



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 04:54 AM GMT

PDB ID : 2ONH
Title : Crystal Structure of of limonene synthase with 2-fluorolinalyl diphosphate(FLPP)
Authors : Hyatt, D.C.; Youn, B.; Croteau, R.; Kang, C.
Deposited on : 2007-01-24
Resolution : 2.70 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

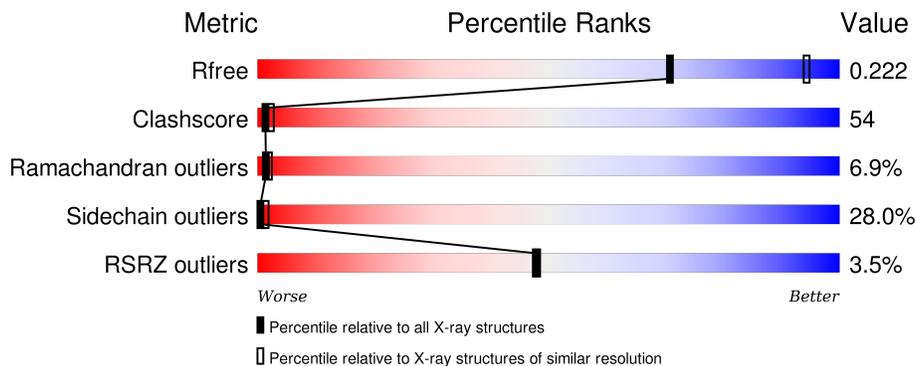
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	2103 (2.70-2.70)
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)
RSRZ outliers	91569	2107 (2.70-2.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	543	 3% 23% 46% 27% .
1	B	543	 3% 28% 42% 24% 6%

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
2	MN	A	602	-	-	-	X
3	F3P	B	1600	-	-	X	-
4	BTB	A	604	-	-	X	X

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 9165 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 4S-limonene synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	543	4497	2873	761	843	20	0	0	0
1	B	543	4497	2873	761	843	20	0	0	0

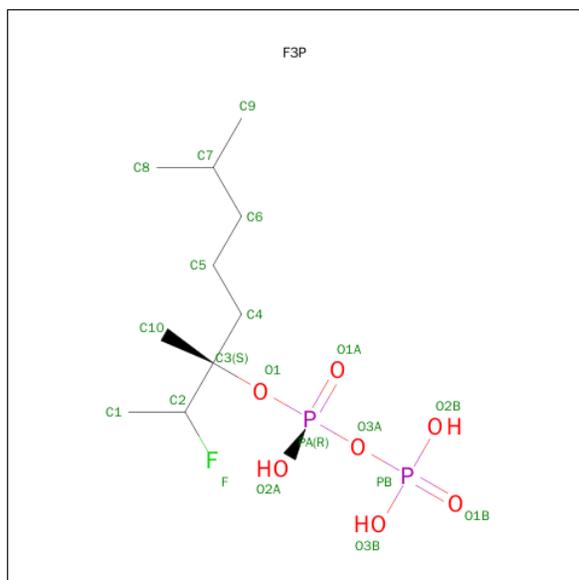
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	57	MET	GLU	ENGINEERED	UNP Q40322
B	57	MET	GLU	ENGINEERED	UNP Q40322

- Molecule 2 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

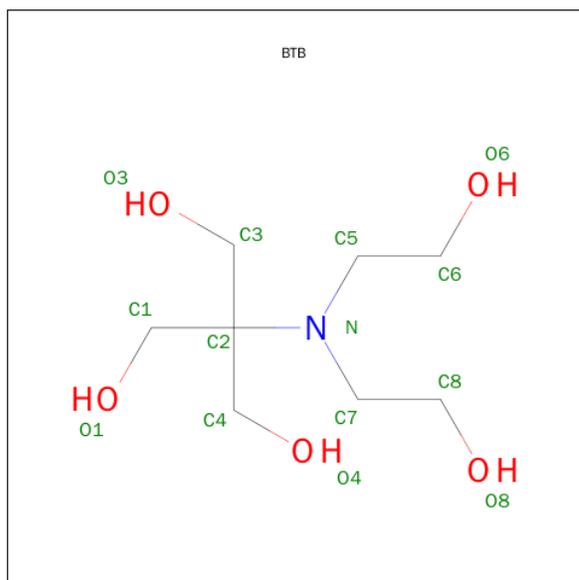
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	3	Total	Mn	0	0
			3	3		
2	A	3	Total	Mn	0	0
			3	3		

- Molecule 3 is (1S)-1-[(1S)-1-FLUOROETHYL]-1,5-DIMETHYLHEXYL TRIHYDROGEN DIPHOSPHATE (three-letter code: F3P) (formula: C₁₀H₂₃FO₇P₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
3	A	1	Total	C	F	O	P	0	0
			20	10	1	7	2		
3	B	1	Total	C	F	O	P	0	0
			20	10	1	7	2		

- Molecule 4 is 2-[BIS-(2-HYDROXY-ETHYL)-AMINO]-2-HYDROXYMETHYL-PROPAN E-1,3-DIOL (three-letter code: BTB) (formula: C₈H₁₉NO₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	
4	A	1	Total	C	N	O	0	0
			14	8	1	5		

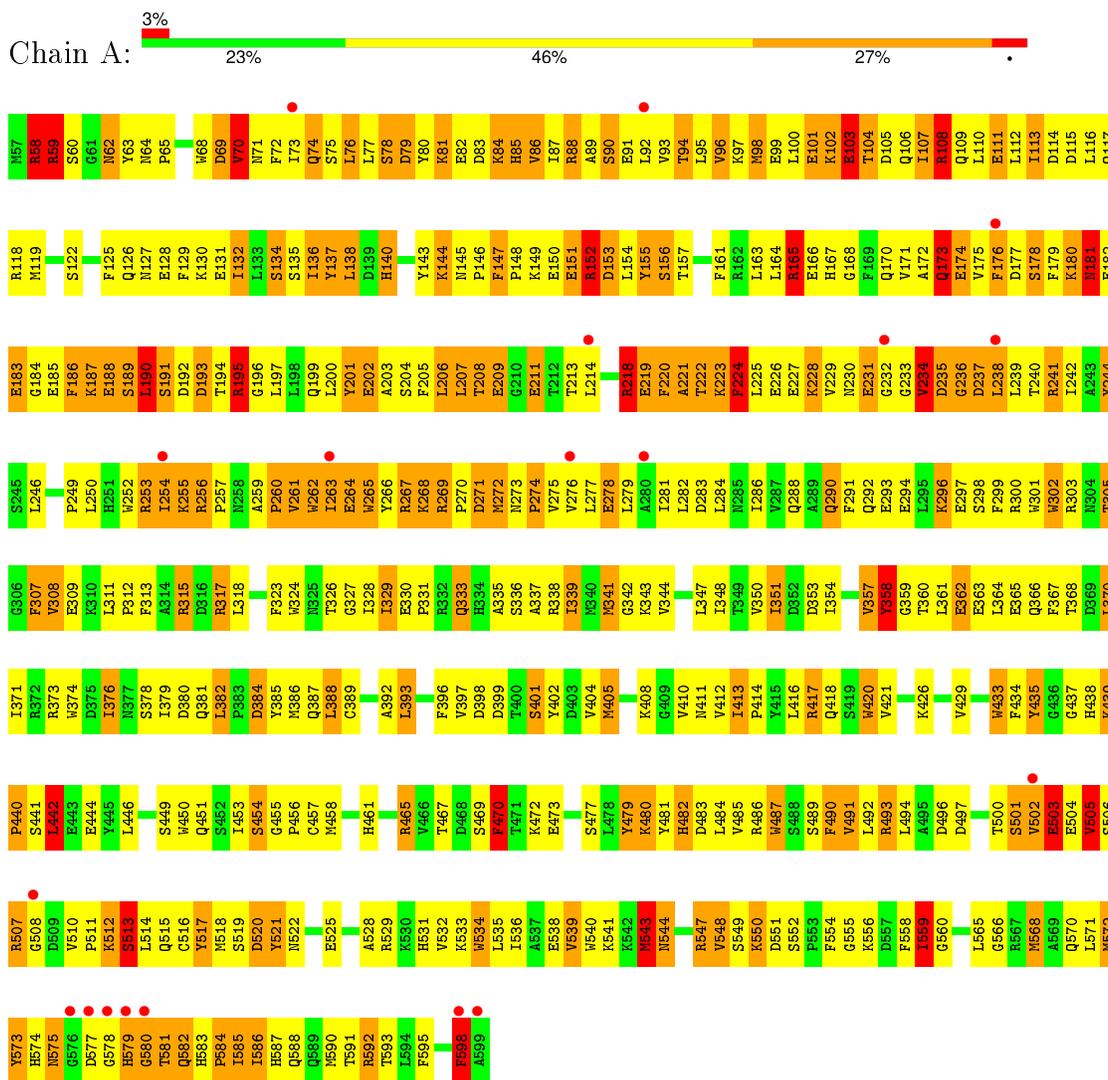
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	58	Total 58	O 58	0	0
5	B	53	Total 53	O 53	0	0

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: 4S-limonene synthase



- Molecule 1: 4S-limonene synthase



4 Data and refinement statistics

Property	Value	Source
Space group	I 4	Depositor
Cell constants a, b, c, α , β , γ	198.66Å 198.66Å 122.67Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.70 46.20 – 2.60	Depositor EDS
% Data completeness (in resolution range)	(Not available) (10.00-2.70) 89.7 (46.20-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.54 (at 2.61Å)	Xtrriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.206 , 0.234 0.213 , 0.222	Depositor DCC
R_{free} test set	2988 reflections (5.39%)	DCC
Wilson B-factor (Å ²)	63.2	Xtrriage
Anisotropy	0.185	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	(Not available) , (Not available)	EDS
Estimated twinning fraction	0.487 for -k,-h,-l	Xtrriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtrriage
Outliers	0 of 65671 reflections	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9165	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.79% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: F3P, MN, BTB

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	1.08	15/4609 (0.3%)	1.48	115/6237 (1.8%)
1	B	1.07	11/4609 (0.2%)	1.49	111/6237 (1.8%)
All	All	1.07	26/9218 (0.3%)	1.48	226/12474 (1.8%)

The worst 5 of 26 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	487	TRP	CG-CD2	-8.29	1.29	1.43
1	B	209	GLU	CG-CD	8.26	1.64	1.51
1	B	219	GLU	CG-CD	7.69	1.63	1.51
1	A	209	GLU	CD-OE2	-7.53	1.17	1.25
1	B	209	GLU	CD-OE1	-7.37	1.17	1.25

The worst 5 of 226 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	266	TYR	CB-CG-CD1	-14.20	112.48	121.00
1	B	358	TYR	CB-CG-CD2	-13.56	112.86	121.00
1	A	266	TYR	CB-CG-CD1	-12.45	113.53	121.00
1	B	420	TRP	CD1-CG-CD2	12.26	116.11	106.30
1	A	374	TRP	CD1-CG-CD2	11.98	115.89	106.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	4497	0	4353	466	0
1	B	4497	0	4353	459	0
2	A	3	0	0	0	0
2	B	3	0	0	0	0
3	A	20	0	18	7	0
3	B	20	0	18	12	0
4	A	14	0	19	29	0
5	A	58	0	0	4	1
5	B	53	0	0	3	1
All	All	9165	0	8761	961	1

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 54.

The worst 5 of 961 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)
4:A:604:BTB:C2	4:A:604:BTB:N	1.68	1.55
3:B:1600:F3P:C6	3:B:1600:F3P:H13	1.62	1.29
3:B:1600:F3P:C1	3:B:1600:F3P:H61	1.62	1.29
4:A:604:BTB:C7	4:A:604:BTB:H32	1.74	1.16
4:A:604:BTB:N	4:A:604:BTB:C3	2.12	1.12

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:737:HOH:O	5:B:707:HOH:O[7_554]	2.12	0.08

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	541/543 (100%)	428 (79%)	78 (14%)	35 (6%)	1	2
1	B	541/543 (100%)	407 (75%)	94 (17%)	40 (7%)	1	1
All	All	1082/1086 (100%)	835 (77%)	172 (16%)	75 (7%)	1	2

5 of 75 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	62	ASN
1	A	76	LEU
1	A	181	ASN
1	A	193	ASP
1	A	235	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	491/492 (100%)	349 (71%)	142 (29%)	0	1
1	B	491/492 (100%)	358 (73%)	133 (27%)	0	1
All	All	982/984 (100%)	707 (72%)	275 (28%)	0	1

5 of 275 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	520	ASP
1	B	86	VAL
1	B	518	MET
1	A	548	VAL
1	A	591	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 32 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	85	HIS
1	B	167	HIS
1	B	575	ASN
1	B	145	ASN
1	B	170	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 9 ligands modelled in this entry, 6 are monoatomic - leaving 3 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	$\# Z > 2$
3	F3P	A	600	2	15,19,19	1.88	4 (26%)	22,29,29	1.26	3 (13%)
4	BTB	A	604	-	12,13,13	2.34	4 (33%)	8,16,16	1.66	2 (25%)
3	F3P	B	1600	2	15,19,19	1.77	3 (20%)	22,29,29	1.24	3 (13%)

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	F3P	A	600	2	-	0/19/25/25	0/0/0/0
4	BTB	A	604	-	-	0/21/21/21	0/0/0/0
3	F3P	B	1600	2	-	0/19/25/25	0/0/0/0

The worst 5 of 11 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	600	F3P	C10-C3	-4.04	1.42	1.52
3	B	1600	F3P	C10-C3	-4.00	1.42	1.52
3	A	600	F3P	PB-O2B	2.04	1.62	1.54
3	B	1600	F3P	PB-O3B	2.21	1.62	1.54
3	A	600	F3P	PB-O3B	2.46	1.63	1.54

The worst 5 of 8 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	604	BTB	C7-N-C5	-3.65	100.89	112.52
3	A	600	F3P	PA-O3A-PB	-3.14	122.12	132.67
3	B	1600	F3P	PA-O3A-PB	-2.96	122.74	132.67
3	B	1600	F3P	O1-C3-C10	2.10	114.80	107.51
3	A	600	F3P	O1-C3-C10	2.13	114.91	107.51

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 48 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	600	F3P	7	0
4	A	604	BTB	29	0
3	B	1600	F3P	12	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 [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	543/543 (100%)	0.24	19 (3%) 48 48	24, 52, 97, 100	0
1	B	543/543 (100%)	0.25	19 (3%) 48 48	23, 52, 98, 100	0
All	All	1086/1086 (100%)	0.25	38 (3%) 48 48	23, 52, 98, 100	0

The worst 5 of 38 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	579	HIS	8.2
1	A	580	GLY	6.6
1	A	577	ASP	6.0
1	A	578	GLY	5.3
1	B	581	THR	4.9

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 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(\AA^2)	Q<0.9
4	BTB	A	604	14/14	0.83	0.42	7.06	84,91,100,100	0
2	MN	A	602	1/1	0.97	0.28	2.33	52,52,52,52	0
3	F3P	B	1600	20/20	0.92	0.26	0.47	45,53,63,66	0
3	F3P	A	600	20/20	0.94	0.24	0.07	45,55,64,67	0
2	MN	B	1603	1/1	0.98	0.26	-	55,55,55,55	0
2	MN	B	1602	1/1	0.94	0.30	-	70,70,70,70	0
2	MN	B	1601	1/1	0.98	0.30	-	49,49,49,49	0
2	MN	A	603	1/1	0.98	0.23	-	54,54,54,54	0
2	MN	A	601	1/1	0.98	0.32	-	61,61,61,61	0

6.5 Other polymers [i](#)

There are no such residues in this entry.