



Full wwPDB X-ray Structure Validation Report

Mar 31, 2014 – 03:37 PM BST

PDB ID : 4L9P
Title : Crystal structure of Aspergillus fumigatus protein farnesyltransferase complexed with the FII analog, FPT-II, and the KCVVM peptide
Authors : Mabanglo, M.F.; Hast, M.A.; Beese, L.S.
Deposited on : 2013-06-18
Resolution : 1.45 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

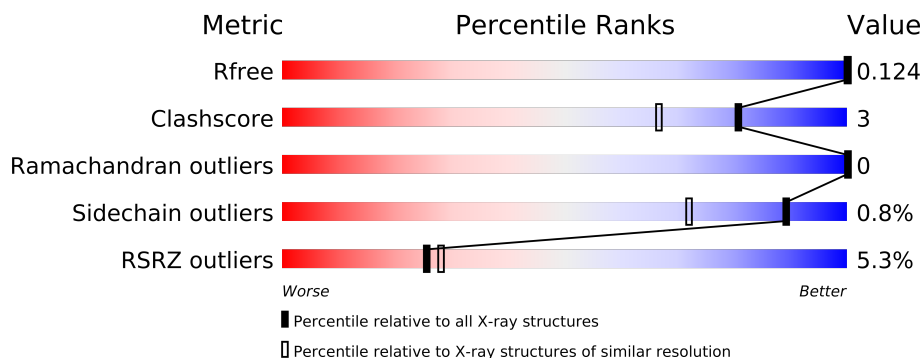
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.15 2013
Xtriage (Phenix)	:	dev-1323
EDS	:	stable23004
Percentile statistics	:	21963
Refmac	:	5.8.0049
CCP4	:	6.3.0 (Settle)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP)	:	stable23004

1 Overall quality at a glance

The reported resolution of this entry is 1.45 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}	66092	2476 (1.50-1.42)
Clashscore	79885	2834 (1.50-1.42)
Ramachandran outliers	78287	2769 (1.50-1.42)
Sidechain outliers	78261	2767 (1.50-1.42)
RSRZ outliers	66119	2477 (1.50-1.42)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	367	
2	B	519	
3	C	5	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
5	EDO	A	402	-	X
5	EDO	A	406	-	X
5	EDO	B	605	-	X
5	EDO	B	606	-	X
5	EDO	B	607	-	X

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 7369 atoms, of which 0 are hydrogen and 0 are deuterium.

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 CaaX farnesyltransferase alpha subunit Ram2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	335	2861	1821	501	529	10	0	17	0

There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-13	MET	-	INITIATING METHIONINE	UNP Q4WP27
A	-12	GLY	-	EXPRESSION TAG	UNP Q4WP27
A	-11	SER	-	EXPRESSION TAG	UNP Q4WP27
A	-10	SER	-	EXPRESSION TAG	UNP Q4WP27
A	-9	HIS	-	EXPRESSION TAG	UNP Q4WP27
A	-8	HIS	-	EXPRESSION TAG	UNP Q4WP27
A	-7	HIS	-	EXPRESSION TAG	UNP Q4WP27
A	-6	HIS	-	EXPRESSION TAG	UNP Q4WP27
A	-5	HIS	-	EXPRESSION TAG	UNP Q4WP27
A	-4	HIS	-	EXPRESSION TAG	UNP Q4WP27
A	-3	SER	-	EXPRESSION TAG	UNP Q4WP27
A	-2	GLN	-	EXPRESSION TAG	UNP Q4WP27
A	-1	ASP	-	EXPRESSION TAG	UNP Q4WP27
A	0	PRO	-	EXPRESSION TAG	UNP Q4WP27
A	146	SER	ASN	ENGINEERED MUTATION	UNP Q4WP27

- Molecule 2 is a protein called CaaX farnesyltransferase beta subunit Ram1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	452	3628	2303	613	687	25	0	21	0

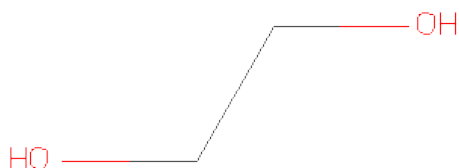
- Molecule 3 is a protein called LYS-CYS-VAL-VAL-MET (CAAX peptide).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	5	Total	C	N	O	S	0	5	0
			38	24	6	6	2			

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	2	Total	Cl	0	0
			2	2		
4	A	1	Total	Cl	0	0
			1	1		

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



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

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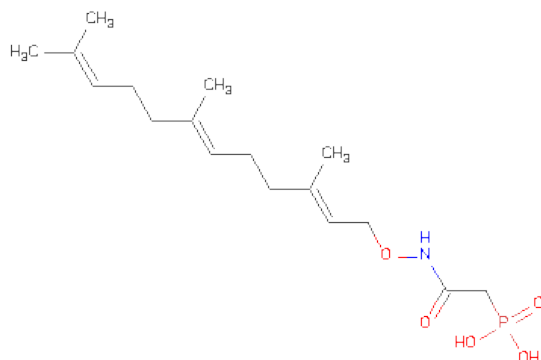
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 6 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 7 is [(3,7,11-TRIMETHYL-DODECA-2,6,10-TRIENYLOXYCARBAMOYL)-METHYL]-PHOSPHONICACID (three-letter code: FII) (formula: C₁₇H₃₀NO₅P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	B	1	Total	C	N	O	P	0	0
			24	17	1	5	1		

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	349	Total	O	0	0
			349	349		
8	B	420	Total	O	0	0
			420	420		

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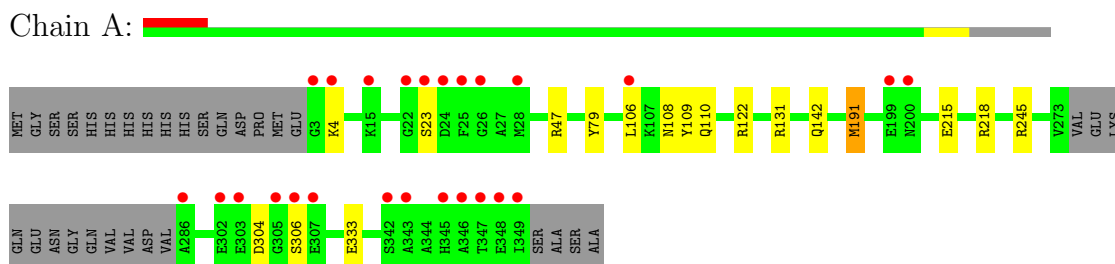
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	C	9	Total	O	0	0
			9	9		

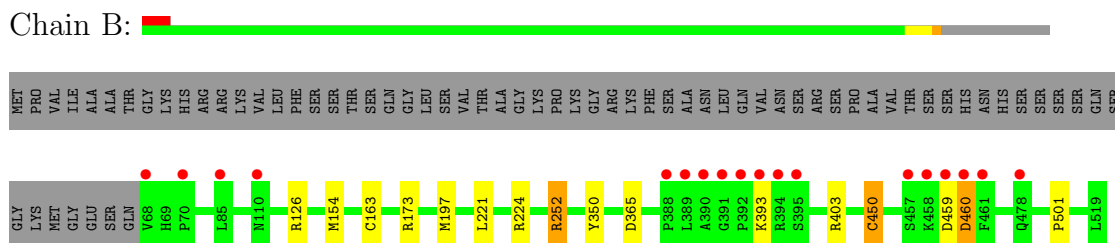
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: CaaX farnesyltransferase alpha subunit Ram2



- Molecule 2: CaaX farnesyltransferase beta subunit Ram1



- Molecule 3: LYS-CYS-VAL-VAL-MET (CAAX peptide)



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	63.24Å 90.33Å 83.01Å 90.00° 111.01° 90.00°	Depositor
Resolution (Å)	22.42 – 1.45 22.42 – 1.45	Depositor EDS
% Data completeness (in resolution range)	98.9 (22.42-1.45) 93.8 (22.42-1.45)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.19 (at 1.44Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8_1069)	Depositor
R, R_{free}	0.125 , 0.152 0.123 , 0.124	Depositor DCC
R_{free} test set	7185 reflections (4.94%)	DCC
Wilson B-factor (Å ²)	12.8	Xtriage
Anisotropy	0.374	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 40.7	EDS
Estimated twinning fraction	0.016 for h,-k,-h-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 153559 reflections	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	7369	wwPDB-VP
Average B, all atoms (Å ²)	14.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.04% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, FII, EDO, CL

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	0.86	1/2972 (0.0%)	0.90	5/4028 (0.1%)
2	B	0.94	2/3752 (0.1%)	0.99	14/5103 (0.3%)
3	C	0.86	0/37	1.21	0/46
All	All	0.91	3/6761 (0.0%)	0.95	19/9177 (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	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	450[A]	CYS	CB-SG	-8.48	1.67	1.82
2	B	450[B]	CYS	CB-SG	-8.48	1.67	1.82
1	A	79	TYR	CD1-CE1	5.08	1.47	1.39

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	224	ARG	NE-CZ-NH2	-13.58	113.51	120.30
2	B	460	ASP	CB-CG-OD1	-9.58	109.68	118.30
2	B	224	ARG	NE-CZ-NH1	8.62	124.61	120.30
2	B	126	ARG	NE-CZ-NH2	-8.17	116.21	120.30
1	A	47	ARG	NE-CZ-NH2	-7.41	116.59	120.30
2	B	350	TYR	CB-CG-CD2	-7.22	116.67	121.00
2	B	403	ARG	NE-CZ-NH2	-6.13	117.24	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	154[A]	MET	CG-SD-CE	-6.04	90.53	100.20
2	B	154[B]	MET	CG-SD-CE	-6.04	90.53	100.20
1	A	245[A]	ARG	NE-CZ-NH2	-5.97	117.32	120.30
1	A	245[B]	ARG	NE-CZ-NH2	-5.97	117.32	120.30
2	B	459	ASP	CB-CG-OD2	-5.85	113.04	118.30
1	A	131	ARG	NE-CZ-NH2	-5.82	117.39	120.30
2	B	365	ASP	CB-CG-OD1	5.75	123.48	118.30
1	A	47	ARG	NE-CZ-NH1	5.73	123.17	120.30
2	B	252	ARG	NE-CZ-NH2	-5.46	117.57	120.30
2	B	173	ARG	NE-CZ-NH2	-5.23	117.68	120.30
2	B	403	ARG	NE-CZ-NH1	5.17	122.88	120.30
2	B	460	ASP	OD1-CG-OD2	5.10	132.99	123.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	306	SER	Peptide

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2861	0	0	16	0
2	B	3628	0	0	3	0
3	C	38	0	0	2	0
4	A	1	0	0	0	0
4	B	2	0	0	0	0
5	A	20	0	0	0	0
5	B	16	0	0	1	0
6	B	1	0	0	0	0
7	B	24	0	0	0	0
8	A	349	0	0	14	0
8	B	420	0	0	2	0
8	C	9	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	7369	0	0	22	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 3.

All (22) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:191:MET:SD	8:A:749:HOH:O	2.40	0.79
1:A:4:LYS:NZ	8:A:789:HOH:O	2.15	0.79
1:A:109[B]:TYR:OH	8:A:750:HOH:O	2.03	0.74
1:A:215[B]:GLU:OE1	8:A:609:HOH:O	2.08	0.72
3:C:7[C]:LYS:N	8:C:107:HOH:O	2.23	0.70
2:B:450[B]:CYS:SG	8:B:991:HOH:O	2.47	0.70
1:A:110[A]:GLN:OE1	8:A:799:HOH:O	2.11	0.68
3:C:11[C]:MET:CB	8:C:109:HOH:O	2.41	0.68
5:B:604:EDO:O1	8:B:1048:HOH:O	2.13	0.65
1:A:142:GLN:NE2	8:A:822:HOH:O	2.30	0.65
8:A:750:HOH:O	2:B:252:ARG:NE	2.31	0.63
1:A:218:ARG:NH2	8:A:657:HOH:O	2.33	0.62
1:A:333[A]:GLU:OE1	8:A:632:HOH:O	2.16	0.61
1:A:106[B]:LEU:O	8:A:771:HOH:O	2.16	0.61
1:A:304:ASP:OD1	8:A:824:HOH:O	2.17	0.60
1:A:106[A]:LEU:N	1:A:106[A]:LEU:CD2	2.71	0.54
1:A:108[A]:ASN:ND2	1:A:110[A]:GLN:OE1	2.41	0.54
1:A:109[B]:TYR:CE2	8:A:750:HOH:O	2.64	0.49
1:A:23:SER:O	1:A:23:SER:OG	2.34	0.45
1:A:110[A]:GLN:CD	8:A:799:HOH:O	2.55	0.43
2:B:163[B]:CYS:SG	2:B:501:PRO:CG	3.07	0.42
1:A:122:ARG:NE	8:A:652:HOH:O	2.52	0.42

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	348/367 (95%)	339 (97%)	9 (3%)	0	100	100
2	B	471/519 (91%)	465 (99%)	6 (1%)	0	100	100
3	C	3/5 (60%)	3 (100%)	0	0	100	100
All	All	822/891 (92%)	807 (98%)	15 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	304/315 (96%)	303 (100%)	1 (0%)	96	88
2	B	393/428 (92%)	389 (99%)	4 (1%)	85	64
3	C	5/5 (100%)	5 (100%)	0	100	100
All	All	702/748 (94%)	697 (99%)	5 (1%)	89	75

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	191	MET
2	B	197	MET
2	B	221	LEU
2	B	393	LYS
2	B	460	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA chains 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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 14 ligands modelled in this entry, 4 are monoatomic - leaving 10 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	$\# Z > 2$
5	EDO	A	402	-	3,3,3	0.63	0	2,2,2	0.51	0
5	EDO	A	403	-	3,3,3	0.72	0	2,2,2	0.54	0
5	EDO	A	404	-	3,3,3	0.43	0	2,2,2	0.15	0
5	EDO	A	405	-	3,3,3	0.65	0	2,2,2	0.56	0
5	EDO	A	406	-	3,3,3	0.44	0	2,2,2	0.54	0
5	EDO	B	604	-	3,3,3	0.25	0	2,2,2	0.72	0
5	EDO	B	605	-	3,3,3	0.47	0	2,2,2	0.86	0
5	EDO	B	606	-	3,3,3	0.55	0	2,2,2	1.04	0
5	EDO	B	607	-	3,3,3	0.62	0	2,2,2	0.12	0
7	FII	B	608	-	23,23,23	3.08	6 (26%)	29,29,29	2.06	9 (31%)

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
5	EDO	A	402	-	-	0/1/1/1	0/0/0/0
5	EDO	A	403	-	-	0/1/1/1	0/0/0/0
5	EDO	A	404	-	-	0/1/1/1	0/0/0/0
5	EDO	A	405	-	-	0/1/1/1	0/0/0/0
5	EDO	A	406	-	-	0/1/1/1	0/0/0/0
5	EDO	B	604	-	-	0/1/1/1	0/0/0/0
5	EDO	B	605	-	-	0/1/1/1	0/0/0/0
5	EDO	B	606	-	-	0/1/1/1	0/0/0/0
5	EDO	B	607	-	-	0/1/1/1	0/0/0/0
7	FII	B	608	-	-	0/24/24/24	0/0/0/0

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	B	608	FII	C22-C15	8.45	1.50	1.32
7	B	608	FII	C34-C27	7.76	1.48	1.32
7	B	608	FII	C43-N42	5.24	1.40	1.33
7	B	608	FII	C10-C2	4.76	1.47	1.32
7	B	608	FII	P46-C45	4.57	1.87	1.79
7	B	608	FII	P46-O50	-2.29	1.50	1.54

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	B	608	FII	C23-C22-C15	-6.19	114.45	127.80
7	B	608	FII	O36-N42-C43	-4.12	112.41	118.31
7	B	608	FII	C24-C27-C34	-3.31	114.72	121.08
7	B	608	FII	C45-C43-N42	-2.95	112.84	115.18
7	B	608	FII	C12-C15-C22	-2.93	115.45	121.08
7	B	608	FII	C18-C15-C12	2.64	119.41	115.39
7	B	608	FII	O51-P46-C45	-2.53	101.66	106.85
7	B	608	FII	O49-P46-C45	-2.15	105.92	110.22
7	B	608	FII	C35-C34-C27	-2.02	122.59	126.19

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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	335/367 (91%)	0.03	25 (7%) 14 15	3, 14, 37, 54	0
2	B	452/519 (87%)	-0.24	18 (3%) 36 40	3, 9, 25, 39	0
3	C	5/5 (100%)	0.81	0 100 100	10, 11, 15, 27	5 (100%)
All	All	792/891 (88%)	-0.12	43 (5%) 25 27	3, 11, 30, 54	5 (0%)

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	25	PHE	17.4
2	B	390	ALA	9.5
2	B	392	PRO	9.3
1	A	24	ASP	9.3
2	B	391	GLY	8.4
2	B	393	LYS	7.6
1	A	307	GLU	7.4
2	B	460	ASP	7.4
2	B	389	LEU	6.8
1	A	305	GLY	6.3
2	B	459	ASP	6.1
2	B	395	SER	5.6
1	A	23	SER	5.2
1	A	26	GLY	4.8
1	A	345	HIS	4.5
1	A	106[A]	LEU	4.3
2	B	68	VAL	4.3
2	B	394	ARG	4.3
1	A	348	GLU	3.8
1	A	303	GLU	3.6
1	A	4	LYS	3.6
1	A	200	ASN	3.5
2	B	388	PRO	3.5

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Mol	Chain	Res	Type	RSRZ
1	A	306	SER	3.4
2	B	457	SER	3.4
1	A	347	THR	3.4
1	A	346	ALA	3.3
1	A	22	GLY	3.0
1	A	286	ALA	3.0
1	A	3	GLY	2.8
2	B	85	LEU	2.7
1	A	349	ILE	2.7
1	A	302	GLU	2.5
1	A	15	LYS	2.5
2	B	458	LYS	2.5
1	A	342	SER	2.5
2	B	478	GLN	2.5
2	B	461	PHE	2.4
2	B	70	PRO	2.4
1	A	28[A]	MET	2.4
1	A	343	ALA	2.3
1	A	199	GLU	2.3
2	B	110	ASN	2.2

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
5	EDO	B	606	4/4	0.17	18.62	23,25,25,26	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
5	EDO	B	607	4/4	0.16	12.23	28,28,28,32	0
5	EDO	A	406	4/4	0.10	6.22	36,36,36,37	0
5	EDO	B	605	4/4	0.09	5.49	18,19,23,23	0
5	EDO	A	402	4/4	0.14	3.26	12,17,20,24	0
5	EDO	A	404	4/4	0.10	1.98	20,22,22,22	0
5	EDO	A	405	4/4	0.09	1.79	22,22,22,23	0
4	CL	B	603	1/1	0.08	1.64	58,58,58,58	0
5	EDO	A	403	4/4	0.08	0.35	14,15,16,17	0
7	FII	B	608	24/24	0.06	-0.14	4,10,14,17	0
5	EDO	B	604	4/4	0.05	-0.15	14,16,17,18	0
4	CL	A	401	1/1	0.12	-0.62	62,62,62,62	0
4	CL	B	602	1/1	0.03	-0.94	19,19,19,19	0
6	ZN	B	601	1/1	0.02	-2.04	5,5,5,5	0

6.5 Other polymers ⓘ

There are no such residues in this entry.