



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 09:14 PM UTC

PDB ID : 2I0V / pdb_00002i0v
Title : c-FMS tyrosine kinase in complex with a quinolone inhibitor
Authors : Schubert, C.; Schalk-Hihi, C.
Deposited on : 2006-08-11
Resolution : 2.80 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

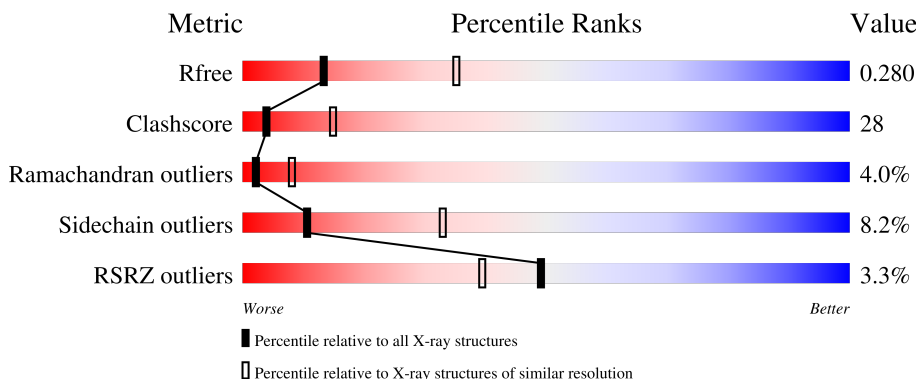
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

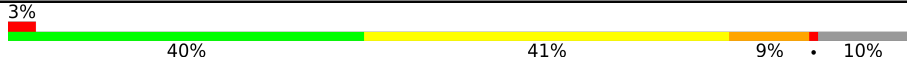
The reported resolution of this entry is 2.80 Å.

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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	335	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 2456 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 cFMS tyrosine kinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	303	2401	1547	409	430	15	0	0	0

There are 24 discrepancies between the modelled and reference sequences:

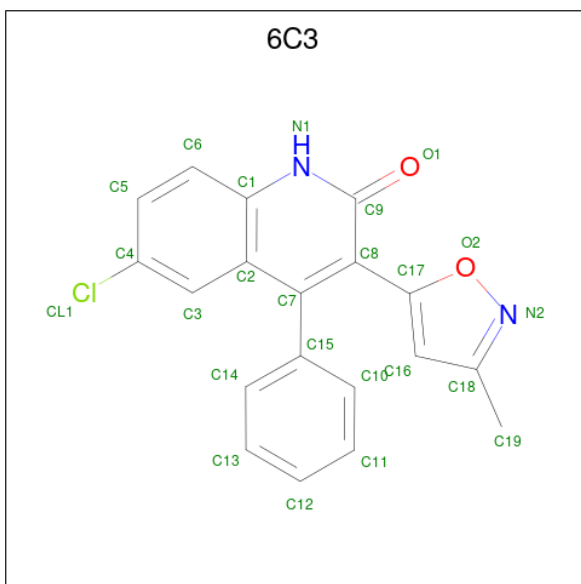
Chain	Residue	Modelled	Actual	Comment	Reference
A	535	GLY	-	expression tag	UNP P07333
A	536	VAL	-	expression tag	UNP P07333
A	537	ASP	-	expression tag	UNP P07333
A	679	ARG	-	SEE REMARK 999	UNP P07333
A	680	PRO	-	SEE REMARK 999	UNP P07333
A	681	PRO	-	SEE REMARK 999	UNP P07333
A	682	GLY	-	SEE REMARK 999	UNP P07333
A	683	LEU	-	SEE REMARK 999	UNP P07333
A	684	GLU	-	SEE REMARK 999	UNP P07333
A	685	TYR	-	SEE REMARK 999	UNP P07333
A	686	SER	-	SEE REMARK 999	UNP P07333
A	687	TYR	-	SEE REMARK 999	UNP P07333
A	688	ASN	-	SEE REMARK 999	UNP P07333
A	689	PRO	-	SEE REMARK 999	UNP P07333
A	690	SER	-	SEE REMARK 999	UNP P07333
A	691	HIS	-	SEE REMARK 999	UNP P07333
A	692	ASN	-	SEE REMARK 999	UNP P07333
A	693	PRO	-	SEE REMARK 999	UNP P07333
A	694	GLU	-	SEE REMARK 999	UNP P07333
A	695	GLU	-	SEE REMARK 999	UNP P07333
A	696	GLN	-	SEE REMARK 999	UNP P07333
A	697	LEU	-	SEE REMARK 999	UNP P07333
A	698	SER	-	SEE REMARK 999	UNP P07333
A	699	SER	-	SEE REMARK 999	UNP P07333

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is 6-CHLORO-3-(3-METHYLISOXAZOL-5-YL)-4-PHENYLQUINOLIN-2(1H)-ONE (CCD ID: 6C3) (formula: $C_{19}H_{13}ClN_2O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	Cl	N	O	
			24	19	1	2	2	

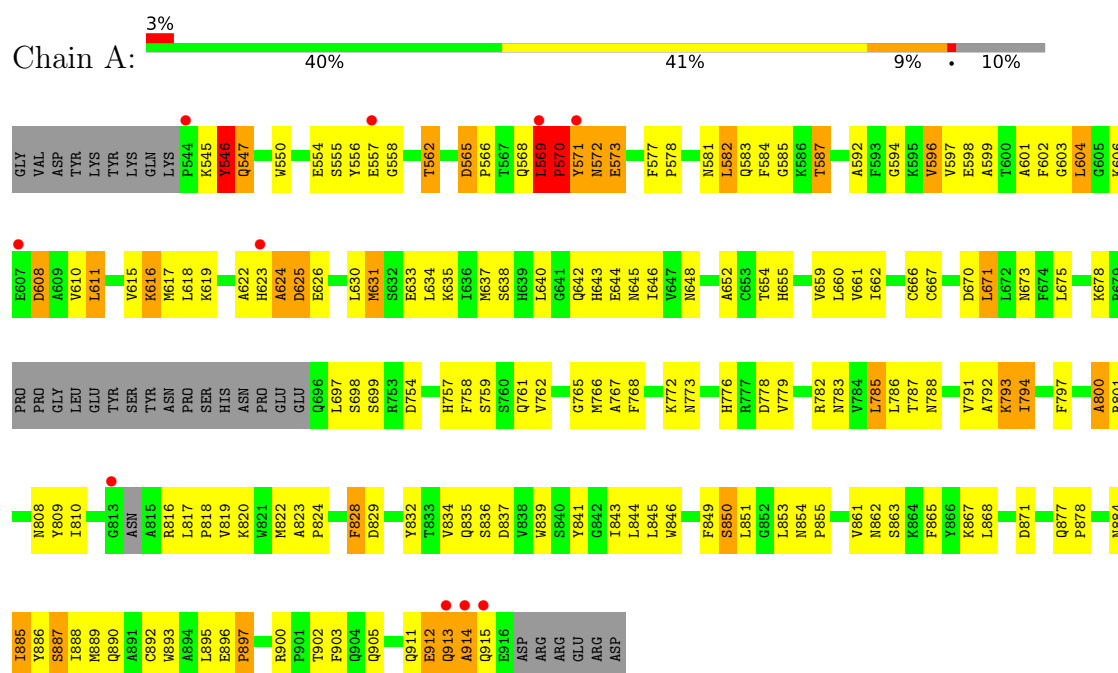
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	26	Total 26	O 26	0	0

3 Residue-property plots

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

• Molecule 1: cFMS tyrosine kinase



4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	82.20Å 82.20Å 143.30Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	26.59 – 2.80 26.59 – 2.80	Depositor EDS
% Data completeness (in resolution range)	98.4 (26.59-2.80) 98.3 (26.59-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.90 (at 2.69Å)	Xtriage
Refinement program	CNX 2005	Depositor
R, R_{free}	0.237 , 0.296 0.229 , 0.280	Depositor DCC
R_{free} test set	895 reflections (9.10%)	wwPDB-VP
Wilson B-factor (Å ²)	37.1	Xtriage
Anisotropy	0.135	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 46.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.044 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	2456	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.75% 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: SO4, 6C3

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	0.54	0/2459	1.12	23/3326 (0.7%)

There are no bond length outliers.

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	667	CYS	N-CA-C	10.02	124.77	112.54
1	A	850	SER	N-CA-C	-8.38	102.78	113.16
1	A	836	SER	N-CA-C	-8.36	102.25	111.36
1	A	819	VAL	N-CA-C	7.92	118.46	110.23
1	A	896	GLU	CA-C-N	7.44	129.14	119.84
1	A	896	GLU	C-N-CA	7.44	129.14	119.84
1	A	611	LEU	N-CA-C	7.36	120.24	110.53
1	A	914	ALA	N-CA-C	-7.30	104.37	113.20
1	A	648	ASN	N-CA-C	6.41	118.80	110.53
1	A	615	VAL	N-CA-C	5.90	116.30	107.51
1	A	642	GLN	N-CA-C	5.79	118.33	108.90
1	A	767	ALA	N-CA-C	-5.76	105.09	111.36
1	A	585	GLY	N-CA-C	-5.68	104.52	112.13
1	A	671	LEU	N-CA-C	5.61	117.39	111.28
1	A	828	PHE	N-CA-C	5.59	118.57	111.69
1	A	565	ASP	CA-C-N	5.55	126.78	119.84
1	A	565	ASP	C-N-CA	5.55	126.78	119.84
1	A	800	ALA	N-CA-C	5.39	119.34	112.87
1	A	616	LYS	N-CA-C	-5.34	102.97	110.50
1	A	913	GLN	N-CA-C	-5.20	106.13	113.20
1	A	888	ILE	N-CA-C	-5.17	105.36	111.00
1	A	569	LEU	CA-C-N	5.09	126.20	119.84
1	A	569	LEU	C-N-CA	5.09	126.20	119.84

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	2401	0	2388	135	0
2	A	5	0	0	0	0
3	A	24	0	13	4	0
4	A	26	0	0	3	0
All	All	2456	0	2401	136	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 28.

All (136) 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:886:TYR:HA	1:A:889:MET:HE3	1.42	1.00
1:A:587:THR:HA	1:A:597:VAL:HG12	1.48	0.94
1:A:654:THR:HA	1:A:659:VAL:HG22	1.56	0.87
1:A:810:ILE:HG13	1:A:861:VAL:HG11	1.55	0.86
1:A:785:LEU:HD13	1:A:797:PHE:CE1	2.22	0.76
1:A:792:ALA:O	1:A:793:LYS:HG3	1.86	0.75
1:A:902:THR:H	1:A:905:GLN:HE21	1.34	0.74
1:A:808:ASN:HD21	1:A:854:ASN:HD21	1.33	0.73
1:A:757:HIS:O	1:A:761:GLN:HG3	1.89	0.71
1:A:633:GLU:O	1:A:637:MET:HG2	1.92	0.70
1:A:666:CYS:SG	1:A:785:LEU:HD23	2.32	0.69
1:A:596:VAL:HG11	1:A:800:ALA:HB2	1.75	0.69
1:A:566:PRO:O	1:A:570:PRO:HD2	1.91	0.69
1:A:596:VAL:CG1	1:A:800:ALA:HB2	2.22	0.69
1:A:778:ASP:O	1:A:783:ASN:ND2	2.27	0.66
1:A:618:LEU:HG	1:A:630:LEU:HD22	1.77	0.65
1:A:785:LEU:HD13	1:A:797:PHE:CZ	2.33	0.64
1:A:606:LYS:HD2	1:A:606:LYS:O	1.97	0.64
1:A:698:SER:N	1:A:754:ASP:OD1	2.32	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:801:ARG:HG3	4:A:11:HOH:O	1.99	0.62
1:A:824:PRO:O	1:A:828:PHE:HD2	1.82	0.62
1:A:877:GLN:HA	1:A:886:TYR:CE1	2.35	0.61
1:A:569:LEU:H	1:A:570:PRO:HD2	1.65	0.60
1:A:646:ILE:HD13	1:A:794:ILE:HG13	1.83	0.60
1:A:577:PHE:CE2	1:A:582:LEU:HD13	2.38	0.59
1:A:573:GLU:OE2	1:A:655:HIS:CD2	2.56	0.59
1:A:845:LEU:O	1:A:849:PHE:HD2	1.85	0.59
1:A:911:GLN:O	1:A:913:GLN:N	2.35	0.59
1:A:602:PHE:HA	1:A:610:VAL:HG13	1.85	0.58
1:A:820:LYS:HE3	1:A:861:VAL:HG23	1.85	0.58
1:A:545:LYS:HG2	1:A:547:GLN:HG3	1.85	0.58
1:A:594:GLY:HA2	1:A:619:LYS:HG3	1.85	0.58
1:A:616:LYS:HB2	1:A:661:VAL:HB	1.86	0.58
1:A:846:TRP:NE1	1:A:878:PRO:HG3	2.19	0.58
1:A:911:GLN:C	1:A:913:GLN:N	2.60	0.58
1:A:892:CYS:O	1:A:900:ARG:HD3	2.04	0.56
1:A:800:ALA:HB1	3:A:1000:6C3:C11	2.35	0.56
1:A:645:ASN:O	1:A:646:ILE:HD13	2.06	0.56
1:A:766:MET:SD	1:A:794:ILE:HD13	2.45	0.56
1:A:583:GLN:HE21	1:A:602:PHE:HZ	1.55	0.55
1:A:631:MET:O	1:A:635:LYS:HG3	2.05	0.55
1:A:809:TYR:HB3	1:A:816:ARG:HB3	1.89	0.55
1:A:911:GLN:C	1:A:913:GLN:H	2.13	0.55
1:A:913:GLN:HA	1:A:913:GLN:OE1	2.06	0.54
1:A:902:THR:H	1:A:905:GLN:NE2	2.03	0.54
1:A:578:PRO:HB2	1:A:581:ASN:OD1	2.08	0.53
1:A:912:GLU:C	1:A:914:ALA:H	2.17	0.52
1:A:644:GLU:O	1:A:793:LYS:NZ	2.42	0.51
1:A:616:LYS:HE3	3:A:1000:6C3:CL1	2.47	0.51
1:A:678:LYS:O	1:A:697:LEU:HD12	2.11	0.51
1:A:577:PHE:CE1	1:A:604:LEU:HG	2.46	0.50
1:A:862:ASN:O	1:A:865:PHE:N	2.45	0.50
1:A:897:PRO:HA	1:A:900:ARG:NH1	2.26	0.50
1:A:583:GLN:O	1:A:599:ALA:HA	2.12	0.50
1:A:766:MET:HE3	1:A:903:PHE:CE2	2.47	0.50
1:A:550:TRP:HE1	1:A:776:HIS:CD2	2.30	0.50
1:A:571:TYR:C	1:A:571:TYR:CD1	2.90	0.49
1:A:758:PHE:O	1:A:762:VAL:HG23	2.12	0.49
1:A:546:TYR:HD2	1:A:546:TYR:C	2.20	0.49
1:A:884:ASN:O	1:A:887:SER:HB2	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:897:PRO:HG3	1:A:900:ARG:HH12	1.78	0.49
1:A:584:PHE:CZ	1:A:617:MET:HE1	2.48	0.48
1:A:546:TYR:C	1:A:546:TYR:CD2	2.91	0.48
1:A:759:SER:OG	1:A:845:LEU:CD2	2.61	0.48
1:A:843:ILE:HD13	1:A:893:TRP:NE1	2.28	0.48
1:A:622:ALA:HA	1:A:626:GLU:OE1	2.14	0.48
1:A:824:PRO:O	1:A:828:PHE:CD2	2.66	0.48
1:A:834:VAL:O	1:A:837:ASP:HB2	2.14	0.47
1:A:854:ASN:O	1:A:855:PRO:C	2.57	0.47
1:A:671:LEU:O	1:A:675:LEU:HG	2.15	0.47
1:A:792:ALA:O	1:A:793:LYS:CG	2.61	0.47
1:A:835:GLN:HG3	4:A:26:HOH:O	2.14	0.47
1:A:587:THR:CA	1:A:597:VAL:HG12	2.32	0.47
1:A:867:LYS:O	1:A:871:ASP:OD2	2.33	0.47
1:A:571:TYR:O	1:A:572:ASN:C	2.57	0.47
1:A:602:PHE:HD2	1:A:610:VAL:CG1	2.27	0.47
1:A:623:HIS:O	1:A:624:ALA:C	2.57	0.47
1:A:678:LYS:HE2	4:A:23:HOH:O	2.15	0.47
1:A:839:TRP:C	1:A:839:TRP:CD1	2.92	0.46
3:A:1000:6C3:C15	3:A:1000:6C3:H16	2.46	0.46
1:A:765:GLY:C	1:A:794:ILE:HD11	2.41	0.46
1:A:603:GLY:HA2	1:A:608:ASP:HA	1.98	0.46
1:A:913:GLN:C	1:A:915:GLN:H	2.23	0.46
1:A:556:TYR:O	1:A:558:GLY:N	2.48	0.46
1:A:782:ARG:NH2	1:A:818:PRO:HB3	2.31	0.46
1:A:787:THR:HG21	1:A:793:LYS:HE2	1.98	0.45
1:A:884:ASN:O	1:A:887:SER:N	2.49	0.45
1:A:550:TRP:NE1	1:A:776:HIS:CD2	2.84	0.45
1:A:850:SER:HB2	1:A:853:LEU:HD12	1.99	0.45
1:A:546:TYR:O	1:A:547:GLN:HB2	2.16	0.45
1:A:555:SER:HB3	1:A:773:ASN:ND2	2.32	0.44
1:A:601:ALA:O	1:A:610:VAL:HA	2.18	0.44
1:A:652:ALA:HA	1:A:660:LEU:O	2.18	0.44
1:A:598:GLU:O	1:A:599:ALA:HB2	2.17	0.44
1:A:841:TYR:O	1:A:844:LEU:HB3	2.17	0.44
1:A:765:GLY:O	1:A:768:PHE:HB3	2.18	0.44
1:A:569:LEU:H	1:A:570:PRO:CD	2.30	0.44
1:A:823:ALA:HB1	1:A:824:PRO:HD2	2.00	0.44
1:A:568:GLN:O	1:A:569:LEU:HD22	2.18	0.43
1:A:785:LEU:O	1:A:786:LEU:HD23	2.17	0.43
1:A:788:ASN:O	1:A:791:VAL:HG23	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:902:THR:N	1:A:905:GLN:NE2	2.65	0.43
1:A:565:ASP:O	1:A:568:GLN:HB2	2.18	0.43
1:A:779:VAL:CG1	1:A:844:LEU:HD12	2.48	0.43
1:A:670:ASP:CG	1:A:673:ASN:HD22	2.27	0.43
1:A:635:LYS:HA	1:A:638:SER:OG	2.19	0.43
1:A:640:LEU:O	1:A:772:LYS:HE2	2.19	0.43
1:A:766:MET:HG2	1:A:794:ILE:HD13	2.01	0.42
1:A:843:ILE:HD13	1:A:893:TRP:HE1	1.83	0.42
1:A:634:LEU:O	1:A:637:MET:HB2	2.20	0.42
1:A:822:MET:HE3	1:A:832:TYR:HE2	1.85	0.42
1:A:787:THR:OG1	1:A:788:ASN:N	2.53	0.42
1:A:562:THR:O	1:A:562:THR:HG22	2.20	0.42
1:A:762:VAL:HG11	1:A:844:LEU:HD11	2.02	0.42
1:A:582:LEU:HD21	1:A:662:ILE:CD1	2.50	0.42
1:A:785:LEU:HD11	3:A:1000:6C3:H14	2.02	0.42
1:A:625:ASP:OD2	1:A:625:ASP:N	2.50	0.41
1:A:556:TYR:C	1:A:558:GLY:N	2.78	0.41
1:A:846:TRP:CD1	1:A:878:PRO:HG3	2.55	0.41
1:A:565:ASP:HA	1:A:566:PRO:HD2	1.80	0.41
1:A:817:LEU:HB3	1:A:822:MET:SD	2.59	0.41
1:A:768:PHE:O	1:A:772:LYS:HG2	2.20	0.41
1:A:845:LEU:HA	1:A:845:LEU:HD23	1.76	0.41
1:A:643:HIS:HB2	1:A:768:PHE:CE2	2.55	0.41
1:A:772:LYS:O	1:A:773:ASN:HB2	2.20	0.41
1:A:849:PHE:C	1:A:851:LEU:H	2.27	0.41
1:A:886:TYR:HE2	1:A:890:GLN:NE2	2.19	0.41
1:A:766:MET:N	1:A:794:ILE:HD11	2.36	0.41
1:A:762:VAL:HG11	1:A:844:LEU:CD1	2.51	0.40
1:A:865:PHE:O	1:A:868:LEU:HB2	2.21	0.40
1:A:654:THR:HA	1:A:659:VAL:CG2	2.38	0.40
1:A:602:PHE:HD2	1:A:610:VAL:HG13	1.85	0.40
1:A:885:ILE:HD12	1:A:885:ILE:HA	1.81	0.40
1:A:587:THR:O	1:A:587:THR:HG22	2.19	0.40
1:A:782:ARG:NH2	1:A:818:PRO:CB	2.84	0.40
1:A:895:LEU:HA	1:A:895:LEU:HD23	1.75	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	297/335 (89%)	254 (86%)	31 (10%)	12 (4%)	2 8

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	557	GLU
1	A	570	PRO
1	A	572	ASN
1	A	624	ALA
1	A	912	GLU
1	A	546	TYR
1	A	547	GLN
1	A	569	LEU
1	A	592	ALA
1	A	793	LYS
1	A	863	SER
1	A	608	ASP

5.3.2 Protein sidechains [i](#)

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	256/287 (89%)	235 (92%)	21 (8%)	10 33

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	546	TYR
1	A	554	GLU
1	A	562	THR
1	A	569	LEU
1	A	570	PRO
1	A	571	TYR
1	A	573	GLU
1	A	582	LEU
1	A	587	THR
1	A	596	VAL
1	A	604	LEU
1	A	611	LEU
1	A	625	ASP
1	A	631	MET
1	A	699	SER
1	A	785	LEU
1	A	794	ILE
1	A	829	ASP
1	A	885	ILE
1	A	887	SER
1	A	897	PRO

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	583	GLN
1	A	655	HIS
1	A	673	ASN
1	A	854	ASN
1	A	874	GLN
1	A	890	GLN
1	A	905	GLN
1	A	911	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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	6C3	A	1000	-	27,27,27	2.10	10 (37%)	34,39,39	1.82	5 (14%)
2	SO4	A	101	-	4,4,4	0.38	0	6,6,6	0.07	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
3	6C3	A	1000	-	-	4/8/8/8	0/4/4/4

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1000	6C3	C8-C9	5.14	1.55	1.44
3	A	1000	6C3	C16-C17	2.74	1.40	1.35
3	A	1000	6C3	C3-C2	2.69	1.44	1.39
3	A	1000	6C3	C8-C17	2.63	1.51	1.46
3	A	1000	6C3	C14-C15	2.57	1.43	1.39
3	A	1000	6C3	C6-C1	2.37	1.43	1.39
3	A	1000	6C3	C10-C15	2.33	1.42	1.39
3	A	1000	6C3	O2-N2	-2.20	1.38	1.42
3	A	1000	6C3	C2-C7	2.19	1.52	1.47
3	A	1000	6C3	C3-C4	2.09	1.41	1.38

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1000	6C3	O2-C17-C16	-8.44	104.28	109.86
3	A	1000	6C3	C17-O2-N2	3.39	112.15	108.38
3	A	1000	6C3	C3-C2-C1	-2.29	117.03	119.09
3	A	1000	6C3	C1-C2-C7	2.12	120.46	118.54
3	A	1000	6C3	C16-C17-C8	-2.00	122.66	131.56

There are no chirality outliers.

All (4) torsion outliers are listed below:

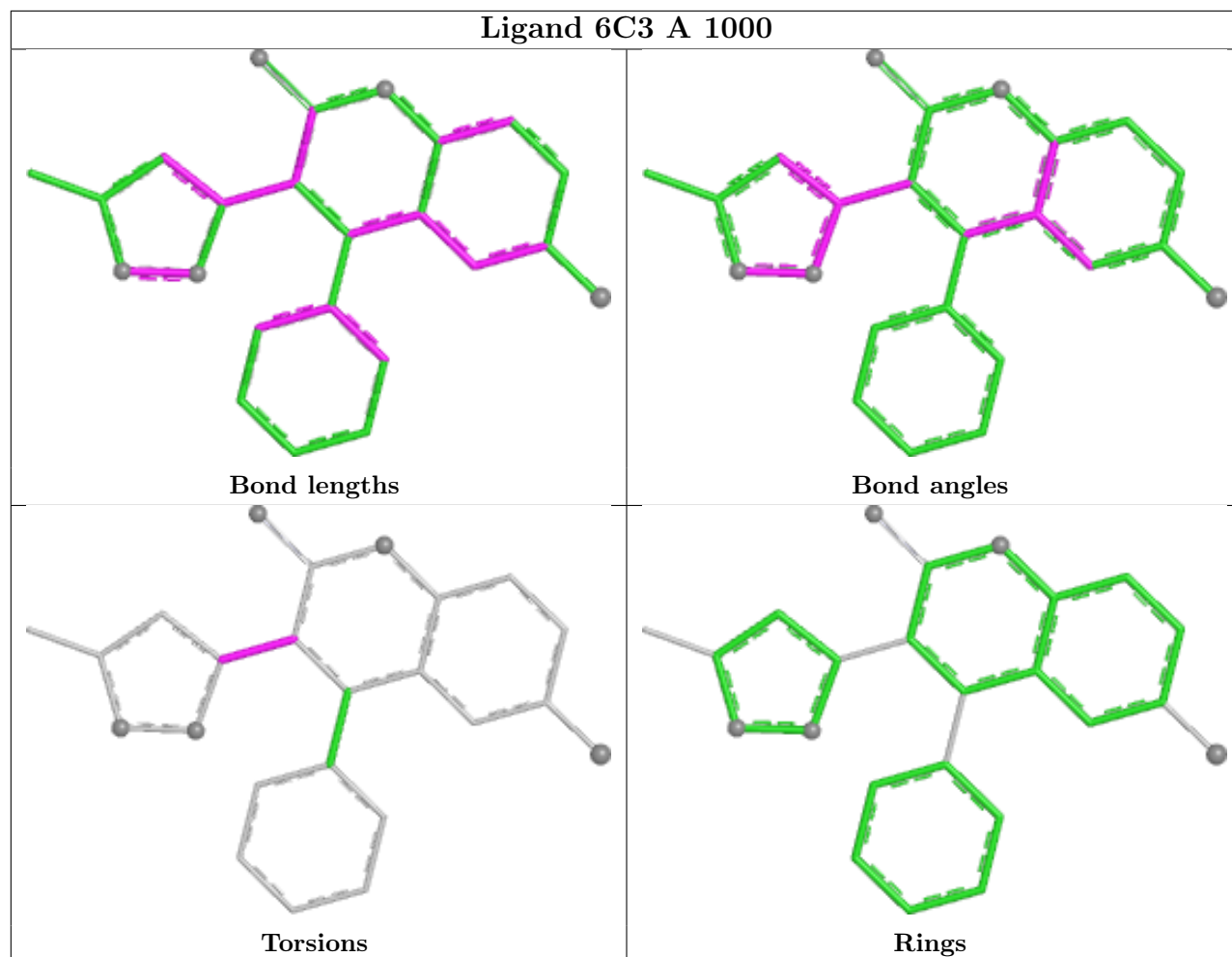
Mol	Chain	Res	Type	Atoms
3	A	1000	6C3	C16-C17-C8-C7
3	A	1000	6C3	O2-C17-C8-C7
3	A	1000	6C3	C16-C17-C8-C9
3	A	1000	6C3	O2-C17-C8-C9

There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1000	6C3	4	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 [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	303/335 (90%)	0.20	10 (3%) 49 39	9, 37, 70, 81	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	571	TYR	3.3
1	A	914	ALA	3.2
1	A	813	GLY	2.8
1	A	569	LEU	2.7
1	A	915	GLN	2.6
1	A	913	GLN	2.5
1	A	544	PRO	2.4
1	A	557	GLU	2.2
1	A	607	GLU	2.2
1	A	623	HIS	2.1

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](#)

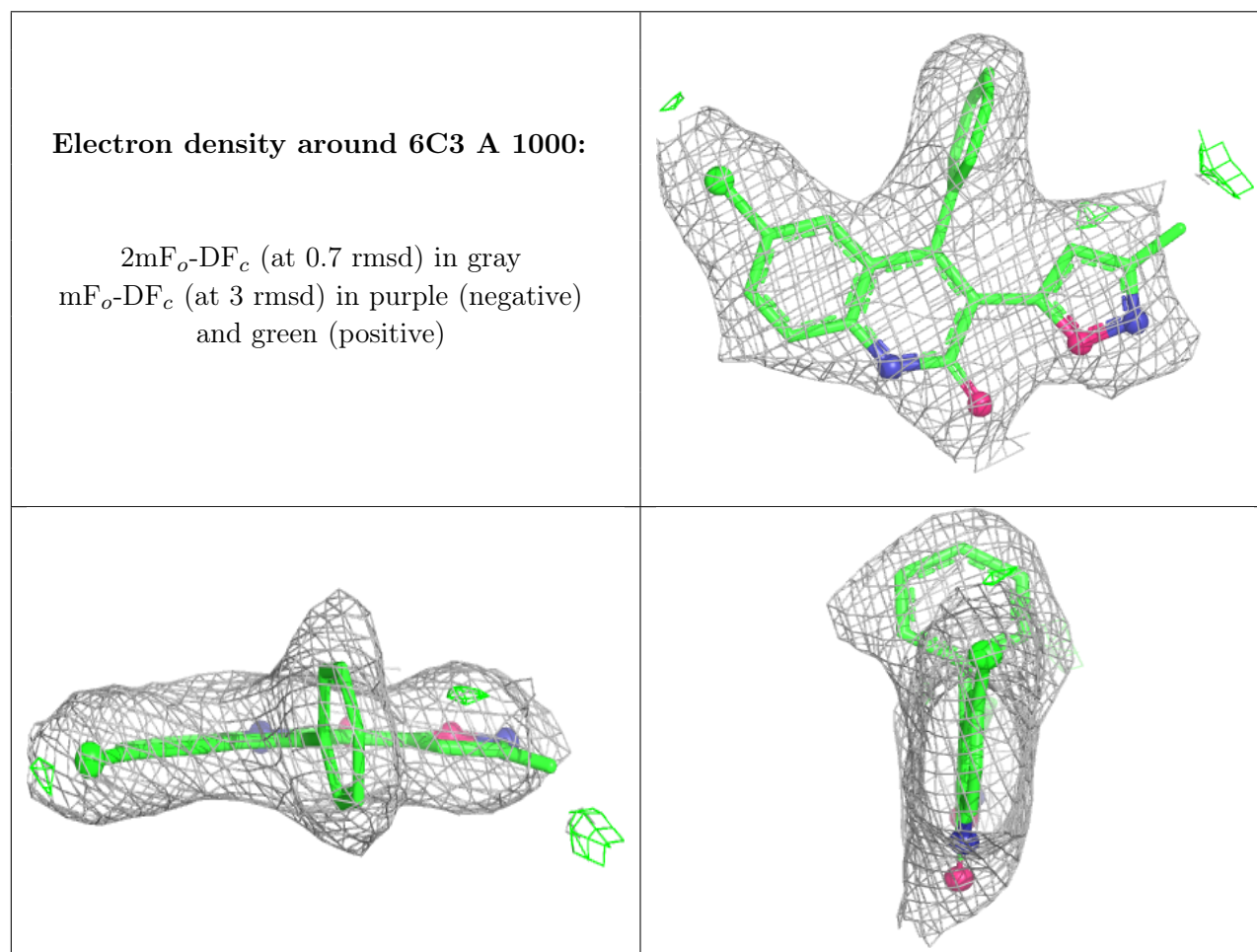
There are no oligosaccharides in this entry.

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	SO4	A	101	5/5	0.83	0.14	75,76,76,76	0
3	6C3	A	1000	24/24	0.91	0.11	23,28,37,40	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.