



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 28, 2026 – 07:21 PM UTC

PDB ID : 1FZW / pdb_00001fzw
Title : THE STRUCTURAL BASIS OF THE CATALYTIC MECHANISM AND
REGULATION OF GLUCOSE-1-PHOSPHATE THYMIDYLYLTRANSFERASE (RMLA). APO ENZYME.
Authors : Blankenfeldt, W.; Asuncion, M.; Lam, J.S.; Naismith, J.H.
Deposited on : 2000-10-04
Resolution : 1.90 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

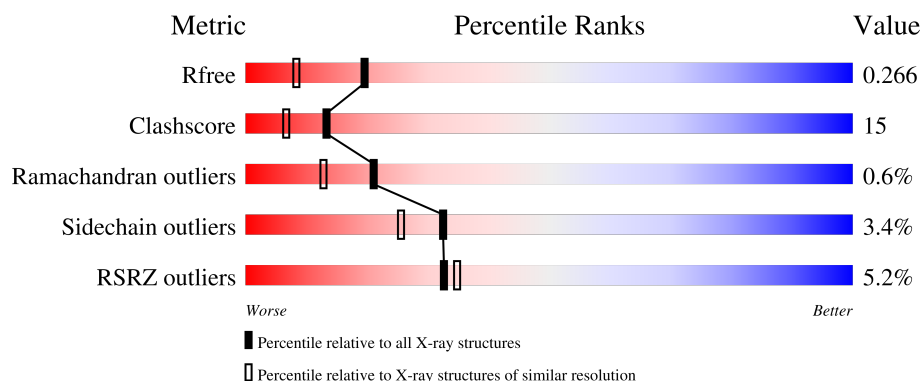
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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}	180053	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. 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	293	4% 63% 32% .
1	B	293	4% 63% 31% 6%
1	C	293	5% 62% 32% 5%
1	D	293	6% 66% 28% 5% .
1	E	293	9% 65% 28% 6%

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Mol	Chain	Length	Quality of chain
1	F	293	
1	G	293	
1	H	293	

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	5000	-	-	X	-
2	SO4	H	6101	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 21094 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 GLUCOSE-1-PHOSPHATE THYMIDYLYLTRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	292	Total	C	N	O	S	0	2	0
			2302	1472	388	437	5			
1	B	292	Total	C	N	O	S	0	5	0
			2322	1484	391	442	5			
1	C	293	Total	C	N	O	S	0	6	0
			2340	1494	394	446	6			
1	D	292	Total	C	N	O	S	0	1	0
			2294	1467	387	436	4			
1	E	292	Total	C	N	O	S	0	7	0
			2339	1491	396	447	5			
1	F	293	Total	C	N	O	S	0	3	0
			2316	1480	389	441	6			
1	G	292	Total	C	N	O	S	0	4	0
			2318	1480	392	441	5			
1	H	292	Total	C	N	O	S	0	0	0
			2285	1462	385	434	4			

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	F	1	Total	O	S	0	0
			5	4	1		
2	G	1	Total	O	S	0	0
			5	4	1		
2	G	1	Total	O	S	0	0
			5	4	1		
2	G	1	Total	O	S	0	0
			5	4	1		
2	H	1	Total	O	S	0	0
			5	4	1		
2	H	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	325	Total	O	0	0
			325	325		
3	B	335	Total	O	0	0
			335	335		
3	C	286	Total	O	0	0
			286	286		
3	D	259	Total	O	0	0
			259	259		
3	E	282	Total	O	0	0
			282	282		
3	F	321	Total	O	0	0
			321	321		
3	G	367	Total	O	0	0
			367	367		
3	H	303	Total	O	0	0
			303	303		

3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

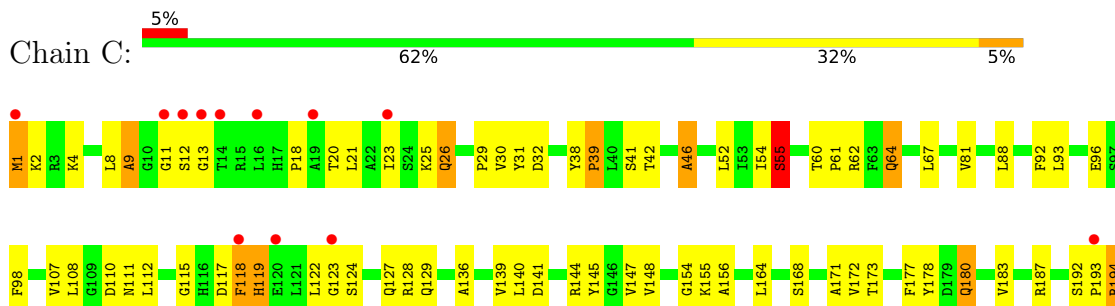
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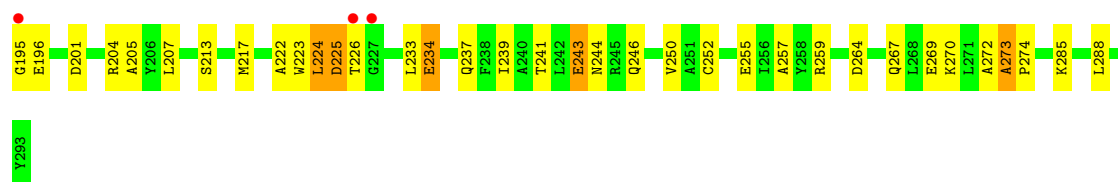


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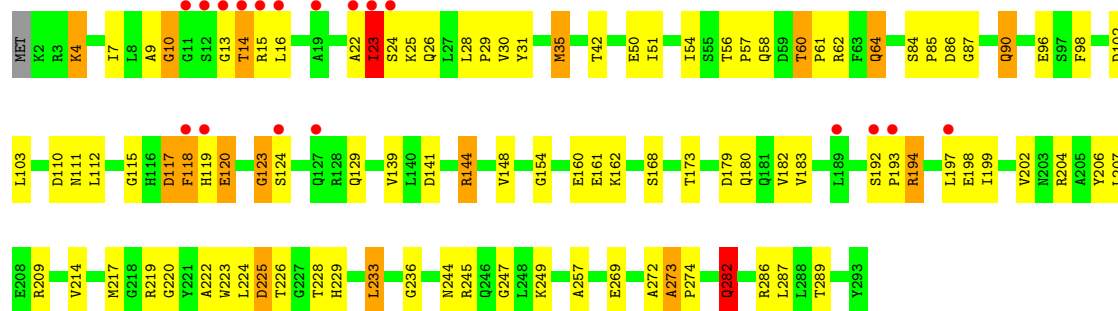


• Molecule 1: GLUCOSE-1-PHOSPHATE THYMIDYLYLTRANSFERASE

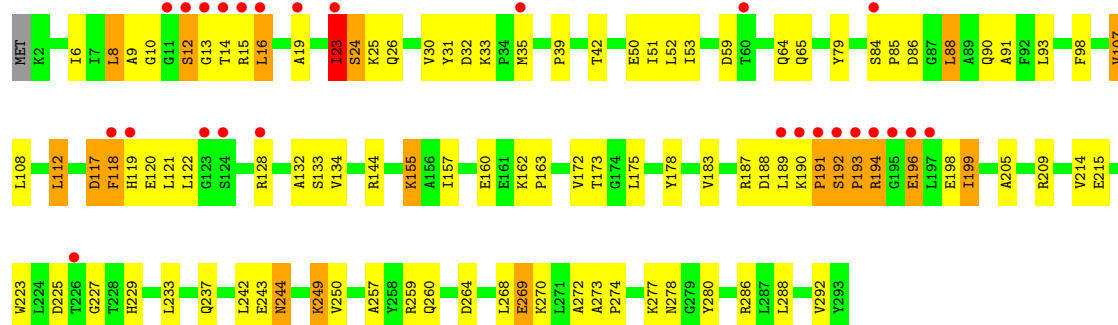




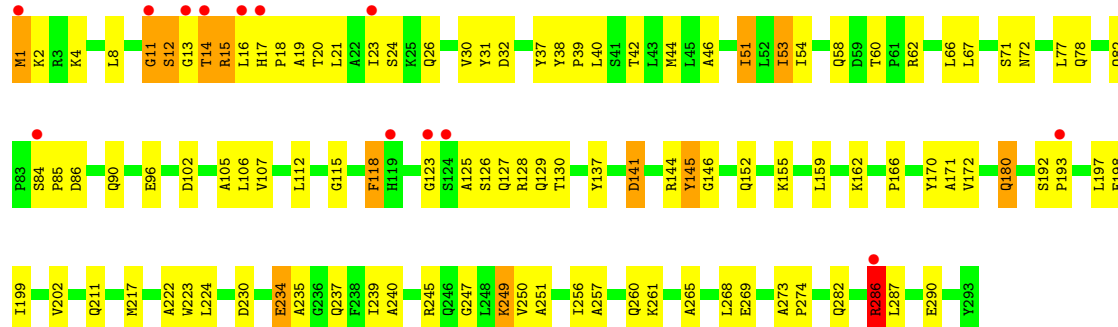
● Molecule 1: GLUCOSE-1-PHOSPHATE THYMIDYLYLTRANSFERASE



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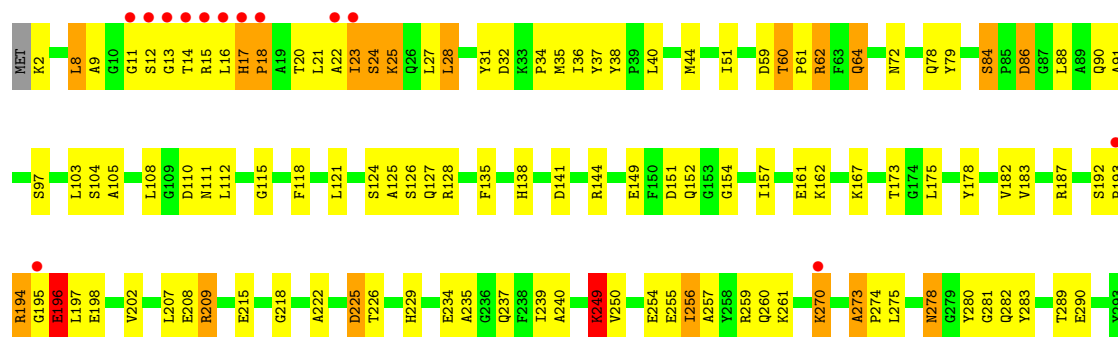


● Molecule 1: GLUCOSE-1-PHOSPHATE THYMIDYLYLTRANSFERASE



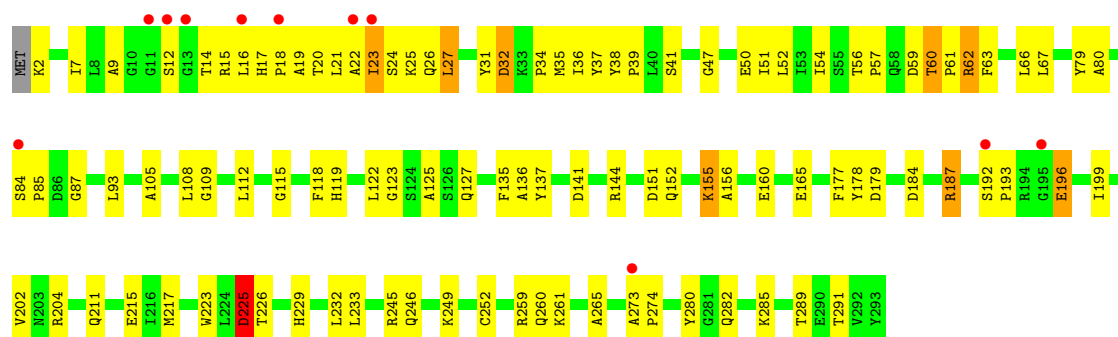
● Molecule 1: GLUCOSE-1-PHOSPHATE THYMIDYLYLTRANSFERASE

Chain G: 4% 59% 33% 6%



• Molecule 1: GLUCOSE-1-PHOSPHATE THYMIDYLYLTRANSFERASE

Chain H: 4% 64% 32%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	71.66Å 73.66Å 134.47Å 89.98° 80.91° 80.91°	Depositor
Resolution (Å)	73.00 – 1.90 73.00 – 1.90	Depositor EDS
% Data completeness (in resolution range)	95.8 (73.00-1.90) 95.1 (73.00-1.90)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.94 (at 1.89Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.176 , 0.251 0.196 , 0.266	Depositor DCC
R_{free} test set	10141 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	21.7	Xtriage
Anisotropy	0.030	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 56.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	21094	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.49% 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 ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: SO4

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	1.98	45/2352 (1.9%)	1.49	17/3190 (0.5%)
1	B	1.93	42/2372 (1.8%)	1.54	22/3217 (0.7%)
1	C	1.91	51/2390 (2.1%)	1.54	26/3239 (0.8%)
1	D	1.90	47/2344 (2.0%)	1.50	17/3180 (0.5%)
1	E	1.79	28/2389 (1.2%)	1.49	13/3240 (0.4%)
1	F	2.04	56/2366 (2.4%)	1.56	18/3208 (0.6%)
1	G	2.12	62/2368 (2.6%)	1.54	16/3212 (0.5%)
1	H	1.89	35/2335 (1.5%)	1.54	19/3168 (0.6%)
All	All	1.95	366/18916 (1.9%)	1.52	148/25654 (0.6%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	1
1	G	0	2
1	H	0	1
All	All	0	4

The worst 5 of 366 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	217	MET	SD-CE	-11.16	1.51	1.79
1	G	44	MET	SD-CE	-10.98	1.52	1.79
1	F	251	ALA	CA-CB	10.56	1.69	1.53
1	F	51	ILE	N-CA	-10.44	1.33	1.46
1	A	128	ARG	CZ-NH2	-10.41	1.20	1.33

The worst 5 of 148 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	118	PHE	N-CA-C	-11.85	98.44	111.36
1	D	118	PHE	N-CA-C	-11.79	98.15	111.71
1	C	154	GLY	N-CA-C	10.15	128.87	115.36
1	C	118	PHE	N-CA-C	-10.11	100.21	112.54
1	B	118	PHE	N-CA-C	-10.01	100.45	111.36

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	119	HIS	Sidechain
1	G	138	HIS	Sidechain
1	G	17	HIS	Sidechain
1	H	17	HIS	Sidechain

5.2 Too-close contacts ⓘ

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	2302	0	2287	89	0
1	B	2322	0	2305	91	0
1	C	2340	0	2324	68	0
1	D	2294	0	2279	67	0
1	E	2339	0	2316	94	0
1	F	2316	0	2301	69	0
1	G	2318	0	2300	61	1
1	H	2285	0	2272	58	0
2	A	10	0	0	0	0
2	B	10	0	0	1	0
2	C	15	0	0	2	0
2	D	15	0	0	0	0
2	E	15	0	0	0	0
2	F	10	0	0	0	0
2	G	15	0	0	1	0
2	H	10	0	0	2	0
3	A	325	0	0	35	1
3	B	335	0	0	24	0
3	C	286	0	0	14	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	259	0	0	17	0
3	E	282	0	0	33	0
3	F	321	0	0	22	4
3	G	367	0	0	22	3
3	H	303	0	0	15	1
All	All	21094	0	18384	550	5

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 15.

The worst 5 of 550 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:A:23:ILE:HG23	1:A:28:LEU:HD21	1.33	1.09
1:E:23:ILE:HG21	1:F:23:ILE:HG12	1.22	1.09
1:C:23:ILE:HG12	1:D:23:ILE:HG21	1.14	1.07
1:A:23:ILE:HB	1:B:23:ILE:HD12	1.11	1.06
1:A:23:ILE:CB	1:B:23:ILE:HD12	1.86	1.06

All (5) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:84:SER:OG	3:F:4852:HOH:O[1_655]	1.25	0.95
3:A:3776:HOH:O	3:F:4954:HOH:O[1_566]	1.99	0.21
3:F:4955:HOH:O	3:G:5148:HOH:O[1_455]	2.06	0.14
3:G:5209:HOH:O	3:H:6340:HOH:O[1_565]	2.09	0.11
3:F:4950:HOH:O	3:G:4974:HOH:O[1_455]	2.18	0.02

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	292/293 (100%)	284 (97%)	7 (2%)	1 (0%)	36	29
1	B	295/293 (101%)	285 (97%)	9 (3%)	1 (0%)	36	29
1	C	297/293 (101%)	290 (98%)	5 (2%)	2 (1%)	18	10
1	D	291/293 (99%)	282 (97%)	7 (2%)	2 (1%)	18	10
1	E	297/293 (101%)	286 (96%)	7 (2%)	4 (1%)	9	3
1	F	294/293 (100%)	285 (97%)	6 (2%)	3 (1%)	12	5
1	G	294/293 (100%)	287 (98%)	6 (2%)	1 (0%)	36	29
1	H	290/293 (99%)	284 (98%)	5 (2%)	1 (0%)	36	29
All	All	2350/2344 (100%)	2283 (97%)	52 (2%)	15 (1%)	21	13

5 of 15 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	14	THR
1	D	14	THR
1	E	31	TYR
1	F	15	ARG
1	F	31	TYR

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	241/240 (100%)	236 (98%)	5 (2%)	47	44
1	B	244/240 (102%)	235 (96%)	9 (4%)	30	22
1	C	246/240 (102%)	237 (96%)	9 (4%)	30	22
1	D	240/240 (100%)	232 (97%)	8 (3%)	33	26
1	E	246/240 (102%)	235 (96%)	11 (4%)	24	17
1	F	243/240 (101%)	235 (97%)	8 (3%)	33	26
1	G	243/240 (101%)	229 (94%)	14 (6%)	18	10
1	H	239/240 (100%)	235 (98%)	4 (2%)	53	52

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	1942/1920 (101%)	1874 (96%)	68 (4%)	32 24

5 of 68 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	G	194[A]	ARG
1	G	196	GLU
1	H	152	GLN
1	D	24	SER
1	D	23	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 47 such sidechains are listed below:

Mol	Chain	Res	Type
1	F	65	GLN
1	G	17	HIS
1	F	78	GLN
1	F	181	GLN
1	G	65	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

20 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	C	3600	-	4,4,4	1.04	0	6,6,6	0.79	0
2	SO4	C	5000	-	4,4,4	0.90	0	6,6,6	0.73	0
2	SO4	E	3607	-	4,4,4	0.80	0	6,6,6	0.74	0
2	SO4	H	6101	-	4,4,4	0.51	0	6,6,6	0.72	0
2	SO4	D	3800	-	4,4,4	0.32	0	6,6,6	1.17	0
2	SO4	C	3702	-	4,4,4	0.32	0	6,6,6	0.66	0
2	SO4	B	3603	-	4,4,4	0.87	0	6,6,6	0.65	0
2	SO4	D	3601	-	4,4,4	1.06	0	6,6,6	0.71	0
2	SO4	D	3703	-	4,4,4	0.52	0	6,6,6	0.61	0
2	SO4	G	3605	-	4,4,4	0.75	0	6,6,6	0.88	0
2	SO4	A	3602	-	4,4,4	0.70	0	6,6,6	0.63	0
2	SO4	H	3604	-	4,4,4	0.88	0	6,6,6	0.59	0
2	SO4	G	3704	-	4,4,4	0.50	0	6,6,6	1.04	0
2	SO4	F	3606	-	4,4,4	0.79	0	6,6,6	0.73	0
2	SO4	A	3700	-	4,4,4	0.28	0	6,6,6	0.60	0
2	SO4	B	3701	-	4,4,4	0.44	0	6,6,6	1.03	0
2	SO4	E	6100	-	4,4,4	0.75	0	6,6,6	0.78	0
2	SO4	F	4700	-	4,4,4	0.51	0	6,6,6	0.84	0
2	SO4	E	5001	-	4,4,4	0.36	0	6,6,6	0.55	0
2	SO4	G	4900	-	4,4,4	0.34	0	6,6,6	0.85	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	5000	SO4	2	0
2	H	6101	SO4	2	0
2	B	3701	SO4	1	0
2	G	4900	SO4	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	292/293 (99%)	0.15	13 (4%)	38	40	8, 20, 41, 60	2 (0%)
1	B	292/293 (99%)	0.06	13 (4%)	38	40	7, 19, 44, 63	5 (1%)
1	C	293/293 (100%)	0.18	15 (5%)	33	36	7, 20, 43, 65	6 (2%)
1	D	292/293 (99%)	0.29	18 (6%)	26	28	7, 21, 47, 72	1 (0%)
1	E	292/293 (99%)	0.44	26 (8%)	15	16	7, 22, 45, 67	7 (2%)
1	F	293/293 (100%)	0.03	13 (4%)	39	41	5, 16, 43, 69	3 (1%)
1	G	292/293 (99%)	-0.02	13 (4%)	38	40	5, 16, 38, 66	4 (1%)
1	H	292/293 (99%)	0.17	11 (3%)	44	47	10, 20, 42, 52	0
All	All	2338/2344 (99%)	0.16	122 (5%)	33	35	5, 19, 43, 72	28 (1%)

The worst 5 of 122 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	13	GLY	8.5
1	F	13	GLY	8.0
1	A	13	GLY	6.5
1	B	13	GLY	6.3
1	B	16	LEU	6.2

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 oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	SO4	H	6101	5/5	0.79	0.11	50,58,59,65	0
2	SO4	C	5000	5/5	0.82	0.11	41,48,52,53	0
2	SO4	G	4900	5/5	0.86	0.10	59,60,61,63	0
2	SO4	D	3703	5/5	0.87	0.12	52,59,62,64	0
2	SO4	E	5001	5/5	0.87	0.10	45,46,50,50	0
2	SO4	F	4700	5/5	0.89	0.10	42,52,58,59	0
2	SO4	B	3701	5/5	0.91	0.10	55,57,58,59	0
2	SO4	E	6100	5/5	0.92	0.09	45,47,49,51	0
2	SO4	C	3702	5/5	0.92	0.09	53,58,58,59	0
2	SO4	G	3704	5/5	0.93	0.09	47,50,52,56	0
2	SO4	D	3800	5/5	0.93	0.10	48,52,56,56	0
2	SO4	A	3700	5/5	0.94	0.08	45,47,52,55	0
2	SO4	F	3606	5/5	0.98	0.07	22,28,31,32	0
2	SO4	C	3600	5/5	0.98	0.08	27,31,35,37	0
2	SO4	E	3607	5/5	0.98	0.06	25,27,31,32	0
2	SO4	D	3601	5/5	0.98	0.07	25,28,30,34	0
2	SO4	H	3604	5/5	0.98	0.07	29,30,30,31	0
2	SO4	A	3602	5/5	0.98	0.07	25,26,29,29	0
2	SO4	B	3603	5/5	0.99	0.06	28,30,32,33	0
2	SO4	G	3605	5/5	0.99	0.07	27,28,31,31	0

6.5 Other polymers [i](#)

There are no such residues in this entry.