



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 10:50 PM GMT

PDB ID : 4ZAS
Title : Crystal structure of sugar aminotransferase CalS13 from *Micromonospora echinospora*
Authors : Wang, F.; Singh, S.; Miller, M.D.; Thorson, J.S.; Phillips Jr., G.N.; Enzyme Discovery for Natural Product Biosynthesis (NatPro)
Deposited on : 2015-04-14
Resolution : 2.47 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

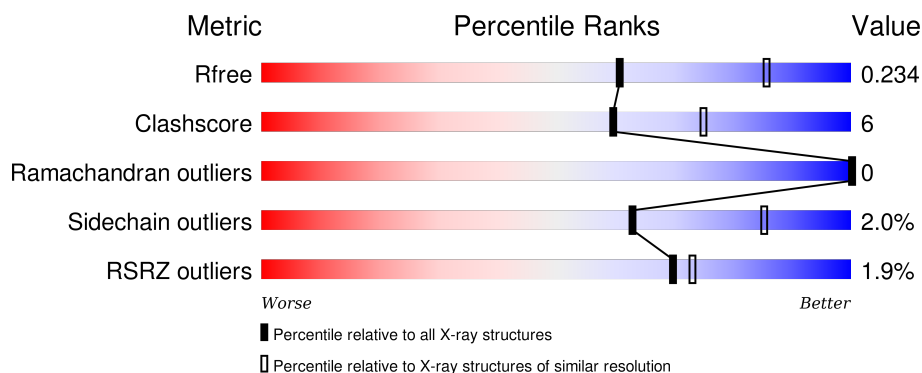
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.47 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	4309 (2.50-2.46)
Clashscore	102246	5050 (2.50-2.46)
Ramachandran outliers	100387	4961 (2.50-2.46)
Sidechain outliers	100360	4963 (2.50-2.46)
RSRZ outliers	91569	4319 (2.50-2.46)

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	404	<div> <div>2%</div> <div>79% 12% 9%</div> </div>
1	B	404	<div> <div></div> <div>79% 12% 9%</div> </div>
1	C	404	<div> <div></div> <div>81% 10% 9%</div> </div>
1	D	404	<div> <div>%</div> <div>79% 12% 9%</div> </div>
1	E	404	<div> <div>%</div> <div>77% 14% 9%</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	404	

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	SO4	C	401	-	-	-	X
3	T46	B	402	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 17830 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 CalS13.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	369	Total	C	N	O	P	S	0	0	0
			2868	1798	526	534	1	9			
1	B	369	Total	C	N	O	P	S	0	0	0
			2868	1798	526	534	1	9			
1	C	369	Total	C	N	O	P	S	0	0	0
			2868	1798	526	534	1	9			
1	D	369	Total	C	N	O	P	S	0	0	0
			2868	1798	526	534	1	9			
1	E	369	Total	C	N	O	P	S	0	0	0
			2868	1798	526	534	1	9			
1	F	369	Total	C	N	O	P	S	0	0	0
			2868	1798	526	534	1	9			

There are 168 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP Q8KND8
A	-18	GLY	-	expression tag	UNP Q8KND8
A	-17	SER	-	expression tag	UNP Q8KND8
A	-16	SER	-	expression tag	UNP Q8KND8
A	-15	HIS	-	expression tag	UNP Q8KND8
A	-14	HIS	-	expression tag	UNP Q8KND8
A	-13	HIS	-	expression tag	UNP Q8KND8
A	-12	HIS	-	expression tag	UNP Q8KND8
A	-11	HIS	-	expression tag	UNP Q8KND8
A	-10	HIS	-	expression tag	UNP Q8KND8
A	-9	SER	-	expression tag	UNP Q8KND8
A	-8	SER	-	expression tag	UNP Q8KND8
A	-7	GLY	-	expression tag	UNP Q8KND8
A	-6	LEU	-	expression tag	UNP Q8KND8
A	-5	VAL	-	expression tag	UNP Q8KND8
A	-4	PRO	-	expression tag	UNP Q8KND8
A	-3	ARG	-	expression tag	UNP Q8KND8

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP Q8KND8
A	-1	SER	-	expression tag	UNP Q8KND8
A	0	HIS	-	expression tag	UNP Q8KND8
A	1	MET	-	expression tag	UNP Q8KND8
A	2	ALA	-	expression tag	UNP Q8KND8
A	3	THR	-	expression tag	UNP Q8KND8
A	4	SER	-	expression tag	UNP Q8KND8
A	5	GLU	-	expression tag	UNP Q8KND8
A	6	ARG	-	expression tag	UNP Q8KND8
A	7	GLY	-	expression tag	UNP Q8KND8
A	8	VAL	-	expression tag	UNP Q8KND8
B	-19	MET	-	expression tag	UNP Q8KND8
B	-18	GLY	-	expression tag	UNP Q8KND8
B	-17	SER	-	expression tag	UNP Q8KND8
B	-16	SER	-	expression tag	UNP Q8KND8
B	-15	HIS	-	expression tag	UNP Q8KND8
B	-14	HIS	-	expression tag	UNP Q8KND8
B	-13	HIS	-	expression tag	UNP Q8KND8
B	-12	HIS	-	expression tag	UNP Q8KND8
B	-11	HIS	-	expression tag	UNP Q8KND8
B	-10	HIS	-	expression tag	UNP Q8KND8
B	-9	SER	-	expression tag	UNP Q8KND8
B	-8	SER	-	expression tag	UNP Q8KND8
B	-7	GLY	-	expression tag	UNP Q8KND8
B	-6	LEU	-	expression tag	UNP Q8KND8
B	-5	VAL	-	expression tag	UNP Q8KND8
B	-4	PRO	-	expression tag	UNP Q8KND8
B	-3	ARG	-	expression tag	UNP Q8KND8
B	-2	GLY	-	expression tag	UNP Q8KND8
B	-1	SER	-	expression tag	UNP Q8KND8
B	0	HIS	-	expression tag	UNP Q8KND8
B	1	MET	-	expression tag	UNP Q8KND8
B	2	ALA	-	expression tag	UNP Q8KND8
B	3	THR	-	expression tag	UNP Q8KND8
B	4	SER	-	expression tag	UNP Q8KND8
B	5	GLU	-	expression tag	UNP Q8KND8
B	6	ARG	-	expression tag	UNP Q8KND8
B	7	GLY	-	expression tag	UNP Q8KND8
B	8	VAL	-	expression tag	UNP Q8KND8
C	-19	MET	-	expression tag	UNP Q8KND8
C	-18	GLY	-	expression tag	UNP Q8KND8
C	-17	SER	-	expression tag	UNP Q8KND8

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-16	SER	-	expression tag	UNP Q8KND8
C	-15	HIS	-	expression tag	UNP Q8KND8
C	-14	HIS	-	expression tag	UNP Q8KND8
C	-13	HIS	-	expression tag	UNP Q8KND8
C	-12	HIS	-	expression tag	UNP Q8KND8
C	-11	HIS	-	expression tag	UNP Q8KND8
C	-10	HIS	-	expression tag	UNP Q8KND8
C	-9	SER	-	expression tag	UNP Q8KND8
C	-8	SER	-	expression tag	UNP Q8KND8
C	-7	GLY	-	expression tag	UNP Q8KND8
C	-6	LEU	-	expression tag	UNP Q8KND8
C	-5	VAL	-	expression tag	UNP Q8KND8
C	-4	PRO	-	expression tag	UNP Q8KND8
C	-3	ARG	-	expression tag	UNP Q8KND8
C	-2	GLY	-	expression tag	UNP Q8KND8
C	-1	SER	-	expression tag	UNP Q8KND8
C	0	HIS	-	expression tag	UNP Q8KND8
C	1	MET	-	expression tag	UNP Q8KND8
C	2	ALA	-	expression tag	UNP Q8KND8
C	3	THR	-	expression tag	UNP Q8KND8
C	4	SER	-	expression tag	UNP Q8KND8
C	5	GLU	-	expression tag	UNP Q8KND8
C	6	ARG	-	expression tag	UNP Q8KND8
C	7	GLY	-	expression tag	UNP Q8KND8
C	8	VAL	-	expression tag	UNP Q8KND8
D	-19	MET	-	expression tag	UNP Q8KND8
D	-18	GLY	-	expression tag	UNP Q8KND8
D	-17	SER	-	expression tag	UNP Q8KND8
D	-16	SER	-	expression tag	UNP Q8KND8
D	-15	HIS	-	expression tag	UNP Q8KND8
D	-14	HIS	-	expression tag	UNP Q8KND8
D	-13	HIS	-	expression tag	UNP Q8KND8
D	-12	HIS	-	expression tag	UNP Q8KND8
D	-11	HIS	-	expression tag	UNP Q8KND8
D	-10	HIS	-	expression tag	UNP Q8KND8
D	-9	SER	-	expression tag	UNP Q8KND8
D	-8	SER	-	expression tag	UNP Q8KND8
D	-7	GLY	-	expression tag	UNP Q8KND8
D	-6	LEU	-	expression tag	UNP Q8KND8
D	-5	VAL	-	expression tag	UNP Q8KND8
D	-4	PRO	-	expression tag	UNP Q8KND8
D	-3	ARG	-	expression tag	UNP Q8KND8

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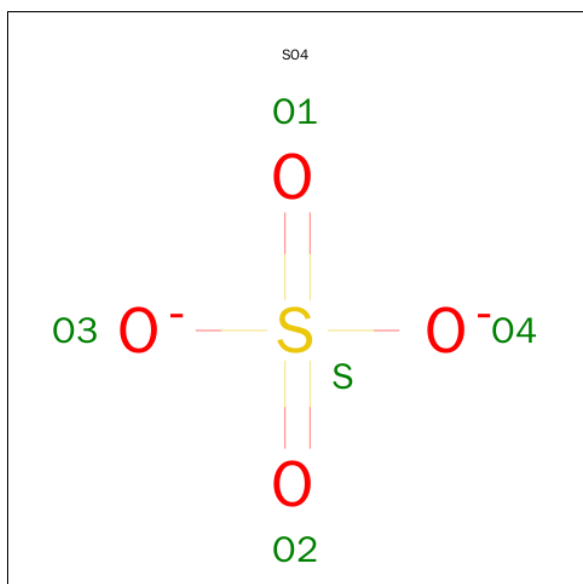
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D	-1	SER	-	expression tag	UNP Q8KND8
D	0	HIS	-	expression tag	UNP Q8KND8
D	1	MET	-	expression tag	UNP Q8KND8
D	2	ALA	-	expression tag	UNP Q8KND8
D	3	THR	-	expression tag	UNP Q8KND8
D	4	SER	-	expression tag	UNP Q8KND8
D	5	GLU	-	expression tag	UNP Q8KND8
D	6	ARG	-	expression tag	UNP Q8KND8
D	7	GLY	-	expression tag	UNP Q8KND8
D	8	VAL	-	expression tag	UNP Q8KND8
E	-19	MET	-	expression tag	UNP Q8KND8
E	-18	GLY	-	expression tag	UNP Q8KND8
E	-17	SER	-	expression tag	UNP Q8KND8
E	-16	SER	-	expression tag	UNP Q8KND8
E	-15	HIS	-	expression tag	UNP Q8KND8
E	-14	HIS	-	expression tag	UNP Q8KND8
E	-13	HIS	-	expression tag	UNP Q8KND8
E	-12	HIS	-	expression tag	UNP Q8KND8
E	-11	HIS	-	expression tag	UNP Q8KND8
E	-10	HIS	-	expression tag	UNP Q8KND8
E	-9	SER	-	expression tag	UNP Q8KND8
E	-8	SER	-	expression tag	UNP Q8KND8
E	-7	GLY	-	expression tag	UNP Q8KND8
E	-6	LEU	-	expression tag	UNP Q8KND8
E	-5	VAL	-	expression tag	UNP Q8KND8
E	-4	PRO	-	expression tag	UNP Q8KND8
E	-3	ARG	-	expression tag	UNP Q8KND8
E	-2	GLY	-	expression tag	UNP Q8KND8
E	-1	SER	-	expression tag	UNP Q8KND8
E	0	HIS	-	expression tag	UNP Q8KND8
E	1	MET	-	expression tag	UNP Q8KND8
E	2	ALA	-	expression tag	UNP Q8KND8
E	3	THR	-	expression tag	UNP Q8KND8
E	4	SER	-	expression tag	UNP Q8KND8
E	5	GLU	-	expression tag	UNP Q8KND8
E	6	ARG	-	expression tag	UNP Q8KND8
E	7	GLY	-	expression tag	UNP Q8KND8
E	8	VAL	-	expression tag	UNP Q8KND8
F	-19	MET	-	expression tag	UNP Q8KND8
F	-18	GLY	-	expression tag	UNP Q8KND8
F	-17	SER	-	expression tag	UNP Q8KND8

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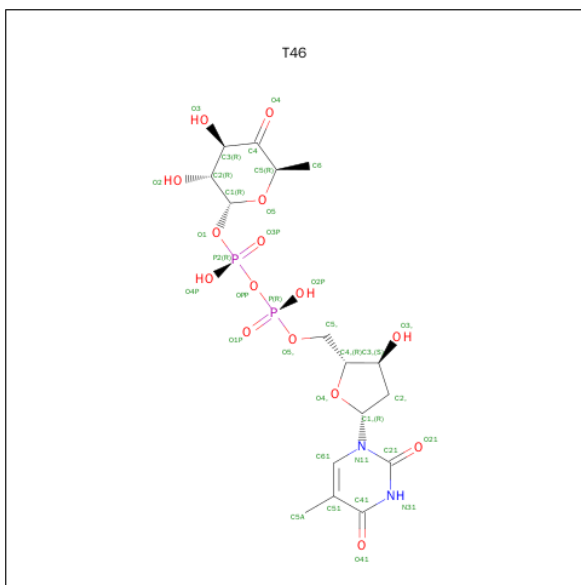
Chain	Residue	Modelled	Actual	Comment	Reference
F	-16	SER	-	expression tag	UNP Q8KND8
F	-15	HIS	-	expression tag	UNP Q8KND8
F	-14	HIS	-	expression tag	UNP Q8KND8
F	-13	HIS	-	expression tag	UNP Q8KND8
F	-12	HIS	-	expression tag	UNP Q8KND8
F	-11	HIS	-	expression tag	UNP Q8KND8
F	-10	HIS	-	expression tag	UNP Q8KND8
F	-9	SER	-	expression tag	UNP Q8KND8
F	-8	SER	-	expression tag	UNP Q8KND8
F	-7	GLY	-	expression tag	UNP Q8KND8
F	-6	LEU	-	expression tag	UNP Q8KND8
F	-5	VAL	-	expression tag	UNP Q8KND8
F	-4	PRO	-	expression tag	UNP Q8KND8
F	-3	ARG	-	expression tag	UNP Q8KND8
F	-2	GLY	-	expression tag	UNP Q8KND8
F	-1	SER	-	expression tag	UNP Q8KND8
F	0	HIS	-	expression tag	UNP Q8KND8
F	1	MET	-	expression tag	UNP Q8KND8
F	2	ALA	-	expression tag	UNP Q8KND8
F	3	THR	-	expression tag	UNP Q8KND8
F	4	SER	-	expression tag	UNP Q8KND8
F	5	GLU	-	expression tag	UNP Q8KND8
F	6	ARG	-	expression tag	UNP Q8KND8
F	7	GLY	-	expression tag	UNP Q8KND8
F	8	VAL	-	expression tag	UNP Q8KND8

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



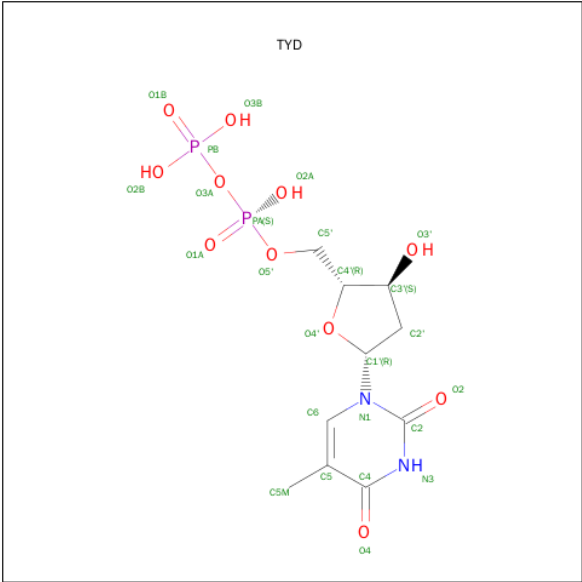
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total 5	O 4	S 1	0	0
2	C	1	Total 5	O 4	S 1	0	0

- Molecule 3 is dTDP-4-keto-6-deoxyglucose (three-letter code: T46) (formula: $\text{C}_{16}\text{H}_{24}\text{N}_2\text{O}_{15}\text{P}_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	B	1	Total	C	N	O	P	0	0
			35	16	2	15	2		

- Molecule 4 is THYMIDINE-5'-DIPHOSPHATE (three-letter code: TYD) (formula: $\text{C}_{10}\text{H}_{16}\text{N}_2\text{O}_{11}\text{P}_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	D	1	Total	C	N	O	P	0	0
			25	10	2	11	2		
4	E	1	Total	C	N	O	P	0	0
			25	10	2	11	2		
4	F	1	Total	C	N	O	P	0	0
			25	10	2	11	2		

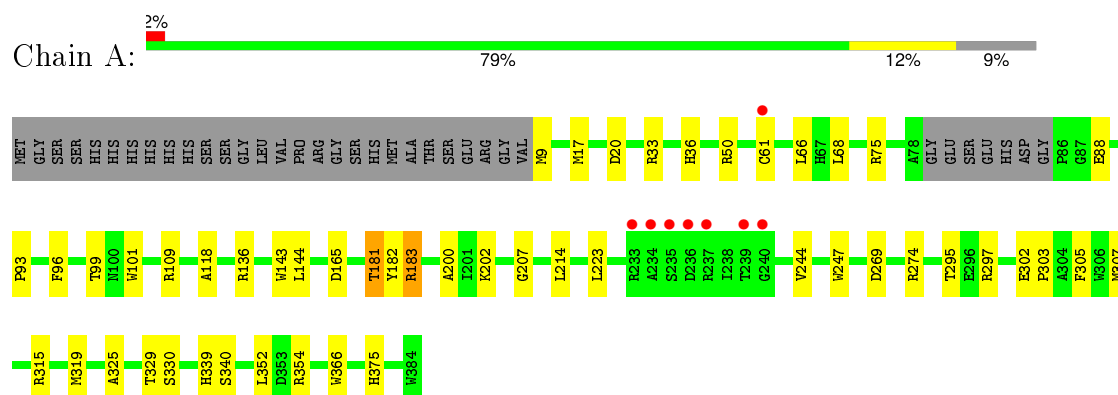
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	95	Total	O	0	0
			95	95		
5	B	112	Total	O	0	0
			112	112		
5	C	69	Total	O	0	0
			69	69		
5	D	101	Total	O	0	0
			101	101		
5	E	67	Total	O	0	0
			67	67		
5	F	58	Total	O	0	0
			58	58		

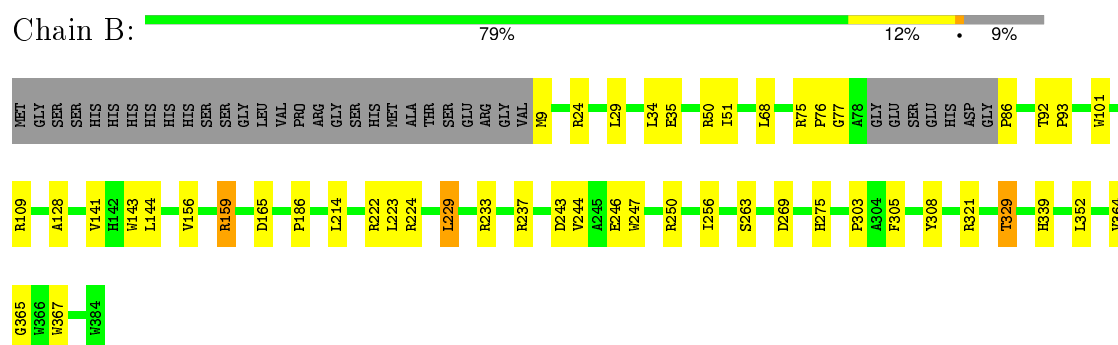
3 Residue-property plots [i](#)

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.

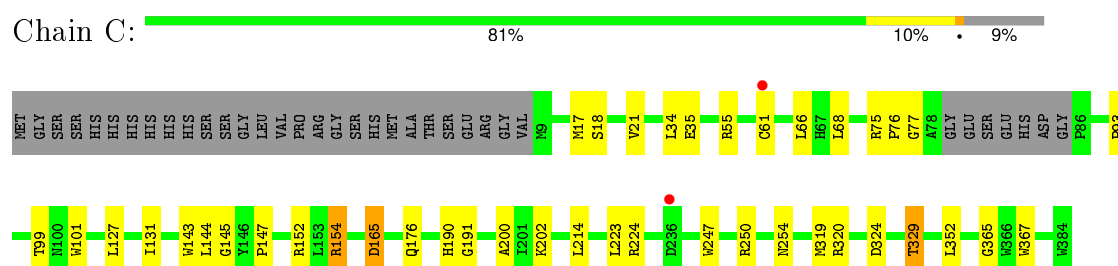
• Molecule 1: CalS13



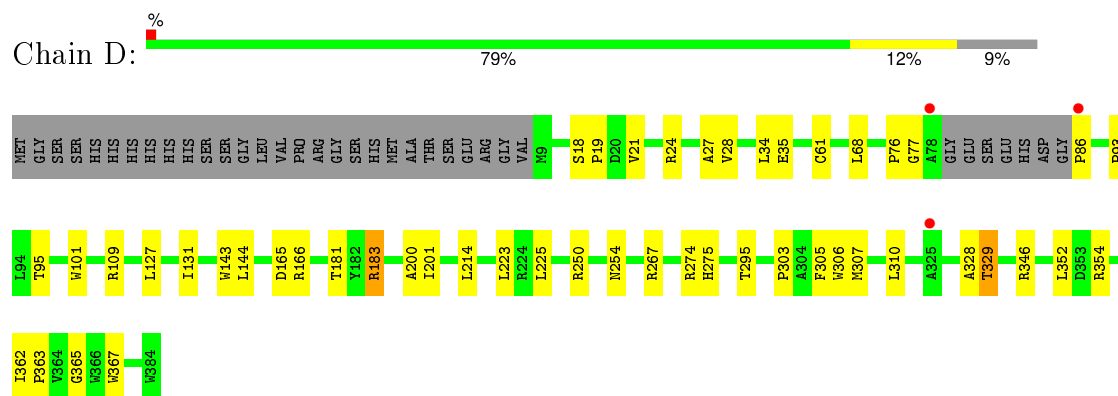
• Molecule 1: CalS13



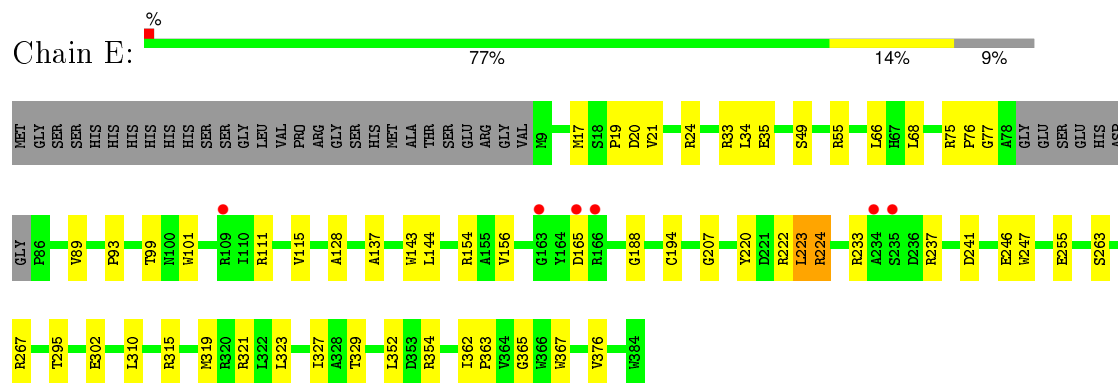
• Molecule 1: CalS13



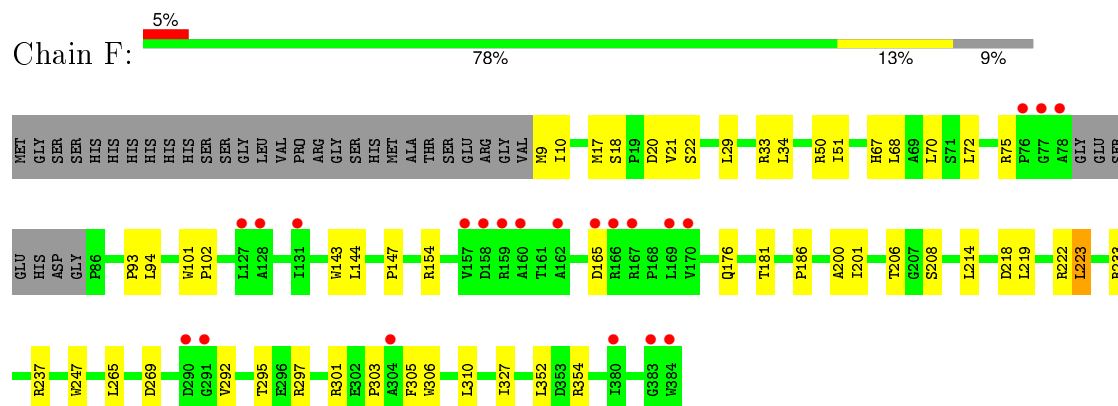
• Molecule 1: CalS13



- Molecule 1: CalS13



- Molecule 1: CalS13



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	71.57Å 88.74Å 189.81Å 90.00° 96.22° 90.00°	Depositor
Resolution (Å)	43.19 – 2.47 44.37 – 2.47	Depositor EDS
% Data completeness (in resolution range)	99.8 (43.19-2.47) 90.9 (44.37-2.47)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.78 (at 2.48Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_1951)	Depositor
R, R_{free}	0.173 , 0.229 0.184 , 0.234	Depositor DCC
R_{free} test set	2006 reflections (2.60%)	DCC
Wilson B-factor (Å ²)	40.3	Xtriage
Anisotropy	0.217	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 36.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	1 of 84859 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	17830	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.74% 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: TYD, LLP, SO4, T46

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	0.44	0/2904	0.62	0/3956
1	B	0.44	0/2904	0.64	1/3956 (0.0%)
1	C	0.40	0/2904	0.59	0/3956
1	D	0.42	0/2904	0.61	0/3956
1	E	0.40	0/2904	0.60	0/3956
1	F	0.37	0/2904	0.57	0/3956
All	All	0.41	0/17424	0.60	1/23736 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	229	LEU	CA-CB-CG	-5.05	103.69	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	2868	0	2841	36	0
1	B	2868	0	2841	36	0
1	C	2868	0	2841	30	0
1	D	2868	0	2841	35	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	2868	0	2841	40	0
1	F	2868	0	2841	36	0
2	B	5	0	0	0	0
2	C	5	0	0	0	0
3	B	35	0	22	2	0
4	D	25	0	13	0	0
4	E	25	0	13	2	0
4	F	25	0	13	0	0
5	A	95	0	0	8	0
5	B	112	0	0	11	0
5	C	69	0	0	4	0
5	D	101	0	0	8	0
5	E	67	0	0	9	0
5	F	58	0	0	7	0
All	All	17830	0	17107	197	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 6.

All (197) 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:E:319:MET:SD	5:E:653:HOH:O	2.34	0.84
1:B:9:MET:N	5:B:501:HOH:O	2.13	0.80
1:A:319:MET:SD	5:A:429:HOH:O	2.38	0.80
1:E:222:ARG:NH1	1:E:246:GLU:OE2	2.25	0.69
1:A:20:ASP:OD1	5:A:401:HOH:O	2.12	0.68
1:F:50:ARG:NH1	1:F:269:ASP:OD1	2.22	0.68
1:C:202:LLP:OP1	5:C:501:HOH:O	2.11	0.68
1:D:295:THR:HG21	5:D:697:HOH:O	1.95	0.67
1:E:302:GLU:HG3	5:E:666:HOH:O	1.94	0.67
1:E:75:ARG:NH2	1:E:247:TRP:O	2.28	0.66
1:C:93:PRO:HG2	1:C:352:LEU:HD22	1.79	0.65
1:B:222:ARG:NH1	1:B:246:GLU:OE2	2.29	0.64
1:F:94:LEU:HG	5:F:646:HOH:O	1.97	0.64
1:C:176:GLN:HE21	1:C:202:LLP:HO3	1.46	0.63
1:E:233:ARG:O	1:E:237:ARG:NH2	2.32	0.63
1:F:143:TRP:CD1	1:F:144:LEU:HG	2.34	0.63
1:D:329:THR:HB	5:D:653:HOH:O	1.99	0.62
1:E:68:LEU:HG	1:E:223:LEU:HG	1.81	0.62
1:E:143:TRP:CD1	1:E:144:LEU:HG	2.35	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:75:ARG:NH2	1:A:247:TRP:O	2.32	0.62
1:B:86:PRO:O	5:B:502:HOH:O	2.16	0.61
1:B:233:ARG:O	1:B:237:ARG:NH2	2.34	0.61
1:C:152:ARG:NH2	5:C:503:HOH:O	2.34	0.61
1:C:329:THR:HB	5:C:516:HOH:O	2.00	0.61
1:B:75:ARG:NH2	1:B:247:TRP:O	2.33	0.60
1:F:201:ILE:HG13	1:F:306:TRP:HZ2	1.67	0.60
1:A:295:THR:HG21	5:A:491:HOH:O	2.01	0.60
1:D:310:LEU:HD12	1:D:362:ILE:HD12	1.84	0.60
1:E:295:THR:HG21	5:E:667:HOH:O	2.01	0.59
1:E:241:ASP:OD1	5:E:602:HOH:O	2.16	0.59
1:F:218:ASP:HB3	5:F:631:HOH:O	2.02	0.59
1:A:136:ARG:NH1	5:A:402:HOH:O	2.22	0.59
1:B:329:THR:HB	5:B:552:HOH:O	2.03	0.59
1:B:128:ALA:HB2	1:B:156:VAL:HG13	1.85	0.59
1:D:28:VAL:HG23	5:D:601:HOH:O	2.01	0.59
1:A:202:LLP:H4'1	3:B:402:T46:O4	2.03	0.59
1:F:295:THR:HB	1:F:297:ARG:HH12	1.68	0.58
1:E:89:VAL:HG22	1:E:137:ALA:HB3	1.85	0.58
1:C:214:LEU:HD11	1:C:223:LEU:HD22	1.86	0.58
1:B:51:ILE:O	1:B:186:PRO:HG2	2.03	0.58
1:A:9:MET:N	5:A:405:HOH:O	2.36	0.58
1:A:68:LEU:HG	1:A:223:LEU:HG	1.86	0.58
1:F:93:PRO:HG2	1:F:352:LEU:HD22	1.87	0.57
1:B:143:TRP:CD1	1:B:144:LEU:HG	2.40	0.57
1:E:111:ARG:HD3	5:E:665:HOH:O	2.02	0.57
1:A:33:ARG:HH12	1:A:36:HIS:HB2	1.69	0.57
1:F:222:ARG:NE	5:F:604:HOH:O	2.30	0.56
1:B:214:LEU:HD11	1:B:223:LEU:HD22	1.86	0.56
1:D:68:LEU:HG	1:D:223:LEU:HG	1.86	0.56
1:B:68:LEU:HG	1:B:223:LEU:HG	1.88	0.56
1:C:18:SER:HB3	1:C:21:VAL:HG23	1.87	0.56
1:C:68:LEU:HG	1:C:223:LEU:HG	1.86	0.56
1:D:214:LEU:HD11	1:D:223:LEU:HD22	1.88	0.56
1:E:21:VAL:HG12	1:F:29:LEU:HD13	1.87	0.56
1:D:18:SER:HB3	1:D:21:VAL:HG23	1.87	0.55
5:B:505:HOH:O	1:D:86:PRO:HB2	2.06	0.55
1:B:93:PRO:HG2	1:B:352:LEU:HD22	1.89	0.55
1:E:93:PRO:HG2	1:E:352:LEU:HD22	1.87	0.55
1:F:75:ARG:NH2	1:F:247:TRP:O	2.39	0.54
1:C:365:GLY:HA3	1:C:367:TRP:CZ3	2.42	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:214:LEU:HD11	1:A:223:LEU:HD22	1.89	0.53
1:E:267:ARG:NH2	5:E:601:HOH:O	2.06	0.53
1:A:66:LEU:HD13	1:A:99:THR:HA	1.91	0.53
1:E:128:ALA:HB2	1:E:156:VAL:HG13	1.91	0.53
1:B:24:ARG:HD2	1:B:263:SER:OG	2.09	0.53
1:A:200:ALA:HB2	1:B:35:GLU:HG3	1.90	0.52
1:D:34:LEU:O	1:D:254:ASN:HB2	2.09	0.52
1:A:93:PRO:HG2	1:A:352:LEU:HD22	1.91	0.52
1:D:143:TRP:CD1	1:D:144:LEU:HG	2.45	0.52
1:B:250:ARG:NH1	5:B:504:HOH:O	2.33	0.52
1:F:72:LEU:HD13	1:F:219:LEU:HD22	1.92	0.51
1:B:76:PRO:HD2	5:B:604:HOH:O	2.11	0.51
1:A:143:TRP:CD1	1:A:144:LEU:HG	2.45	0.51
1:F:20:ASP:OD2	5:F:601:HOH:O	2.18	0.51
3:B:402:T46:O5,	3:B:402:T46:H9	2.10	0.51
1:B:76:PRO:N	1:B:77:GLY:HA3	2.24	0.51
1:E:354:ARG:NH2	5:E:603:HOH:O	2.22	0.51
1:C:200:ALA:HB2	1:D:35:GLU:HG3	1.92	0.51
1:D:354:ARG:HD3	5:D:694:HOH:O	2.09	0.51
1:F:22:SER:N	5:F:611:HOH:O	2.43	0.51
1:F:50:ARG:NH2	1:F:265:LEU:O	2.44	0.51
1:E:365:GLY:HA3	1:E:367:TRP:CH2	2.46	0.50
1:C:165:ASP:N	5:C:505:HOH:O	2.44	0.50
1:B:29:LEU:HD21	1:B:256:ILE:HD11	1.93	0.50
1:C:143:TRP:CD1	1:C:144:LEU:HG	2.46	0.50
1:E:188:GLY:HA2	1:E:194:CYS:SG	2.52	0.50
1:F:354:ARG:HG2	1:F:354:ARG:O	2.12	0.50
1:F:10:ILE:HD12	1:F:327:ILE:HG12	1.94	0.49
1:C:154:ARG:NH1	1:C:190:HIS:O	2.45	0.49
1:C:154:ARG:HH22	1:C:191:GLY:HA3	1.77	0.49
1:F:147:PRO:HB2	1:F:301:ARG:HD3	1.94	0.49
1:E:66:LEU:HD13	1:E:99:THR:HA	1.93	0.49
1:B:109:ARG:HD3	5:B:609:HOH:O	2.13	0.49
1:D:144:LEU:HD22	1:D:307:MET:HB2	1.95	0.48
1:D:365:GLY:HA3	1:D:367:TRP:CZ3	2.47	0.48
1:B:50:ARG:NH1	1:B:269:ASP:OD1	2.46	0.48
1:E:33:ARG:NH2	4:E:500:TYD:H2'2	2.28	0.48
1:D:76:PRO:N	1:D:77:GLY:HA3	2.29	0.48
1:D:27:ALA:HB3	5:D:601:HOH:O	2.13	0.48
1:E:365:GLY:HA3	1:E:367:TRP:CZ3	2.49	0.48
1:D:201:ILE:HG13	1:D:306:TRP:HZ2	1.78	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:166:ARG:NH1	5:D:605:HOH:O	2.46	0.48
1:D:93:PRO:HG2	1:D:352:LEU:HD22	1.95	0.48
1:B:303:PRO:HB2	1:B:305:PHE:CE2	2.48	0.48
1:F:214:LEU:HD11	1:F:223:LEU:HD22	1.95	0.48
1:A:274:ARG:NH2	5:A:406:HOH:O	2.37	0.48
1:F:67:HIS:HA	1:F:102:PRO:HB3	1.95	0.48
1:B:92:THR:HG22	5:B:503:HOH:O	2.13	0.47
1:D:346:ARG:HD3	5:D:698:HOH:O	2.14	0.47
1:C:61:CYS:SG	1:D:250:ARG:HG3	2.54	0.47
1:B:275:HIS:CE1	1:B:365:GLY:HA2	2.49	0.47
1:E:327:ILE:HD13	1:E:376:VAL:HA	1.97	0.47
1:C:75:ARG:NH2	1:C:247:TRP:O	2.46	0.47
1:F:51:ILE:O	1:F:186:PRO:HG2	2.15	0.47
1:D:275:HIS:CE1	1:D:365:GLY:HA2	2.50	0.47
1:D:19:PRO:HD2	1:D:267:ARG:HE	1.80	0.47
1:A:17:MET:CE	1:B:34:LEU:HD11	2.46	0.46
1:A:61:CYS:SG	1:B:250:ARG:HG3	2.56	0.46
1:C:127:LEU:O	1:C:131:ILE:HG13	2.15	0.46
1:C:176:GLN:NE2	1:C:202:LLP:O3	2.39	0.46
1:C:76:PRO:N	1:C:77:GLY:HA3	2.30	0.46
1:A:182:TYR:CE2	1:A:183:ARG:HG3	2.50	0.46
1:C:320:ARG:NH1	1:C:324:ASP:OD1	2.49	0.46
1:B:75:ARG:HD3	5:B:590:HOH:O	2.16	0.46
1:E:35:GLU:HG3	1:F:200:ALA:HB2	1.97	0.46
1:F:10:ILE:HA	5:F:602:HOH:O	2.15	0.45
1:F:233:ARG:HA	1:F:237:ARG:HH21	1.81	0.45
1:F:292:VAL:HG13	1:F:310:LEU:HD22	1.98	0.45
1:D:303:PRO:HB2	1:D:305:PHE:CE2	2.51	0.45
1:E:49:SER:HB3	5:E:641:HOH:O	2.16	0.45
1:E:207:GLY:HA2	1:F:34:LEU:HD13	1.97	0.45
1:A:207:GLY:HA2	1:B:34:LEU:HD13	1.97	0.45
1:D:274:ARG:NH2	5:D:604:HOH:O	2.41	0.45
1:E:24:ARG:HD2	1:E:263:SER:OG	2.17	0.44
1:D:183:ARG:NH2	5:E:601:HOH:O	2.50	0.44
1:B:365:GLY:HA3	1:B:367:TRP:CZ3	2.52	0.44
1:C:34:LEU:O	1:C:254:ASN:HB2	2.18	0.44
1:A:50:ARG:NH1	1:A:269:ASP:OD1	2.43	0.44
1:E:223:LEU:HD12	1:E:223:LEU:HA	1.64	0.44
1:D:183:ARG:HH11	1:E:20:ASP:CG	2.21	0.44
1:A:274:ARG:NH1	1:A:366:TRP:O	2.50	0.44
1:B:141:VAL:HG23	5:B:503:HOH:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:33:ARG:O	1:E:255:GLU:HG3	2.17	0.44
1:A:315:ARG:O	1:A:319:MET:HG3	2.17	0.44
1:A:340:SER:OG	1:B:243:ASP:OD2	2.29	0.44
1:E:19:PRO:HD2	1:E:267:ARG:HE	1.82	0.44
1:A:181:THR:HG22	1:A:302:GLU:HB2	1.99	0.43
1:B:159:ARG:NH2	5:B:508:HOH:O	2.47	0.43
1:B:365:GLY:HA3	1:B:367:TRP:CH2	2.53	0.43
1:F:206:THR:OG1	1:F:208:SER:O	2.23	0.43
1:C:66:LEU:HD13	1:C:99:THR:HA	2.01	0.43
1:B:308:TYR:HB2	1:B:364:VAL:HB	2.00	0.43
1:D:127:LEU:O	1:D:131:ILE:HG13	2.19	0.43
1:C:365:GLY:HA3	1:C:367:TRP:CH2	2.54	0.43
1:F:68:LEU:HG	1:F:223:LEU:HG	2.00	0.43
1:A:244:VAL:HG23	1:B:339:HIS:CE1	2.54	0.43
1:C:145:GLY:O	1:C:147:PRO:HD3	2.18	0.43
1:D:95:THR:HA	1:D:143:TRP:CZ3	2.54	0.43
1:E:17:MET:CE	1:F:34:LEU:HD11	2.49	0.43
1:E:315:ARG:O	1:E:319:MET:HG3	2.18	0.43
1:F:144:LEU:N	1:F:176:GLN:OE1	2.45	0.42
4:E:500:TYD:O5'	4:E:500:TYD:H6	2.19	0.42
1:E:76:PRO:N	1:E:77:GLY:HA3	2.33	0.42
1:D:223:LEU:HD12	1:D:223:LEU:HA	1.85	0.42
1:F:303:PRO:HB2	1:F:305:PHE:CE2	2.55	0.42
1:C:319:MET:HG2	1:C:329:THR:HG21	2.01	0.42
1:A:136:ARG:HD3	5:A:402:HOH:O	2.19	0.42
1:F:9:MET:O	5:F:602:HOH:O	2.22	0.42
1:A:144:LEU:HD22	1:A:307:MET:HB2	2.02	0.42
1:F:18:SER:O	1:F:21:VAL:HG23	2.19	0.42
1:E:34:LEU:HD11	1:F:17:MET:CE	2.50	0.42
1:A:330:SER:HA	5:A:429:HOH:O	2.20	0.41
1:D:365:GLY:HA3	1:D:367:TRP:CH2	2.55	0.41
1:C:55:ARG:HD2	1:C:55:ARG:HH11	1.65	0.41
1:D:328:ALA:O	1:D:363:PRO:HD3	2.20	0.41
1:F:147:PRO:HG2	1:F:301:ARG:HB2	2.02	0.41
1:E:220:TYR:O	1:E:224:ARG:HB2	2.21	0.41
1:A:118:ALA:O	1:A:295:THR:HG23	2.20	0.41
1:E:310:LEU:HD12	1:E:362:ILE:HD12	2.03	0.41
1:A:354:ARG:HA	1:A:354:ARG:HD2	1.75	0.41
1:E:323:LEU:HA	1:E:323:LEU:HD12	1.84	0.41
1:A:339:HIS:CE1	1:B:244:VAL:HG23	2.55	0.41
1:F:201:ILE:HG13	1:F:306:TRP:CZ2	2.52	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:17:MET:SD	1:C:21:VAL:HG21	2.60	0.41
1:F:70:LEU:HD12	1:F:102:PRO:HB2	2.03	0.41
1:E:55:ARG:HH11	1:E:55:ARG:HD2	1.72	0.41
1:A:96:PHE:CZ	1:B:229:LEU:HD13	2.56	0.41
1:C:250:ARG:HG3	1:D:61:CYS:SG	2.61	0.41
1:C:35:GLU:HG3	1:D:200:ALA:HB2	2.02	0.41
1:C:223:LEU:HA	1:C:223:LEU:HD12	1.80	0.40
1:A:303:PRO:HB2	1:A:305:PHE:CE2	2.56	0.40
1:A:295:THR:HB	1:A:297:ARG:HH12	1.86	0.40
1:D:225:LEU:HA	1:D:225:LEU:HD23	1.85	0.40
1:A:88:GLU:HG2	1:A:109:ARG:HB2	2.03	0.40
1:E:115:VAL:HG21	1:E:352:LEU:HD13	2.03	0.40
1:E:362:ILE:HA	1:E:363:PRO:HD3	1.98	0.40
1:A:325:ALA:HB1	1:A:375:HIS:HE2	1.86	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	364/404 (90%)	355 (98%)	9 (2%)	0	100	100
1	B	364/404 (90%)	354 (97%)	10 (3%)	0	100	100
1	C	364/404 (90%)	355 (98%)	9 (2%)	0	100	100
1	D	364/404 (90%)	355 (98%)	9 (2%)	0	100	100
1	E	364/404 (90%)	355 (98%)	9 (2%)	0	100	100
1	F	364/404 (90%)	356 (98%)	8 (2%)	0	100	100
All	All	2184/2424 (90%)	2130 (98%)	54 (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	297/325 (91%)	292 (98%)	5 (2%)	68	88
1	B	297/325 (91%)	291 (98%)	6 (2%)	63	85
1	C	297/325 (91%)	292 (98%)	5 (2%)	68	88
1	D	297/325 (91%)	290 (98%)	7 (2%)	57	81
1	E	297/325 (91%)	290 (98%)	7 (2%)	57	81
1	F	297/325 (91%)	291 (98%)	6 (2%)	63	85
All	All	1782/1950 (91%)	1746 (98%)	36 (2%)	63	85

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	101	TRP
1	A	165	ASP
1	A	181	THR
1	A	183	ARG
1	A	329	THR
1	B	101	TRP
1	B	159	ARG
1	B	165	ASP
1	B	224	ARG
1	B	321	ARG
1	B	329	THR
1	C	101	TRP
1	C	154	ARG
1	C	165	ASP
1	C	224	ARG
1	C	329	THR
1	D	24	ARG
1	D	101	TRP
1	D	109	ARG
1	D	165	ASP
1	D	181	THR
1	D	183	ARG

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Mol	Chain	Res	Type
1	D	329	THR
1	E	101	TRP
1	E	154	ARG
1	E	165	ASP
1	E	223	LEU
1	E	224	ARG
1	E	321	ARG
1	E	329	THR
1	F	33	ARG
1	F	101	TRP
1	F	154	ARG
1	F	165	ASP
1	F	181	THR
1	F	223	LEU

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 molecules in this entry.

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

6 non-standard protein/DNA/RNA residues are 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
1	LLP	A	202	1	23,24,25	2.79	7 (30%)	28,32,34	1.77	6 (21%)
1	LLP	B	202	1	23,24,25	2.98	7 (30%)	28,32,34	1.12	2 (7%)
1	LLP	C	202	1	23,24,25	2.87	9 (39%)	28,32,34	1.42	4 (14%)
1	LLP	D	202	1	23,24,25	2.89	10 (43%)	28,32,34	1.29	3 (10%)
1	LLP	E	202	1	23,24,25	2.83	7 (30%)	28,32,34	1.31	5 (17%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	LLP	F	202	1	23,24,25	2.81	8 (34%)	28,32,34	1.32	5 (17%)

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
1	LLP	A	202	1	-	0/15/17/19	0/1/1/1
1	LLP	B	202	1	-	0/15/17/19	0/1/1/1
1	LLP	C	202	1	-	0/15/17/19	0/1/1/1
1	LLP	D	202	1	-	0/15/17/19	0/1/1/1
1	LLP	E	202	1	-	0/15/17/19	0/1/1/1
1	LLP	F	202	1	-	0/15/17/19	0/1/1/1

All (48) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	202	LLP	C4-C5	-3.66	1.37	1.42
1	F	202	LLP	C4-C5	-3.58	1.37	1.42
1	B	202	LLP	C4-C5	-3.30	1.37	1.42
1	D	202	LLP	C4-C5	-3.20	1.37	1.42
1	F	202	LLP	C4-C3	-3.02	1.37	1.40
1	C	202	LLP	C4-C3	-2.98	1.37	1.40
1	E	202	LLP	C4-C5	-2.88	1.38	1.42
1	D	202	LLP	CB-CA	-2.72	1.51	1.53
1	E	202	LLP	C4-C3	-2.68	1.37	1.40
1	F	202	LLP	CB-CA	-2.62	1.51	1.53
1	D	202	LLP	C4-C3	-2.54	1.37	1.40
1	A	202	LLP	P-OP2	-2.48	1.45	1.54
1	B	202	LLP	C4-C3	-2.32	1.37	1.40
1	A	202	LLP	C4-C5	-2.19	1.39	1.42
1	C	202	LLP	P-OP4	-2.06	1.53	1.60
1	D	202	LLP	P-OP2	-2.06	1.47	1.54
1	C	202	LLP	CB-CA	-2.05	1.51	1.53
1	D	202	LLP	C5'-C5	2.00	1.56	1.50
1	E	202	LLP	C3-C2	2.05	1.42	1.40
1	F	202	LLP	C3-C2	2.17	1.42	1.40
1	A	202	LLP	C6-N1	2.32	1.39	1.34
1	E	202	LLP	C6-N1	2.47	1.39	1.34
1	D	202	LLP	C6-N1	2.56	1.39	1.34
1	B	202	LLP	C3-C2	2.63	1.42	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	202	LLP	C3-C2	2.68	1.42	1.40
1	C	202	LLP	C3-C2	2.74	1.42	1.40
1	A	202	LLP	C3-C2	2.82	1.42	1.40
1	C	202	LLP	C2'-C2	2.83	1.56	1.50
1	E	202	LLP	C2'-C2	2.87	1.56	1.50
1	F	202	LLP	C6-N1	2.88	1.40	1.34
1	F	202	LLP	C2'-C2	2.89	1.56	1.50
1	D	202	LLP	C2'-C2	3.02	1.56	1.50
1	A	202	LLP	C2'-C2	3.06	1.56	1.50
1	C	202	LLP	C6-N1	3.14	1.41	1.34
1	B	202	LLP	C6-N1	3.16	1.41	1.34
1	B	202	LLP	C2'-C2	3.27	1.56	1.50
1	C	202	LLP	C4'-NZ	5.75	1.44	1.27
1	F	202	LLP	C4'-NZ	5.77	1.44	1.27
1	A	202	LLP	C4'-NZ	5.84	1.44	1.27
1	D	202	LLP	C4'-NZ	5.95	1.45	1.27
1	E	202	LLP	C4'-NZ	6.11	1.45	1.27
1	B	202	LLP	C4'-NZ	6.38	1.46	1.27
1	F	202	LLP	C4-C4'	9.07	1.62	1.46
1	C	202	LLP	C4-C4'	9.37	1.63	1.46
1	D	202	LLP	C4-C4'	9.37	1.63	1.46
1	A	202	LLP	C4-C4'	9.49	1.63	1.46
1	E	202	LLP	C4-C4'	9.62	1.63	1.46
1	B	202	LLP	C4-C4'	9.75	1.63	1.46

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	202	LLP	C4-C4'-NZ	-3.37	106.30	125.06
1	F	202	LLP	C4-C4'-NZ	-3.19	107.29	125.06
1	A	202	LLP	C5'-C5-C6	-3.14	113.34	119.28
1	A	202	LLP	C3-C4-C4'	-3.10	116.15	120.16
1	E	202	LLP	C3-C4-C4'	-2.72	116.64	120.16
1	D	202	LLP	C4-C4'-NZ	-2.65	110.32	125.06
1	D	202	LLP	C3-C4-C4'	-2.59	116.80	120.16
1	E	202	LLP	C4-C4'-NZ	-2.48	111.27	125.06
1	C	202	LLP	C4-C4'-NZ	-2.47	111.31	125.06
1	C	202	LLP	C5-C6-N1	-2.42	119.66	123.86
1	B	202	LLP	C4-C4'-NZ	-2.35	111.96	125.06
1	C	202	LLP	O-C-CA	-2.26	119.60	125.49
1	F	202	LLP	C5-C6-N1	-2.13	120.17	123.86
1	E	202	LLP	C5-C6-N1	-2.06	120.28	123.86

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	202	LLP	O-C-CA	-2.06	120.13	125.49
1	F	202	LLP	O-C-CA	-2.06	120.13	125.49
1	F	202	LLP	C2'-C2-C3	-2.00	118.62	121.04
1	B	202	LLP	OP4-C5'-C5	2.19	112.61	108.99
1	D	202	LLP	C5-C4-C4'	2.21	124.69	121.52
1	F	202	LLP	OP4-C5'-C5	2.52	113.16	108.99
1	E	202	LLP	C5-C4-C4'	2.64	125.32	121.52
1	C	202	LLP	OP4-C5'-C5	3.09	114.11	108.99
1	A	202	LLP	C5-C4-C4'	3.20	126.12	121.52
1	A	202	LLP	C5'-C5-C4	3.24	126.92	121.47
1	A	202	LLP	OP4-C5'-C5	4.00	115.60	108.99

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	202	LLP	1	0
1	C	202	LLP	3	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

6 ligands are 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	SO4	B	401	-	4,4,4	0.32	0	6,6,6	0.24	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	T46	B	402	-	29,37,37	3.71	15 (51%)	39,57,57	2.03	10 (25%)
2	SO4	C	401	-	4,4,4	0.50	0	6,6,6	0.16	0
4	TYD	D	500	-	19,26,26	2.87	7 (36%)	27,40,40	2.10	6 (22%)
4	TYD	E	500	-	19,26,26	2.85	7 (36%)	27,40,40	2.03	5 (18%)
4	TYD	F	500	-	19,26,26	2.98	5 (26%)	27,40,40	2.03	6 (22%)

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	SO4	B	401	-	-	0/0/0/0	0/0/0/0
3	T46	B	402	-	-	0/17/53/53	0/3/3/3
2	SO4	C	401	-	-	0/0/0/0	0/0/0/0
4	TYD	D	500	-	-	0/12/28/28	0/2/2/2
4	TYD	E	500	-	-	0/12/28/28	0/2/2/2
4	TYD	F	500	-	-	0/12/28/28	0/2/2/2

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	402	T46	C3,-C4,	-8.68	1.28	1.53
3	B	402	T46	C2,-C1,	-5.91	1.35	1.52
3	B	402	T46	C2-C3	-4.12	1.48	1.53
4	F	500	TYD	O4-C4	-2.66	1.18	1.24
3	B	402	T46	O41-C41	-2.57	1.18	1.24
4	E	500	TYD	O4-C4	-2.50	1.18	1.24
4	E	500	TYD	PB-O3B	-2.26	1.46	1.54
4	D	500	TYD	PB-O2B	-2.20	1.46	1.54
4	E	500	TYD	C2'-C3'	-2.16	1.47	1.52
4	F	500	TYD	PB-O2B	-2.14	1.47	1.54
4	E	500	TYD	PB-O2B	-2.09	1.47	1.54
4	D	500	TYD	PB-O3B	-2.08	1.47	1.54
4	D	500	TYD	PA-O2A	-2.08	1.46	1.54
3	B	402	T46	O4-C4	-2.05	1.18	1.21
4	D	500	TYD	O4-C4	-2.02	1.19	1.24
3	B	402	T46	O3-C3	2.11	1.46	1.42
3	B	402	T46	O5-C1	2.29	1.47	1.41
3	B	402	T46	C5-C4	2.39	1.56	1.51
4	E	500	TYD	O4'-C1'	2.46	1.48	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	F	500	TYD	O4'-C1'	2.59	1.48	1.42
3	B	402	T46	C5A-C51	2.70	1.56	1.51
4	D	500	TYD	O4'-C1'	2.73	1.48	1.42
3	B	402	T46	C61-C51	3.10	1.48	1.40
3	B	402	T46	O4,-C1,	3.85	1.51	1.42
3	B	402	T46	C2,-C3,	4.32	1.64	1.52
4	F	500	TYD	C4-N3	4.86	1.42	1.33
4	E	500	TYD	C4-N3	4.90	1.42	1.33
4	D	500	TYD	C4-N3	5.18	1.42	1.33
3	B	402	T46	C41-N31	7.13	1.46	1.33
3	B	402	T46	O4,-C4,	7.83	1.63	1.45
3	B	402	T46	C61-N11	8.14	1.47	1.35
4	E	500	TYD	C6-N1	9.35	1.49	1.35
4	D	500	TYD	C6-N1	9.45	1.49	1.35
4	F	500	TYD	C6-N1	10.19	1.50	1.35

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	500	TYD	C5-C4-N3	-5.35	119.19	125.14
4	F	500	TYD	C5-C4-N3	-4.57	120.05	125.14
4	E	500	TYD	C5-C4-N3	-4.52	120.10	125.14
3	B	402	T46	C51-C41-N31	-4.03	120.64	125.14
3	B	402	T46	O5-C1-O1	-3.00	107.41	111.36
4	E	500	TYD	C2'-C1'-N1	-2.86	107.19	114.16
4	F	500	TYD	O2B-PB-O1B	-2.56	102.32	110.58
3	B	402	T46	O1-C1-C2	-2.35	104.00	108.39
3	B	402	T46	C2,-C1,-N11	-2.20	108.82	114.16
3	B	402	T46	P2-OPP-P	-2.16	126.66	132.73
4	D	500	TYD	PA-O3A-PB	-2.16	125.44	132.67
4	D	500	TYD	O2B-PB-O1B	-2.04	104.00	110.58
4	E	500	TYD	O2B-PB-O3A	2.13	114.74	105.09
4	F	500	TYD	O3B-PB-O3A	2.17	114.95	105.09
3	B	402	T46	C3,-C2,-C1,	2.24	107.78	102.40
4	F	500	TYD	C2'-C3'-C4'	2.43	107.82	102.77
4	D	500	TYD	O3A-PA-O5'	2.50	109.58	102.94
3	B	402	T46	C2,-C3,-C4,	2.55	108.06	102.77
3	B	402	T46	OPP-P2-O1	2.64	111.24	103.63
4	E	500	TYD	C2'-C3'-C4'	2.69	108.36	102.77
4	D	500	TYD	C2'-C3'-C4'	3.01	109.02	102.77
4	F	500	TYD	O3A-PA-O5'	3.13	111.24	102.94
3	B	402	T46	O4,-C1,-N11	3.25	113.35	107.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	E	500	TYD	C4-N3-C2	5.74	120.21	115.25
4	D	500	TYD	C4-N3-C2	5.81	120.27	115.25
4	F	500	TYD	C4-N3-C2	6.19	120.59	115.25
3	B	402	T46	C41-N31-C21	8.50	122.59	115.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	402	T46	2	0
4	E	500	TYD	2	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	368/404 (91%)	-0.20	8 (2%) 65 69	32, 47, 76, 140	0
1	B	368/404 (91%)	-0.36	0 100 100	29, 47, 69, 88	0
1	C	368/404 (91%)	-0.05	2 (0%) 91 92	38, 57, 88, 119	0
1	D	368/404 (91%)	-0.26	3 (0%) 87 89	35, 50, 79, 111	0
1	E	368/404 (91%)	-0.16	6 (1%) 74 77	38, 59, 87, 125	0
1	F	368/404 (91%)	0.30	22 (5%) 25 28	40, 73, 107, 127	0
All	All	2208/2424 (91%)	-0.12	41 (1%) 70 73	29, 54, 92, 140	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	235	SER	4.8
1	A	236	ASP	4.6
1	F	127	LEU	4.0
1	F	77	GLY	3.9
1	F	128	ALA	3.9
1	F	162	ALA	3.8
1	F	380	ILE	3.7
1	A	239	THR	3.4
1	F	157	VAL	3.4
1	A	240	GLY	3.2
1	F	384	TRP	3.2
1	F	159	ARG	3.2
1	F	291	GLY	3.1
1	F	165	ASP	3.1
1	D	78	ALA	3.0
1	A	233	ARG	3.0
1	E	234	ALA	3.0
1	C	236	ASP	3.0
1	E	109	ARG	2.9

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Mol	Chain	Res	Type	RSRZ
1	F	167	ARG	2.9
1	F	169	LEU	2.9
1	F	166	ARG	2.8
1	F	78	ALA	2.7
1	A	234	ALA	2.5
1	D	325	ALA	2.5
1	F	290	ASP	2.5
1	F	170	VAL	2.4
1	F	160	ALA	2.4
1	F	76	PRO	2.4
1	F	131	ILE	2.4
1	A	61	CYS	2.3
1	E	165	ASP	2.3
1	D	86	PRO	2.2
1	F	158	ASP	2.1
1	E	163	GLY	2.1
1	E	235	SER	2.1
1	A	237	ARG	2.1
1	C	61	CYS	2.1
1	E	166	ARG	2.1
1	F	304	ALA	2.1
1	F	383	GLY	2.0

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

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
1	LLP	E	202	24/25	0.97	0.21	-	40,53,61,64	0
1	LLP	C	202	24/25	0.97	0.15	-	37,51,63,66	0
1	LLP	A	202	24/25	0.97	0.19	-	31,44,49,54	0
1	LLP	F	202	24/25	0.97	0.20	-	51,65,78,80	0
1	LLP	B	202	24/25	0.97	0.17	-	27,47,65,72	0
1	LLP	D	202	24/25	0.97	0.17	-	41,50,62,69	0

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
2	SO4	C	401	5/5	0.93	0.35	5.76	127,130,131,133	0
3	T46	B	402	35/35	0.88	0.22	2.94	50,64,88,95	35
4	TYD	F	500	25/25	0.85	0.21	1.23	47,92,103,105	25
4	TYD	E	500	25/25	0.87	0.20	1.05	58,85,101,107	25
4	TYD	D	500	25/25	0.91	0.17	0.60	52,70,101,108	0
2	SO4	B	401	5/5	0.97	0.13	-	73,76,80,90	0

6.5 Other polymers [i](#)

There are no such residues in this entry.