W O R L D W I D E

wwPDB

P R O T E I N   D A T A   B A N K

wwPDB X-ray Structure Validation Summary Report ⓘ

May 13, 2020 – 04:41 am BST

PDB ID	:	3FIW
Title	:	Structure of SCO0253, a Tetr-family transcriptional regulator from Streptomyces coelicolor
Authors	:	Singer, A.U.; Xu, X.; Chang, C.; Gu, J.; Edwards, A.M.; Joachimiak, A.; Savchenko, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on	:	2008-12-12
Resolution	:	2.20 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
https://www.wwpdb.org/validation/2017/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	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

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 2.20 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.
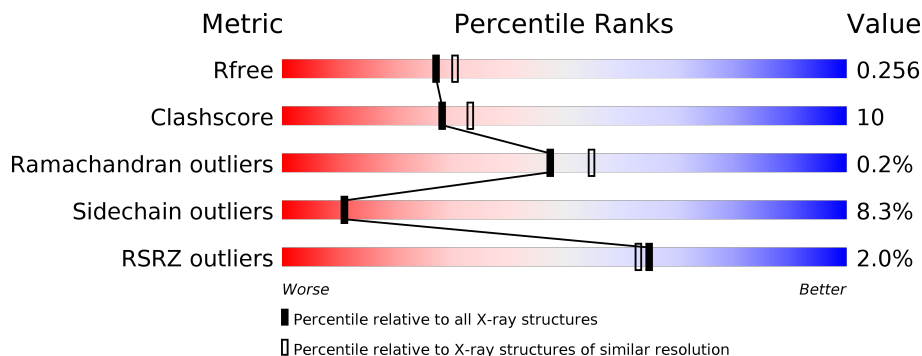
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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 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 <=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	211	% 67% 12% • 18%
1	B	211	% 61% 15% • 21%
1	C	211	64% 16% • 18%
1	D	211	4% 64% 14% • 20%

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	PEG	A	191	-	-	X	-

2   Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 5833 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 Putative tetR-family transcriptional regulator.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	173	Total	C	N	O	S	Se	0	0	0
			1363	847	261	248	1	6			
1	B	167	Total	C	N	O	S	Se	0	1	0
			1323	825	253	238	1	6			
1	C	173	Total	C	N	O	S	Se	0	1	0
			1374	853	265	249	1	6			
1	D	169	Total	C	N	O	S	Se	0	1	0
			1340	833	257	244	1	5			

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-20	MSE	-	expression tag	UNP Q9S1N8
A	-19	GLY	-	expression tag	UNP Q9S1N8
A	-18	SER	-	expression tag	UNP Q9S1N8
A	-17	SER	-	expression tag	UNP Q9S1N8
A	-16	HIS	-	expression tag	UNP Q9S1N8
A	-15	HIS	-	expression tag	UNP Q9S1N8
A	-14	HIS	-	expression tag	UNP Q9S1N8
A	-13	HIS	-	expression tag	UNP Q9S1N8
A	-12	HIS	-	expression tag	UNP Q9S1N8
A	-11	HIS	-	expression tag	UNP Q9S1N8
A	-10	SER	-	expression tag	UNP Q9S1N8
A	-9	SER	-	expression tag	UNP Q9S1N8
A	-8	GLY	-	expression tag	UNP Q9S1N8
A	-7	ARG	-	expression tag	UNP Q9S1N8
A	-6	GLU	-	expression tag	UNP Q9S1N8
A	-5	ASN	-	expression tag	UNP Q9S1N8
A	-4	LEU	-	expression tag	UNP Q9S1N8
A	-3	TYR	-	expression tag	UNP Q9S1N8
A	-2	PHE	-	expression tag	UNP Q9S1N8
A	-1	GLN	-	expression tag	UNP Q9S1N8
A	0	GLY	-	expression tag	UNP Q9S1N8

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-20	MSE	-	expression tag	UNP Q9S1N8
B	-19	GLY	-	expression tag	UNP Q9S1N8
B	-18	SER	-	expression tag	UNP Q9S1N8
B	-17	SER	-	expression tag	UNP Q9S1N8
B	-16	HIS	-	expression tag	UNP Q9S1N8
B	-15	HIS	-	expression tag	UNP Q9S1N8
B	-14	HIS	-	expression tag	UNP Q9S1N8
B	-13	HIS	-	expression tag	UNP Q9S1N8
B	-12	HIS	-	expression tag	UNP Q9S1N8
B	-11	HIS	-	expression tag	UNP Q9S1N8
B	-10	SER	-	expression tag	UNP Q9S1N8
B	-9	SER	-	expression tag	UNP Q9S1N8
B	-8	GLY	-	expression tag	UNP Q9S1N8
B	-7	ARG	-	expression tag	UNP Q9S1N8
B	-6	GLU	-	expression tag	UNP Q9S1N8
B	-5	ASN	-	expression tag	UNP Q9S1N8
B	-4	LEU	-	expression tag	UNP Q9S1N8
B	-3	TYR	-	expression tag	UNP Q9S1N8
B	-2	PHE	-	expression tag	UNP Q9S1N8
B	-1	GLN	-	expression tag	UNP Q9S1N8
B	0	GLY	-	expression tag	UNP Q9S1N8
C	-20	MSE	-	expression tag	UNP Q9S1N8
C	-19	GLY	-	expression tag	UNP Q9S1N8
C	-18	SER	-	expression tag	UNP Q9S1N8
C	-17	SER	-	expression tag	UNP Q9S1N8
C	-16	HIS	-	expression tag	UNP Q9S1N8
C	-15	HIS	-	expression tag	UNP Q9S1N8
C	-14	HIS	-	expression tag	UNP Q9S1N8
C	-13	HIS	-	expression tag	UNP Q9S1N8
C	-12	HIS	-	expression tag	UNP Q9S1N8
C	-11	HIS	-	expression tag	UNP Q9S1N8
C	-10	SER	-	expression tag	UNP Q9S1N8
C	-9	SER	-	expression tag	UNP Q9S1N8
C	-8	GLY	-	expression tag	UNP Q9S1N8
C	-7	ARG	-	expression tag	UNP Q9S1N8
C	-6	GLU	-	expression tag	UNP Q9S1N8
C	-5	ASN	-	expression tag	UNP Q9S1N8
C	-4	LEU	-	expression tag	UNP Q9S1N8
C	-3	TYR	-	expression tag	UNP Q9S1N8
C	-2	PHE	-	expression tag	UNP Q9S1N8
C	-1	GLN	-	expression tag	UNP Q9S1N8
C	0	GLY	-	expression tag	UNP Q9S1N8

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-20	MSE	-	expression tag	UNP Q9S1N8
D	-19	GLY	-	expression tag	UNP Q9S1N8
D	-18	SER	-	expression tag	UNP Q9S1N8
D	-17	SER	-	expression tag	UNP Q9S1N8
D	-16	HIS	-	expression tag	UNP Q9S1N8
D	-15	HIS	-	expression tag	UNP Q9S1N8
D	-14	HIS	-	expression tag	UNP Q9S1N8
D	-13	HIS	-	expression tag	UNP Q9S1N8
D	-12	HIS	-	expression tag	UNP Q9S1N8
D	-11	HIS	-	expression tag	UNP Q9S1N8
D	-10	SER	-	expression tag	UNP Q9S1N8
D	-9	SER	-	expression tag	UNP Q9S1N8
D	-8	GLY	-	expression tag	UNP Q9S1N8
D	-7	ARG	-	expression tag	UNP Q9S1N8
D	-6	GLU	-	expression tag	UNP Q9S1N8
D	-5	ASN	-	expression tag	UNP Q9S1N8
D	-4	LEU	-	expression tag	UNP Q9S1N8
D	-3	TYR	-	expression tag	UNP Q9S1N8
D	-2	PHE	-	expression tag	UNP Q9S1N8
D	-1	GLN	-	expression tag	UNP Q9S1N8
D	0	GLY	-	expression tag	UNP Q9S1N8

- Molecule 2 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$).
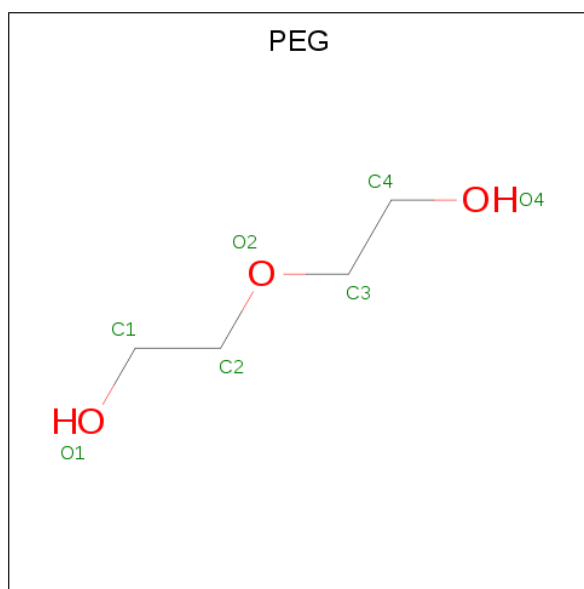
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total 7	C 4	O 3	0	0

• Molecule 3 is water.

| Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	98	Total 98	O 98	0	0
3	B	97	Total 97	O 97	0	0
3	C	108	Total 108	O 108	0	0
3	D	123	Total 123	O 123	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 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.
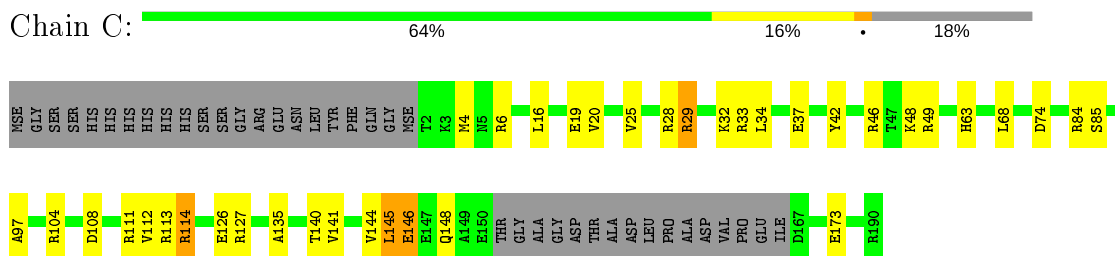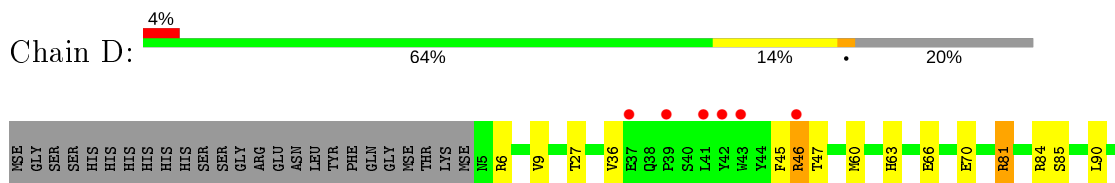
- Molecule 1: Putative tetR-family transcriptional regulator
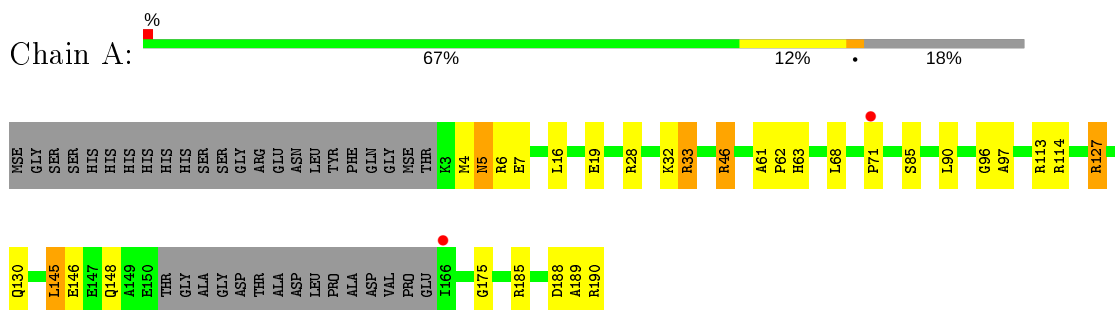
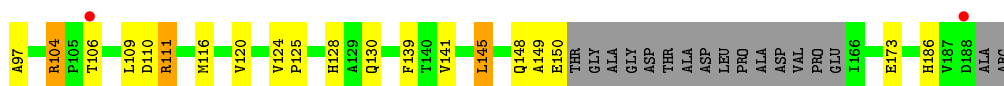- Molecule 1: Putative tetR-family transcriptional regulator
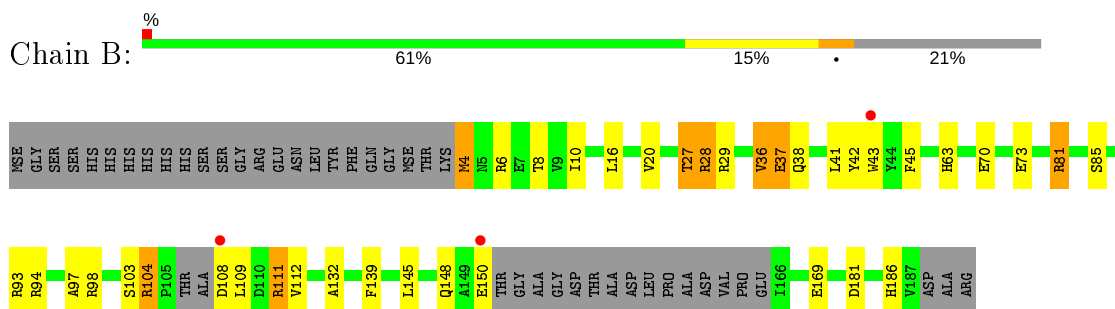
- Molecule 1: Putative tetR-family transcriptional regulator

- Molecule 1: Putative tetR-family transcriptional regulator

4   Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	64.49Å 114.35Å 67.45Å 90.00° 90.14° 90.00°	Depositor
Resolution (Å)	42.80 – 2.20 42.78 – 2.20	Depositor EDS
% Data completeness (in resolution range)	97.5 (42.80-2.20) 97.2 (42.78-2.20)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$<I/\sigma(I)>$ [1]	4.58 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.189 , 0.256 0.189 , 0.256	Depositor DCC
R_{free} test set	2462 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å^2)	32.4	Xtriage
Anisotropy	0.026	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 35.6	EDS
L-test for twinning[2]	$<|L|> = 0.49$, $<L^2> = 0.32$	Xtriage
Estimated twinning fraction	0.000 for l,k,-h 0.010 for h,-k,-l 0.000 for l,-k,h	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	5833	wwPDB-VP
Average B, all atoms (Å^2)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 32.24 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 9.7003e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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[1] Intensities estimated from amplitudes.

[2] Theoretical values of $<|L|>$, $<L^2>$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 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: PEG

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	0.82	0/1382	0.85	1/1854 (0.1%)
1	B	1.02	1/1342 (0.1%)	0.99	7/1802 (0.4%)
1	C	0.80	1/1393 (0.1%)	0.87	0/1868
1	D	0.90	0/1360	0.93	2/1829 (0.1%)
All	All	0.89	2/5477 (0.0%)	0.91	10/7353 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	146	GLU	CG-CD	5.65	1.60	1.51
1	B	132	ALA	CA-CB	-5.04	1.41	1.52

The worst 5 of 10 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	81	ARG	NE-CZ-NH1	8.36	124.48	120.30
1	D	111	ARG	NE-CZ-NH2	-8.11	116.25	120.30
1	D	111	ARG	NE-CZ-NH1	7.40	124.00	120.30
1	B	81	ARG	NE-CZ-NH2	-6.66	116.97	120.30
1	A	145	LEU	CA-CB-CG	6.63	130.55	115.30

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	1363	0	1350	33	0
1	B	1323	0	1307	26	1
1	C	1374	0	1362	29	0
1	D	1340	0	1322	25	0
2	A	7	0	10	11	0
3	A	98	0	0	4	0
3	B	97	0	0	4	0
3	C	108	0	0	5	0
3	D	123	0	0	4	1
All	All	5833	0	5351	105	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 10.

The worst 5 of 105 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:108:ASP:HA	1:B:111:ARG:HB3	1.33	1.06
1:A:146:GLU:HG2	1:B:104:ARG:NH2	1.74	1.00
1:A:5:ASN:HD21	1:A:7:GLU:HG2	1.39	0.88
1:D:60:MSE:HE1	1:D:90:LEU:HD23	1.53	0.88
1:C:25:VAL:O	1:C:48:LYS:HE3	1.77	0.84

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)
1:B:111:ARG:NH2	3:D:310:HOH:O[1_654]	2.19	0.01

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	169/211 (80%)	162 (96%)	6 (4%)	1 (1%)	25	26
1	B	162/211 (77%)	158 (98%)	4 (2%)	0	100	100
1	C	170/211 (81%)	166 (98%)	4 (2%)	0	100	100
1	D	166/211 (79%)	163 (98%)	3 (2%)	0	100	100
All	All	667/844 (79%)	649 (97%)	17 (2%)	1 (0%)	47	60

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	189	ALA

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	137/160 (86%)	126 (92%)	11 (8%)	12	12
1	B	133/160 (83%)	119 (90%)	14 (10%)	7	6
1	C	138/160 (86%)	128 (93%)	10 (7%)	14	15
1	D	135/160 (84%)	124 (92%)	11 (8%)	11	12
All	All	543/640 (85%)	497 (92%)	46 (8%)	11	10

5 of 46 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	103	SER
1	C	29	ARG
1	D	104	ARG
1	B	104	ARG
1	B	150	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 18 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	186	HIS
1	C	63	HIS
1	D	63	HIS
1	B	100	HIS
1	B	148	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)

1 ligand is 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	$\#|Z| > 2$
2	PEG	A	191	-	6,6,6	0.83	0	5,5,5	0.71	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	PEG	A	191	-	-	3/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	191	PEG	C4-C3-O2-C2
2	A	191	PEG	O1-C1-C2-O2
2	A	191	PEG	O2-C3-C4-O4

There are no ring outliers.

1 monomer is involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	191	PEG	11	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	167/211 (79%)	-0.19	2 (1%)	79	77	21, 33, 50, 58	0
1	B	161/211 (76%)	-0.07	3 (1%)	66	65	19, 29, 49, 71	0
1	C	167/211 (79%)	-0.18	0	100	100	27, 39, 52, 59	0
1	D	164/211 (77%)	-0.41	8 (4%)	29	28	19, 30, 70, 77	0
All	All	659/844 (78%)	-0.21	13 (1%)	65	63	19, 34, 56, 77	0

The worst 5 of 13 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	39	PRO	4.9
1	B	43	TRP	4.1
1	D	188	ASP	4.1
1	D	43	TRP	3.7
1	D	46	ARG	3.3

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. 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(Å^2)	Q<0.9
2	PEG	A	191	7/7	0.60	0.36	60,62,65,65	0

6.5 Other polymers (i)

There are no such residues in this entry.