



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2026 – 11:09 PM UTC

PDB ID : 3SFY / pdb_00003sfy
Title : Cryptococcus neoformans protein farnesyltransferase in complex with FPT-II and ethylenediamine inhibitor 2
Authors : Hast, M.A.; Beese, L.S.
Deposited on : 2011-06-14
Resolution : 2.10 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

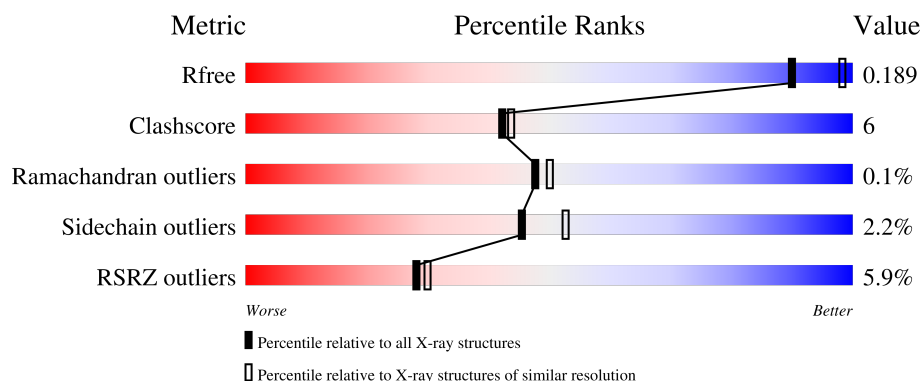
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	349	
2	B	520	

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 7079 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 *Cryptococcus neoformans* protein farnesyltransferase alpha subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	315	Total	C	N	O	S	0	0	0
			2621	1691	446	473	11			

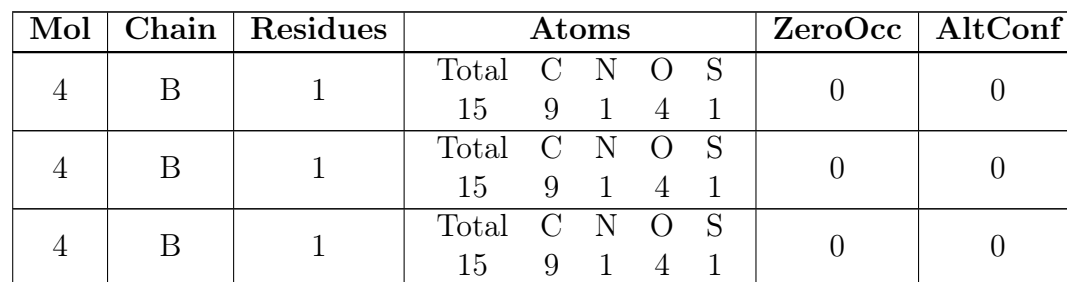
- Molecule 2 is a protein called *Cryptococcus neoformans* protein farnesyltransferase beta subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	482	Total	C	N	O	S	0	1	0
			3702	2345	646	696	15			

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Zn	0	0
			1	1		

- Molecule 4 is (2R)-3-(cyclohexylamino)-2-hydroxypropane-1-sulfonic acid (CCD ID: 3FX) (formula: C₉H₁₉NO₄S).

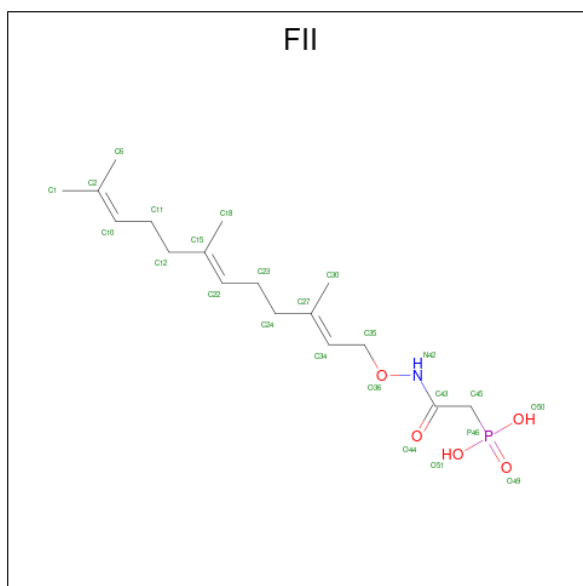


- Molecule 5 is N-(2-{(4-bromophenyl)[(1-methyl-1H-imidazol-5-yl)methyl]amino}ethyl)-1-methyl-N-(2-methylbenzyl)-1H-imidazole-4-sulfonamide (CCD ID: 3FY) (formula: $\text{C}_{25}\text{H}_{29}\text{BrN}_6\text{O}_2\text{S}$).



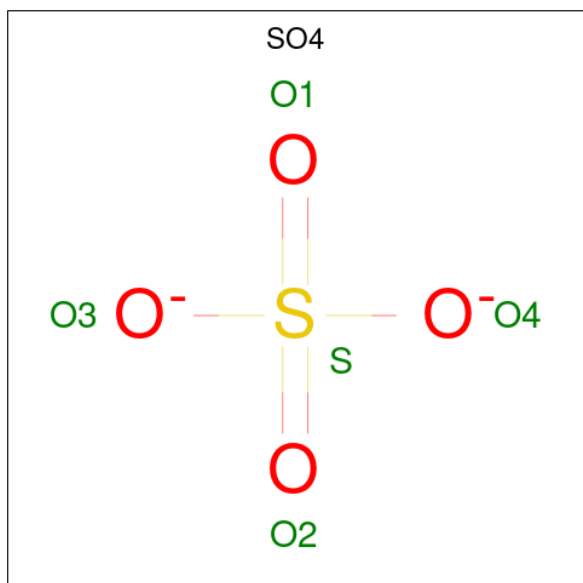
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	B	1	Total	Br	C	N	O	S	
			35	1	25	6	2	1	

- Molecule 6 is [(3,7,11-TRIMETHYL-DODECA-2,6,10-TRIENYLOXYCARBAMOYL)-METHYL]-PHOSPHONIC ACID (CCD ID: FII) (formula: $C_{17}H_{30}NO_5P$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	B	1	Total	C	N	O	P		
			24	17	1	5	1		

- Molecule 7 is SULFATE ION (CCD ID: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	B	1	Total	O	S	0	0
			5	4	1		

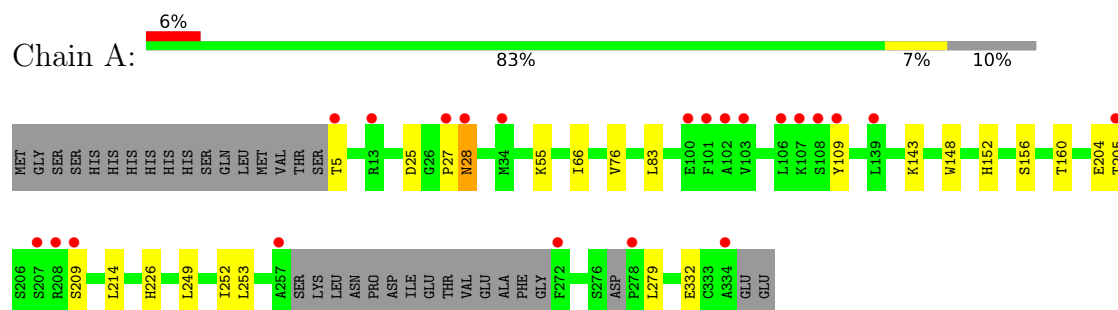
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	228	Total	O	0	0
			228	228		
8	B	418	Total	O	0	0
			418	418		

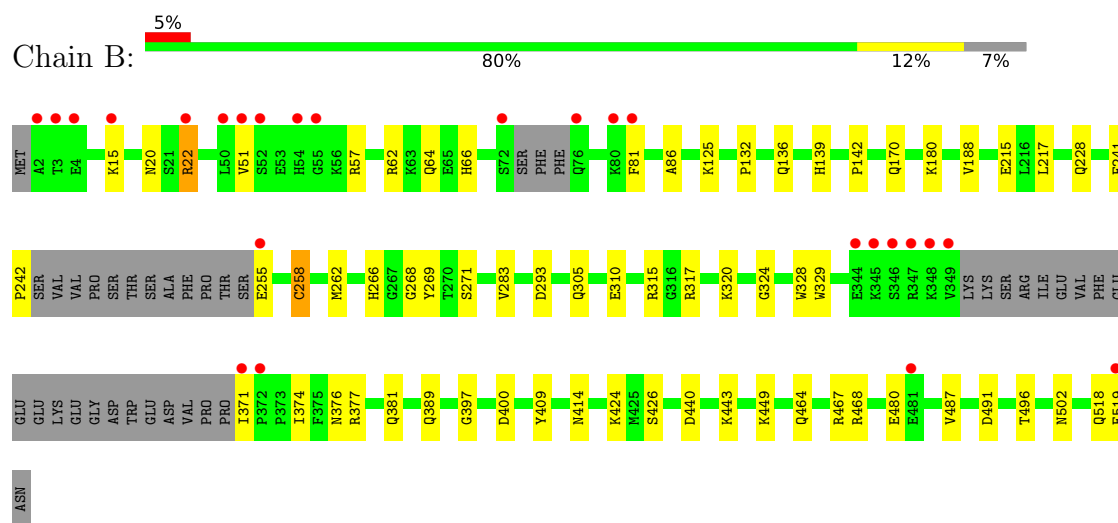
3 Residue-property plots [i](#)

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: *Cryptococcus neoformans* protein farnesyltransferase alpha subunit



- Molecule 2: *Cryptococcus neoformans* protein farnesyltransferase beta subunit



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	141.16Å 141.16Å 129.63Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.57 – 2.10 46.57 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.9 (46.57-2.10) 100.0 (46.57-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.55 (at 2.05Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.178 , 0.203 0.188 , 0.189	Depositor DCC
R_{free} test set	4123 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	27.0	Xtriage
Anisotropy	0.468	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 34.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7079	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.45% 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: 3FY, SO4, ZN, 3FX, FII

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	0.64	0/2703	0.82	0/3680
2	B	0.70	0/3796	0.84	0/5155
All	All	0.67	0/6499	0.83	0/8835

There are no bond length outliers.

There are no bond angle outliers.

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	2621	0	2544	19	0
2	B	3702	0	3633	62	0
3	B	1	0	0	0	0
4	B	45	0	57	4	0
5	B	35	0	29	5	0
6	B	24	0	28	5	0
7	B	5	0	0	1	0
8	A	228	0	0	3	0
8	B	418	0	0	17	0
All	All	7079	0	6291	80	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 (80) 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:28:ASN:HB3	8:B:2794:HOH:O	1.56	1.03
2:B:258:CYS:HB2	8:B:2540:HOH:O	1.63	0.97
2:B:496:THR:HG23	8:B:2053:HOH:O	1.70	0.92
2:B:440:ASP:H	2:B:518:GLN:HE22	1.10	0.90
2:B:228:GLN:HE22	2:B:293:ASP:H	1.19	0.89
2:B:464:GLN:HE22	2:B:467:ARG:HH11	1.15	0.88
2:B:62:ARG:H	2:B:389:GLN:HE22	1.27	0.82
2:B:170:GLN:HE22	2:B:449:LYS:H	1.22	0.82
1:A:143:LYS:HD3	2:B:320:LYS:HD3	1.61	0.80
1:A:109:TYR:HA	8:A:2596:HOH:O	1.81	0.79
2:B:305:GLN:HE22	2:B:376:ASN:H	1.33	0.77
2:B:255:GLU:HA	8:B:2757:HOH:O	1.86	0.74
2:B:266:HIS:HD2	2:B:268:GLY:H	1.37	0.73
2:B:440:ASP:H	2:B:518:GLN:NE2	1.85	0.72
2:B:66:HIS:HD1	2:B:397:GLY:H	1.39	0.70
2:B:170:GLN:NE2	2:B:449:LYS:H	1.93	0.67
1:A:148:TRP:O	1:A:152:HIS:HD2	1.78	0.66
1:A:226:HIS:H	2:B:20:ASN:HD21	1.42	0.66
2:B:262[B]:MET:HA	2:B:262[B]:MET:HE2	1.79	0.65
2:B:496:THR:HG21	8:B:2600:HOH:O	1.97	0.63
1:A:143:LYS:CD	2:B:320:LYS:HD3	2.30	0.61
4:B:522:3FX:HAJA	8:B:1900:HOH:O	2.03	0.58
1:A:28:ASN:HA	8:B:2137:HOH:O	2.04	0.57
2:B:180:LYS:HE2	2:B:215:GLU:O	2.04	0.57
2:B:266:HIS:CD2	2:B:268:GLY:H	2.21	0.57
1:A:156:SER:O	1:A:160:THR:HG23	2.05	0.57
5:B:525:3FY:HAC	8:B:2751:HOH:O	2.05	0.56
2:B:22:ARG:HB2	8:B:2748:HOH:O	2.06	0.55
2:B:468:ARG:HD2	8:B:2150:HOH:O	2.05	0.55
2:B:57:ARG:HD2	8:B:1608:HOH:O	2.08	0.53
2:B:62:ARG:H	2:B:389:GLN:NE2	2.02	0.53
2:B:266:HIS:HE1	6:B:526:FII:O49	1.91	0.53
2:B:271:SER:HB2	2:B:329:TRP:O	2.09	0.53
2:B:64:GLN:HB2	4:B:523:3FX:HAK	1.92	0.52
2:B:381:GLN:HE22	2:B:487:VAL:H	1.55	0.52
2:B:170:GLN:HE22	2:B:449:LYS:N	2.00	0.52
2:B:86:ALA:HA	2:B:136:GLN:HE22	1.74	0.52
1:A:249:LEU:HG	1:A:253:LEU:HD13	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:125:LYS:H	4:B:524:3FX:HOAB	1.57	0.51
2:B:305:GLN:HE22	2:B:376:ASN:N	2.05	0.51
2:B:464:GLN:HE22	2:B:467:ARG:NH1	1.97	0.50
1:A:226:HIS:H	2:B:20:ASN:ND2	2.08	0.50
2:B:440:ASP:N	2:B:518:GLN:HE22	1.93	0.50
5:B:525:3FY:HAAA	8:B:2916:HOH:O	2.12	0.50
2:B:139:HIS:CE1	2:B:188:VAL:HB	2.46	0.49
2:B:139:HIS:HB3	2:B:142:PRO:HG2	1.93	0.49
1:A:152:HIS:HE1	8:A:1787:HOH:O	1.96	0.48
2:B:317:ARG:HG3	2:B:320:LYS:HG3	1.96	0.48
2:B:518:GLN:O	2:B:519:GLU:C	2.57	0.47
2:B:320:LYS:HE3	6:B:526:FII:O50	2.14	0.47
2:B:258:CYS:CB	8:B:2540:HOH:O	2.39	0.47
2:B:305:GLN:NE2	2:B:376:ASN:H	2.07	0.47
2:B:328:TRP:CZ2	2:B:502:ASN:HB2	2.49	0.47
2:B:242:PRO:C	8:B:2634:HOH:O	2.58	0.47
2:B:255:GLU:CA	8:B:2757:HOH:O	2.55	0.47
2:B:142:PRO:HD3	7:B:527:SO4:O2	2.15	0.47
1:A:109:TYR:CE2	5:B:525:3FY:HAR	2.51	0.46
1:A:27:PRO:HA	8:A:2714:HOH:O	2.14	0.46
2:B:426:SER:OG	2:B:480:GLU:OE1	2.30	0.46
2:B:324:GLY:HA3	2:B:414:ASN:HD21	1.82	0.45
2:B:377:ARG:NH1	2:B:424:LYS:HD2	2.31	0.45
2:B:269:TYR:CE2	6:B:526:FII:H302	2.52	0.45
1:A:55:LYS:HD3	1:A:83:LEU:HD23	1.98	0.44
2:B:381:GLN:NE2	2:B:487:VAL:H	2.14	0.44
2:B:409:TYR:CE2	5:B:525:3FY:HAO	2.53	0.44
2:B:310:GLU:OE1	2:B:315:ARG:NH2	2.43	0.43
6:B:526:FII:H111	6:B:526:FII:H221	1.67	0.43
2:B:491:ASP:OD2	4:B:522:3FX:NAL	2.52	0.43
2:B:496:THR:HG23	2:B:496:THR:O	2.18	0.43
1:A:55:LYS:HA	1:A:83:LEU:CD2	2.49	0.43
2:B:139:HIS:HB3	2:B:142:PRO:CG	2.48	0.42
2:B:255:GLU:HB3	8:B:2569:HOH:O	2.19	0.42
6:B:526:FII:H232	6:B:526:FII:H181	1.80	0.42
1:A:160:THR:HG22	2:B:241:PHE:HD2	1.84	0.42
1:A:109:TYR:CD2	5:B:525:3FY:HAR	2.55	0.42
2:B:81:PHE:HZ	8:B:2746:HOH:O	2.03	0.41
2:B:400:ASP:C	2:B:400:ASP:OD1	2.63	0.41
1:A:66:ILE:HG21	1:A:76:VAL:HG21	2.02	0.41
2:B:217:LEU:HD12	2:B:283:VAL:HG11	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:25:ASP:HB3	2:B:132:PRO:HG3	2.04	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	309/349 (88%)	298 (96%)	11 (4%)	0	100	100
2	B	475/520 (91%)	466 (98%)	8 (2%)	1 (0%)	43	44
All	All	784/869 (90%)	764 (97%)	19 (2%)	1 (0%)	48	50

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	51	VAL

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	285/316 (90%)	276 (97%)	9 (3%)	34	38
2	B	401/436 (92%)	395 (98%)	6 (2%)	57	65
All	All	686/752 (91%)	671 (98%)	15 (2%)	45	53

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	THR
1	A	28	ASN
1	A	204	GLU
1	A	205	THR
1	A	209	SER
1	A	214	LEU
1	A	252	ILE
1	A	279	LEU
1	A	332	GLU
2	B	15	LYS
2	B	22	ARG
2	B	258	CYS
2	B	371	ILE
2	B	374	ILE
2	B	443	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	104	GLN
1	A	105	ASN
1	A	110	GLN
1	A	152	HIS
1	A	168	GLN
1	A	320	GLN
2	B	20	ASN
2	B	76	GLN
2	B	108	GLN
2	B	136	GLN
2	B	170	GLN
2	B	219	ASN
2	B	228	GLN
2	B	266	HIS
2	B	304	GLN
2	B	305	GLN
2	B	381	GLN
2	B	389	GLN
2	B	413	ASN
2	B	414	ASN
2	B	464	GLN
2	B	495	ASN

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Mol	Chain	Res	Type
2	B	518	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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 7 ligands modelled in this entry, 1 is monoatomic - leaving 6 for Mogul analysis.

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$
4	3FX	B	522	-	14,15,15	0.87	0	16,20,20	1.10	1 (6%)
6	FII	B	526	-	23,23,23	2.36	9 (39%)	27,29,29	1.46	6 (22%)
4	3FX	B	524	-	14,15,15	0.97	0	16,20,20	1.09	2 (12%)
7	SO4	B	527	-	4,4,4	0.28	0	6,6,6	0.37	0
5	3FY	B	525	3	36,38,38	2.11	8 (22%)	44,54,54	3.55	11 (25%)
4	3FX	B	523	-	14,15,15	0.95	0	16,20,20	0.98	1 (6%)

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	3FX	B	522	-	-	2/10/18/18	0/1/1/1
6	FII	B	526	-	-	5/23/24/24	-
4	3FX	B	524	-	-	6/10/18/18	0/1/1/1
5	3FY	B	525	3	-	6/27/29/29	0/4/4/4
4	3FX	B	523	-	-	7/10/18/18	0/1/1/1

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	525	3FY	OAE-SBI	7.29	1.50	1.43
5	B	525	3FY	OAD-SBI	6.78	1.50	1.43
6	B	526	FII	P46-O49	5.99	1.62	1.50
6	B	526	FII	C23-C22	-3.64	1.39	1.50
6	B	526	FII	C11-C10	-3.62	1.39	1.50
6	B	526	FII	C35-C34	-3.43	1.39	1.49
6	B	526	FII	P46-O50	3.32	1.62	1.55
6	B	526	FII	P46-O51	-2.91	1.48	1.55
6	B	526	FII	C34-C27	2.76	1.39	1.33
5	B	525	3FY	CAO-CBC	-2.76	1.33	1.36
6	B	526	FII	C22-C15	2.74	1.39	1.33
5	B	525	3FY	CBC-NBH	-2.69	1.33	1.38
5	B	525	3FY	CAV-CBC	2.60	1.53	1.50
5	B	525	3FY	CAC-NBH	2.57	1.51	1.46
6	B	526	FII	C10-C2	2.33	1.39	1.32
5	B	525	3FY	CAO-NAW	-2.18	1.33	1.37
5	B	525	3FY	CAR-NBG	-2.18	1.33	1.37

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	525	3FY	NBG-CAQ-NAX	-13.74	107.40	112.66
5	B	525	3FY	NBH-CAP-NAW	-10.71	106.56	112.94
5	B	525	3FY	OAE-SBI-OAD	-8.15	107.99	119.69
5	B	525	3FY	CAR-NBG-CAQ	7.80	109.79	106.71
5	B	525	3FY	CAU-NBF-SBI	-4.32	107.28	117.47
5	B	525	3FY	CAQ-NAX-CBD	4.25	107.30	103.39
5	B	525	3FY	CAS-CAT-NBF	-3.72	106.09	112.43
5	B	525	3FY	CAT-NBF-SBI	-3.48	110.74	117.88
6	B	526	FII	O51-P46-C45	3.11	113.37	106.84
5	B	525	3FY	CAU-NBF-CAT	-3.05	111.72	116.98
4	B	524	3FX	OAD-SAO-CAK	2.78	110.91	106.76
6	B	526	FII	C30-C27-C24	2.75	120.01	115.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	523	3FX	OAB-SAO-CAK	2.75	111.29	105.97
6	B	526	FII	C18-C15-C12	2.74	119.98	115.23
6	B	526	FII	O50-P46-O49	-2.73	105.32	112.39
5	B	525	3FY	CAO-NAW-CAP	2.73	108.81	105.24
4	B	522	3FX	OAB-SAO-CAK	2.55	110.91	105.97
6	B	526	FII	O49-P46-C45	-2.49	105.35	111.10
6	B	526	FII	C6-C2-C1	2.36	120.02	114.59
5	B	525	3FY	CAK-CAZ-CAJ	-2.27	117.89	121.33
4	B	524	3FX	OAA-SAO-CAK	2.26	110.13	106.76

There are no chirality outliers.

All (26) torsion outliers are listed below:

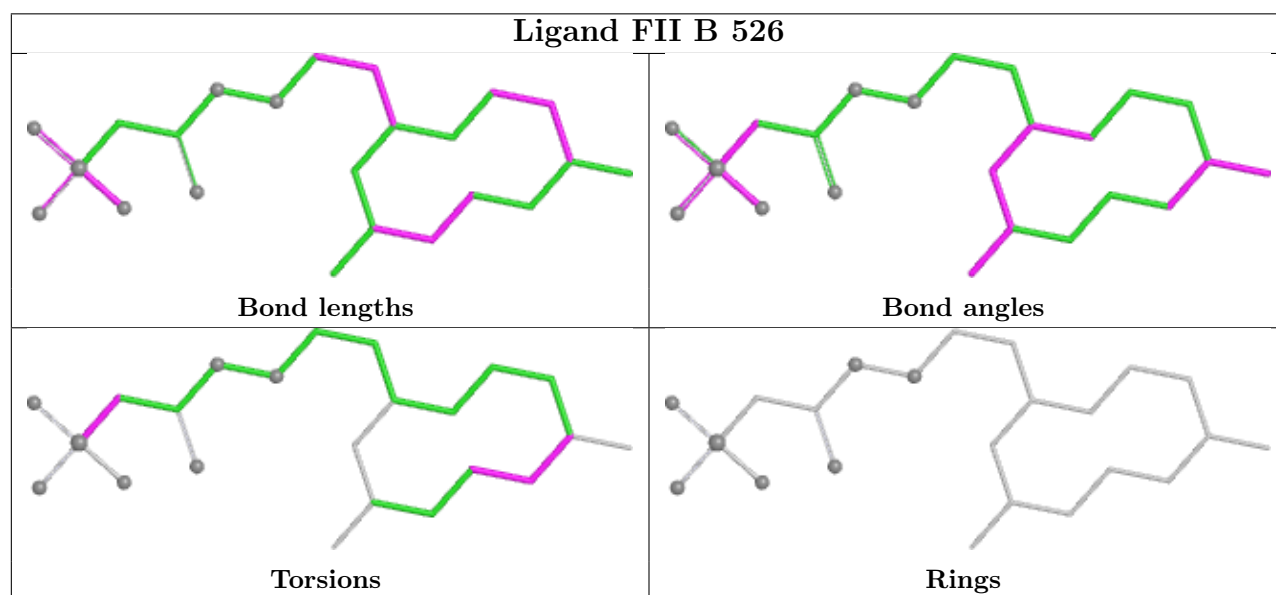
Mol	Chain	Res	Type	Atoms
4	B	523	3FX	NAL-CAJ-CAM-OAC
4	B	523	3FX	NAL-CAJ-CAM-CAK
4	B	523	3FX	CAM-CAK-SAO-OAA
4	B	523	3FX	CAM-CAK-SAO-OAB
4	B	523	3FX	CAM-CAK-SAO-OAD
4	B	524	3FX	NAL-CAJ-CAM-OAC
4	B	524	3FX	NAL-CAJ-CAM-CAK
5	B	525	3FY	CAR-CBD-SBI-OAE
5	B	525	3FY	NAX-CBD-SBI-OAE
6	B	526	FII	C10-C11-C12-C15
5	B	525	3FY	CAU-NBF-SBI-OAD
4	B	523	3FX	CAH-CAN-NAL-CAJ
4	B	524	3FX	CAI-CAN-NAL-CAJ
5	B	525	3FY	CAU-NBF-SBI-CBD
6	B	526	FII	C11-C12-C15-C22
6	B	526	FII	C11-C12-C15-C18
4	B	522	3FX	SAO-CAK-CAM-OAC
4	B	524	3FX	SAO-CAK-CAM-OAC
6	B	526	FII	C43-C45-P46-O51
5	B	525	3FY	CAT-NBF-SBI-OAE
4	B	522	3FX	SAO-CAK-CAM-CAJ
4	B	524	3FX	SAO-CAK-CAM-CAJ
5	B	525	3FY	CAS-CAT-NBF-CAU
4	B	523	3FX	CAI-CAN-NAL-CAJ
4	B	524	3FX	CAH-CAN-NAL-CAJ
6	B	526	FII	C43-C45-P46-O49

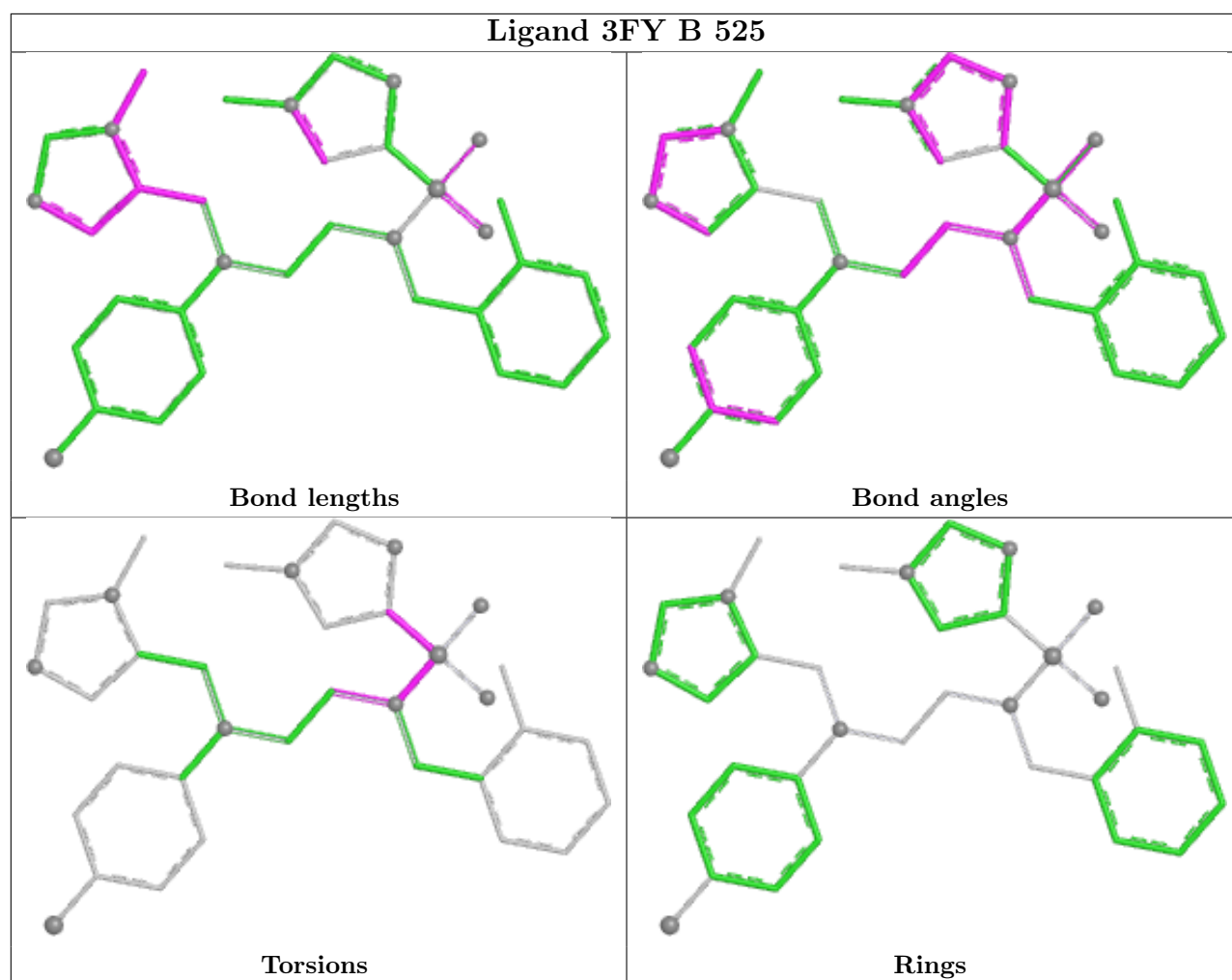
There are no ring outliers.

6 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	522	3FX	2	0
6	B	526	FII	5	0
4	B	524	3FX	1	0
7	B	527	SO4	1	0
5	B	525	3FY	5	0
4	B	523	3FX	1	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	315/349 (90%)	0.32	22 (6%) 22 24	24, 39, 62, 73	0
2	B	482/520 (92%)	-0.12	25 (5%) 33 35	17, 30, 51, 82	1 (0%)
All	All	797/869 (91%)	0.05	47 (5%) 28 30	17, 34, 59, 82	1 (0%)

All (47) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	2	ALA	6.7
2	B	51	VAL	6.3
1	A	109	TYR	5.3
1	A	107	LYS	4.9
2	B	371	ILE	4.8
1	A	103	VAL	4.6
1	A	334	ALA	4.4
2	B	349	VAL	4.2
1	A	106	LEU	4.0
1	A	27	PRO	3.9
2	B	255	GLU	3.8
2	B	50	LEU	3.8
1	A	28	ASN	3.7
2	B	76	GLN	3.7
1	A	139	LEU	3.6
2	B	481	GLU	3.5
1	A	13	ARG	3.4
1	A	108	SER	3.3
1	A	272	PHE	3.3
2	B	519	GLU	3.3
1	A	102	ALA	3.2
1	A	5	THR	3.2
2	B	81	PHE	3.2
2	B	372	PRO	3.0

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Mol	Chain	Res	Type	RSRZ
2	B	346	SER	3.0
1	A	278	PRO	2.9
1	A	208	ARG	2.8
2	B	52	SER	2.7
1	A	101	PHE	2.7
2	B	345	LYS	2.5
1	A	209	SER	2.5
1	A	34	MET	2.5
1	A	205	THR	2.5
2	B	348	LYS	2.5
1	A	207	SER	2.4
2	B	72	SER	2.4
2	B	347	ARG	2.4
2	B	22	ARG	2.3
2	B	4	GLU	2.3
2	B	3	THR	2.3
2	B	54	HIS	2.2
1	A	257	ALA	2.1
2	B	15	LYS	2.1
2	B	55	GLY	2.1
2	B	80	LYS	2.1
1	A	100	GLU	2.0
2	B	344	GLU	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.

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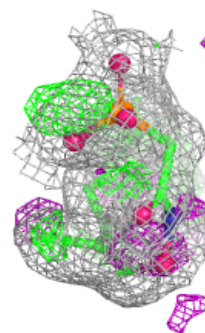
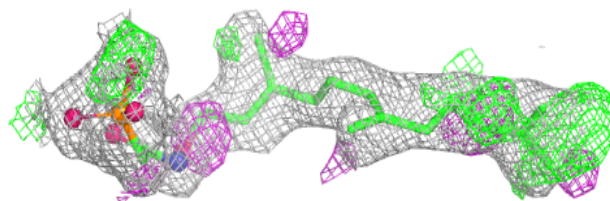
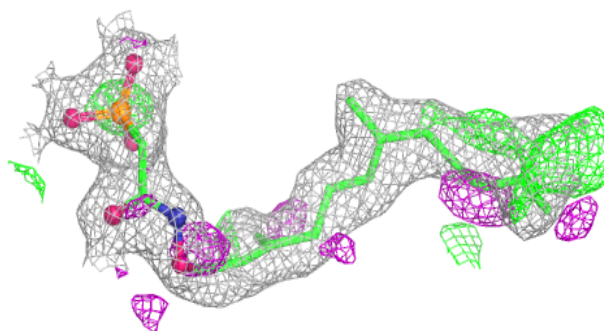
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	3FX	B	524	15/15	0.83	0.19	55,62,72,72	0
7	SO4	B	527	5/5	0.90	0.24	51,53,56,56	0
6	FII	B	526	24/24	0.91	0.16	28,33,35,38	0
5	3FY	B	525	35/35	0.91	0.21	37,54,62,62	35
4	3FX	B	523	15/15	0.95	0.13	40,52,58,58	0
4	3FX	B	522	15/15	0.97	0.07	28,29,32,33	0
3	ZN	B	521	1/1	1.00	0.03	28,28,28,28	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.

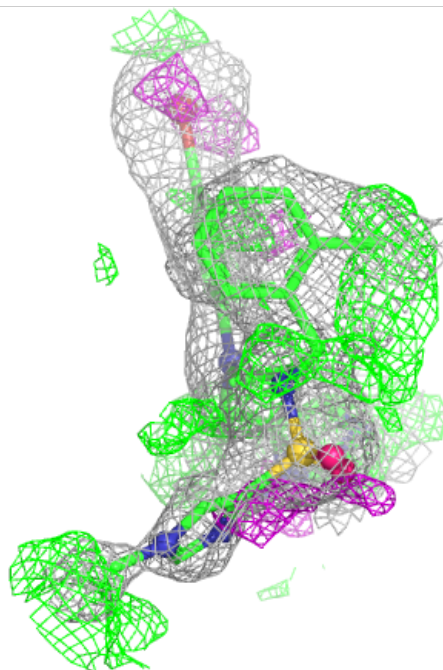
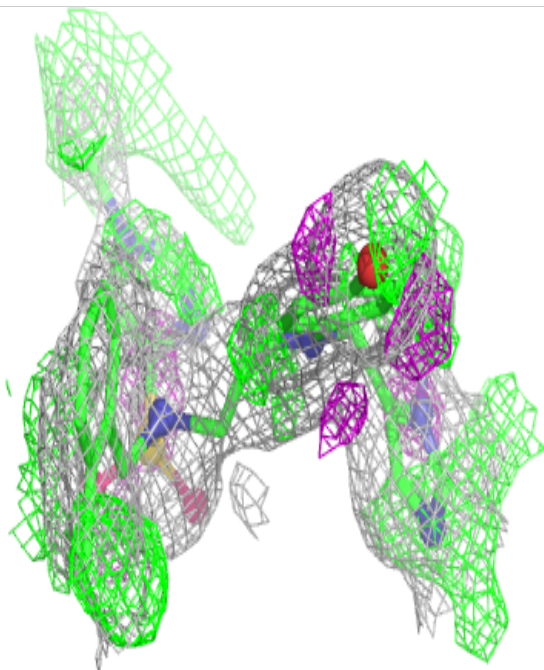
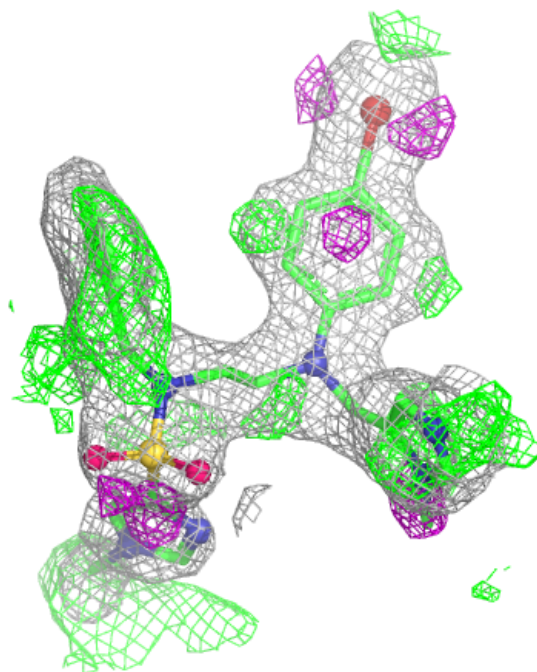
Electron density around FII B 526:

2mF_o-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around 3FY B 525:

$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers ⓘ

There are no such residues in this entry.