



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 01:35 AM UTC

PDB ID : 4LRC / pdb_00004lrc
Title : Phosphopentomutase V158L variant
Authors : Birmingham, W.A.; Starbird, C.A.; Panosian, T.D.; Nannemann, D.P.; Iverson, T.M.; Bachmann, B.O.
Deposited on : 2013-07-19
Resolution : 1.89 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

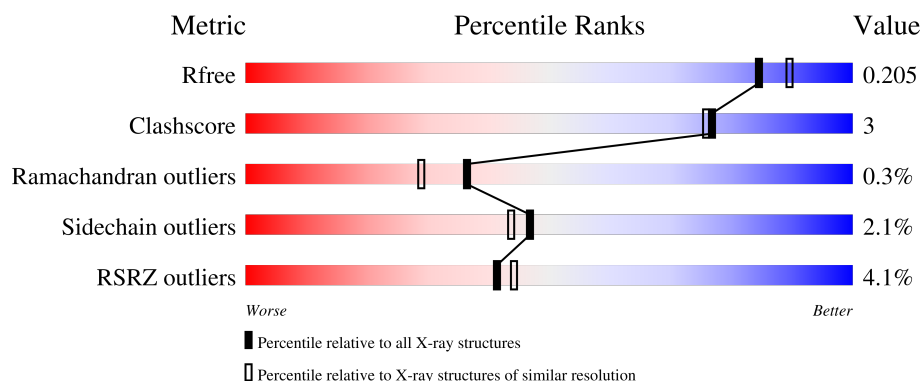
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.89 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	416	11% 88% 5% 6%
1	B	416	85% 8% 6%
1	C	416	82% 10% 6%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 9965 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 Phosphopentomutase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	391	Total	C	N	O	P	S	0	0	0
			3069	1940	505	606	1	17			
1	B	390	Total	C	N	O	P	S	0	0	0
			3053	1932	503	600	1	17			
1	C	390	Total	C	N	O	P	S	2	2	0
			3067	1941	504	603	1	18			

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-21	MET	-	expression tag	UNP Q818Z9
A	-20	GLY	-	expression tag	UNP Q818Z9
A	-19	SER	-	expression tag	UNP Q818Z9
A	-18	SER	-	expression tag	UNP Q818Z9
A	-17	HIS	-	expression tag	UNP Q818Z9
A	-16	HIS	-	expression tag	UNP Q818Z9
A	-15	HIS	-	expression tag	UNP Q818Z9
A	-14	HIS	-	expression tag	UNP Q818Z9
A	-13	HIS	-	expression tag	UNP Q818Z9
A	-12	HIS	-	expression tag	UNP Q818Z9
A	-11	SER	-	expression tag	UNP Q818Z9
A	-10	SER	-	expression tag	UNP Q818Z9
A	-9	GLY	-	expression tag	UNP Q818Z9
A	-8	LEU	-	expression tag	UNP Q818Z9
A	-7	VAL	-	expression tag	UNP Q818Z9
A	-6	PRO	-	expression tag	UNP Q818Z9
A	-5	ARG	-	expression tag	UNP Q818Z9
A	-4	GLY	-	expression tag	UNP Q818Z9
A	-3	SER	-	expression tag	UNP Q818Z9
A	-2	HIS	-	expression tag	UNP Q818Z9
A	-1	MET	-	expression tag	UNP Q818Z9
A	0	ALA	-	expression tag	UNP Q818Z9
A	1	SER	-	expression tag	UNP Q818Z9

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Chain	Residue	Modelled	Actual	Comment	Reference
A	158	LEU	VAL	engineered mutation	UNP Q818Z9
B	-21	MET	-	expression tag	UNP Q818Z9
B	-20	GLY	-	expression tag	UNP Q818Z9
B	-19	SER	-	expression tag	UNP Q818Z9
B	-18	SER	-	expression tag	UNP Q818Z9
B	-17	HIS	-	expression tag	UNP Q818Z9
B	-16	HIS	-	expression tag	UNP Q818Z9
B	-15	HIS	-	expression tag	UNP Q818Z9
B	-14	HIS	-	expression tag	UNP Q818Z9
B	-13	HIS	-	expression tag	UNP Q818Z9
B	-12	HIS	-	expression tag	UNP Q818Z9
B	-11	SER	-	expression tag	UNP Q818Z9
B	-10	SER	-	expression tag	UNP Q818Z9
B	-9	GLY	-	expression tag	UNP Q818Z9
B	-8	LEU	-	expression tag	UNP Q818Z9
B	-7	VAL	-	expression tag	UNP Q818Z9
B	-6	PRO	-	expression tag	UNP Q818Z9
B	-5	ARG	-	expression tag	UNP Q818Z9
B	-4	GLY	-	expression tag	UNP Q818Z9
B	-3	SER	-	expression tag	UNP Q818Z9
B	-2	HIS	-	expression tag	UNP Q818Z9
B	-1	MET	-	expression tag	UNP Q818Z9
B	0	ALA	-	expression tag	UNP Q818Z9
B	1	SER	-	expression tag	UNP Q818Z9
B	158	LEU	VAL	engineered mutation	UNP Q818Z9
C	-21	MET	-	expression tag	UNP Q818Z9
C	-20	GLY	-	expression tag	UNP Q818Z9
C	-19	SER	-	expression tag	UNP Q818Z9
C	-18	SER	-	expression tag	UNP Q818Z9
C	-17	HIS	-	expression tag	UNP Q818Z9
C	-16	HIS	-	expression tag	UNP Q818Z9
C	-15	HIS	-	expression tag	UNP Q818Z9
C	-14	HIS	-	expression tag	UNP Q818Z9
C	-13	HIS	-	expression tag	UNP Q818Z9
C	-12	HIS	-	expression tag	UNP Q818Z9
C	-11	SER	-	expression tag	UNP Q818Z9
C	-10	SER	-	expression tag	UNP Q818Z9
C	-9	GLY	-	expression tag	UNP Q818Z9
C	-8	LEU	-	expression tag	UNP Q818Z9
C	-7	VAL	-	expression tag	UNP Q818Z9
C	-6	PRO	-	expression tag	UNP Q818Z9
C	-5	ARG	-	expression tag	UNP Q818Z9

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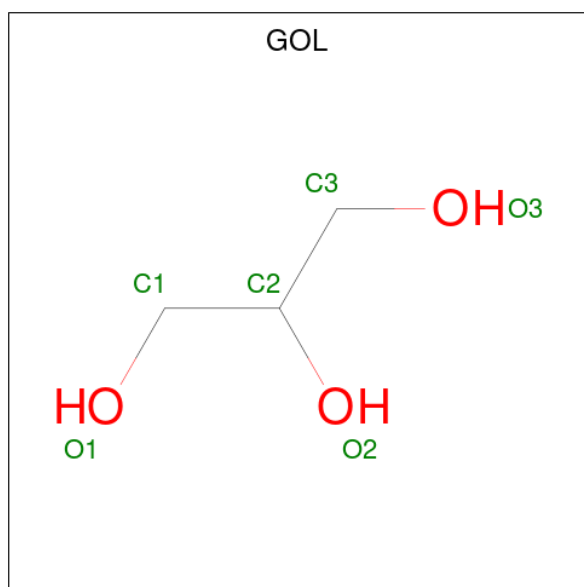
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Chain	Residue	Modelled	Actual	Comment	Reference
C	-4	GLY	-	expression tag	UNP Q818Z9
C	-3	SER	-	expression tag	UNP Q818Z9
C	-2	HIS	-	expression tag	UNP Q818Z9
C	-1	MET	-	expression tag	UNP Q818Z9
C	0	ALA	-	expression tag	UNP Q818Z9
C	1	SER	-	expression tag	UNP Q818Z9
C	158	LEU	VAL	engineered mutation	UNP Q818Z9

- Molecule 2 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	3	Total Mn 3 3	0	0
2	B	2	Total Mn 2 2	0	0
2	C	3	Total Mn 3 3	0	0

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



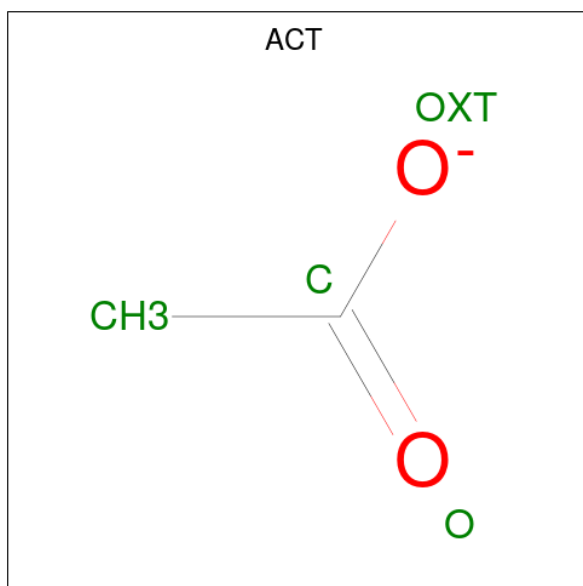
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 6 3 3	0	0
3	A	1	Total C O 6 3 3	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is ACETATE ION (CCD ID: ACT) (formula: $\text{C}_2\text{H}_3\text{O}_2$).



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

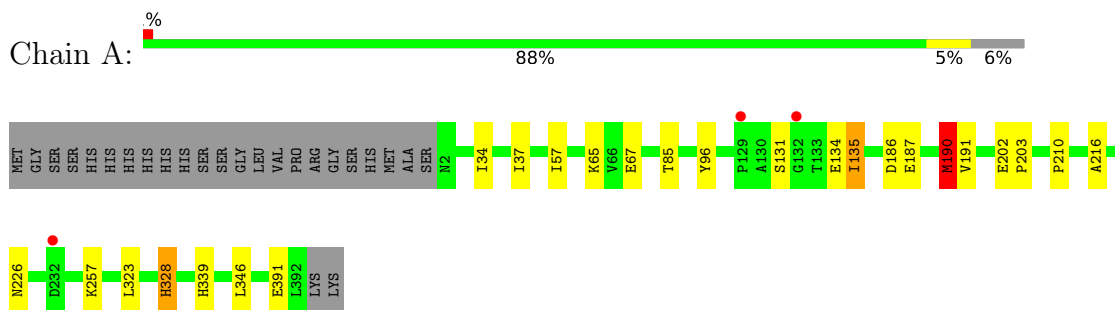
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	283	Total	O	0	0
			283	283		
5	B	308	Total	O	0	0
			308	308		
5	C	155	Total	O	0	0
			155	155		

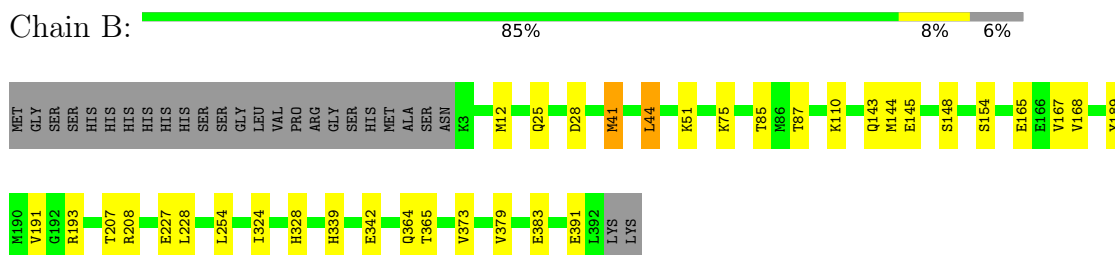
3 Residue-property plots

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.

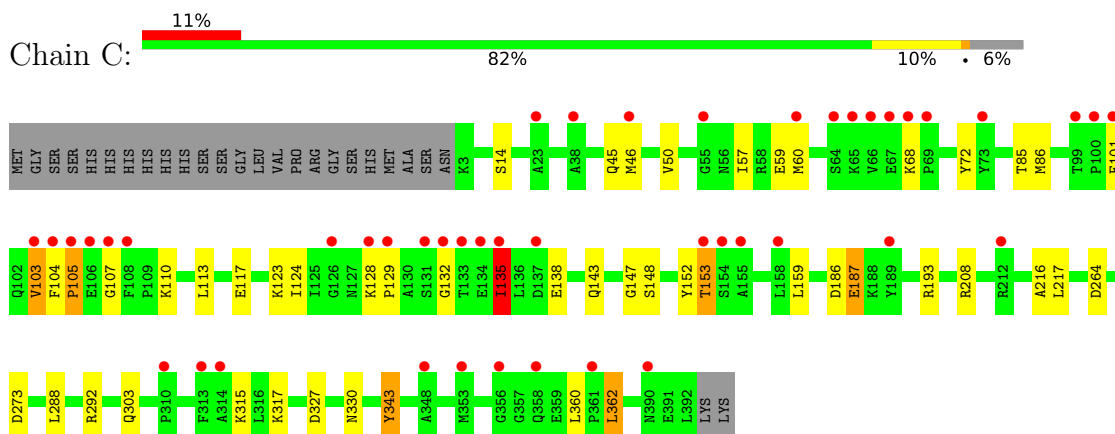
• Molecule 1: Phosphopentomutase



• Molecule 1: Phosphopentomutase



• Molecule 1: Phosphopentomutase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	91.72Å 76.80Å 107.17Å 90.00° 108.70° 90.00°	Depositor
Resolution (Å)	38.78 – 1.89 38.78 – 1.89	Depositor EDS
% Data completeness (in resolution range)	91.3 (38.78-1.89) 91.3 (38.78-1.89)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.45 (at 1.89Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.155 , 0.197 0.167 , 0.205	Depositor DCC
R_{free} test set	4956 reflections (3.81%)	wwPDB-VP
Wilson B-factor (Å ²)	19.6	Xtriage
Anisotropy	0.644	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 44.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	9965	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 5.06% 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: TPO, GOL, MN, ACT

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	1.34	6/3118 (0.2%)	1.18	3/4209 (0.1%)
1	B	1.32	6/3102 (0.2%)	1.15	3/4188 (0.1%)
1	C	1.28	5/3119 (0.2%)	1.14	4/4210 (0.1%)
All	All	1.31	17/9339 (0.2%)	1.16	10/12607 (0.1%)

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	186	ASP	CG-OD2	18.00	1.59	1.25
1	C	186	ASP	CG-OD1	7.20	1.39	1.25
1	B	207	THR	C-O	-6.57	1.16	1.24
1	C	135	ILE	CA-CB	6.41	1.61	1.54
1	A	67	GLU	C-O	-6.00	1.16	1.24
1	C	68	LYS	CA-C	5.86	1.60	1.52
1	B	342	GLU	CA-C	-5.84	1.45	1.52
1	A	37	ILE	CA-CB	5.64	1.61	1.54
1	A	186	ASP	CA-C	-5.64	1.46	1.52
1	C	327	ASP	CA-C	5.63	1.59	1.52
1	B	373	VAL	N-CA	-5.38	1.40	1.46
1	B	228	LEU	N-CA	5.33	1.52	1.46
1	B	87	THR	N-CA	-5.26	1.40	1.46
1	A	346	LEU	N-CA	-5.25	1.39	1.46
1	A	323	LEU	CA-C	-5.24	1.46	1.52
1	A	226	ASN	N-CA	5.23	1.52	1.46
1	B	254	LEU	C-O	-5.08	1.18	1.24

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	190	MET	CG-SD-CE	-8.52	82.15	100.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	57	ILE	N-CA-C	-6.49	104.43	110.53
1	C	147	GLY	N-CA-C	-5.79	107.27	115.43
1	B	168	VAL	CA-C-N	-5.57	114.51	120.03
1	B	168	VAL	C-N-CA	-5.57	114.51	120.03
1	B	41	MET	N-CA-C	-5.55	106.46	113.18
1	C	57	ILE	N-CA-C	-5.22	105.62	110.53
1	A	34	ILE	N-CA-CB	5.18	116.61	110.55
1	C	343	TYR	N-CA-C	-5.17	103.71	110.53
1	C	103	VAL	N-CA-C	-5.14	100.48	108.86

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	3069	0	3009	11	0
1	B	3053	0	2995	19	0
1	C	3067	0	3009	26	0
2	A	3	0	0	0	0
2	B	2	0	0	0	0
2	C	3	0	0	0	0
3	A	12	0	16	1	0
3	B	6	0	8	0	0
4	A	4	0	3	0	0
5	A	283	0	0	1	0
5	B	308	0	0	9	0
5	C	155	0	0	2	0
All	All	9965	0	9040	53	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 3.

All (53) 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:B:51:LYS:NZ	5:B:746:HOH:O	2.11	0.84
1:C:187:GLU:H	1:C:187:GLU:CD	1.98	0.71
1:C:117[B]:GLU:HG2	1:C:123:LYS:HA	1.73	0.70
1:B:28:ASP:OD1	5:B:542:HOH:O	2.10	0.70
1:C:264:ASP:OD1	5:C:655:HOH:O	2.12	0.68
1:B:383:GLU:OE1	5:B:807:HOH:O	2.14	0.64
1:C:72:TYR:HB3	1:C:360:LEU:HD21	1.82	0.62
1:C:86[A]:MET:HE2	5:C:564:HOH:O	2.00	0.61
1:A:216:ALA:HB1	5:A:729:HOH:O	2.01	0.58
1:C:193:ARG:NH1	1:C:216:ALA:HB2	2.17	0.58
1:C:45:GLN:H	1:C:303:GLN:HE21	1.52	0.57
1:C:187:GLU:CD	1:C:187:GLU:N	2.65	0.55
1:A:187:GLU:HA	1:A:190:MET:HE3	1.89	0.53
1:A:131:SER:OG	1:A:134:GLU:HG3	2.09	0.52
1:C:113:LEU:HD13	1:C:124:ILE:HG21	1.90	0.52
1:C:288:LEU:O	1:C:292:ARG:HD2	2.09	0.52
1:B:143:GLN:HG2	1:B:148:SER:O	2.11	0.50
1:C:143:GLN:HG2	1:C:148:SER:O	2.11	0.50
1:C:72:TYR:HB3	1:C:360:LEU:CD2	2.43	0.49
1:A:257:LYS:HD2	1:B:145:GLU:OE1	2.13	0.48
1:B:75:LYS:NZ	5:B:694:HOH:O	2.45	0.48
1:B:227:GLU:OE2	5:B:808:HOH:O	2.19	0.48
1:B:193:ARG:NH1	5:B:724:HOH:O	2.38	0.47
1:C:101:PHE:HE2	1:C:217:LEU:HD12	1.79	0.46
1:A:328:HIS:CE1	1:A:339:HIS:CE1	3.03	0.46
1:C:46:MET:O	1:C:50:VAL:HG23	2.16	0.46
1:C:86[B]:MET:CE	1:C:193:ARG:HH21	2.28	0.46
1:C:343:TYR:CE1	1:C:362:LEU:HD11	2.52	0.45
1:C:14:SER:HB2	1:C:330:ASN:HB2	1.98	0.44
1:C:128:LYS:O	1:C:129:PRO:C	2.59	0.44
1:A:202:GLU:HB2	1:A:203:PRO:HD2	1.99	0.44
1:C:128:LYS:NZ	1:C:138:GLU:OE1	2.36	0.44
1:C:129:PRO:CA	1:C:153:THR:HG22	2.48	0.44
1:B:12:MET:HE2	1:B:324:ILE:HG22	1.98	0.44
1:C:103:VAL:HG23	1:C:105:PRO:HD2	2.01	0.43
1:B:208:ARG:NH1	5:B:804:HOH:O	2.50	0.43
1:A:202:GLU:HB2	1:A:203:PRO:CD	2.49	0.43
1:C:59:GLU:C	1:C:60:MET:HG3	2.44	0.42
1:C:132:GLY:HA2	1:C:135:ILE:HG22	2.01	0.42
1:A:190:MET:HE3	1:A:190:MET:HB2	1.23	0.42
1:B:154:SER:HB2	5:B:760:HOH:O	2.19	0.42
1:A:135:ILE:HD12	1:A:135:ILE:HA	1.87	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:210:PRO:HG2	1:B:167:VAL:HG13	2.01	0.42
1:B:144:MET:HA	1:B:167:VAL:HG21	2.02	0.42
1:B:41:MET:HG3	1:B:44:LEU:HD13	2.02	0.42
1:B:165:GLU:OE2	1:C:273:ASP:OD2	2.37	0.42
1:B:227:GLU:HG3	1:B:379:VAL:HB	2.02	0.42
1:B:227:GLU:CG	1:B:379:VAL:HB	2.49	0.41
1:C:152:TYR:OH	1:C:208:ARG:NH2	2.47	0.41
1:A:96:TYR:CE1	3:A:404:GOL:H12	2.55	0.41
1:B:25:GLN:NE2	5:B:709:HOH:O	2.53	0.40
1:B:208:ARG:HH11	1:B:208:ARG:HG3	1.86	0.40
1:C:104:PHE:HB3	1:C:107:GLY:O	2.22	0.40

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	388/416 (93%)	372 (96%)	15 (4%)	1 (0%)	36	29
1	B	387/416 (93%)	371 (96%)	14 (4%)	2 (0%)	24	16
1	C	389/416 (94%)	365 (94%)	24 (6%)	0	100	100
All	All	1164/1248 (93%)	1108 (95%)	53 (5%)	3 (0%)	36	29

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	328	HIS
1	B	328	HIS
1	B	339	HIS

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	332/353 (94%)	327 (98%)	5 (2%)	57	56
1	B	329/353 (93%)	322 (98%)	7 (2%)	47	44
1	C	331/353 (94%)	322 (97%)	9 (3%)	39	34
All	All	992/1059 (94%)	971 (98%)	21 (2%)	47	44

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	65	LYS
1	A	135	ILE
1	A	190	MET
1	A	191	VAL
1	A	391	GLU
1	B	44	LEU
1	B	110	LYS
1	B	189	TYR
1	B	191	VAL
1	B	364	GLN
1	B	365	THR
1	B	391	GLU
1	C	105	PRO
1	C	110	LYS
1	C	135	ILE
1	C	153	THR
1	C	159	LEU
1	C	187	GLU
1	C	315	LYS
1	C	317	LYS
1	C	362	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	2	ASN
1	A	45	GLN
1	A	211	ASN
1	A	281	ASN
1	B	205	ASN
1	C	102	GLN
1	C	211	ASN
1	C	281	ASN
1	C	303	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

3 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	TPO	A	85	2,1	8,10,11	1.05	1 (12%)	10,14,16	1.00	1 (10%)
1	TPO	B	85	2,1	8,10,11	2.02	1 (12%)	10,14,16	1.28	0
1	TPO	C	85	2,1	8,10,11	1.01	0	10,14,16	1.27	1 (10%)

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	TPO	A	85	2,1	-	2/9/11/13	-
1	TPO	B	85	2,1	-	2/9/11/13	-
1	TPO	C	85	2,1	-	2/9/11/13	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	85	TPO	P-OG1	5.30	1.68	1.59
1	A	85	TPO	P-OG1	2.66	1.64	1.59

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	85	TPO	OG1-P-O1P	-2.41	100.74	109.33
1	A	85	TPO	O-C-CA	-2.13	119.29	124.77

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	85	TPO	O-C-CA-CB
1	B	85	TPO	O-C-CA-CB
1	B	85	TPO	C-CA-CB-CG2
1	A	85	TPO	C-CA-CB-CG2
1	C	85	TPO	C-CA-CB-CG2
1	C	85	TPO	O-C-CA-CB

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 12 ligands modelled in this entry, 8 are monoatomic - leaving 4 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
3	GOL	A	405	-	5,5,5	0.64	0	5,5,5	0.41	0
3	GOL	B	403	-	5,5,5	0.21	0	5,5,5	0.83	0
4	ACT	A	406	-	3,3,3	0.64	0	3,3,3	1.14	0
3	GOL	A	404	-	5,5,5	0.88	0	5,5,5	1.51	1 (20%)

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
3	GOL	A	405	-	-	2/4/4/4	-
3	GOL	B	403	-	-	0/4/4/4	-
3	GOL	A	404	-	-	1/4/4/4	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	404	GOL	C3-C2-C1	2.71	121.73	111.80

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	405	GOL	O1-C1-C2-C3
3	A	405	GOL	O1-C1-C2-O2
3	A	404	GOL	O2-C2-C3-O3

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	404	GOL	1	0

5.7 Other polymers [i](#)

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	390/416 (93%)	-0.31	3 (0%) 82 85	10, 18, 39, 60	14 (3%)
1	B	389/416 (93%)	-0.31	0 100 100	10, 19, 36, 51	10 (2%)
1	C	389/416 (93%)	0.63	45 (11%) 9 10	14, 30, 56, 86	11 (2%)
All	All	1168/1248 (93%)	0.00	48 (4%) 41 44	10, 22, 45, 86	35 (2%)

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	105	PRO	4.8
1	C	390	ASN	4.7
1	C	154	SER	4.7
1	C	155	ALA	4.1
1	C	133	THR	4.0
1	C	128	LYS	3.7
1	C	189	TYR	3.6
1	C	129	PRO	3.4
1	C	153	THR	3.4
1	C	310	PRO	3.3
1	A	132	GLY	3.1
1	C	23	ALA	3.0
1	C	73	TYR	2.9
1	C	55	GLY	2.9
1	C	158	LEU	2.9
1	C	356	GLY	2.9
1	A	129	PRO	2.8
1	C	135	ILE	2.8
1	C	132	GLY	2.8
1	C	67	GLU	2.8
1	C	103	VAL	2.8
1	C	137	ASP	2.7
1	C	100	PRO	2.7

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Mol	Chain	Res	Type	RSRZ
1	C	313	PHE	2.7
1	C	99	THR	2.6
1	C	64	SER	2.6
1	C	134	GLU	2.5
1	C	361	PRO	2.5
1	C	65	LYS	2.5
1	C	108	PHE	2.5
1	C	46	MET	2.5
1	A	232	ASP	2.4
1	C	104	PHE	2.4
1	C	353	MET	2.3
1	C	212	ARG	2.2
1	C	106	GLU	2.2
1	C	66	VAL	2.2
1	C	314	ALA	2.2
1	C	131	SER	2.1
1	C	107	GLY	2.1
1	C	60	MET	2.1
1	C	358	GLN	2.1
1	C	126	GLY	2.1
1	C	69	PRO	2.1
1	C	348	ALA	2.1
1	C	68	LYS	2.0
1	C	101	PHE	2.0
1	C	38	ALA	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. 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(Å ²)	Q<0.9
1	TPO	C	85	11/12	0.95	0.14	18,27,67,71	0
1	TPO	A	85	11/12	0.96	0.11	12,15,55,58	0
1	TPO	B	85	11/12	0.97	0.12	12,18,61,67	0

6.3 Carbohydrates ⓘ

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
4	ACT	A	406	4/4	0.82	0.17	46,55,56,61	0
3	GOL	A	404	6/6	0.83	0.13	33,35,41,42	0
3	GOL	A	405	6/6	0.95	0.11	19,42,45,46	0
3	GOL	B	403	6/6	0.96	0.07	26,31,36,36	0
2	MN	A	403	1/1	0.99	0.08	38,38,38,38	0
2	MN	C	401	1/1	0.99	0.03	19,19,19,19	0
2	MN	C	402	1/1	0.99	0.02	22,22,22,22	0
2	MN	C	403	1/1	0.99	0.05	41,41,41,41	0
2	MN	B	402	1/1	1.00	0.02	13,13,13,13	0
2	MN	A	402	1/1	1.00	0.01	12,12,12,12	0
2	MN	A	401	1/1	1.00	0.01	12,12,12,12	0
2	MN	B	401	1/1	1.00	0.01	14,14,14,14	0

6.5 Other polymers [i](#)

There are no such residues in this entry.