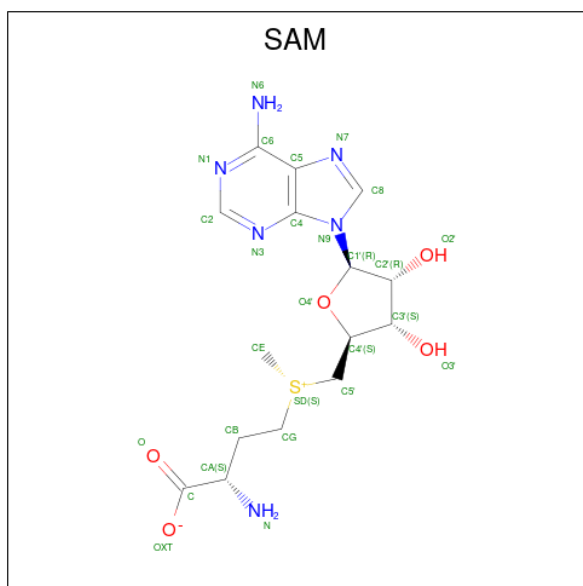
- Molecule 1 is a protein called WBDD.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	501	4047	2588	705	742	12	2	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	234	GLY	ASP	conflict	UNP A0A446F868

- Molecule 2 is S-ADENOSYLMETHIONINE (CCD ID: SAM) (formula: $C_{15}H_{22}N_6O_5S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
2	A	1	27	15	6	5	1	0	0

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: WBDD



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 3	Depositor
Cell constants a, b, c, α , β , γ	181.29Å 181.29Å 181.29Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	128.19 – 3.87 128.19 – 3.87	Depositor EDS
% Data completeness (in resolution range)	98.4 (128.19-3.87) 98.4 (128.19-3.87)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.32 (at 3.89Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.268 , 0.341 0.277 , 0.353	Depositor DCC
R_{free} test set	443 reflections (4.78%)	wwPDB-VP
Wilson B-factor (Å ²)	167.8	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 198.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.032 for -l,-k,-h	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	4074	wwPDB-VP
Average B, all atoms (Å ²)	171.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.24% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SAM

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	0.59	2/4156 (0.0%)	1.08	10/5651 (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	5

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	17	GLN	CD-OE1	18.14	1.58	1.23
1	A	17	GLN	CG-CD	-5.79	1.37	1.52

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	17	GLN	OE1-CD-NE2	30.25	152.85	122.60
1	A	17	GLN	CG-CD-OE1	-24.53	71.74	120.80
1	A	420	GLN	CA-C-N	-9.29	108.23	119.84
1	A	420	GLN	C-N-CA	-9.29	108.23	119.84
1	A	376	THR	C-N-CD	-8.76	89.07	125.00
1	A	17	GLN	CG-CD-NE2	-8.28	103.98	116.40
1	A	329	GLU	N-CA-C	7.56	119.21	110.97
1	A	410	SER	N-CA-C	6.37	121.40	112.93
1	A	399	ALA	N-CA-C	5.56	117.95	108.67
1	A	429	VAL	CB-CA-C	-5.05	105.33	112.14

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	17	GLN	Sidechain
1	A	261	GLY	Peptide
1	A	30	ALA	Peptide
1	A	409	ARG	Peptide
1	A	420	GLN	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	4047	0	3918	256	0
2	A	27	0	22	5	0
All	All	4074	0	3940	256	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 32.

All (256) 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:31:ARG:NH1	1:A:197:HIS:ND1	1.67	1.42
1:A:376:THR:CG2	1:A:377:PRO:HD3	1.69	1.22
1:A:31:ARG:CZ	1:A:197:HIS:ND1	2.10	1.14
1:A:342:LEU:HD21	1:A:347:PHE:HB2	1.12	1.11
1:A:376:THR:CB	1:A:377:PRO:HD3	1.69	1.07
1:A:342:LEU:CD2	1:A:347:PHE:HB2	1.83	1.07
1:A:376:THR:HG22	1:A:377:PRO:HD3	1.27	1.07
1:A:342:LEU:HD21	1:A:347:PHE:CB	1.86	1.06
1:A:31:ARG:HH12	1:A:197:HIS:CG	1.75	1.03
1:A:376:THR:CB	1:A:377:PRO:CD	2.38	1.02
1:A:31:ARG:NH2	1:A:197:HIS:ND1	2.12	0.98
1:A:331:ILE:HD11	1:A:397:LEU:HD21	1.42	0.97
1:A:31:ARG:NH2	1:A:197:HIS:CE1	2.34	0.96
1:A:477:GLN:O	1:A:480:VAL:HG22	1.67	0.94

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:31:ARG:NH1	1:A:197:HIS:CE1	2.36	0.93
1:A:182:ILE:HD11	1:A:208:VAL:CG1	2.01	0.90
1:A:31:ARG:CZ	1:A:197:HIS:CE1	2.55	0.89
1:A:31:ARG:HH22	1:A:197:HIS:CE1	1.89	0.89
1:A:376:THR:HG22	1:A:377:PRO:CD	2.04	0.88
1:A:31:ARG:HH12	1:A:197:HIS:CE1	1.92	0.88
1:A:376:THR:HB	1:A:377:PRO:CD	2.05	0.86
1:A:166:GLU:OE2	1:A:196:THR:HG21	1.75	0.85
1:A:465:TRP:CZ2	1:A:469:GLN:HG3	2.11	0.84
1:A:425:TRP:CZ2	1:A:446:PHE:HB2	2.12	0.84
1:A:424:ASN:HA	1:A:455:ALA:HB2	1.60	0.82
1:A:382:TRP:HB3	1:A:383:PRO:HD3	1.60	0.81
1:A:342:LEU:HD11	1:A:347:PHE:C	2.06	0.81
1:A:376:THR:HB	1:A:377:PRO:HD2	1.61	0.81
1:A:459:ARG:NH2	1:A:464:GLN:OE1	2.14	0.80
1:A:413:VAL:O	1:A:414:HIS:CG	2.36	0.79
1:A:155:GLN:NE2	1:A:212:ARG:HD2	1.98	0.78
1:A:83:PHE:HD1	1:A:84:GLN:H	1.32	0.77
1:A:458:GLN:OE1	1:A:458:GLN:N	2.18	0.77
1:A:36:ARG:NH1	1:A:160:GLU:OE1	2.19	0.76
1:A:477:GLN:O	1:A:480:VAL:CG2	2.35	0.75
1:A:182:ILE:HD11	1:A:208:VAL:HG12	1.68	0.75
1:A:413:VAL:O	1:A:414:HIS:CD2	2.40	0.74
1:A:215:ILE:HG22	1:A:216:ASN:N	2.02	0.74
1:A:331:ILE:HD11	1:A:397:LEU:CD2	2.17	0.74
1:A:15:ILE:HB	1:A:29:ALA:HB3	1.69	0.74
1:A:83:PHE:HD1	1:A:84:GLN:N	1.86	0.73
1:A:126:ILE:HG23	1:A:128:LEU:HG	1.70	0.73
1:A:398:PHE:HZ	1:A:422:TRP:CD2	2.06	0.73
1:A:72:ALA:HB2	1:A:78:ILE:HD13	1.70	0.73
1:A:170:TRP:C	1:A:174:GLN:HE22	1.97	0.73
1:A:5:LEU:O	1:A:9:VAL:HG23	1.90	0.72
1:A:215:ILE:HG22	1:A:216:ASN:H	1.54	0.72
1:A:404:TRP:CE3	1:A:474:LEU:HB3	2.25	0.70
1:A:162:ALA:HA	1:A:203:ARG:HG2	1.73	0.69
1:A:244:PHE:HB2	1:A:251:CYS:HB3	1.73	0.69
1:A:424:ASN:OD1	1:A:453:PRO:O	2.11	0.68
1:A:170:TRP:O	1:A:174:GLN:NE2	2.27	0.68
1:A:342:LEU:HD13	1:A:342:LEU:C	2.18	0.68
1:A:289:PHE:CE1	1:A:291:ALA:HB3	2.29	0.67
1:A:425:TRP:CE3	1:A:426:LEU:CD2	2.78	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:415:PRO:HB3	1:A:426:LEU:HB2	1.76	0.67
1:A:376:THR:CG2	1:A:377:PRO:CD	2.59	0.67
1:A:339:LEU:HD21	1:A:439:PHE:CD1	2.31	0.66
1:A:166:GLU:CG	1:A:167:PRO:HD2	2.26	0.66
1:A:155:GLN:HE22	1:A:212:ARG:HD2	1.58	0.66
1:A:425:TRP:CE3	1:A:426:LEU:HD21	2.31	0.65
1:A:413:VAL:HG12	1:A:470:GLU:OE2	1.97	0.65
1:A:477:GLN:C	1:A:480:VAL:HG22	2.21	0.65
1:A:38:ASP:O	1:A:41:THR:OG1	2.14	0.64
1:A:174:GLN:HB3	1:A:175:PRO:HD2	1.80	0.64
1:A:391:PHE:CD1	1:A:426:LEU:HD12	2.33	0.64
1:A:212:ARG:HA	1:A:221:PRO:HA	1.78	0.63
1:A:118:GLU:HB2	1:A:121:GLU:HB2	1.81	0.62
1:A:424:ASN:OD1	1:A:454:SER:C	2.43	0.62
1:A:83:PHE:CD1	1:A:84:GLN:N	2.66	0.62
1:A:342:LEU:HD11	1:A:348:TRP:N	2.15	0.61
1:A:391:PHE:CE1	1:A:426:LEU:HD12	2.36	0.61
1:A:289:PHE:CE1	1:A:291:ALA:CB	2.83	0.61
1:A:339:LEU:HD21	1:A:439:PHE:CE1	2.35	0.61
1:A:415:PRO:CB	1:A:426:LEU:HB2	2.31	0.61
1:A:316:LEU:HB3	1:A:357:VAL:CG1	2.31	0.61
1:A:382:TRP:O	1:A:383:PRO:C	2.44	0.60
1:A:9:VAL:O	1:A:12:LEU:HB3	2.02	0.60
1:A:350:ASP:HB2	1:A:380:CYS:HA	1.84	0.59
1:A:409:ARG:HB3	1:A:410:SER:HB2	1.84	0.59
1:A:306:LEU:HD23	1:A:307:VAL:N	2.18	0.59
1:A:342:LEU:HD11	1:A:348:TRP:HA	1.84	0.59
1:A:166:GLU:HG3	1:A:167:PRO:HD2	1.85	0.59
1:A:462:THR:OG1	1:A:463:GLU:N	2.32	0.59
1:A:382:TRP:HB3	1:A:383:PRO:CD	2.32	0.59
1:A:461:ALA:O	1:A:463:GLU:N	2.36	0.58
1:A:457:GLN:HB2	1:A:458:GLN:OE1	2.04	0.58
1:A:477:GLN:HA	1:A:480:VAL:HG22	1.86	0.58
1:A:12:LEU:HD12	1:A:90:VAL:HG21	1.85	0.58
1:A:163:VAL:HG23	1:A:178:PRO:HD3	1.86	0.57
1:A:407:PHE:O	1:A:408:TRP:CG	2.57	0.57
1:A:415:PRO:HB3	1:A:426:LEU:CB	2.34	0.57
1:A:350:ASP:O	1:A:380:CYS:SG	2.60	0.57
1:A:332:LEU:HD21	1:A:397:LEU:CD2	2.35	0.57
1:A:154:THR:O	1:A:155:GLN:CG	2.53	0.57
1:A:199:SER:OG	1:A:200:PRO:HD2	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:387:VAL:HG21	1:A:437:TRP:HB3	1.86	0.56
1:A:328:ARG:NH1	1:A:397:LEU:O	2.39	0.56
1:A:404:TRP:HZ3	1:A:475:GLU:CA	2.17	0.56
1:A:215:ILE:CG2	1:A:216:ASN:H	2.18	0.56
1:A:460:GLY:O	1:A:461:ALA:O	2.25	0.56
1:A:443:LEU:O	1:A:446:PHE:HB3	2.07	0.55
1:A:289:PHE:HE1	1:A:291:ALA:HB3	1.72	0.55
1:A:407:PHE:O	1:A:408:TRP:CB	2.54	0.55
1:A:415:PRO:O	1:A:418:LEU:N	2.40	0.55
1:A:398:PHE:CD2	1:A:418:LEU:HD22	2.42	0.55
1:A:404:TRP:CE3	1:A:474:LEU:CB	2.89	0.55
1:A:37:LEU:HA	1:A:40:ILE:HD12	1.89	0.55
1:A:425:TRP:NE1	1:A:446:PHE:N	2.54	0.54
1:A:477:GLN:CA	1:A:480:VAL:HG22	2.37	0.54
1:A:469:GLN:OE1	1:A:469:GLN:HA	2.07	0.54
1:A:83:PHE:HE1	2:A:1506:SAM:HO2'	1.54	0.54
1:A:46:ASN:O	1:A:49:ARG:HG2	2.07	0.54
1:A:456:GLU:HA	1:A:459:ARG:HD3	1.88	0.54
1:A:425:TRP:O	1:A:429:VAL:HG23	2.07	0.54
1:A:383:PRO:HG2	1:A:389:SER:HA	1.90	0.53
1:A:256:TYR:CE1	1:A:272:LYS:HG3	2.43	0.53
1:A:404:TRP:HZ3	1:A:475:GLU:N	2.05	0.53
1:A:110:ILE:O	1:A:113:VAL:HG22	2.09	0.53
1:A:127:GLY:HA3	1:A:159:LEU:CD2	2.38	0.53
1:A:135:VAL:CG1	1:A:174:GLN:HG3	2.38	0.53
1:A:215:ILE:CG2	1:A:216:ASN:N	2.69	0.53
1:A:316:LEU:CB	1:A:357:VAL:CG1	2.87	0.53
1:A:465:TRP:CE3	1:A:466:ILE:N	2.76	0.53
1:A:253:PHE:HE2	1:A:255:TYR:OH	1.91	0.53
1:A:376:THR:HG22	1:A:377:PRO:CG	2.39	0.53
1:A:465:TRP:CH2	1:A:469:GLN:HG3	2.42	0.53
1:A:37:LEU:O	1:A:41:THR:HG23	2.09	0.52
1:A:454:SER:O	1:A:456:GLU:N	2.42	0.52
1:A:72:ALA:HB2	1:A:78:ILE:CD1	2.37	0.52
1:A:88:ILE:HG13	1:A:107:VAL:HG12	1.90	0.52
1:A:353:ARG:HH12	1:A:392:VAL:CG1	2.22	0.52
1:A:378:GLN:HB2	1:A:386:LEU:HG	1.92	0.52
1:A:182:ILE:CD1	1:A:208:VAL:CG1	2.82	0.51
1:A:425:TRP:HE1	1:A:446:PHE:N	2.08	0.51
1:A:311:LEU:HB2	1:A:312:PRO:HD2	1.92	0.51
1:A:68:SER:HB3	1:A:78:ILE:HG21	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:353:ARG:HH12	1:A:392:VAL:HG11	1.76	0.50
1:A:289:PHE:HD1	1:A:291:ALA:N	2.10	0.50
1:A:415:PRO:HA	1:A:418:LEU:HG	1.93	0.50
1:A:342:LEU:HD11	1:A:348:TRP:CA	2.42	0.50
1:A:407:PHE:O	1:A:408:TRP:HB2	2.12	0.50
1:A:44:TYR:CD1	1:A:44:TYR:C	2.90	0.50
1:A:342:LEU:HD22	1:A:347:PHE:HB2	1.88	0.50
1:A:386:LEU:O	1:A:389:SER:N	2.45	0.50
1:A:416:PHE:CD1	1:A:459:ARG:HD2	2.47	0.50
1:A:443:LEU:HD12	1:A:443:LEU:C	2.37	0.50
1:A:9:VAL:O	1:A:12:LEU:CB	2.60	0.49
1:A:61:GLY:O	2:A:1506:SAM:N	2.44	0.49
1:A:465:TRP:CE3	1:A:466:ILE:CA	2.95	0.49
1:A:104:GLU:O	1:A:105:PHE:HB2	2.12	0.49
1:A:282:LEU:HD21	1:A:370:PHE:CD2	2.47	0.49
1:A:395:ASN:HB3	1:A:411:ALA:HB1	1.95	0.49
1:A:477:GLN:HA	1:A:480:VAL:CG2	2.42	0.49
1:A:174:GLN:NE2	1:A:174:GLN:H	2.10	0.49
1:A:176:ASP:OD1	1:A:180:GLU:OE2	2.31	0.48
1:A:356:ASN:HA	1:A:368:ILE:HG22	1.95	0.48
1:A:31:ARG:O	1:A:33:CYS:SG	2.67	0.48
1:A:157:VAL:HB	1:A:208:VAL:CG2	2.44	0.48
1:A:90:VAL:O	1:A:94:LEU:HG	2.13	0.48
1:A:164:LYS:HG3	1:A:171:GLY:O	2.13	0.48
1:A:253:PHE:CE2	1:A:255:TYR:CZ	3.02	0.48
1:A:348:TRP:O	1:A:374:VAL:N	2.46	0.48
1:A:46:ASN:HA	1:A:49:ARG:HD3	1.96	0.48
1:A:398:PHE:HZ	1:A:422:TRP:CE3	2.32	0.48
1:A:462:THR:O	1:A:465:TRP:N	2.46	0.48
1:A:131:PHE:HE2	1:A:159:LEU:HD22	1.79	0.47
1:A:452:LEU:HB3	1:A:453:PRO:HD2	1.96	0.47
1:A:253:PHE:CE2	1:A:255:TYR:CE1	3.02	0.47
1:A:391:PHE:CE1	1:A:426:LEU:CD1	2.97	0.47
1:A:135:VAL:HG13	1:A:174:GLN:HG3	1.96	0.47
1:A:253:PHE:CD1	1:A:305:TRP:HB3	2.49	0.47
1:A:408:TRP:O	1:A:409:ARG:O	2.33	0.47
1:A:391:PHE:HB2	1:A:413:VAL:HG23	1.96	0.47
1:A:465:TRP:CE3	1:A:466:ILE:HA	2.50	0.47
1:A:422:TRP:O	1:A:423:SER:C	2.58	0.47
1:A:425:TRP:CE2	1:A:446:PHE:HA	2.50	0.47
1:A:463:GLU:O	1:A:467:ILE:HG13	2.14	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:127:GLY:HA3	1:A:159:LEU:HD23	1.97	0.46
1:A:117:LEU:HD13	1:A:153:VAL:HG21	1.96	0.46
1:A:258:MET:HA	1:A:258:MET:HE2	1.98	0.46
1:A:289:PHE:CD1	1:A:291:ALA:HB2	2.51	0.46
1:A:17:GLN:HB2	1:A:64:GLN:HB2	1.97	0.46
1:A:404:TRP:CZ3	1:A:475:GLU:N	2.83	0.46
1:A:347:PHE:CD2	1:A:375:THR:HG22	2.50	0.46
1:A:110:ILE:HA	1:A:113:VAL:HG22	1.98	0.46
1:A:342:LEU:CD1	1:A:348:TRP:HA	2.46	0.46
1:A:353:ARG:HB3	1:A:354:PRO:HD2	1.98	0.46
1:A:342:LEU:HD21	1:A:347:PHE:HB3	1.89	0.45
1:A:215:ILE:CG2	1:A:305:TRP:CZ3	2.99	0.45
1:A:421:PRO:HD2	1:A:422:TRP:CE3	2.51	0.45
1:A:57:VAL:HB	1:A:78:ILE:HG13	1.97	0.45
1:A:149:ARG:O	1:A:153:VAL:HG13	2.16	0.45
1:A:256:TYR:CD1	1:A:272:LYS:HG3	2.52	0.45
1:A:289:PHE:HD1	1:A:291:ALA:H	1.65	0.45
1:A:131:PHE:CE2	1:A:159:LEU:HD22	2.52	0.45
1:A:368:ILE:HG12	1:A:369:ASP:H	1.81	0.45
1:A:12:LEU:HG	1:A:13:PRO:HD2	1.99	0.44
1:A:398:PHE:CZ	1:A:422:TRP:CE3	3.05	0.44
1:A:311:LEU:C	1:A:311:LEU:HD12	2.43	0.44
1:A:157:VAL:O	1:A:208:VAL:HG22	2.18	0.44
1:A:457:GLN:C	1:A:459:ARG:H	2.23	0.44
1:A:496:THR:O	1:A:500:GLN:HG3	2.17	0.44
1:A:135:VAL:HB	1:A:143:VAL:HG21	1.98	0.44
1:A:386:LEU:O	1:A:387:VAL:C	2.60	0.44
1:A:413:VAL:N	1:A:470:GLU:OE2	2.51	0.44
1:A:109:ARG:O	1:A:112:GLU:HG2	2.18	0.43
1:A:12:LEU:HG	1:A:13:PRO:CD	2.48	0.43
1:A:141:ASP:HA	1:A:144:LYS:HE2	2.00	0.43
1:A:490:LEU:O	1:A:494:ILE:HG13	2.18	0.43
1:A:128:LEU:O	2:A:1506:SAM:N	2.52	0.43
1:A:147:LEU:HD12	1:A:181:LEU:O	2.19	0.43
1:A:357:VAL:HG22	1:A:358:MET:N	2.33	0.43
1:A:40:ILE:HG23	1:A:158:ILE:HD11	2.00	0.43
1:A:443:LEU:HD12	1:A:444:ALA:N	2.34	0.43
1:A:179:ARG:O	1:A:182:ILE:HG22	2.18	0.43
1:A:339:LEU:CD2	1:A:439:PHE:CD1	3.00	0.43
1:A:379:ASP:O	1:A:380:CYS:CB	2.65	0.43
1:A:390:PHE:O	1:A:394:VAL:HG23	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:12:LEU:CD1	1:A:90:VAL:HG21	2.49	0.42
1:A:153:VAL:HG23	1:A:154:THR:HG23	2.01	0.42
1:A:470:GLU:O	1:A:474:LEU:HD13	2.19	0.42
1:A:128:LEU:HB2	2:A:1506:SAM:HN1	1.84	0.42
1:A:427:TYR:O	1:A:431:GLN:HG3	2.19	0.42
1:A:331:ILE:HD12	1:A:332:LEU:N	2.35	0.42
1:A:461:ALA:O	1:A:462:THR:OG1	2.32	0.42
1:A:73:SER:HA	1:A:101:PHE:CE1	2.54	0.42
1:A:316:LEU:CB	1:A:357:VAL:HG11	2.50	0.42
1:A:140:ILE:HG22	1:A:181:LEU:HD21	2.02	0.42
1:A:349:HIS:CD2	1:A:367:LEU:HD23	2.55	0.42
1:A:143:VAL:HA	1:A:146:LEU:HD23	2.02	0.41
1:A:335:LEU:HD23	1:A:335:LEU:C	2.44	0.41
1:A:12:LEU:CD2	1:A:13:PRO:HD2	2.50	0.41
1:A:188:TYR:HA	1:A:207:LEU:O	2.21	0.41
1:A:504:LEU:HD12	1:A:504:LEU:HA	1.91	0.41
1:A:169:TYR:CG	1:A:170:TRP:N	2.88	0.41
1:A:364:HIS:O	1:A:364:HIS:ND1	2.51	0.41
1:A:182:ILE:HG13	1:A:208:VAL:HG11	2.03	0.41
1:A:264:THR:O	1:A:265:ALA:C	2.64	0.41
1:A:424:ASN:OD1	1:A:455:ALA:CB	2.69	0.41
1:A:425:TRP:NE1	1:A:445:LEU:C	2.78	0.41
1:A:132:HIS:HA	1:A:135:VAL:HG12	2.03	0.41
1:A:316:LEU:CD1	1:A:326:ILE:HD13	2.51	0.41
1:A:83:PHE:HE1	2:A:1506:SAM:O2'	2.04	0.40
1:A:382:TRP:CG	1:A:383:PRO:N	2.87	0.40
1:A:425:TRP:HE1	1:A:445:LEU:CB	2.34	0.40
1:A:461:ALA:C	1:A:462:THR:HG1	2.25	0.40
1:A:382:TRP:CB	1:A:383:PRO:CD	2.98	0.40
1:A:223:GLN:OE1	1:A:223:GLN:N	2.55	0.40
1:A:465:TRP:CD2	1:A:465:TRP:C	2.99	0.40
1:A:38:ASP:C	1:A:41:THR:HG1	2.24	0.40
1:A:289:PHE:CD1	1:A:291:ALA:CB	3.04	0.40
1:A:15:ILE:HG22	1:A:17:GLN:O	2.22	0.40
1:A:378:GLN:O	1:A:379:ASP:CG	2.64	0.40
1:A:425:TRP:HE1	1:A:445:LEU:C	2.30	0.40
1:A:476:LEU:HD23	1:A:476:LEU:HA	1.68	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	499/505 (99%)	414 (83%)	67 (13%)	18 (4%)	2	23

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	28	ASP
1	A	155	GLN
1	A	382	TRP
1	A	386	LEU
1	A	408	TRP
1	A	409	ARG
1	A	421	PRO
1	A	379	ASP
1	A	459	ARG
1	A	461	ALA
1	A	13	PRO
1	A	380	CYS
1	A	455	ALA
1	A	462	THR
1	A	377	PRO
1	A	383	PRO
1	A	460	GLY
1	A	61	GLY

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	428/432 (99%)	415 (97%)	13 (3%)	36 57

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14	GLU
1	A	83	PHE
1	A	155	GLN
1	A	165	GLU
1	A	166	GLU
1	A	174	GLN
1	A	203	ARG
1	A	262	ILE
1	A	267	GLU
1	A	348	TRP
1	A	355	TRP
1	A	380	CYS
1	A	504	LEU

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

Mol	Chain	Res	Type
1	A	56	ASN
1	A	155	GLN
1	A	174	GLN
1	A	227	ASN

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

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	SAM	A	1506	-	27,29,29	1.53	4 (14%)	34,42,42	2.09	11 (32%)

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	SAM	A	1506	-	-	2/17/33/33	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1506	SAM	C5-C4	4.91	1.47	1.39
2	A	1506	SAM	C5-C6	2.66	1.48	1.41
2	A	1506	SAM	C8-N7	2.47	1.36	1.31
2	A	1506	SAM	C5-N7	-2.44	1.34	1.39

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1506	SAM	C5-C4-N3	-5.68	118.89	126.72
2	A	1506	SAM	N3-C4-N9	4.36	134.58	127.17
2	A	1506	SAM	C2-N3-C4	3.60	120.62	111.83
2	A	1506	SAM	N3-C2-N1	-3.47	123.33	128.58
2	A	1506	SAM	C4-C5-N7	-3.43	106.66	110.58
2	A	1506	SAM	CG-SD-C5'	3.28	111.45	103.43
2	A	1506	SAM	C5-N7-C8	2.63	107.58	103.45
2	A	1506	SAM	C2-N1-C6	2.29	122.50	118.73
2	A	1506	SAM	C3'-C2'-C1'	2.19	105.60	101.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1506	SAM	C4-N9-C8	2.18	108.03	105.74
2	A	1506	SAM	O4'-C1'-N9	2.09	112.10	108.09

There are no chirality outliers.

All (2) torsion outliers are listed below:

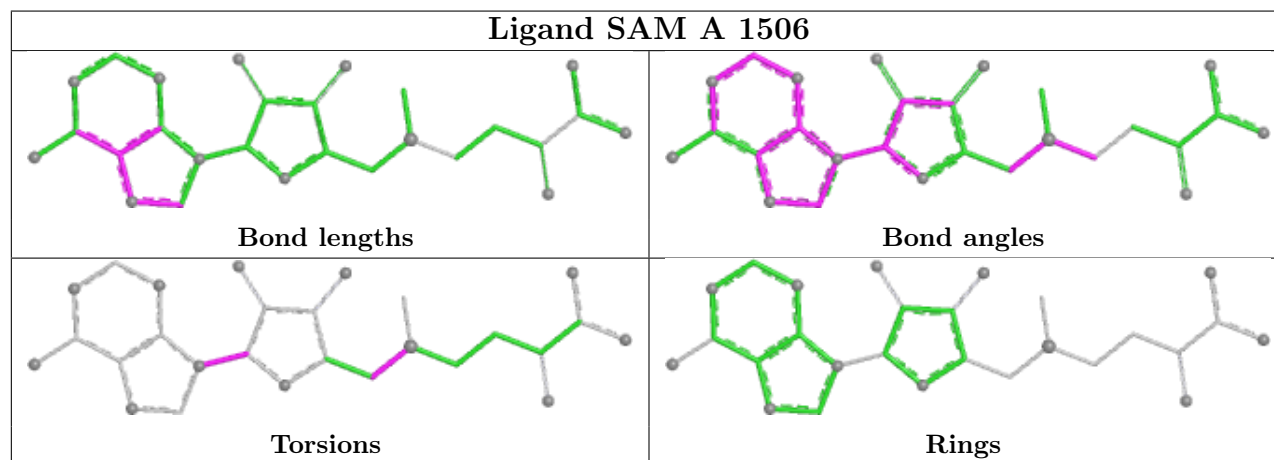
Mol	Chain	Res	Type	Atoms
2	A	1506	SAM	C4'-C5'-SD-CE
2	A	1506	SAM	C2'-C1'-N9-C8

There are no ring outliers.

1 monomer is involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1506	SAM	5	0

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



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

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	501/505 (99%)	0.59	34 (6%) 23 20	124, 168, 215, 247	1 (0%)

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	494	ILE	6.3
1	A	490	LEU	5.6
1	A	85	GLN	4.7
1	A	367	LEU	4.1
1	A	439	PHE	3.8
1	A	483	GLU	3.7
1	A	167	PRO	3.7
1	A	497	LEU	3.6
1	A	165	GLU	3.6
1	A	505	GLN	3.6
1	A	477	GLN	3.5
1	A	493	GLN	3.4
1	A	368	ILE	3.2
1	A	352	VAL	3.1
1	A	411	ALA	3.1
1	A	480	VAL	3.0
1	A	384	THR	3.0
1	A	487	SER	2.9
1	A	398	PHE	2.9
1	A	200	PRO	2.6
1	A	84	GLN	2.6
1	A	390	PHE	2.4
1	A	396	GLU	2.4
1	A	504	LEU	2.4
1	A	498	GLU	2.4
1	A	198	LEU	2.2
1	A	346	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	94	LEU	2.2
1	A	169	TYR	2.1
1	A	397	LEU	2.1
1	A	393	PHE	2.1
1	A	124	LEU	2.0
1	A	168	LEU	2.0
1	A	217	ASP	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](#)

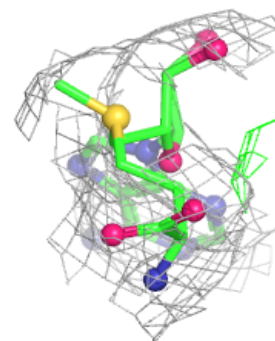
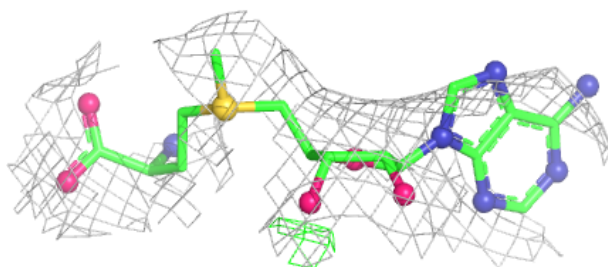
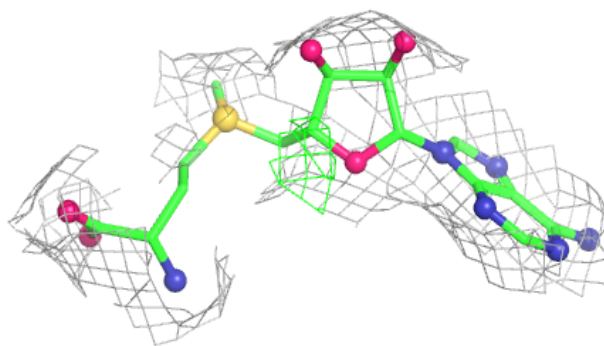
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
2	SAM	A	1506	27/27	0.89	0.10	117,126,132,132	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 SAM A 1506:

$2mF_o-DF_c$ (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.