



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 14, 2018 – 05:55 PM EDT

PDB ID : 6BYF
Title : Crystal structure of the core catalytic domain of PP-IP phosphatase SIW14 from *S. cerevisiae* in complex with citrate
Authors : Wang, H.; Shears, S.B.
Deposited on : 2017-12-20
Resolution : 2.35 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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20030736
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
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) : rb-20030736

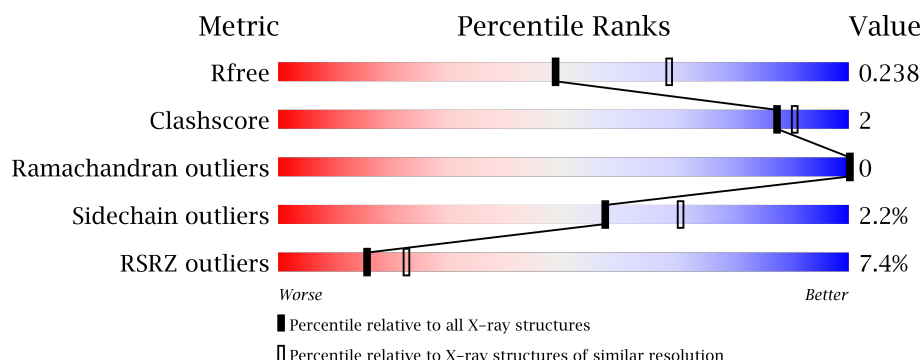
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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}	100719	1522 (2.38-2.34)
Clashscore	112137	1626 (2.38-2.34)
Ramachandran outliers	110173	1605 (2.38-2.34)
Sidechain outliers	110143	1606 (2.38-2.34)
RSRZ outliers	101464	1528 (2.38-2.34)

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	170	<div> <div>4%</div> <div>85% 11% . .</div> </div>
1	B	170	<div> <div>8%</div> <div>91% 5% .</div> </div>
1	C	170	<div> <div>5%</div> <div>89% 6% . .</div> </div>
1	D	170	<div> <div>4%</div> <div>89% 8% .</div> </div>
1	E	170	<div> <div>6%</div> <div>89% 7% .</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	170	<div><div></div><div>14%</div><div>88%</div><div>7%</div><div>5%</div></div>
1	G	170	<div><div></div><div>7%</div><div>95%</div><div></div><div></div></div>
1	H	170	<div><div></div><div>8%</div><div>85%</div><div>11%</div><div></div></div>
1	I	170	<div><div></div><div>8%</div><div>88%</div><div>8%</div><div></div></div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 12692 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 Tyrosine-protein phosphatase SIW14.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	165	Total	C	N	O	S	Se	0	0	0
			1355	874	238	237	3	3			
1	B	164	Total	C	N	O	S	Se	0	0	0
			1356	878	237	235	3	3			
1	C	163	Total	C	N	O	S	Se	0	0	0
			1345	869	236	234	3	3			
1	D	165	Total	C	N	O	S	Se	0	0	0
			1355	874	238	237	3	3			
1	E	164	Total	C	N	O	S	Se	0	0	0
			1356	878	237	235	3	3			
1	F	162	Total	C	N	O	S	Se	0	1	0
			1345	870	235	233	3	4			
1	G	164	Total	C	N	O	S	Se	0	0	0
			1347	870	236	235	3	3			
1	H	163	Total	C	N	O	S	Se	0	0	0
			1345	869	236	234	3	3			
1	I	163	Total	C	N	O	S	Se	0	0	0
			1345	869	236	234	3	3			

There are 36 discrepancies between the modelled and reference sequences:

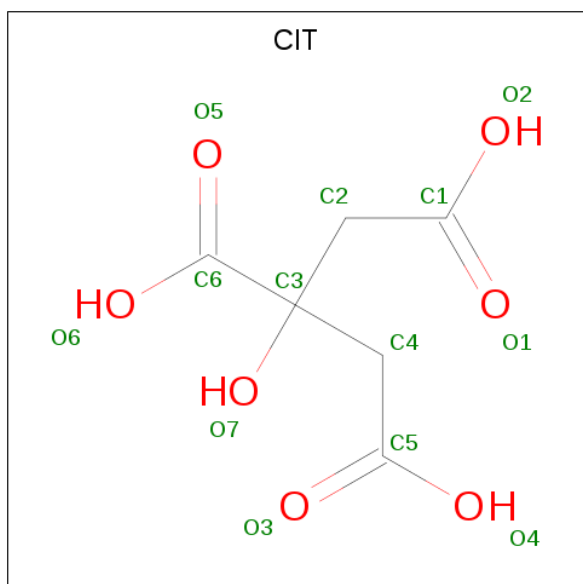
Chain	Residue	Modelled	Actual	Comment	Reference
A	112	GLY	-	expression tag	UNP P53965
A	113	SER	-	expression tag	UNP P53965
A	114	GLY	-	expression tag	UNP P53965
A	115	SER	-	expression tag	UNP P53965
B	112	GLY	-	expression tag	UNP P53965
B	113	SER	-	expression tag	UNP P53965
B	114	GLY	-	expression tag	UNP P53965
B	115	SER	-	expression tag	UNP P53965
C	112	GLY	-	expression tag	UNP P53965
C	113	SER	-	expression tag	UNP P53965
C	114	GLY	-	expression tag	UNP P53965

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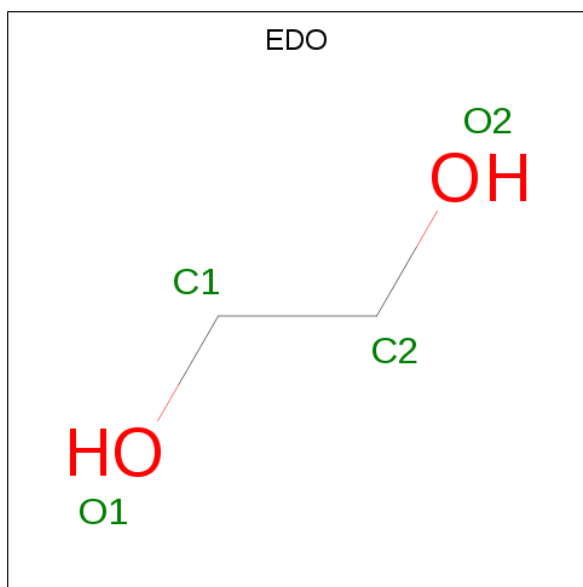
Chain	Residue	Modelled	Actual	Comment	Reference
C	115	SER	-	expression tag	UNP P53965
D	112	GLY	-	expression tag	UNP P53965
D	113	SER	-	expression tag	UNP P53965
D	114	GLY	-	expression tag	UNP P53965
D	115	SER	-	expression tag	UNP P53965
E	112	GLY	-	expression tag	UNP P53965
E	113	SER	-	expression tag	UNP P53965
E	114	GLY	-	expression tag	UNP P53965
E	115	SER	-	expression tag	UNP P53965
F	112	GLY	-	expression tag	UNP P53965
F	113	SER	-	expression tag	UNP P53965
F	114	GLY	-	expression tag	UNP P53965
F	115	SER	-	expression tag	UNP P53965
G	112	GLY	-	expression tag	UNP P53965
G	113	SER	-	expression tag	UNP P53965
G	114	GLY	-	expression tag	UNP P53965
G	115	SER	-	expression tag	UNP P53965
H	112	GLY	-	expression tag	UNP P53965
H	113	SER	-	expression tag	UNP P53965
H	114	GLY	-	expression tag	UNP P53965
H	115	SER	-	expression tag	UNP P53965
I	112	GLY	-	expression tag	UNP P53965
I	113	SER	-	expression tag	UNP P53965
I	114	GLY	-	expression tag	UNP P53965
I	115	SER	-	expression tag	UNP P53965

- Molecule 2 is CITRIC ACID (three-letter code: CIT) (formula: C₆H₈O₇).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			13	6	7		
2	B	1	Total	C	O	0	0
			13	6	7		
2	C	1	Total	C	O	0	0
			13	6	7		
2	D	1	Total	C	O	0	0
			13	6	7		
2	E	1	Total	C	O	0	0
			13	6	7		
2	F	1	Total	C	O	0	0
			13	6	7		
2	G	1	Total	C	O	0	0
			13	6	7		
2	H	1	Total	C	O	0	0
			13	6	7		
2	I	1	Total	C	O	0	0
			13	6	7		

- Molecule 3 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	C	1	Total	C	O	0	0
			4	2	2		
3	D	1	Total	C	O	0	0
			4	2	2		

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	G	1	Total 1	Cl 1	0	0
4	D	1	Total 1	Cl 1	0	0

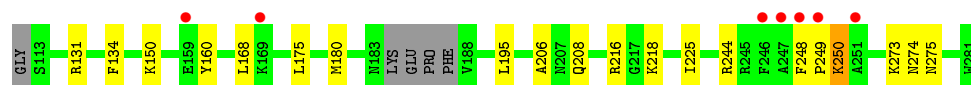
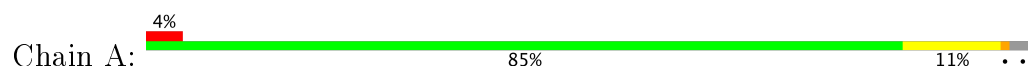
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	51	Total 51	O 51	0	0
5	B	47	Total 47	O 47	0	0
5	C	46	Total 46	O 46	0	0
5	D	48	Total 48	O 48	0	0
5	E	59	Total 59	O 59	0	0
5	F	36	Total 36	O 36	0	0
5	G	46	Total 46	O 46	0	0
5	H	41	Total 41	O 41	0	0
5	I	38	Total 38	O 38	0	0

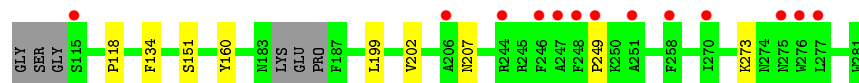
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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.

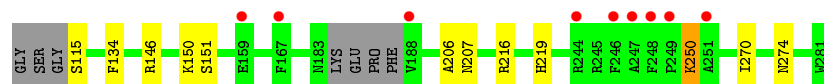
- Molecule 1: Tyrosine-protein phosphatase SIW14



- Molecule 1: Tyrosine-protein phosphatase SIW14



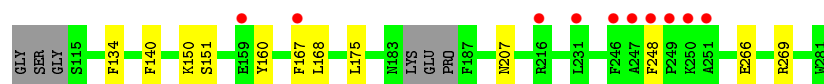
- Molecule 1: Tyrosine-protein phosphatase SIW14



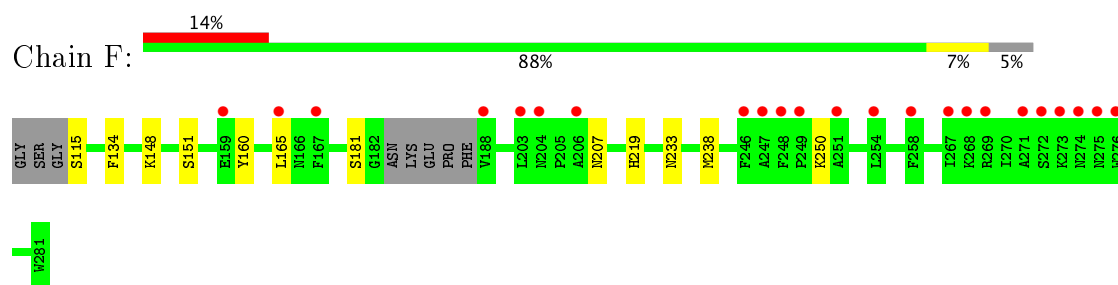
- Molecule 1: Tyrosine-protein phosphatase SIW14



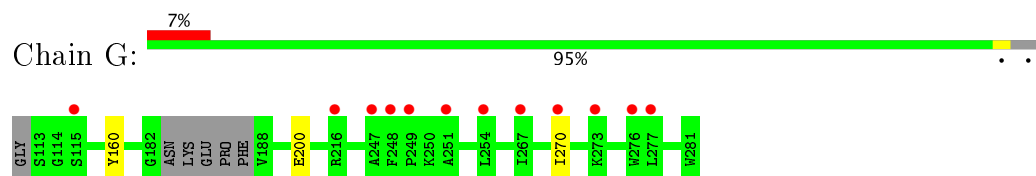
- Molecule 1: Tyrosine-protein phosphatase SIW14



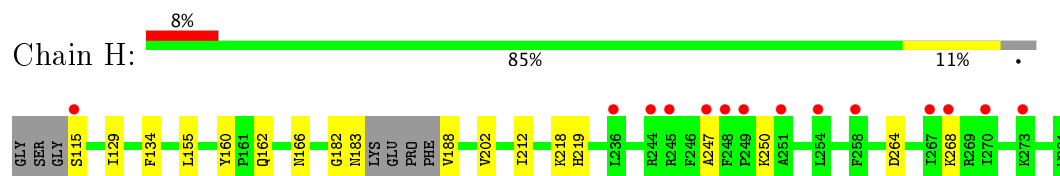
- Molecule 1: Tyrosine-protein phosphatase SIW14



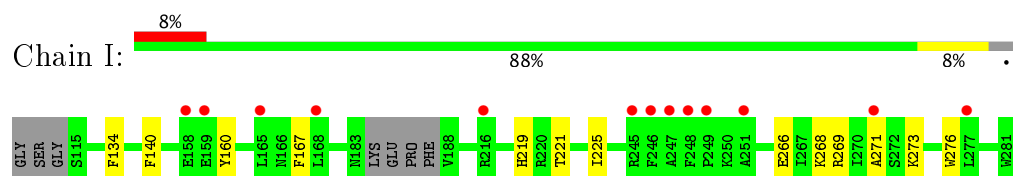
- Molecule 1: Tyrosine-protein phosphatase SIW14



- Molecule 1: Tyrosine-protein phosphatase SIW14



- Molecule 1: Tyrosine-protein phosphatase SIW14



4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	92.97Å 92.97Å 814.72Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.45 – 2.35 49.45 – 2.34	Depositor EDS
% Data completeness (in resolution range)	98.0 (49.45-2.35) 92.2 (49.45-2.34)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.30 (at 2.34Å)	Xtriage
Refinement program	PHENIX (1.12_2829: ???)	Depositor
R, R_{free}	0.206 , 0.229 0.225 , 0.238	Depositor DCC
R_{free} test set	1901 reflections (2.24%)	DCC
Wilson B-factor (Å ²)	37.8	Xtriage
Anisotropy	0.386	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 43.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	12692	wwPDB-VP
Average B, all atoms (Å ²)	46.0	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 49.83 % 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 7.0559e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.30	0/1385	0.48	0/1867
1	B	0.26	0/1387	0.45	0/1870
1	C	0.40	0/1375	0.47	0/1854
1	D	0.40	0/1385	0.49	0/1867
1	E	0.30	0/1387	0.46	0/1870
1	F	0.30	0/1375	0.48	0/1853
1	G	0.26	0/1377	0.47	0/1856
1	H	0.36	0/1375	0.48	0/1854
1	I	0.36	0/1375	0.47	0/1854
All	All	0.33	0/12421	0.47	0/16745

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 [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	1355	0	1368	15	0
1	B	1356	0	1369	4	0
1	C	1345	0	1360	9	0
1	D	1355	0	1368	11	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	1356	0	1369	5	0
1	F	1345	0	1362	3	0
1	G	1347	0	1362	1	0
1	H	1345	0	1360	9	0
1	I	1345	0	1360	5	0
2	A	13	0	5	1	0
2	B	13	0	5	0	0
2	C	13	0	5	1	0
2	D	13	0	5	1	0
2	E	13	0	5	0	0
2	F	13	0	5	0	0
2	G	13	0	5	1	0
2	H	13	0	5	0	0
2	I	13	0	5	1	0
3	A	4	0	6	0	0
3	C	4	0	6	0	0
3	D	4	0	6	0	0
4	D	1	0	0	0	0
4	G	1	0	0	0	0
5	A	51	0	0	0	0
5	B	47	0	0	0	0
5	C	46	0	0	0	0
5	D	48	0	0	3	0
5	E	59	0	0	0	0
5	F	36	0	0	0	0
5	G	46	0	0	0	0
5	H	41	0	0	1	0
5	I	38	0	0	0	0
All	All	12692	0	12341	59	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 2.

The worst 5 of 59 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:C:216:ARG:CZ	1:C:250:LYS:HE3	2.07	0.84
1:D:216:ARG:NH1	2:D:301:CIT:O2	2.24	0.70
1:C:216:ARG:NH1	1:C:250:LYS:HE3	2.07	0.69
1:A:216:ARG:NH1	2:A:301:CIT:O7	2.26	0.68
1:D:244:ARG:NE	5:D:401:HOH:O	2.33	0.62

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	161/170 (95%)	155 (96%)	6 (4%)	0	100	100
1	B	160/170 (94%)	155 (97%)	5 (3%)	0	100	100
1	C	159/170 (94%)	155 (98%)	4 (2%)	0	100	100
1	D	161/170 (95%)	157 (98%)	4 (2%)	0	100	100
1	E	160/170 (94%)	155 (97%)	5 (3%)	0	100	100
1	F	159/170 (94%)	154 (97%)	5 (3%)	0	100	100
1	G	160/170 (94%)	154 (96%)	6 (4%)	0	100	100
1	H	159/170 (94%)	154 (97%)	5 (3%)	0	100	100
1	I	159/170 (94%)	153 (96%)	6 (4%)	0	100	100
All	All	1438/1530 (94%)	1392 (97%)	46 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	151/152 (99%)	147 (97%)	4 (3%)	51	64
1	B	151/152 (99%)	148 (98%)	3 (2%)	60	73
1	C	150/152 (99%)	147 (98%)	3 (2%)	60	73

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	151/152 (99%)	149 (99%)	2 (1%)	73	84
1	E	151/152 (99%)	148 (98%)	3 (2%)	60	73
1	F	150/152 (99%)	143 (95%)	7 (5%)	30	38
1	G	150/152 (99%)	149 (99%)	1 (1%)	87	93
1	H	150/152 (99%)	147 (98%)	3 (2%)	60	73
1	I	150/152 (99%)	146 (97%)	4 (3%)	50	63
All	All	1354/1368 (99%)	1324 (98%)	30 (2%)	57	70

5 of 30 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	150	LYS
1	F	134	PHE
1	I	160	TYR
1	F	115	SER
1	F	148	LYS

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

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 14 ligands modelled in this entry, 2 are monoatomic - leaving 12 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	CIT	A	301	-	3,12,12	1.27	0	3,17,17	1.49	1 (33%)
3	EDO	A	302	-	3,3,3	0.44	0	2,2,2	0.33	0
2	CIT	B	301	-	3,12,12	1.32	0	3,17,17	1.85	1 (33%)
2	CIT	C	301	-	3,12,12	1.27	0	3,17,17	1.97	2 (66%)
3	EDO	C	302	-	3,3,3	0.48	0	2,2,2	0.30	0
2	CIT	D	301	-	3,12,12	1.29	0	3,17,17	2.22	2 (66%)
3	EDO	D	302	-	3,3,3	0.45	0	2,2,2	0.32	0
2	CIT	E	301	-	3,12,12	1.09	0	3,17,17	1.97	2 (66%)
2	CIT	F	301	-	3,12,12	1.25	0	3,17,17	1.75	1 (33%)
2	CIT	G	301	-	3,12,12	1.18	0	3,17,17	2.81	1 (33%)
2	CIT	H	301	-	3,12,12	1.39	0	3,17,17	2.23	2 (66%)
2	CIT	I	301	-	3,12,12	1.20	0	3,17,17	0.61	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	CIT	A	301	-	-	0/6/16/16	0/0/0/0
3	EDO	A	302	-	-	0/1/1/1	0/0/0/0
2	CIT	B	301	-	-	0/6/16/16	0/0/0/0
2	CIT	C	301	-	-	0/6/16/16	0/0/0/0
3	EDO	C	302	-	-	0/1/1/1	0/0/0/0
2	CIT	D	301	-	-	0/6/16/16	0/0/0/0
3	EDO	D	302	-	-	0/1/1/1	0/0/0/0
2	CIT	E	301	-	-	0/6/16/16	0/0/0/0
2	CIT	F	301	-	-	0/6/16/16	0/0/0/0
2	CIT	G	301	-	-	0/6/16/16	0/0/0/0
2	CIT	H	301	-	-	0/6/16/16	0/0/0/0
2	CIT	I	301	-	-	0/6/16/16	0/0/0/0

There are no bond length outliers.

The worst 5 of 12 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	G	301	CIT	C3-C4-C5	-4.80	107.45	114.95
2	H	301	CIT	C3-C2-C1	-3.25	109.87	114.95
2	D	301	CIT	C3-C4-C5	-2.66	110.80	114.95
2	D	301	CIT	C3-C2-C1	-2.65	110.81	114.95
2	B	301	CIT	C3-C2-C1	-2.63	110.85	114.95

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	301	CIT	1	0
2	C	301	CIT	1	0
2	D	301	CIT	1	0
2	G	301	CIT	1	0
2	I	301	CIT	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 [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	162/170 (95%)	0.41	7 (4%) 36 48	34, 43, 63, 74	0
1	B	161/170 (94%)	0.49	13 (8%) 13 18	34, 44, 64, 75	0
1	C	160/170 (94%)	0.41	9 (5%) 25 36	32, 43, 63, 79	0
1	D	162/170 (95%)	0.33	6 (3%) 42 54	30, 41, 61, 72	0
1	E	161/170 (94%)	0.45	10 (6%) 21 30	34, 43, 64, 75	0
1	F	159/170 (93%)	0.76	23 (14%) 3 5	36, 51, 79, 96	0
1	G	161/170 (94%)	0.56	12 (7%) 15 22	31, 43, 65, 76	0
1	H	160/170 (94%)	0.57	14 (8%) 11 16	33, 49, 68, 86	0
1	I	160/170 (94%)	0.41	13 (8%) 13 18	36, 47, 66, 79	0
All	All	1446/1530 (94%)	0.49	107 (7%) 15 22	30, 44, 67, 96	0

The worst 5 of 107 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	249	PRO	6.1
1	A	248	PHE	5.6
1	I	248	PHE	5.5
1	A	249	PRO	5.5
1	D	249	PRO	5.4

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(Å ²)	Q<0.9
3	EDO	A	302	4/4	0.91	0.27	1.92	32,34,34,40	0
2	CIT	B	301	13/13	0.78	0.23	0.53	41,60,70,75	0
3	EDO	D	302	4/4	0.93	0.20	0.23	36,37,38,38	0
2	CIT	C	301	13/13	0.84	0.19	0.12	42,58,79,80	0
2	CIT	G	301	13/13	0.75	0.21	-0.05	42,59,71,73	0
3	EDO	C	302	4/4	0.96	0.18	-0.38	31,32,32,35	0
2	CIT	F	301	13/13	0.87	0.17	-0.45	44,58,81,85	0
2	CIT	E	301	13/13	0.90	0.15	-0.56	37,56,64,68	0
2	CIT	I	301	13/13	0.89	0.15	-0.66	38,49,66,74	0
2	CIT	H	301	13/13	0.85	0.16	-0.73	44,55,69,71	0
2	CIT	D	301	13/13	0.88	0.16	-0.80	36,54,67,67	0
2	CIT	A	301	13/13	0.89	0.13	-0.85	42,55,69,69	0
4	CL	G	302	1/1	0.96	0.09	-2.45	41,41,41,41	0
4	CL	D	303	1/1	0.96	0.05	-3.13	44,44,44,44	0

6.5 Other polymers [i](#)

There are no such residues in this entry.