



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 07:17 AM UTC

PDB ID : 5JNB / pdb_00005jnb
Title : structure of GLD-2/RNP-8 complex
Authors : Nakel, K.; Bonneau, F.; Basquin, C.; Eckmann, C.R.; Conti, E.
Deposited on : 2016-04-29
Resolution : 2.49 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

<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:

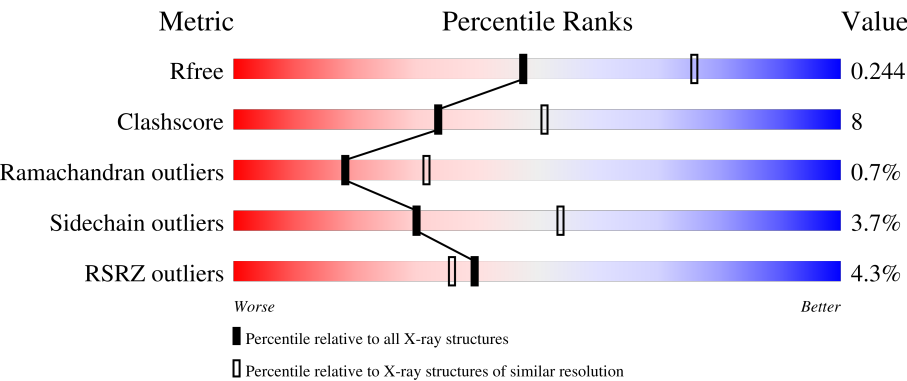
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry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

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.49 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	7589 (2.50-2.46)
Clashscore	190562	8295 (2.50-2.46)
Ramachandran outliers	187476	8164 (2.50-2.46)
Sidechain outliers	187428	8166 (2.50-2.46)
RSRZ outliers	180081	7593 (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 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	338	% 74% 22% ..
1	B	338	% 78% 17% ..
1	C	338	4% 72% 11% 14%
1	D	338	9% 64% 12% 22%
2	E	74	% 50% 11% 38%

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Mol	Chain	Length	Quality of chain
2	F	74	
2	G	74	
2	H	74	

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
5	EDO	A	1003	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 10776 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 Poly(A) RNA polymerase gld-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	329	Total	C	N	O	S	0	0	0
			2623	1669	451	489	14			
1	B	324	Total	C	N	O	S	0	0	0
			2590	1650	444	481	15			
1	C	289	Total	C	N	O	S	0	0	0
			2259	1440	380	425	14			
1	D	263	Total	C	N	O	S	0	0	0
			2004	1287	330	378	9			

There are 180 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	544	GLY	-	expression tag	UNP O17087
A	545	ALA	-	expression tag	UNP O17087
A	668	ALA	ASP	engineered mutation	UNP O17087
A	?	-	ALA	deletion	UNP O17087
A	?	-	LEU	deletion	UNP O17087
A	?	-	ALA	deletion	UNP O17087
A	?	-	VAL	deletion	UNP O17087
A	?	-	ARG	deletion	UNP O17087
A	?	-	PRO	deletion	UNP O17087
A	?	-	LYS	deletion	UNP O17087
A	?	-	ILE	deletion	UNP O17087
A	?	-	HIS	deletion	UNP O17087
A	?	-	SER	deletion	UNP O17087
A	?	-	ASN	deletion	UNP O17087
A	?	-	SER	deletion	UNP O17087
A	?	-	GLU	deletion	UNP O17087
A	?	-	GLY	deletion	UNP O17087
A	?	-	ASP	deletion	UNP O17087
A	?	-	LYS	deletion	UNP O17087
A	?	-	GLU	deletion	UNP O17087
A	?	-	THR	deletion	UNP O17087

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Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	PRO	deletion	UNP O17087
A	?	-	PRO	deletion	UNP O17087
A	?	-	PRO	deletion	UNP O17087
A	?	-	SER	deletion	UNP O17087
A	?	-	SER	deletion	UNP O17087
A	?	-	SER	deletion	UNP O17087
A	?	-	ALA	deletion	UNP O17087
A	?	-	SER	deletion	UNP O17087
A	?	-	THR	deletion	UNP O17087
A	?	-	SER	deletion	UNP O17087
A	?	-	SER	deletion	UNP O17087
A	?	-	ILE	deletion	UNP O17087
A	?	-	HIS	deletion	UNP O17087
A	?	-	ASN	deletion	UNP O17087
A	?	-	GLY	deletion	UNP O17087
A	?	-	GLY	deletion	UNP O17087
A	?	-	THR	deletion	UNP O17087
A	?	-	PRO	deletion	UNP O17087
A	?	-	GLY	deletion	UNP O17087
A	?	-	ILE	deletion	UNP O17087
A	?	-	PRO	deletion	UNP O17087
A	?	-	MET	deletion	UNP O17087
A	?	-	HIS	deletion	UNP O17087
A	?	-	HIS	deletion	UNP O17087
B	544	GLY	-	expression tag	UNP O17087
B	545	ALA	-	expression tag	UNP O17087
B	668	ALA	ASP	engineered mutation	UNP O17087
B	?	-	ALA	deletion	UNP O17087
B	?	-	LEU	deletion	UNP O17087
B	?	-	ALA	deletion	UNP O17087
B	?	-	VAL	deletion	UNP O17087
B	?	-	ARG	deletion	UNP O17087
B	?	-	PRO	deletion	UNP O17087
B	?	-	LYS	deletion	UNP O17087
B	?	-	ILE	deletion	UNP O17087
B	?	-	HIS	deletion	UNP O17087
B	?	-	SER	deletion	UNP O17087
B	?	-	ASN	deletion	UNP O17087
B	?	-	SER	deletion	UNP O17087
B	?	-	GLU	deletion	UNP O17087
B	?	-	GLY	deletion	UNP O17087
B	?	-	ASP	deletion	UNP O17087

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Chain	Residue	Modelled	Actual	Comment	Reference
B	?	-	LYS	deletion	UNP O17087
B	?	-	GLU	deletion	UNP O17087
B	?	-	THR	deletion	UNP O17087
B	?	-	PRO	deletion	UNP O17087
B	?	-	PRO	deletion	UNP O17087
B	?	-	PRO	deletion	UNP O17087
B	?	-	SER	deletion	UNP O17087
B	?	-	SER	deletion	UNP O17087
B	?	-	SER	deletion	UNP O17087
B	?	-	ALA	deletion	UNP O17087
B	?	-	SER	deletion	UNP O17087
B	?	-	THR	deletion	UNP O17087
B	?	-	SER	deletion	UNP O17087
B	?	-	SER	deletion	UNP O17087
B	?	-	ILE	deletion	UNP O17087
B	?	-	HIS	deletion	UNP O17087
B	?	-	ASN	deletion	UNP O17087
B	?	-	GLY	deletion	UNP O17087
B	?	-	GLY	deletion	UNP O17087
B	?	-	THR	deletion	UNP O17087
B	?	-	PRO	deletion	UNP O17087
B	?	-	GLY	deletion	UNP O17087
B	?	-	ILE	deletion	UNP O17087
B	?	-	PRO	deletion	UNP O17087
B	?	-	MET	deletion	UNP O17087
B	?	-	HIS	deletion	UNP O17087
B	?	-	HIS	deletion	UNP O17087
C	544	GLY	-	expression tag	UNP O17087
C	545	ALA	-	expression tag	UNP O17087
C	668	ALA	ASP	engineered mutation	UNP O17087
C	?	-	ALA	deletion	UNP O17087
C	?	-	LEU	deletion	UNP O17087
C	?	-	ALA	deletion	UNP O17087
C	?	-	VAL	deletion	UNP O17087
C	?	-	ARG	deletion	UNP O17087
C	?	-	PRO	deletion	UNP O17087
C	?	-	LYS	deletion	UNP O17087
C	?	-	ILE	deletion	UNP O17087
C	?	-	HIS	deletion	UNP O17087
C	?	-	SER	deletion	UNP O17087
C	?	-	ASN	deletion	UNP O17087
C	?	-	SER	deletion	UNP O17087

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Chain	Residue	Modelled	Actual	Comment	Reference
C	?	-	GLU	deletion	UNP O17087
C	?	-	GLY	deletion	UNP O17087
C	?	-	ASP	deletion	UNP O17087
C	?	-	LYS	deletion	UNP O17087
C	?	-	GLU	deletion	UNP O17087
C	?	-	THR	deletion	UNP O17087
C	?	-	PRO	deletion	UNP O17087
C	?	-	PRO	deletion	UNP O17087
C	?	-	PRO	deletion	UNP O17087
C	?	-	SER	deletion	UNP O17087
C	?	-	SER	deletion	UNP O17087
C	?	-	SER	deletion	UNP O17087
C	?	-	ALA	deletion	UNP O17087
C	?	-	SER	deletion	UNP O17087
C	?	-	THR	deletion	UNP O17087
C	?	-	SER	deletion	UNP O17087
C	?	-	SER	deletion	UNP O17087
C	?	-	ILE	deletion	UNP O17087
C	?	-	HIS	deletion	UNP O17087
C	?	-	ASN	deletion	UNP O17087
C	?	-	GLY	deletion	UNP O17087
C	?	-	GLY	deletion	UNP O17087
C	?	-	THR	deletion	UNP O17087
C	?	-	PRO	deletion	UNP O17087
C	?	-	GLY	deletion	UNP O17087
C	?	-	ILE	deletion	UNP O17087
C	?	-	PRO	deletion	UNP O17087
C	?	-	MET	deletion	UNP O17087
C	?	-	HIS	deletion	UNP O17087
C	?	-	HIS	deletion	UNP O17087
D	544	GLY	-	expression tag	UNP O17087
D	545	ALA	-	expression tag	UNP O17087
D	668	ALA	ASP	engineered mutation	UNP O17087
D	?	-	ALA	deletion	UNP O17087
D	?	-	LEU	deletion	UNP O17087
D	?	-	ALA	deletion	UNP O17087
D	?	-	VAL	deletion	UNP O17087
D	?	-	ARG	deletion	UNP O17087
D	?	-	PRO	deletion	UNP O17087
D	?	-	LYS	deletion	UNP O17087
D	?	-	ILE	deletion	UNP O17087
D	?	-	HIS	deletion	UNP O17087

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Chain	Residue	Modelled	Actual	Comment	Reference
D	?	-	SER	deletion	UNP O17087
D	?	-	ASN	deletion	UNP O17087
D	?	-	SER	deletion	UNP O17087
D	?	-	GLU	deletion	UNP O17087
D	?	-	GLY	deletion	UNP O17087
D	?	-	ASP	deletion	UNP O17087
D	?	-	LYS	deletion	UNP O17087
D	?	-	GLU	deletion	UNP O17087
D	?	-	THR	deletion	UNP O17087
D	?	-	PRO	deletion	UNP O17087
D	?	-	PRO	deletion	UNP O17087
D	?	-	PRO	deletion	UNP O17087
D	?	-	SER	deletion	UNP O17087
D	?	-	SER	deletion	UNP O17087
D	?	-	SER	deletion	UNP O17087
D	?	-	ALA	deletion	UNP O17087
D	?	-	SER	deletion	UNP O17087
D	?	-	THR	deletion	UNP O17087
D	?	-	SER	deletion	UNP O17087
D	?	-	SER	deletion	UNP O17087
D	?	-	ILE	deletion	UNP O17087
D	?	-	HIS	deletion	UNP O17087
D	?	-	ASN	deletion	UNP O17087
D	?	-	GLY	deletion	UNP O17087
D	?	-	GLY	deletion	UNP O17087
D	?	-	THR	deletion	UNP O17087
D	?	-	PRO	deletion	UNP O17087
D	?	-	GLY	deletion	UNP O17087
D	?	-	ILE	deletion	UNP O17087
D	?	-	PRO	deletion	UNP O17087
D	?	-	MET	deletion	UNP O17087
D	?	-	HIS	deletion	UNP O17087
D	?	-	HIS	deletion	UNP O17087

- Molecule 2 is a protein called RNP (RRM RNA binding domain) containing.

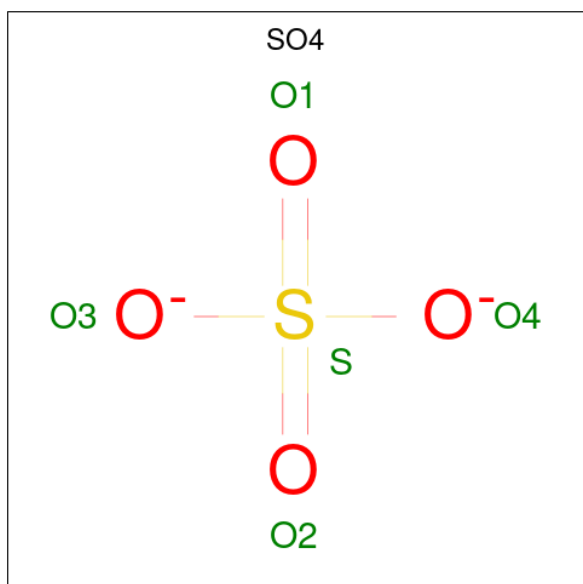
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	46	Total	C	N	O	0	0	0
			360	230	60	70			
2	F	46	Total	C	N	O	0	0	0
			364	232	60	72			
2	G	28	Total	C	N	O	0	0	0
			201	128	30	43			

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	H	25	Total	C	N	O	0	0	0
			187	121	28	38			

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

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

- Molecule 5 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



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

- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	47	Total O 47 47	0	0
6	B	37	Total O 37 37	0	0
6	C	20	Total O 20 20	0	0

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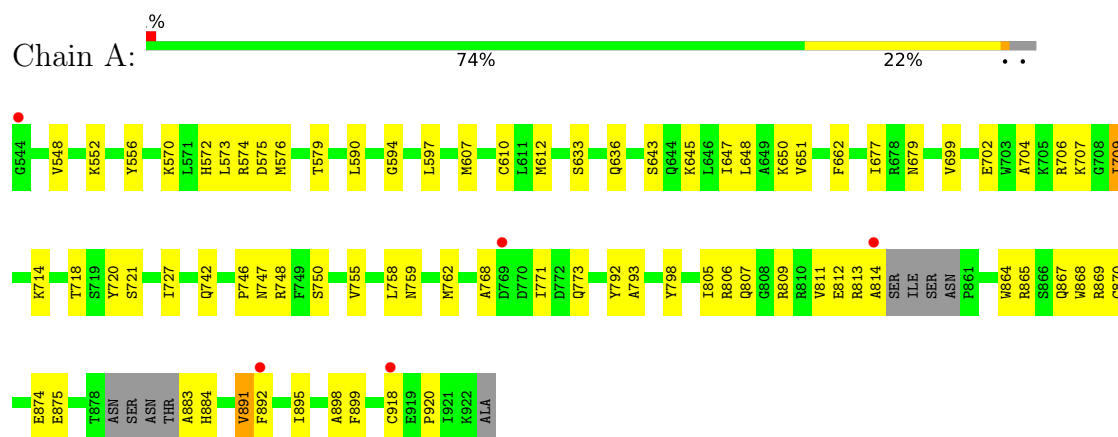
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	D	9	Total 9	O 9	0	0
6	E	9	Total 9	O 9	0	0
6	F	7	Total 7	O 7	0	0
6	G	1	Total 1	O 1	0	0

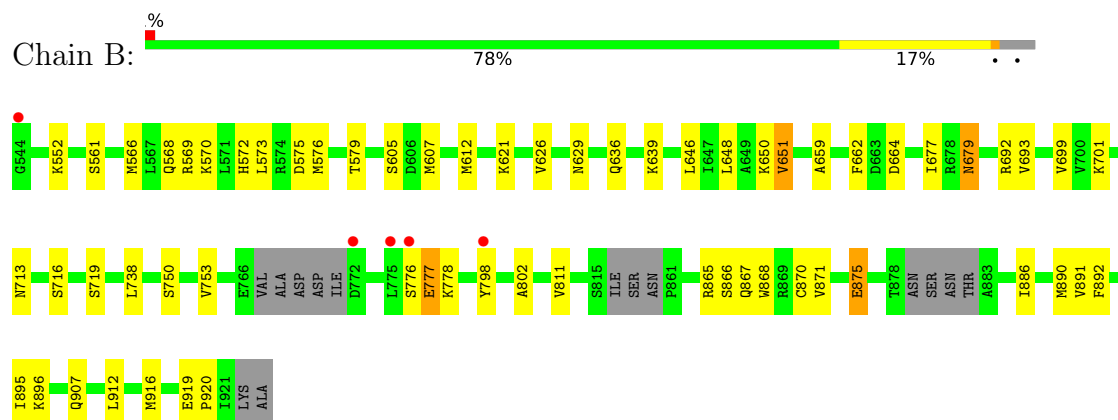
3 Residue-property plots [i](#)

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.

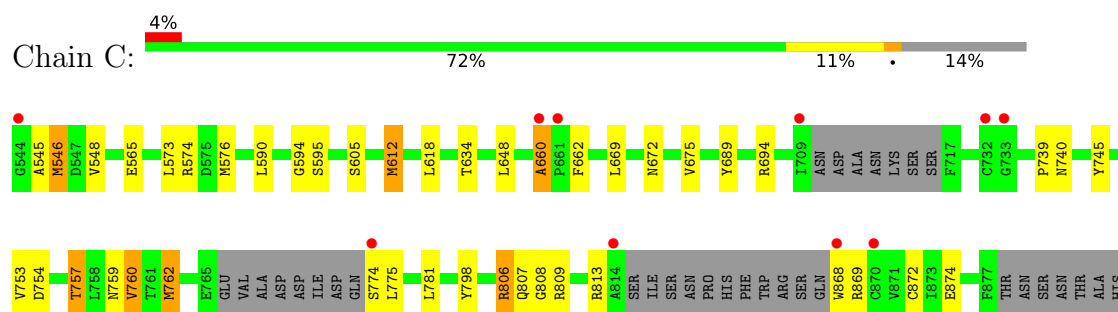
- Molecule 1: Poly(A) RNA polymerase gld-2

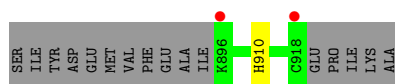


- Molecule 1: Poly(A) RNA polymerase gld-2

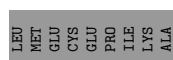
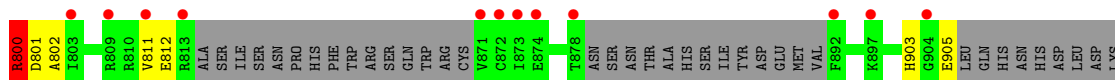
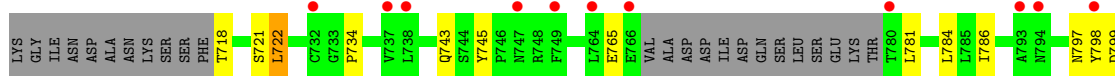
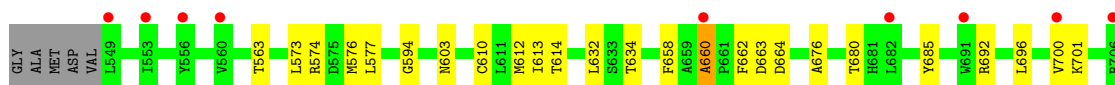


- Molecule 1: Poly(A) RNA polymerase gld-2

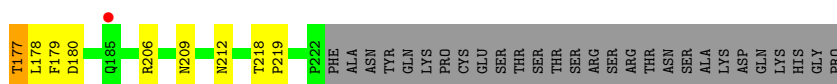




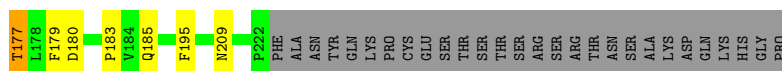
- Molecule 1: Poly(A) RNA polymerase gld-2



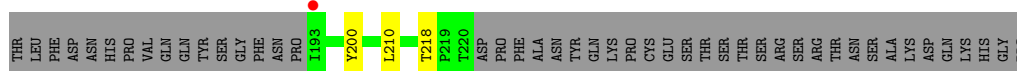
- Molecule 2: RNP (RRM RNA binding domain) containing



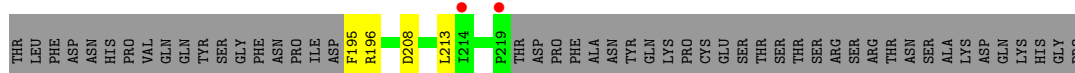
- Molecule 2: RNP (RRM RNA binding domain) containing



- Molecule 2: RNP (RRM RNA binding domain) containing



- Molecule 2: RNP (RRM RNA binding domain) containing



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	65.35Å 86.44Å 91.91Å 89.88° 93.13° 90.22°	Depositor
Resolution (Å)	46.25 – 2.49 46.25 – 2.49	Depositor EDS
% Data completeness (in resolution range)	95.6 (46.25-2.49) 95.5 (46.25-2.49)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.45 (at 2.48Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.192 , 0.242 0.198 , 0.244	Depositor DCC
R_{free} test set	3383 reflections (4.88%)	wwPDB-VP
Wilson B-factor (Å ²)	50.3	Xtriage
Anisotropy	0.358	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 38.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.025 for -h,k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10776	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.58% 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: EDO, MG, 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	0.60	0/2677	0.90	1/3630 (0.0%)
1	B	0.59	0/2643	0.86	6/3580 (0.2%)
1	C	0.56	0/2300	0.92	4/3123 (0.1%)
1	D	0.46	0/2041	0.90	3/2785 (0.1%)
2	E	0.62	0/371	0.85	0/508
2	F	0.61	0/375	0.85	0/513
2	G	0.54	0/205	0.87	0/281
2	H	0.41	0/191	0.88	0/259
All	All	0.56	0/10803	0.89	14/14679 (0.1%)

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	660	ALA	CA-C-N	-13.21	103.33	119.84
1	C	660	ALA	C-N-CA	-13.21	103.33	119.84
1	D	660	ALA	CA-C-N	-11.53	106.94	120.89
1	D	660	ALA	C-N-CA	-11.53	106.94	120.89
1	A	807	GLN	N-CA-C	-7.49	100.64	110.53
1	B	651	VAL	CB-CA-C	6.17	116.55	109.89
1	B	778	LYS	N-CA-C	-5.72	106.91	112.97
1	C	745	TYR	CA-C-N	5.64	125.17	119.19
1	C	745	TYR	C-N-CA	5.64	125.17	119.19
1	D	800	ARG	N-CA-C	-5.59	107.10	114.31
1	B	875	GLU	CA-C-N	5.10	125.14	119.32
1	B	875	GLU	C-N-CA	5.10	125.14	119.32
1	B	919	GLU	CA-C-N	5.03	126.12	119.84
1	B	919	GLU	C-N-CA	5.03	126.12	119.84

There are no chirality outliers.

There are no planarity outliers.

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	2623	0	2570	57	0
1	B	2590	0	2547	41	0
1	C	2259	0	2191	29	0
1	D	2004	0	1908	30	0
2	E	360	0	316	6	0
2	F	364	0	320	6	0
2	G	201	0	161	5	0
2	H	187	0	158	5	0
3	A	5	0	0	0	0
3	B	5	0	0	1	0
3	C	5	0	0	1	0
3	D	5	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	12	0	18	6	0
5	B	12	0	18	6	0
5	C	8	0	12	1	0
5	D	4	0	6	1	0
6	A	47	0	0	2	0
6	B	37	0	0	2	0
6	C	20	0	0	0	0
6	D	9	0	0	1	0
6	E	9	0	0	2	0
6	F	7	0	0	1	0
6	G	1	0	0	0	0
All	All	10776	0	10225	161	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 8.

All (161) 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)
2:E:177:THR:HG23	2:E:179:PHE:H	1.40	0.86
2:F:177:THR:HG23	2:F:179:PHE:H	1.41	0.85
1:B:692:ARG:NH2	1:B:777:GLU:OE1	2.10	0.85
1:A:573:LEU:HA	1:A:576:MET:HE3	1.64	0.80
1:A:574:ARG:NE	6:A:1101:HOH:O	2.16	0.79
1:D:594:GLY:H	1:D:680:THR:HG21	1.48	0.79
1:A:610:CYS:SG	1:A:612:MET:HE1	2.24	0.78
1:B:659:ALA:O	6:B:1101:HOH:O	2.00	0.78
2:E:206:ARG:NH1	6:E:401:HOH:O	2.18	0.76
1:D:573:LEU:HA	1:D:576:MET:HE3	1.71	0.72
1:B:679:ASN:HD22	5:B:1003:EDO:H12	1.54	0.71
1:A:884:HIS:HD1	1:A:892:PHE:HE2	1.41	0.67
1:C:740:ASN:ND2	1:C:808:GLY:H	1.93	0.67
1:D:718:THR:HG23	1:D:721:SER:HB2	1.77	0.66
1:C:798:TYR:O	1:C:813:ARG:NH2	2.28	0.66
1:B:573:LEU:HA	1:B:576:MET:HE3	1.78	0.65
1:C:576:MET:HE1	1:C:662:PHE:CE2	2.31	0.65
1:A:792:TYR:O	1:A:798:TYR:OH	2.12	0.65
1:D:797:ASN:OD1	1:D:798:TYR:N	2.30	0.64
2:E:209:ASN:OD1	2:E:212:ASN:ND2	2.30	0.64
1:A:870:CYS:HB3	1:A:892:PHE:CE1	2.33	0.63
1:A:643:SER:HB2	1:B:713:ASN:HB3	1.80	0.63
1:A:594:GLY:HA2	5:A:1005:EDO:H22	1.80	0.63
1:B:650:LYS:NZ	2:E:180:ASP:OD1	2.32	0.63
1:A:771:ILE:HG22	1:A:773:GLN:HG2	1.81	0.63
1:B:629:ASN:HB3	5:B:1002:EDO:H22	1.81	0.62
1:B:626:VAL:HA	5:B:1002:EDO:H11	1.81	0.61
1:D:563:THR:HG23	1:D:603:ASN:HD21	1.66	0.61
1:A:574:ARG:HH11	1:A:574:ARG:HG2	1.66	0.60
1:C:573:LEU:HA	1:C:576:MET:HE3	1.85	0.59
1:C:648:LEU:HD12	2:F:195:PHE:HE2	1.69	0.57
1:D:676:ALA:O	1:D:680:THR:HG23	2.04	0.57
1:D:594:GLY:H	1:D:680:THR:CG2	2.18	0.57
1:A:748:ARG:NH2	1:A:762:MET:O	2.39	0.56
1:D:612:MET:HE3	6:D:1109:HOH:O	2.05	0.56
1:B:870:CYS:HB3	1:B:892:PHE:CZ	2.42	0.55
1:A:576:MET:HE1	1:A:662:PHE:CZ	2.41	0.55
1:A:721:SER:OG	1:A:865:ARG:NH1	2.39	0.55
1:A:812:GLU:O	1:A:814:ALA:N	2.39	0.55
1:A:891:VAL:O	1:A:895:ILE:HG12	2.07	0.54
1:D:799:ASP:N	1:D:799:ASP:OD1	2.38	0.54
1:A:650:LYS:NZ	2:F:180:ASP:OD2	2.39	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:866:SER:H	5:B:1004:EDO:H11	1.72	0.54
1:C:759:ASN:CG	1:C:762:MET:HB3	2.32	0.54
1:A:714:LYS:HB3	5:A:1003:EDO:H21	1.91	0.53
1:D:800:ARG:O	1:D:812:GLU:HB2	2.09	0.53
1:A:707:LYS:HD3	1:A:898:ALA:HB1	1.90	0.53
1:B:575:ASP:O	1:B:579:THR:HG23	2.09	0.53
1:D:802:ALA:N	1:D:811:VAL:O	2.41	0.52
1:A:806:ARG:HH12	1:A:874:GLU:CD	2.17	0.52
1:B:891:VAL:O	1:B:895:ILE:HG12	2.09	0.52
1:C:605:SER:OG	3:C:1001:SO4:O2	2.26	0.52
1:B:750:SER:HB3	1:B:753:VAL:HG23	1.92	0.51
1:B:802:ALA:HB3	1:B:811:VAL:HG23	1.92	0.51
1:A:575:ASP:O	1:A:579:THR:HG23	2.10	0.51
1:D:696:LEU:O	1:D:700:VAL:HG23	2.10	0.51
1:A:755:VAL:HA	1:A:758:LEU:HG	1.93	0.51
1:A:867:GLN:HG3	1:A:868:TRP:CD2	2.46	0.51
1:A:759:ASN:HD21	1:A:762:MET:HB2	1.75	0.50
1:A:576:MET:HE1	1:A:662:PHE:CE2	2.47	0.50
1:B:605:SER:OG	3:B:1001:SO4:O3	2.27	0.49
5:A:1003:EDO:C2	1:B:636:GLN:HB2	2.42	0.49
1:A:648:LEU:HB2	1:B:646:LEU:HB3	1.95	0.49
1:D:610:CYS:SG	1:D:612:MET:HE1	2.52	0.49
1:B:867:GLN:HG3	1:B:868:TRP:CD2	2.48	0.49
1:D:701:LYS:HA	1:D:722:LEU:HD23	1.93	0.49
1:D:574:ARG:HH12	2:H:208:ASP:CG	2.20	0.49
1:A:869:ARG:NH2	2:F:185:GLN:OE1	2.45	0.48
1:A:870:CYS:HB3	1:A:892:PHE:CZ	2.49	0.48
1:D:685:TYR:CE1	2:H:213:LEU:HD11	2.49	0.48
1:D:614:THR:HG22	2:H:196:ARG:O	2.14	0.48
5:A:1003:EDO:H22	1:B:636:GLN:HB2	1.97	0.47
1:B:865:ARG:HD2	1:B:875:GLU:OE2	2.13	0.47
1:C:694:ARG:HH12	2:G:218:THR:HG22	1.79	0.47
1:B:679:ASN:ND2	5:B:1003:EDO:H12	2.24	0.47
1:C:806:ARG:HD3	1:C:874:GLU:CD	2.40	0.47
1:B:798:TYR:CD2	1:B:896:LYS:HE3	2.50	0.47
1:D:797:ASN:O	1:D:801:ASP:HB2	2.15	0.46
1:A:574:ARG:HD2	1:A:590:LEU:O	2.15	0.46
1:A:709:ILE:HD13	1:A:709:ILE:HA	1.72	0.46
2:E:206:ARG:NH2	6:E:402:HOH:O	2.35	0.46
1:A:570:LYS:HB3	1:A:607:MET:SD	2.56	0.46
1:A:718:THR:HG22	1:A:720:TYR:N	2.31	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:612:MET:HG2	2:G:200:TYR:CE1	2.51	0.46
1:D:811:VAL:HG22	1:D:812:GLU:H	1.81	0.46
1:A:793:ALA:HB2	1:A:899:PHE:HB3	1.98	0.46
1:B:576:MET:HE1	1:B:662:PHE:CZ	2.50	0.46
1:B:646:LEU:HD12	2:F:183:PRO:HG2	1.98	0.46
1:B:870:CYS:HB3	1:B:892:PHE:CE1	2.50	0.46
1:C:807:GLN:O	1:C:809:ARG:N	2.49	0.46
1:A:718:THR:HG22	1:A:720:TYR:H	1.81	0.45
1:B:867:GLN:HG3	1:B:868:TRP:CE3	2.51	0.45
1:C:694:ARG:HH12	2:G:218:THR:CG2	2.28	0.45
1:C:694:ARG:NH1	2:G:218:THR:HG22	2.32	0.45
1:A:702:GLU:O	1:A:706:ARG:HG3	2.16	0.45
1:C:740:ASN:HD22	1:C:808:GLY:H	1.60	0.45
1:D:685:TYR:CZ	2:H:213:LEU:HD11	2.52	0.45
1:A:556:TYR:OH	1:A:702:GLU:HG3	2.16	0.45
1:A:612:MET:HE3	1:A:677:ILE:HD11	1.98	0.45
1:A:727:ILE:HG22	1:A:805:ILE:HD12	1.99	0.45
1:B:693:VAL:HG21	1:B:738:LEU:HD21	1.98	0.44
1:A:643:SER:CB	1:B:713:ASN:HB3	2.47	0.44
1:B:566:MET:HE3	1:B:570:LYS:HE3	1.99	0.44
1:B:576:MET:HE1	1:B:662:PHE:CE2	2.52	0.44
1:A:720:TYR:HD2	1:A:865:ARG:CZ	2.30	0.44
1:B:569:ARG:NH2	1:B:664:ASP:OD2	2.51	0.44
1:D:573:LEU:HD21	1:D:658:PHE:CE2	2.52	0.44
1:D:576:MET:HE1	1:D:662:PHE:CZ	2.53	0.44
1:A:574:ARG:NH1	1:A:575:ASP:OD1	2.46	0.44
1:A:883:ALA:HA	1:A:884:HIS:HD2	1.83	0.44
1:A:884:HIS:N	1:A:884:HIS:CD2	2.86	0.44
1:C:868:TRP:O	1:C:869:ARG:HG3	2.18	0.44
1:A:645:LYS:HD2	1:A:647:ILE:HD11	1.99	0.43
1:D:786:ILE:H	1:D:786:ILE:HG13	1.68	0.43
1:B:912:LEU:O	1:B:916:MET:HG2	2.18	0.43
1:A:875:GLU:HB3	6:A:1131:HOH:O	2.17	0.43
1:C:594:GLY:HA2	5:C:1002:EDO:H12	2.00	0.43
1:C:760:VAL:HG22	2:G:210:LEU:HG	2.00	0.43
1:C:618:LEU:HD12	1:C:618:LEU:HA	1.83	0.43
1:A:747:ASN:ND2	1:A:748:ARG:HG2	2.34	0.43
1:B:662:PHE:HB2	6:B:1101:HOH:O	2.18	0.43
1:B:701:LYS:HE2	1:B:719:SER:HB2	2.00	0.43
1:B:612:MET:HE3	1:B:677:ILE:HD11	2.01	0.43
1:C:813:ARG:NH1	1:C:872:CYS:HB2	2.34	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:633:SER:O	1:A:636:GLN:HG2	2.19	0.43
1:B:798:TYR:HD1	1:B:871:VAL:HG22	1.84	0.43
1:D:577:LEU:HD23	1:D:577:LEU:HA	1.89	0.43
1:D:664:ASP:OD1	1:D:664:ASP:N	2.48	0.43
1:D:692:ARG:HE	1:D:784:LEU:HD11	1.84	0.43
1:A:865:ARG:HA	5:A:1004:EDO:H11	2.01	0.43
1:C:774:SER:OG	1:C:775:LEU:N	2.52	0.42
2:F:209:ASN:ND2	6:F:301:HOH:O	2.45	0.42
1:A:714:LYS:CB	5:A:1003:EDO:H21	2.49	0.42
1:A:809:ARG:NH1	1:A:811:VAL:HG12	2.34	0.42
1:B:886:ILE:HG21	2:H:195:PHE:CE1	2.55	0.42
1:C:574:ARG:O	1:C:574:ARG:HG2	2.19	0.42
1:A:864:TRP:CZ3	1:A:874:GLU:HB2	2.54	0.42
1:C:689:TYR:CE2	1:C:739:PRO:HD3	2.55	0.42
1:A:805:ILE:HD13	1:A:805:ILE:HA	1.90	0.41
1:A:572:HIS:CE1	1:A:576:MET:HE2	2.55	0.41
1:D:745:TYR:OH	1:D:765:GLU:O	2.21	0.41
1:B:798:TYR:CD1	1:B:871:VAL:HG22	2.56	0.41
1:A:809:ARG:CZ	1:A:811:VAL:HG12	2.50	0.41
1:A:704:ALA:HB1	1:A:709:ILE:HG22	2.02	0.41
1:C:590:LEU:HD12	1:C:590:LEU:HA	1.94	0.41
1:D:680:THR:HG22	5:D:1002:EDO:O1	2.21	0.41
1:B:679:ASN:HB3	5:B:1003:EDO:H12	2.01	0.41
1:C:669:LEU:C	1:C:669:LEU:HD23	2.45	0.41
1:A:707:LYS:HG3	1:A:918:CYS:SG	2.60	0.41
1:A:746:PRO:O	1:A:750:SER:HB2	2.20	0.41
1:C:754:ASP:O	1:C:757:THR:HB	2.20	0.41
1:C:781:LEU:HD23	1:C:781:LEU:HA	1.86	0.41
2:E:218:THR:HA	2:E:219:PRO:HD2	1.93	0.41
1:D:613:ILE:HG22	1:D:614:THR:HG23	2.02	0.40
1:A:870:CYS:HB3	1:A:892:PHE:CD1	2.56	0.40
1:C:545:ALA:O	1:C:548:VAL:HB	2.21	0.40
1:B:570:LYS:HB3	1:B:607:MET:SD	2.61	0.40
1:C:546:MET:HG3	1:C:910:HIS:CD2	2.56	0.40
1:C:813:ARG:O	1:C:813:ARG:HG3	2.21	0.40
1:D:903:HIS:C	1:D:905:GLU:H	2.29	0.40
1:B:572:HIS:ND1	1:B:576:MET:HE2	2.37	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	323/338 (96%)	309 (96%)	11 (3%)	3 (1%)	14	25
1	B	316/338 (94%)	306 (97%)	9 (3%)	1 (0%)	36	53
1	C	279/338 (82%)	265 (95%)	13 (5%)	1 (0%)	30	46
1	D	253/338 (75%)	243 (96%)	7 (3%)	3 (1%)	10	18
2	E	44/74 (60%)	40 (91%)	3 (7%)	1 (2%)	5	7
2	F	44/74 (60%)	41 (93%)	3 (7%)	0	100	100
2	G	26/74 (35%)	24 (92%)	2 (8%)	0	100	100
2	H	23/74 (31%)	23 (100%)	0	0	100	100
All	All	1308/1648 (79%)	1251 (96%)	48 (4%)	9 (1%)	18	32

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	813	ARG
1	C	660	ALA
1	D	660	ALA
1	A	920	PRO
2	E	178	LEU
1	D	734	PRO
1	D	781	LEU
1	B	920	PRO
1	A	768	ALA

5.3.2 Protein sidechains

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was

analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	289/306 (94%)	280 (97%)	9 (3%)	35	60
1	B	287/306 (94%)	273 (95%)	14 (5%)	22	42
1	C	247/306 (81%)	235 (95%)	12 (5%)	22	42
1	D	212/306 (69%)	206 (97%)	6 (3%)	38	63
2	E	38/67 (57%)	37 (97%)	1 (3%)	40	65
2	F	39/67 (58%)	38 (97%)	1 (3%)	40	65
2	G	19/67 (28%)	19 (100%)	0	100	100
2	H	18/67 (27%)	18 (100%)	0	100	100
All	All	1149/1492 (77%)	1106 (96%)	43 (4%)	30	54

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	548	VAL
1	A	552	LYS
1	A	597	LEU
1	A	651	VAL
1	A	679	ASN
1	A	699	VAL
1	A	709	ILE
1	A	742	GLN
1	A	891	VAL
1	B	552	LYS
1	B	561	SER
1	B	568	GLN
1	B	621	LYS
1	B	639	LYS
1	B	648	LEU
1	B	651	VAL
1	B	679	ASN
1	B	699	VAL
1	B	716	SER
1	B	776	SER
1	B	777	GLU
1	B	890	MET
1	B	907	GLN
1	C	546	MET
1	C	565	GLU

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Mol	Chain	Res	Type
1	C	595	SER
1	C	612	MET
1	C	634	THR
1	C	672	ASN
1	C	675	VAL
1	C	753	VAL
1	C	757	THR
1	C	760	VAL
1	C	762	MET
1	C	806	ARG
1	D	632	LEU
1	D	634	THR
1	D	663	ASP
1	D	722	LEU
1	D	743	GLN
1	D	800	ARG
2	E	177	THR
2	F	177	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	636	GLN
1	A	731	GLN
1	A	807	GLN
1	B	679	ASN
1	B	742	GLN
1	C	740	ASN
1	C	807	GLN
1	C	903	HIS
1	C	909	ASN
1	D	562	GLN
1	D	728	HIS
1	D	731	GLN
2	E	209	ASN
2	E	212	ASN

5.3.3 RNA ⓘ

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 15 ligands modelled in this entry, 2 are monoatomic - leaving 13 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	$\# Z > 2$
5	EDO	B	1004	-	3,3,3	0.67	0	2,2,2	0.16	0
3	SO4	A	1001	-	4,4,4	0.18	0	6,6,6	0.32	0
5	EDO	B	1003	-	3,3,3	0.36	0	2,2,2	0.75	0
5	EDO	C	1002	-	3,3,3	0.49	0	2,2,2	0.31	0
5	EDO	A	1005	-	3,3,3	0.35	0	2,2,2	0.67	0
5	EDO	A	1003	-	3,3,3	0.70	0	2,2,2	0.50	0
5	EDO	A	1004	-	3,3,3	0.46	0	2,2,2	0.42	0
3	SO4	C	1001	-	4,4,4	0.25	0	6,6,6	0.24	0
3	SO4	B	1001	-	4,4,4	0.20	0	6,6,6	0.34	0
5	EDO	C	1003	-	3,3,3	0.34	0	2,2,2	0.49	0
5	EDO	B	1002	-	3,3,3	0.51	0	2,2,2	0.28	0
5	EDO	D	1002	-	3,3,3	0.47	0	2,2,2	0.38	0
3	SO4	D	1001	-	4,4,4	0.23	0	6,6,6	0.20	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
5	EDO	C	1002	-	-	1/1/1/1	-
5	EDO	B	1003	-	-	1/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	EDO	A	1005	-	-	0/1/1/1	-
5	EDO	A	1004	-	-	0/1/1/1	-
5	EDO	A	1003	-	-	1/1/1/1	-
5	EDO	C	1003	-	-	0/1/1/1	-
5	EDO	B	1002	-	-	1/1/1/1	-
5	EDO	D	1002	-	-	0/1/1/1	-
5	EDO	B	1004	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	1003	EDO	O1-C1-C2-O2
5	A	1003	EDO	O1-C1-C2-O2
5	C	1002	EDO	O1-C1-C2-O2
5	B	1002	EDO	O1-C1-C2-O2

There are no ring outliers.

10 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	1004	EDO	1	0
5	B	1003	EDO	3	0
5	C	1002	EDO	1	0
5	A	1005	EDO	1	0
5	A	1003	EDO	4	0
5	A	1004	EDO	1	0
3	C	1001	SO4	1	0
3	B	1001	SO4	1	0
5	B	1002	EDO	2	0
5	D	1002	EDO	1	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	329/338 (97%)	0.08	5 (1%) 72 69	29, 45, 69, 90	0
1	B	324/338 (95%)	0.05	5 (1%) 72 69	30, 45, 68, 89	0
1	C	289/338 (85%)	0.12	12 (4%) 40 36	28, 48, 73, 84	0
1	D	263/338 (77%)	0.83	32 (12%) 8 7	37, 72, 112, 118	0
2	E	46/74 (62%)	0.13	1 (2%) 62 59	33, 44, 69, 78	0
2	F	46/74 (62%)	0.15	0 100 100	34, 44, 62, 77	0
2	G	28/74 (37%)	0.30	1 (3%) 46 42	37, 52, 79, 87	0
2	H	25/74 (33%)	0.77	2 (8%) 18 16	53, 71, 94, 101	0
All	All	1350/1648 (81%)	0.25	58 (4%) 40 36	28, 49, 89, 118	0

All (58) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	814	ALA	5.0
1	D	549	LEU	4.8
1	C	868	TRP	3.8
1	C	918	CYS	3.6
1	D	560	VAL	3.5
2	H	219	PRO	3.4
1	D	766	GLU	3.3
1	D	892	PHE	3.3
1	D	747	ASN	3.2
1	C	814	ALA	3.1
1	C	870	CYS	3.1
1	D	780	THR	3.1
1	B	775	LEU	3.1
1	D	682	LEU	3.1
1	C	774	SER	3.0
1	D	691	TRP	3.0

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Mol	Chain	Res	Type	RSRZ
1	D	871	VAL	2.9
1	A	544	GLY	2.9
1	C	544	GLY	2.8
1	B	776	SER	2.7
1	D	873	ILE	2.7
1	D	878	THR	2.6
1	D	749	PHE	2.6
1	C	733	GLY	2.6
1	C	709	ILE	2.6
1	D	794	ASN	2.6
1	D	811	VAL	2.6
1	A	918	CYS	2.5
1	B	544	GLY	2.5
1	B	772	ASP	2.5
1	D	813	ARG	2.5
2	G	193	ILE	2.4
1	D	803	ILE	2.4
2	H	214	ILE	2.4
1	D	556	TYR	2.4
1	D	738	LEU	2.4
1	D	809	ARG	2.4
1	A	769	ASP	2.4
1	A	892	PHE	2.4
1	D	553	ILE	2.3
1	D	706	ARG	2.3
1	D	732	CYS	2.3
1	D	793	ALA	2.3
1	D	904	GLY	2.3
1	D	660	ALA	2.2
1	C	896	LYS	2.2
1	C	660	ALA	2.2
1	B	798	TYR	2.2
2	E	185	GLN	2.2
1	D	872	CYS	2.2
1	C	732	CYS	2.1
1	D	700	VAL	2.1
1	D	798	TYR	2.1
1	D	897	LYS	2.1
1	D	737	VAL	2.1
1	D	764	LEU	2.1
1	D	874	GLU	2.1
1	C	661	PRO	2.0

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
5	EDO	B	1004	4/4	0.78	0.24	48,50,51,58	0
4	MG	A	1002	1/1	0.80	0.15	58,58,58,58	0
5	EDO	A	1003	4/4	0.82	0.18	49,51,53,59	0
3	SO4	D	1001	5/5	0.87	0.09	79,87,91,93	0
5	EDO	B	1002	4/4	0.90	0.16	44,49,53,57	0
5	EDO	B	1003	4/4	0.90	0.19	31,36,38,48	0
5	EDO	A	1005	4/4	0.90	0.16	32,32,36,44	0
5	EDO	C	1003	4/4	0.90	0.18	35,46,48,54	0
5	EDO	A	1004	4/4	0.92	0.12	48,49,52,55	0
5	EDO	D	1002	4/4	0.92	0.14	57,59,61,71	0
3	SO4	C	1001	5/5	0.94	0.12	59,62,72,73	0
5	EDO	C	1002	4/4	0.94	0.10	41,43,48,48	0
4	MG	B	1005	1/1	0.96	0.14	51,51,51,51	0
3	SO4	B	1001	5/5	0.97	0.14	43,44,53,60	0
3	SO4	A	1001	5/5	0.98	0.07	46,46,54,55	0

6.5 Other polymers [i](#)

There are no such residues in this entry.