



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 15, 2017 – 02:34 am GMT

PDB ID : 4ACB
Title : CRYSTAL STRUCTURE OF TRANSLATION ELONGATION FACTOR
SELB FROM METHANOCOCCUS MARIPALUDIS IN COMPLEX WITH
THE GTP ANALOGUE GPPNHP
Authors : Leibundgut, M.; Frick, C.; Thanbichler, M.; Boeck, A.; Ban, N.
Deposited on : 2011-12-14
Resolution : 3.34 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 : trunk28620
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) : recalc28949

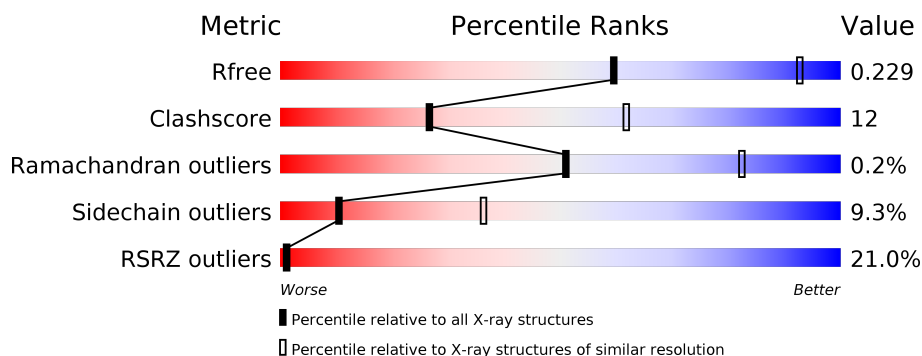
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.34 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	1167 (3.40-3.28)
Clashscore	112137	1239 (3.40-3.28)
Ramachandran outliers	110173	1219 (3.40-3.28)
Sidechain outliers	110143	1218 (3.40-3.28)
RSRZ outliers	101464	1176 (3.40-3.28)

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	482	
1	B	482	
1	C	482	
1	D	482	

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
1	CMH	D	340	-	-	X	-
5	SO4	B	1472	-	-	-	X

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 14623 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 TRANSLATION ELONGATION FACTOR SELB.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	452	Total	C	Hg	N	O	S	0	0	0
			3506	2240	4	597	651	14			
1	B	456	Total	C	Hg	N	O	S	0	0	0
			3533	2257	4	603	655	14			
1	C	471	Total	C	Hg	N	O	S	0	0	0
			3651	2327	4	627	679	14			
1	D	467	Total	C	Hg	N	O	S	0	0	0
			3615	2305	4	618	675	13			

There are 56 discrepancies between the modelled and reference sequences:

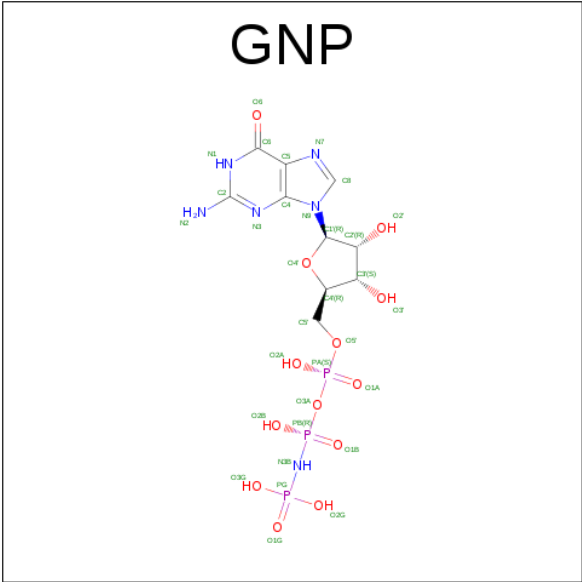
Chain	Residue	Modelled	Actual	Comment	Reference
A	-13	MET	-	EXPRESSION TAG	UNP Q8J307
A	-12	HIS	-	EXPRESSION TAG	UNP Q8J307
A	-11	HIS	-	EXPRESSION TAG	UNP Q8J307
A	-10	HIS	-	EXPRESSION TAG	UNP Q8J307
A	-9	HIS	-	EXPRESSION TAG	UNP Q8J307
A	-8	HIS	-	EXPRESSION TAG	UNP Q8J307
A	-7	HIS	-	EXPRESSION TAG	UNP Q8J307
A	-6	SER	-	EXPRESSION TAG	UNP Q8J307
A	-5	ILE	-	EXPRESSION TAG	UNP Q8J307
A	-4	GLU	-	EXPRESSION TAG	UNP Q8J307
A	-3	GLY	-	EXPRESSION TAG	UNP Q8J307
A	-2	ARG	-	EXPRESSION TAG	UNP Q8J307
A	-1	PRO	-	EXPRESSION TAG	UNP Q8J307
A	0	HIS	-	EXPRESSION TAG	UNP Q8J307
B	-13	MET	-	EXPRESSION TAG	UNP Q8J307
B	-12	HIS	-	EXPRESSION TAG	UNP Q8J307
B	-11	HIS	-	EXPRESSION TAG	UNP Q8J307
B	-10	HIS	-	EXPRESSION TAG	UNP Q8J307
B	-9	HIS	-	EXPRESSION TAG	UNP Q8J307
B	-8	HIS	-	EXPRESSION TAG	UNP Q8J307
B	-7	HIS	-	EXPRESSION TAG	UNP Q8J307

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-6	SER	-	EXPRESSION TAG	UNP Q8J307
B	-5	ILE	-	EXPRESSION TAG	UNP Q8J307
B	-4	GLU	-	EXPRESSION TAG	UNP Q8J307
B	-3	GLY	-	EXPRESSION TAG	UNP Q8J307
B	-2	ARG	-	EXPRESSION TAG	UNP Q8J307
B	-1	PRO	-	EXPRESSION TAG	UNP Q8J307
B	0	HIS	-	EXPRESSION TAG	UNP Q8J307
C	-13	MET	-	EXPRESSION TAG	UNP Q8J307
C	-12	HIS	-	EXPRESSION TAG	UNP Q8J307
C	-11	HIS	-	EXPRESSION TAG	UNP Q8J307
C	-10	HIS	-	EXPRESSION TAG	UNP Q8J307
C	-9	HIS	-	EXPRESSION TAG	UNP Q8J307
C	-8	HIS	-	EXPRESSION TAG	UNP Q8J307
C	-7	HIS	-	EXPRESSION TAG	UNP Q8J307
C	-6	SER	-	EXPRESSION TAG	UNP Q8J307
C	-5	ILE	-	EXPRESSION TAG	UNP Q8J307
C	-4	GLU	-	EXPRESSION TAG	UNP Q8J307
C	-3	GLY	-	EXPRESSION TAG	UNP Q8J307
C	-2	ARG	-	EXPRESSION TAG	UNP Q8J307
C	-1	PRO	-	EXPRESSION TAG	UNP Q8J307
C	0	HIS	-	EXPRESSION TAG	UNP Q8J307
D	-13	MET	-	EXPRESSION TAG	UNP Q8J307
D	-12	HIS	-	EXPRESSION TAG	UNP Q8J307
D	-11	HIS	-	EXPRESSION TAG	UNP Q8J307
D	-10	HIS	-	EXPRESSION TAG	UNP Q8J307
D	-9	HIS	-	EXPRESSION TAG	UNP Q8J307
D	-8	HIS	-	EXPRESSION TAG	UNP Q8J307
D	-7	HIS	-	EXPRESSION TAG	UNP Q8J307
D	-6	SER	-	EXPRESSION TAG	UNP Q8J307
D	-5	ILE	-	EXPRESSION TAG	UNP Q8J307
D	-4	GLU	-	EXPRESSION TAG	UNP Q8J307
D	-3	GLY	-	EXPRESSION TAG	UNP Q8J307
D	-2	ARG	-	EXPRESSION TAG	UNP Q8J307
D	-1	PRO	-	EXPRESSION TAG	UNP Q8J307
D	0	HIS	-	EXPRESSION TAG	UNP Q8J307

- Molecule 2 is PHOSPHOAMINOPHOSPHONIC ACID-GUANYLATE ESTER (three-letter code: GNP) (formula: C₁₀H₁₇N₆O₁₃P₃).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			32	10	6	13	3		

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

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

- Molecule 4 is GUANOSINE-5'-DIPHOSPHATE (three-letter code: GDP) (formula: C₁₀H₁₅N₅O₁₁P₂).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	B	1	Total 28	C 10	N 5	O 11	P 2	0	0

- Molecule 5 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



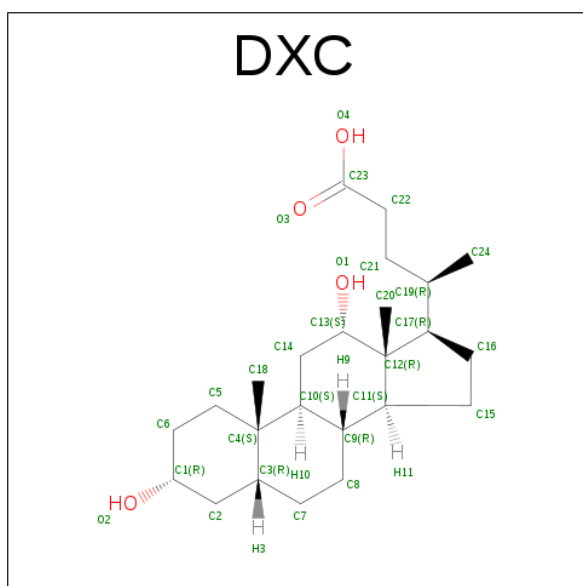
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total 5	O 4	S 1	0	0
5	B	1	Total 5	O 4	S 1	0	0
5	C	1	Total 5	O 4	S 1	0	0

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

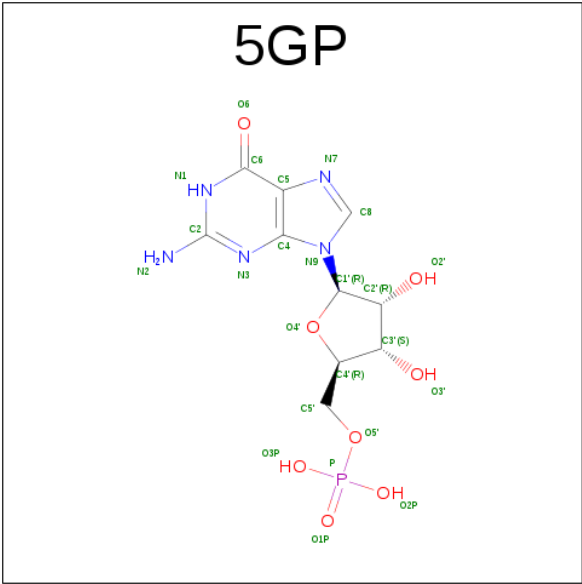
- Molecule 6 is (3ALPHA,5BETA,12ALPHA)-3,12-DIHYDROXYCHOLAN-24-OIC ACID (three-letter code: DXC) (formula: C₂₄H₄₀O₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	B	1	Total	C	O	0	0
			28	24	4		
6	C	1	Total	C	O	0	0
			28	24	4		
6	C	1	Total	C	O	0	0
			28	24	4		
6	C	1	Total	C	O	0	0
			28	24	4		
6	C	1	Total	C	O	0	0
			28	24	4		
6	C	1	Total	C	O	0	0
			28	24	4		

- Molecule 7 is GUANOSINE-5'-MONOPHOSPHATE (three-letter code: 5GP) (formula:

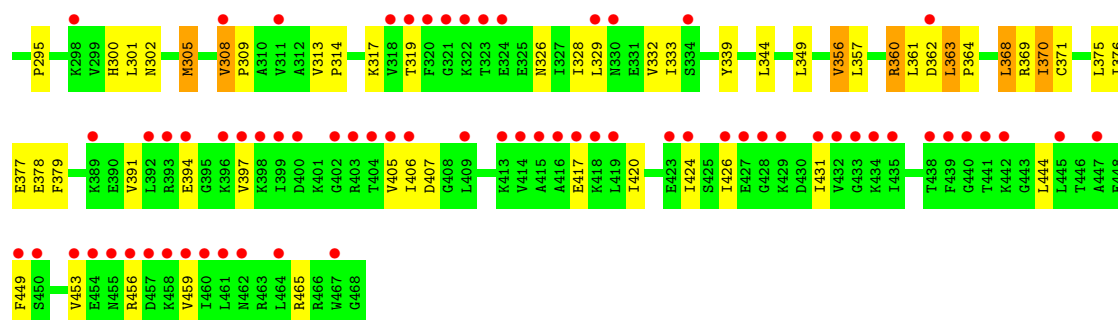
C₁₀H₁₄N₅O₈P).



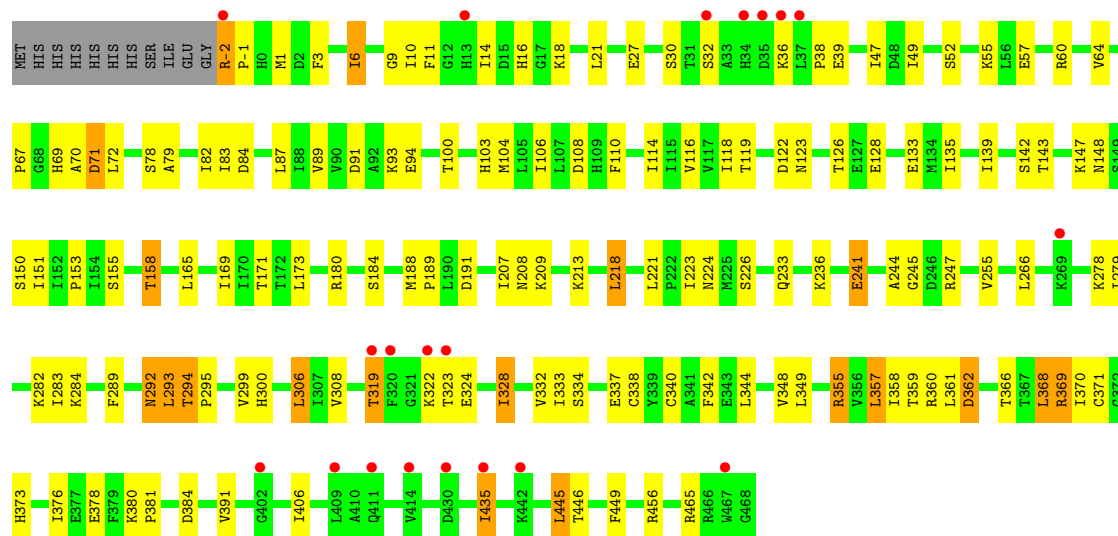
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	B	1	Total	C	N	O	P	0	0
			24	10	5	8	1		

- Molecule 8 is water.

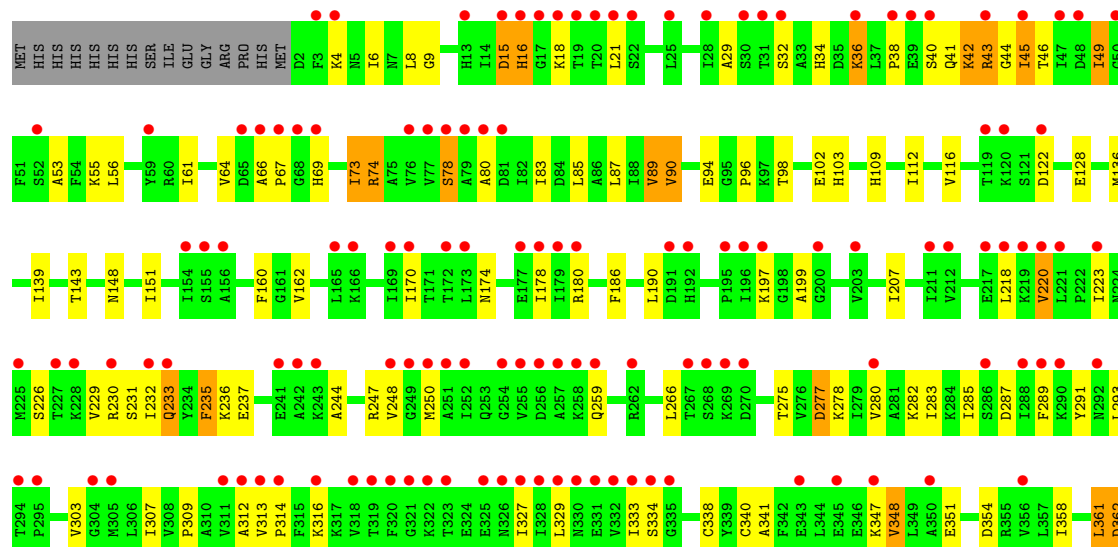
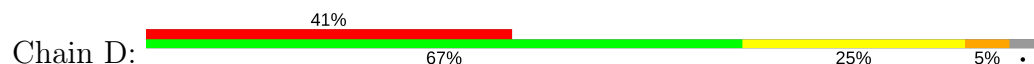
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	2	Total	O	0
			2	2	
8	B	4	Total	O	0
			4	4	

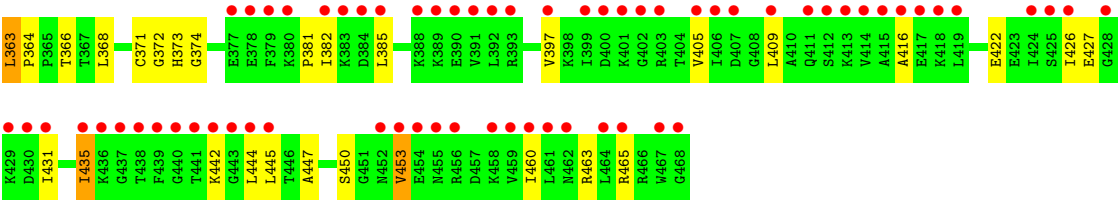


• Molecule 1: TRANSLATION ELONGATION FACTOR SELB



• Molecule 1: TRANSLATION ELONGATION FACTOR SELB





4 Data and refinement statistics

Property	Value	Source
Space group	P 31 1 2	Depositor
Cell constants a, b, c, α , β , γ	146.63Å 146.63Å 297.22Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.94 – 3.34 34.48 – 3.34	Depositor EDS
% Data completeness (in resolution range)	82.8 (19.94-3.34) 88.5 (34.48-3.34)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.66 (at 3.32Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE: 1.8_1069)	Depositor
R, R_{free}	0.179 , 0.223 0.186 , 0.229	Depositor DCC
R_{free} test set	1989 reflections (4.19%)	DCC
Wilson B-factor (Å ²)	110.2	Xtriage
Anisotropy	0.481	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 153.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.043 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	14623	wwPDB-VP
Average B, all atoms (Å ²)	198.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.99% 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, 0.333 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: GDP, GNP, CMH, MG, SO4, 5GP, DXC

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	0.31	0/3515	0.55	0/4727
1	B	0.35	0/3541	0.57	0/4760
1	C	0.40	0/3664	0.63	0/4929
1	D	0.28	0/3626	0.53	0/4878
All	All	0.34	0/14346	0.57	0/19294

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	3506	0	3687	73	0
1	B	3533	0	3726	84	0
1	C	3651	0	3838	82	0
1	D	3615	0	3801	107	0
2	A	32	0	13	2	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	B	28	0	12	2	0
5	B	10	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	C	20	0	0	2	0
6	B	28	0	39	4	0
6	C	168	0	234	15	0
7	B	24	0	12	2	0
8	A	2	0	0	0	0
8	B	4	0	0	0	0
All	All	14623	0	15362	349	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 12.

The worst 5 of 349 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:6:ILE:HD12	1:C:173:LEU:HD13	1.54	0.90
1:C:338:CMH:HB3	1:C:340:CMH:CM	2.05	0.86
1:D:43:ARG:HD3	1:D:45:ILE:HD11	1.57	0.85
1:D:283:ILE:HG13	1:D:340:CMH:HB3	1.61	0.82
1:D:223:ILE:HG21	1:D:259:GLN:HB3	1.61	0.80

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	444/482 (92%)	420 (95%)	24 (5%)	0	100	100
1	B	448/482 (93%)	429 (96%)	19 (4%)	0	100	100
1	C	465/482 (96%)	435 (94%)	28 (6%)	2 (0%)	38	73
1	D	461/482 (96%)	435 (94%)	25 (5%)	1 (0%)	51	82
All	All	1818/1928 (94%)	1719 (95%)	96 (5%)	3 (0%)	51	82

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	42	LYS
1	C	57	GLU
1	C	47	ILE

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	386/412 (94%)	352 (91%)	34 (9%)	12	40
1	B	388/412 (94%)	354 (91%)	34 (9%)	12	40
1	C	402/412 (98%)	358 (89%)	44 (11%)	7	29
1	D	398/412 (97%)	364 (92%)	34 (8%)	12	42
All	All	1574/1648 (96%)	1428 (91%)	146 (9%)	10	36

5 of 146 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	394	GLU
1	C	171	THR
1	D	235	PHE
1	B	453	VAL
1	C	114	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 6 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	233	GLN
1	D	373	HIS
1	D	69	HIS
1	A	300	HIS
1	D	233	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

16 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	CMH	A	264	1	6,7,8	1.15	1 (16%)	2,7,9	1.10	0
1	CMH	A	338	1	6,7,8	0.83	0	2,7,9	1.25	0
1	CMH	A	340	1	6,7,8	0.99	1 (16%)	2,7,9	1.31	0
1	CMH	A	371	1	6,7,8	1.17	1 (16%)	2,7,9	1.36	0
1	CMH	B	264	1	6,7,8	1.35	1 (16%)	2,7,9	1.34	0
1	CMH	B	338	1	6,7,8	0.65	0	2,7,9	1.02	0
1	CMH	B	340	1	6,7,8	0.86	0	2,7,9	1.21	0
1	CMH	B	371	1	6,7,8	0.77	0	2,7,9	1.49	0
1	CMH	C	264	1	6,7,8	1.21	1 (16%)	2,7,9	1.16	0
1	CMH	C	338	1	6,7,8	0.66	0	2,7,9	1.07	0
1	CMH	C	340	1	6,7,8	1.02	0	2,7,9	1.15	0
1	CMH	C	371	1	6,7,8	0.65	0	2,7,9	1.11	0
1	CMH	D	264	1	6,7,8	0.87	0	2,7,9	1.06	0
1	CMH	D	338	1	6,7,8	1.15	1 (16%)	2,7,9	1.42	1 (50%)
1	CMH	D	340	1	6,7,8	1.55	1 (16%)	2,7,9	1.09	0
1	CMH	D	371	1	6,7,8	0.96	0	2,7,9	1.24	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
1	CMH	A	264	1	-	0/0/6/8	0/0/0/0
1	CMH	A	338	1	-	0/0/6/8	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	CMH	A	340	1	-	0/0/6/8	0/0/0/0
1	CMH	A	371	1	-	0/0/6/8	0/0/0/0
1	CMH	B	264	1	-	0/0/6/8	0/0/0/0
1	CMH	B	338	1	-	0/0/6/8	0/0/0/0
1	CMH	B	340	1	-	0/0/6/8	0/0/0/0
1	CMH	B	371	1	-	0/0/6/8	0/0/0/0
1	CMH	C	264	1	-	0/0/6/8	0/0/0/0
1	CMH	C	338	1	-	0/0/6/8	0/0/0/0
1	CMH	C	340	1	-	0/0/6/8	0/0/0/0
1	CMH	C	371	1	-	0/0/6/8	0/0/0/0
1	CMH	D	264	1	-	0/0/6/8	0/0/0/0
1	CMH	D	338	1	-	0/0/6/8	0/0/0/0
1	CMH	D	340	1	-	0/0/6/8	0/0/0/0
1	CMH	D	371	1	-	0/0/6/8	0/0/0/0

The worst 5 of 7 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	340	CMH	CA-C	2.04	1.52	1.50
1	D	338	CMH	CA-C	2.50	1.53	1.50
1	A	264	CMH	CA-C	2.53	1.53	1.50
1	A	371	CMH	CA-C	2.60	1.53	1.50
1	C	264	CMH	CA-C	2.62	1.53	1.50

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	338	CMH	O-C-CA	-2.01	119.46	125.02

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

9 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	338	CMH	1	0
1	A	340	CMH	2	0
1	B	371	CMH	2	0
1	C	338	CMH	1	0
1	C	340	CMH	2	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	C	371	CMH	3	0
1	D	338	CMH	2	0
1	D	340	CMH	4	0
1	D	371	CMH	2	0

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 18 ligands modelled in this entry, 2 are monoatomic - leaving 16 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	GNP	A	1469	3	27,34,34	2.68	7 (25%)	26,54,54	1.66	5 (19%)
4	GDP	B	1469	3	25,30,30	1.21	2 (8%)	26,47,47	2.04	6 (23%)
5	SO4	B	1471	-	4,4,4	0.20	0	6,6,6	0.08	0
5	SO4	B	1472	-	4,4,4	0.16	0	6,6,6	0.16	0
6	DXC	B	1473	-	28,31,31	1.32	3 (10%)	45,49,49	2.13	12 (26%)
7	5GP	B	1474	-	22,26,26	0.53	0	26,40,40	1.78	8 (30%)
5	SO4	C	1471	-	4,4,4	0.34	0	6,6,6	0.68	0
5	SO4	C	1472	-	4,4,4	0.31	0	6,6,6	0.19	0
5	SO4	C	1473	-	4,4,4	0.19	0	6,6,6	0.12	0
6	DXC	C	1475	-	28,31,31	1.60	5 (17%)	45,49,49	2.27	18 (40%)
6	DXC	C	1476	-	28,31,31	1.49	5 (17%)	45,49,49	1.87	13 (28%)
6	DXC	C	1477	-	28,31,31	1.60	6 (21%)	45,49,49	1.87	12 (26%)
6	DXC	C	1478	-	28,31,31	1.35	5 (17%)	45,49,49	1.89	15 (33%)
6	DXC	C	1479	-	28,31,31	1.54	7 (25%)	45,49,49	2.17	17 (37%)
6	DXC	C	1480	-	28,31,31	1.60	7 (25%)	45,49,49	2.10	15 (33%)
5	SO4	C	1481	-	4,4,4	0.22	0	6,6,6	0.12	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	GNP	A	1469	3	-	0/16/38/38	0/3/3/3
4	GDP	B	1469	3	-	0/12/32/32	0/3/3/3
5	SO4	B	1471	-	-	0/0/0/0	0/0/0/0
5	SO4	B	1472	-	-	0/0/0/0	0/0/0/0
6	DXC	B	1473	-	-	0/7/71/71	0/4/4/4
7	5GP	B	1474	-	-	0/6/26/26	0/3/3/3
5	SO4	C	1471	-	-	0/0/0/0	0/0/0/0
5	SO4	C	1472	-	-	0/0/0/0	0/0/0/0
5	SO4	C	1473	-	-	0/0/0/0	0/0/0/0
6	DXC	C	1475	-	-	0/7/71/71	0/4/4/4
6	DXC	C	1476	-	-	0/7/71/71	0/4/4/4
6	DXC	C	1477	-	-	0/7/71/71	0/4/4/4
6	DXC	C	1478	-	-	0/7/71/71	0/4/4/4
6	DXC	C	1479	-	-	0/7/71/71	0/4/4/4
6	DXC	C	1480	-	-	0/7/71/71	0/4/4/4
5	SO4	C	1481	-	-	0/0/0/0	0/0/0/0

The worst 5 of 47 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1469	GNP	C4-N9	-7.09	1.38	1.47
2	A	1469	GNP	C5-C6	-6.13	1.41	1.53
2	A	1469	GNP	PB-O3A	-4.87	1.53	1.59
6	C	1475	DXC	C12-C13	-4.55	1.47	1.54
6	C	1477	DXC	C12-C13	-3.80	1.48	1.54

The worst 5 of 121 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	C	1478	DXC	C10-C14-C13	-5.64	106.89	114.32
6	C	1475	DXC	C15-C11-C12	-5.30	98.29	103.57
6	C	1479	DXC	C18-C4-C3	-5.02	101.64	110.30
6	C	1475	DXC	C3-C2-C1	-4.99	105.55	112.87
6	C	1479	DXC	C6-C5-C4	-4.88	104.30	112.80

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

14 monomers are involved in 29 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1469	GNP	2	0
4	B	1469	GDP	2	0
5	B	1471	SO4	1	0
5	B	1472	SO4	1	0
6	B	1473	DXC	4	0
7	B	1474	5GP	2	0
5	C	1471	SO4	1	0
5	C	1473	SO4	1	0
6	C	1475	DXC	5	0
6	C	1476	DXC	2	0
6	C	1477	DXC	3	0
6	C	1478	DXC	2	0
6	C	1479	DXC	1	0
6	C	1480	DXC	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 [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	448/482 (92%)	1.05	97 (21%) 1 1	107, 205, 299, 372	0
1	B	452/482 (93%)	0.75	71 (15%) 2 2	80, 152, 347, 414	0
1	C	467/482 (96%)	0.24	20 (4%) 36 36	74, 123, 227, 308	0
1	D	463/482 (96%)	2.14	196 (42%) 0 0	140, 274, 369, 416	0
All	All	1830/1928 (94%)	1.04	384 (20%) 1 1	74, 188, 338, 416	0

The worst 5 of 384 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	453	VAL	12.7
1	D	401	LYS	11.9
1	D	78	SER	10.9
1	B	404	THR	10.9
1	D	438	THR	10.8

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	CMH	D	264	8/9	0.92	0.26	-	241,255,280,281	2
1	CMH	C	264	8/9	0.99	0.16	-	70,108,137,210	0
1	CMH	B	338	8/9	0.92	0.28	-	151,200,232,276	2
1	CMH	A	371	8/9	0.88	0.36	-	168,213,225,228	2
1	CMH	C	371	8/9	0.99	0.25	-	78,109,123,156	2

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
1	CMH	C	340	8/9	0.98	0.24	-	69,99,124,131	2
1	CMH	B	340	8/9	0.98	0.25	-	160,194,264,332	0
1	CMH	A	264	8/9	0.97	0.17	-	171,215,224,235	2
1	CMH	D	340	8/9	0.90	0.21	-	207,306,312,317	2
1	CMH	A	340	8/9	0.98	0.19	-	169,213,226,240	2
1	CMH	B	264	8/9	1.00	0.17	-	73,89,99,112	2
1	CMH	D	371	8/9	0.98	0.19	-	145,181,227,234	2
1	CMH	D	338	8/9	0.81	0.11	-	180,242,266,268	2
1	CMH	C	338	8/9	0.96	0.18	-	74,109,174,331	2
1	CMH	B	371	8/9	0.93	0.30	-	174,197,251,332	2
1	CMH	A	338	8/9	0.95	0.14	-	174,190,268,307	2

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
5	SO4	B	1472	5/5	0.81	0.89	9.91	139,226,336,344	0
7	5GP	B	1474	24/24	0.82	0.38	1.55	242,294,308,311	0
6	DXC	C	1479	28/28	0.93	0.29	1.17	82,104,174,196	28
6	DXC	C	1477	28/28	0.94	0.31	0.98	91,111,162,182	0
6	DXC	C	1476	28/28	0.92	0.33	0.64	159,168,217,253	0
6	DXC	C	1478	28/28	0.96	0.24	0.57	93,118,158,196	0
5	SO4	C	1471	5/5	0.97	0.21	0.44	101,149,159,220	0
6	DXC	B	1473	28/28	0.95	0.28	0.28	68,97,134,143	0
6	DXC	C	1475	28/28	0.96	0.27	0.28	61,95,139,146	0
4	GDP	B	1469	28/28	0.92	0.21	-0.12	126,197,212,235	0
6	DXC	C	1480	28/28	0.97	0.23	-0.46	63,85,106,127	0
2	GNP	A	1469	32/32	0.97	0.12	-1.01	120,165,202,249	0
5	SO4	B	1471	5/5	0.78	0.18	-	205,213,271,283	0
3	MG	B	1470	1/1	0.98	0.20	-	342,342,342,342	0
5	SO4	C	1473	5/5	0.83	0.14	-	164,166,220,230	5

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
5	SO4	C	1472	5/5	0.93	0.19	-	133,154,189,210	0
3	MG	A	1470	1/1	0.98	0.07	-	244,244,244,244	0
5	SO4	C	1481	5/5	0.90	1.65	-	172,210,221,235	5

6.5 Other polymers [i](#)

There are no such residues in this entry.