



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:58 AM GMT

PDB ID : 2JI8
Title : X-RAY STRUCTURE OF OXALYL-COA DECARBOXYLASE IN COMPLEX WITH FORMYL-COA
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Deposited on : 2007-02-26
Resolution : 2.15 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

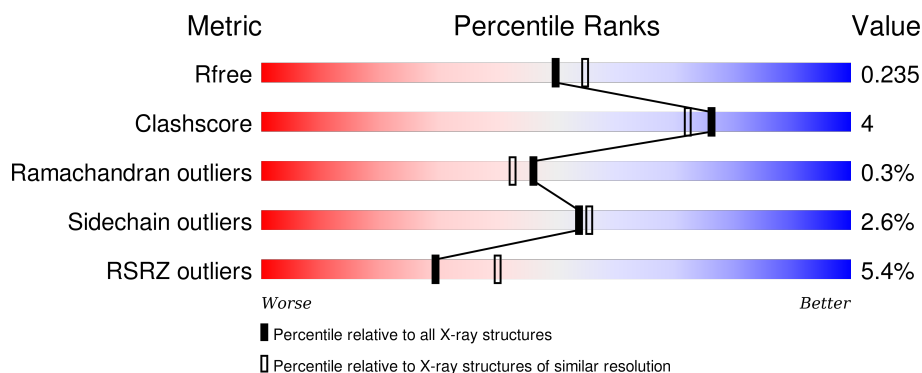
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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}	91344	1045 (2.16-2.16)
Clashscore	102246	1152 (2.16-2.16)
Ramachandran outliers	100387	1131 (2.16-2.16)
Sidechain outliers	100360	1131 (2.16-2.16)
RSRZ outliers	91569	1050 (2.16-2.16)

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. 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	568	<div> <div>5%</div> <div>90%</div> <div>8% ..</div> </div>
1	B	568	<div> <div>6%</div> <div>89%</div> <div>8% ..</div> </div>

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
6	PGE	B	1567	-	-	-	X

2 Entry composition [i](#)

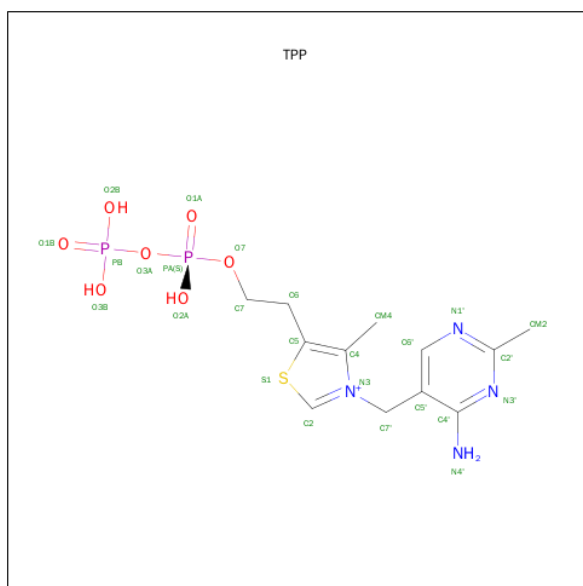
There are 7 unique types of molecules in this entry. The entry contains 9206 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 OXALYL-COA DECARBOXYLASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	559	Total	C	N	O	S	0	4	0
			4194	2641	727	789	37			
1	B	556	Total	C	N	O	S	0	4	0
			4170	2625	723	786	36			

- Molecule 2 is THIAMINE DIPHOSPHATE (three-letter code: TPP) (formula: C₁₂H₁₉N₄O₇P₂S).

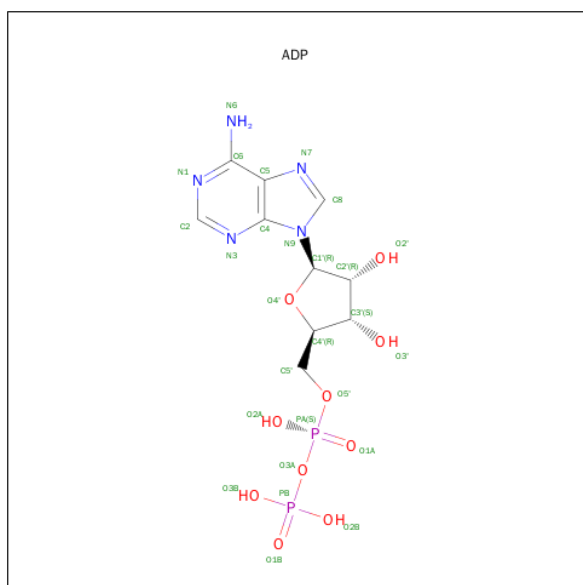


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P S	0	0
			26	12	4	7	2 1		
2	B	1	Total	C	N	O	P S	0	0
			26	12	4	7	2 1		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

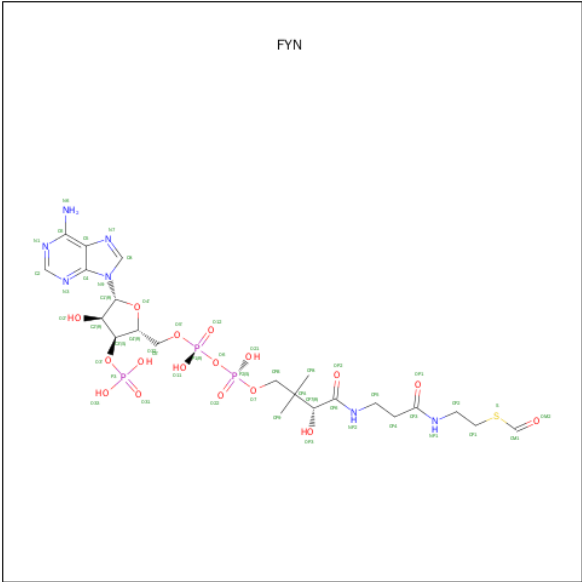
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total Mg 1 1	0	0
3	A	1	Total Mg 1 1	0	0

- Molecule 4 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



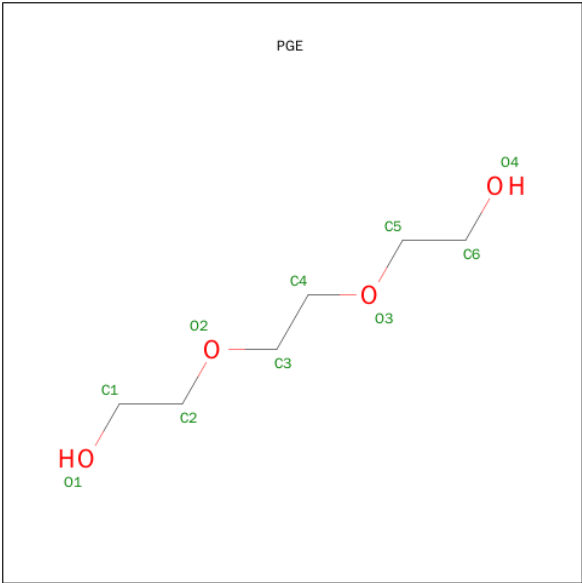
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C N O P 27 10 5 10 2	0	0
4	B	1	Total C N O P 27 10 5 10 2	0	0

- Molecule 5 is S-{(9R,13S,15R)-17-[(2R,3S,4R,5R)-5-(6-AMINO-9H-PURIN-9-YL)-4-HYDROXY-3-(PHOSPHONOXY)TETRAHYDROFURAN-2-YL]-9,13,15-TRIHYDROXY-10,10-DIMETHYL-13,15-DIOXIDO-4,8-DIOXO-12,14,16-TRIOXA-3,7-DIAZA-13,15-DIPHOSPHAHEPTADEC-1-YL} THIOFORMATE (three-letter code: FYN) (formula: $C_{22}H_{36}N_7O_{17}P_3S$).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	S	0	0
			50	22	7	17	3	1		
5	B	1	Total	C	N	O	P	S	0	0
			50	22	7	17	3	1		

- Molecule 6 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: C₆H₁₄O₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			10	6	4		
6	B	1	Total	C	O	0	0
			10	6	4		

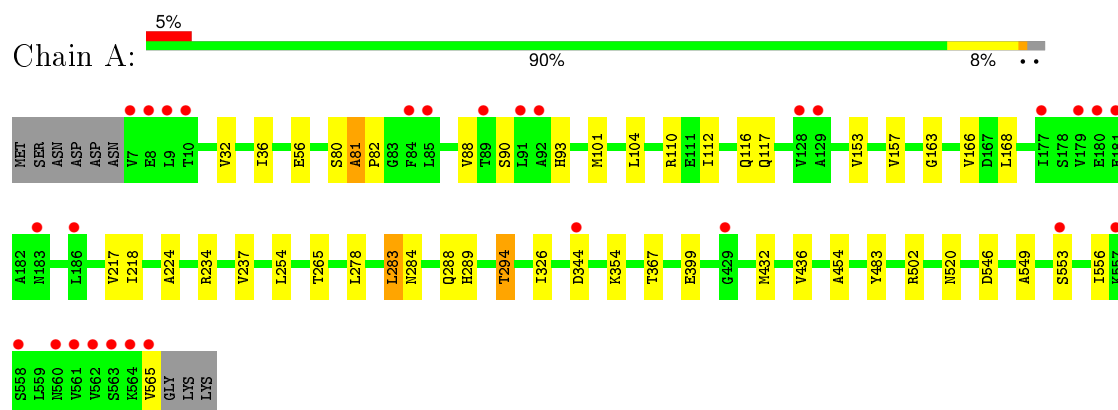
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	316	Total 316	O 316	0	0
7	B	298	Total 298	O 298	0	0

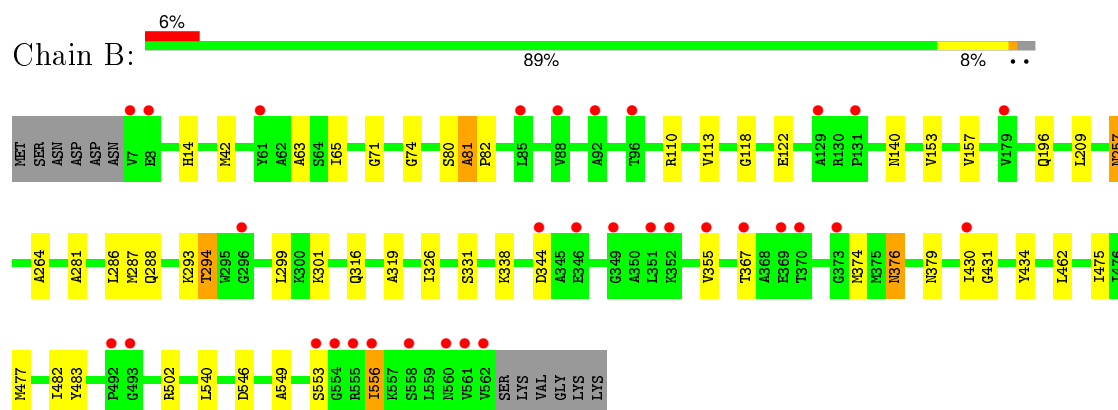
3 Residue-property plots [i](#)

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: OXALYL-COA DECARBOXYLASE



• Molecule 1: OXALYL-COA DECARBOXYLASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	127.57Å 127.57Å 152.03Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	25.00 – 2.15 46.06 – 1.95	Depositor EDS
% Data completeness (in resolution range)	98.8 (25.00-2.15) 98.6 (46.06-1.95)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.15 (at 1.95Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.189 , 0.231 0.193 , 0.235	Depositor DCC
R_{free} test set	3658 reflections (4.98%)	DCC
Wilson B-factor (Å ²)	22.6	Xtriage
Anisotropy	1.120	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 39.3	EDS
Estimated twinning fraction	0.021 for -h,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	5 of 102861 reflections (0.005%)	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9206	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.83% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: MG, PGE, FYN, ADP, TPP

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	0.46	0/4280	0.60	0/5793
1	B	0.45	0/4256	0.61	0/5762
All	All	0.46	0/8536	0.60	0/11555

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	4194	0	4246	30	0
1	B	4170	0	4215	36	0
2	A	26	0	16	0	0
2	B	26	0	16	2	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	27	0	12	0	0
4	B	27	0	12	1	0
5	A	50	0	32	1	0
5	B	50	0	32	2	0
6	A	10	0	14	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	10	0	14	1	0
7	A	316	0	0	2	0
7	B	298	0	0	1	0
All	All	9206	0	8609	65	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 4.

All (65) 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:B:153:VAL:O	1:B:157:VAL:HG22	1.89	0.73
1:B:288:GLN:HG2	1:B:294:THR:HG21	1.72	0.71
1:B:113:VAL:HG22	1:B:122:GLU:HG3	1.79	0.64
1:B:14:HIS:CE1	1:B:42:MET:CE	2.81	0.63
1:A:153:VAL:O	1:A:157:VAL:HG22	1.99	0.63
1:B:376:ASN:HD22	1:B:376:ASN:C	2.03	0.62
1:B:81:ALA:HB3	1:B:82:PRO:CD	2.30	0.61
1:B:299:LEU:HD13	1:B:316:GLN:NE2	2.15	0.61
1:A:278:LEU:CD1	1:A:283:LEU:HD13	2.34	0.57
1:A:556:ILE:HD11	5:A:1569:FYN:S	2.44	0.57
1:A:283:LEU:O	1:A:289:HIS:HA	2.04	0.57
1:B:376:ASN:HD21	1:B:379:ASN:H	1.53	0.57
1:A:81:ALA:HB3	1:A:82:PRO:CD	2.36	0.55
1:B:556:ILE:HD11	5:B:1566:FYN:S	2.48	0.54
1:B:140:ASN:ND2	7:B:2110:HOH:O	2.40	0.54
1:B:257:ASN:C	1:B:257:ASN:HD22	2.12	0.53
1:B:81:ALA:CB	1:B:82:PRO:CD	2.87	0.52
1:A:112:ILE:HG23	1:A:117:GLN:HB2	1.90	0.52
1:B:63:ALA:O	1:B:74:GLY:HA3	2.11	0.50
1:A:56:GLU:OE2	2:B:1563:TPP:N1'	2.45	0.50
1:B:430:ILE:HD12	1:B:431:GLY:N	2.26	0.50
1:A:278:LEU:HD11	1:A:283:LEU:HD13	1.93	0.49
1:A:104:LEU:HD23	1:A:166:VAL:HB	1.93	0.49
1:B:326:ILE:HG13	4:B:1565:ADP:C2	2.46	0.49
1:A:104:LEU:HD22	1:A:168:LEU:HD11	1.95	0.49
1:B:546:ASP:HB3	1:B:549:ALA:HB2	1.92	0.49
1:B:209:LEU:HD21	1:B:301:LYS:HD3	1.95	0.49
1:A:288:GLN:OE1	1:A:294:THR:HG21	2.12	0.49
1:A:81:ALA:HB3	1:A:82:PRO:HD3	1.95	0.48
1:B:81:ALA:HB3	1:B:82:PRO:HD3	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:288:GLN:HG2	1:B:294:THR:CG2	2.41	0.47
1:B:14:HIS:CE1	1:B:42:MET:HE1	2.48	0.47
1:B:376:ASN:ND2	1:B:379:ASN:H	2.12	0.47
1:A:483:TYR:CE1	1:A:553:SER:HA	2.49	0.47
1:A:278:LEU:HD13	1:A:283:LEU:HD13	1.97	0.47
1:B:264:ALA:HB3	1:B:355:VAL:HG13	1.96	0.46
1:A:546:ASP:HB3	1:A:549:ALA:HB2	1.96	0.46
1:B:14:HIS:CE1	1:B:42:MET:HE3	2.50	0.46
1:B:430:ILE:HD12	1:B:430:ILE:C	2.36	0.46
1:A:278:LEU:HD11	1:A:283:LEU:CD1	2.46	0.45
1:A:432:MET:O	1:A:436:VAL:HG23	2.17	0.45
1:A:224:ALA:HA	1:A:326:ILE:HD13	1.98	0.45
1:A:116:GLN:NE2	7:A:2059:HOH:O	2.49	0.45
1:B:475:ILE:HG22	1:B:477:MET:CE	2.46	0.45
1:A:32:VAL:HA	1:A:36:ILE:HD11	1.97	0.45
1:B:483:TYR:CE1	1:B:553:SER:HA	2.52	0.44
1:A:520:ASN:ND2	7:A:2285:HOH:O	2.51	0.44
1:A:218:ILE:HD12	1:A:237:VAL:HG22	1.99	0.44
1:A:565:VAL:O	1:A:565:VAL:HG12	2.18	0.43
1:A:234:ARG:HA	1:A:254:LEU:HD11	2.00	0.43
1:B:71:GLY:HA3	6:B:1567:PGE:H2	2.00	0.43
1:A:81:ALA:CB	1:A:82:PRO:CD	2.96	0.43
1:B:286:LEU:O	5:B:1566:FYN:HP83	2.18	0.43
1:B:209:LEU:HD11	1:B:319:ALA:HB3	1.99	0.43
1:B:540:LEU:C	1:B:540:LEU:HD23	2.39	0.43
1:B:196:GLN:HE21	1:B:196:GLN:HA	1.83	0.43
1:A:399:GLU:HG3	1:A:454:ALA:HB1	2.01	0.42
1:A:101:MET:O	1:A:163:GLY:HA2	2.20	0.42
1:B:281:ALA:HB1	1:B:287:MET:CE	2.50	0.41
1:A:90:SER:O	1:A:93:HIS:HB3	2.20	0.41
1:A:284:ASN:HB2	1:B:118:GLY:O	2.20	0.41
1:B:482:ILE:HG23	2:B:1563:TPP:H62	2.02	0.41
1:B:65:ILE:HD13	1:B:434:TYR:CZ	2.56	0.41
1:A:265:THR:O	1:A:265:THR:HG23	2.21	0.41
1:B:475:ILE:HG22	1:B:477:MET:HE2	2.03	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	561/568 (99%)	540 (96%)	20 (4%)	1 (0%)	52	51
1	B	558/568 (98%)	541 (97%)	15 (3%)	2 (0%)	39	34
All	All	1119/1136 (98%)	1081 (97%)	35 (3%)	3 (0%)	46	42

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	81	ALA
1	B	81	ALA
1	B	556	ILE

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	441/445 (99%)	431 (98%)	10 (2%)	58	62
1	B	438/445 (98%)	424 (97%)	14 (3%)	46	45
All	All	879/890 (99%)	855 (97%)	24 (3%)	54	53

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	80	SER
1	A	88	VAL
1	A	110	ARG

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Mol	Chain	Res	Type
1	A	217	VAL
1	A	283	LEU
1	A	294	THR
1	A	344	ASP
1	A	354	LYS
1	A	367	THR
1	A	502	ARG
1	B	80	SER
1	B	110	ARG
1	B	257	ASN
1	B	293	LYS
1	B	294	THR
1	B	331[A]	SER
1	B	331[B]	SER
1	B	338	LYS
1	B	344	ASP
1	B	367	THR
1	B	374	MET
1	B	376	ASN
1	B	462	LEU
1	B	502	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	116	GLN
1	A	228	GLN
1	A	520	ASN
1	B	140	ASN
1	B	183	ASN
1	B	196	GLN
1	B	228	GLN
1	B	257	ASN
1	B	304	GLN
1	B	316	GLN
1	B	376	ASN
1	B	491	GLN
1	B	520	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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 10 ligands modelled in this entry, 2 are monoatomic - leaving 8 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$
2	TPP	A	1566	3	20,27,27	1.63	5 (25%)	31,40,40	1.55	8 (25%)
4	ADP	A	1568	-	22,29,29	1.01	1 (4%)	27,45,45	1.96	6 (22%)
5	FYN	A	1569	-	41,52,52	1.01	1 (2%)	52,77,77	1.96	8 (15%)
6	PGE	A	1570	-	9,9,9	0.55	0	8,8,8	0.29	0
2	TPP	B	1563	3	20,27,27	1.48	5 (25%)	31,40,40	1.56	7 (22%)
4	ADP	B	1565	-	22,29,29	1.03	1 (4%)	27,45,45	2.02	5 (18%)
5	FYN	B	1566	-	41,52,52	1.02	2 (4%)	52,77,77	1.92	8 (15%)
6	PGE	B	1567	-	9,9,9	0.56	0	8,8,8	0.33	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
2	TPP	A	1566	3	-	0/16/17/17	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	ADP	A	1568	-	-	0/12/32/32	0/3/3/3
5	FYN	A	1569	-	-	0/45/66/66	0/3/3/3
6	PGE	A	1570	-	-	0/7/7/7	0/0/0/0
2	TPP	B	1563	3	-	0/16/17/17	0/2/2/2
4	ADP	B	1565	-	-	0/12/32/32	0/3/3/3
5	FYN	B	1566	-	-	0/45/66/66	0/3/3/3
6	PGE	B	1567	-	-	0/7/7/7	0/0/0/0

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	1569	FYN	CP1-S	-5.25	1.74	1.81
5	B	1566	FYN	CP1-S	-5.10	1.75	1.81
2	A	1566	TPP	C4-N3	-4.58	1.35	1.39
2	B	1563	TPP	C4-N3	-3.81	1.36	1.39
2	A	1566	TPP	C7'-N3	-2.56	1.44	1.48
2	B	1563	TPP	C7'-N3	-2.05	1.45	1.48
2	B	1563	TPP	C6'-N1'	2.06	1.38	1.34
5	B	1566	FYN	O4'-C1'	2.08	1.43	1.41
2	B	1563	TPP	C2'-N1'	2.14	1.38	1.34
2	B	1563	TPP	C4'-N3'	2.16	1.38	1.35
2	A	1566	TPP	C2'-N3'	2.34	1.38	1.34
2	A	1566	TPP	C2'-N1'	2.34	1.38	1.34
2	A	1566	TPP	C4'-N3'	2.36	1.38	1.35
4	B	1565	ADP	C5-C4	3.02	1.47	1.40
4	A	1568	ADP	C5-C4	3.22	1.47	1.40

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1569	FYN	N3-C2-N1	-10.46	120.89	128.89
5	B	1566	FYN	N3-C2-N1	-10.06	121.19	128.89
4	B	1565	ADP	N3-C2-N1	-7.98	122.78	128.89
4	A	1568	ADP	N3-C2-N1	-7.32	123.29	128.89
5	A	1569	FYN	CP1-CP2-NP1	-4.71	102.94	112.36
5	B	1566	FYN	CP1-CP2-NP1	-4.33	103.69	112.36
5	B	1566	FYN	CP4-CP5-NP2	-3.61	103.97	111.88
4	A	1568	ADP	C4-C5-N7	-3.27	106.47	109.48
4	A	1568	ADP	C1'-N9-C4	-3.01	122.40	126.94
5	A	1569	FYN	CP4-CP5-NP2	-2.93	105.46	111.88
4	B	1565	ADP	C4-C5-N7	-2.89	106.82	109.48
5	B	1566	FYN	C1'-N9-C4	-2.83	122.67	126.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1563	TPP	N1'-C2'-N3'	-2.79	120.43	125.60
2	B	1563	TPP	CM4-C4-C5	-2.77	122.67	128.90
2	A	1566	TPP	N1'-C2'-N3'	-2.65	120.69	125.60
4	B	1565	ADP	C1'-N9-C4	-2.64	122.95	126.94
5	B	1566	FYN	C4-C5-N7	-2.58	107.10	109.48
2	B	1563	TPP	PA-O3A-PB	-2.56	124.07	132.67
5	A	1569	FYN	C1'-N9-C4	-2.56	123.08	126.94
2	A	1566	TPP	C6-C5-S1	-2.52	116.70	120.24
5	A	1569	FYN	C4-C5-N7	-2.24	107.41	109.48
2	A	1566	TPP	C5'-C6'-N1'	-2.20	120.03	123.86
5	B	1566	FYN	P2-O6-P1	-2.14	126.73	132.73
4	B	1565	ADP	PA-O3A-PB	-2.12	125.56	132.67
2	A	1566	TPP	PA-O3A-PB	-2.07	125.73	132.67
4	A	1568	ADP	PA-O3A-PB	-2.06	125.76	132.67
5	A	1569	FYN	O33-P3-O32	2.00	115.00	107.38
2	B	1563	TPP	C6'-N1'-C2'	2.09	119.42	115.77
4	A	1568	ADP	C2-N1-C6	2.13	122.58	118.77
5	A	1569	FYN	CP5-NP2-CP6	2.15	126.78	122.53
2	A	1566	TPP	N4'-C4'-N3'	2.20	120.14	116.95
4	B	1565	ADP	C2'-C1'-N9	2.20	117.66	114.29
5	B	1566	FYN	CP5-NP2-CP6	2.21	126.91	122.53
2	A	1566	TPP	C6'-N1'-C2'	2.29	119.78	115.77
4	A	1568	ADP	O3B-PB-O2B	2.35	116.33	107.38
2	B	1563	TPP	C6'-C5'-C4'	2.53	119.35	115.72
2	B	1563	TPP	CM2-C2'-N1'	2.60	120.15	117.03
2	A	1566	TPP	CM4-C4-N3	2.61	126.07	122.59
2	A	1566	TPP	C6'-C5'-C4'	3.02	120.06	115.72
5	B	1566	FYN	CP2-NP1-CP3	3.12	128.92	122.79
5	A	1569	FYN	CP2-NP1-CP3	3.31	129.31	122.79
2	B	1563	TPP	CM4-C4-N3	3.59	127.38	122.59

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	1569	FYN	1	0
2	B	1563	TPP	2	0
4	B	1565	ADP	1	0
5	B	1566	FYN	2	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	B	1567	PGE	1	0

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	559/568 (98%)	0.29	28 (5%)	32 43	21, 29, 37, 63	0
1	B	556/568 (97%)	0.32	32 (5%)	26 36	22, 29, 38, 64	0
All	All	1115/1136 (98%)	0.31	60 (5%)	29 40	21, 29, 38, 64	0

All (60) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	565	VAL	8.6
1	A	7	VAL	7.2
1	B	7	VAL	5.9
1	B	562	VAL	5.3
1	A	563	SER	5.2
1	A	561	VAL	4.8
1	A	564	LYS	4.5
1	B	492	PRO	4.4
1	A	562	VAL	4.3
1	B	367	THR	4.1
1	B	355	VAL	4.0
1	A	85	LEU	3.9
1	A	9	LEU	3.9
1	B	561	VAL	3.8
1	B	560	ASN	3.6
1	B	296	GLY	3.6
1	A	558	SER	3.5
1	B	558	SER	3.4
1	B	344	ASP	3.4
1	B	493	GLY	3.4
1	A	8	GLU	3.4
1	B	8	GLU	3.2
1	A	128	VAL	3.2
1	B	349	GLY	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	557	LYS	3.2
1	B	430	ILE	3.1
1	B	369	GLU	3.1
1	B	555	ARG	3.0
1	A	553	SER	3.0
1	A	177	ILE	2.9
1	A	560	ASN	2.8
1	B	88	VAL	2.7
1	B	554	GLY	2.7
1	B	179	VAL	2.7
1	B	351	LEU	2.6
1	B	556	ILE	2.6
1	A	181	GLU	2.5
1	B	553	SER	2.5
1	B	373	GLY	2.5
1	A	129	ALA	2.5
1	B	370	THR	2.5
1	A	179	VAL	2.5
1	B	92	ALA	2.4
1	B	346	GLU	2.3
1	A	183	ASN	2.3
1	B	61	TYR	2.3
1	A	91	LEU	2.3
1	A	84	PHE	2.3
1	A	89	THR	2.3
1	A	10	THR	2.2
1	B	85	LEU	2.2
1	A	92	ALA	2.2
1	B	131	PRO	2.2
1	B	129	ALA	2.2
1	A	186	LEU	2.1
1	B	96	THR	2.1
1	A	344	ASP	2.1
1	A	180	GLU	2.0
1	B	352	LYS	2.0
1	A	429	GLY	2.0

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 [i](#)

There are no carbohydrates 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. 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	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
6	PGE	B	1567	10/10	0.73	0.23	5.92	49,50,50,50	0
6	PGE	A	1570	10/10	0.81	0.14	1.82	54,55,55,56	0
5	FYN	B	1566	50/50	0.79	0.26	1.74	32,40,45,46	50
5	FYN	A	1569	50/50	0.88	0.22	1.61	32,40,42,46	50
2	TPP	A	1566	26/26	0.95	0.11	-0.63	19,27,28,29	0
2	TPP	B	1563	26/26	0.95	0.11	-1.00	21,28,30,32	0
4	ADP	B	1565	27/27	0.95	0.10	-1.08	20,22,26,28	0
4	ADP	A	1568	27/27	0.96	0.09	-1.33	19,22,24,24	0
3	MG	A	1567	1/1	0.89	0.07	-1.52	26,26,26,26	0
3	MG	B	1564	1/1	0.91	0.08	-1.90	28,28,28,28	0

6.5 Other polymers [i](#)

There are no such residues in this entry.