



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 08:19 AM UTC

PDB ID : 6B65 / pdb_00006b65
Title : IMPase (AF2372) R92Q/K164E
Authors : Goldstein, R.I.; Roberts, M.
Deposited on : 2017-10-01
Resolution : 2.70 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

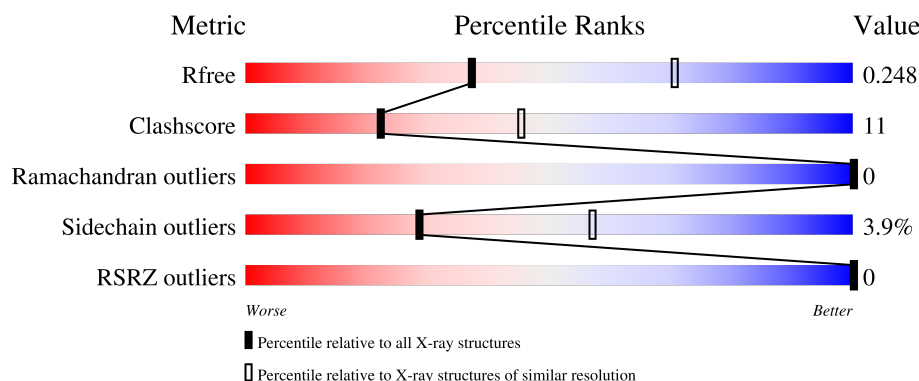
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	252	 75% 22% .
1	B	252	 85% 13% .

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	IPA	B	301	-	-	X	-
5	IPA	B	307	-	-	X	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 4060 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.

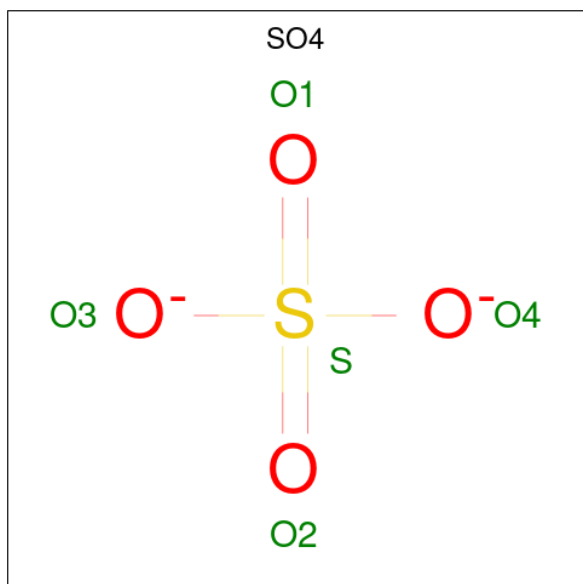
- Molecule 1 is a protein called Fructose-1,6-bisphosphatase/inositol-1-monophosphatase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	252	Total	C	N	O	S	0	2	0
			1978	1250	333	385	10			
1	B	252	Total	C	N	O	S	0	1	0
			1968	1244	331	383	10			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	92	GLN	ARG	engineered mutation	UNP O30298
A	164	GLU	LYS	engineered mutation	UNP O30298
B	92	GLN	ARG	engineered mutation	UNP O30298
B	164	GLU	LYS	engineered mutation	UNP O30298

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

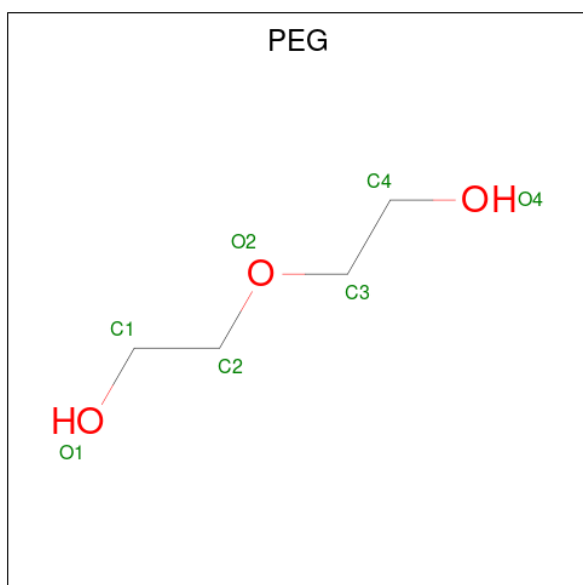


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

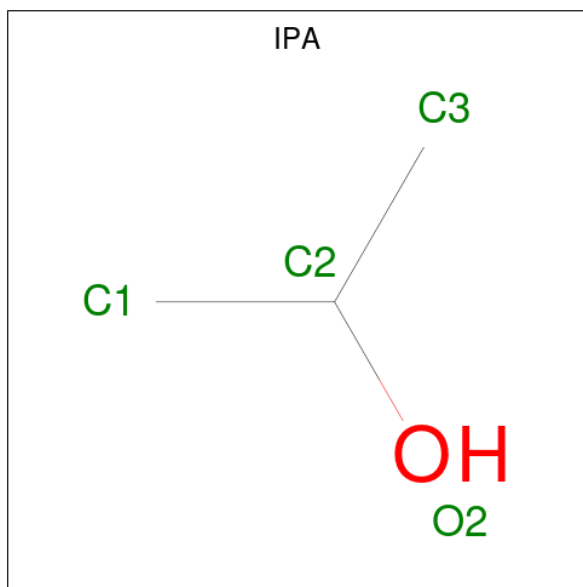
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Mg	0	0
			2	2		
3	B	2	Total	Mg	0	0
			2	2		

- Molecule 4 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: C₄H₁₀O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			7	4	3		
4	B	1	Total	C	O	0	0
			7	4	3		
4	B	1	Total	C	O	0	0
			7	4	3		

- Molecule 5 is ISOPROPYL ALCOHOL (CCD ID: IPA) (formula: C₃H₈O).



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

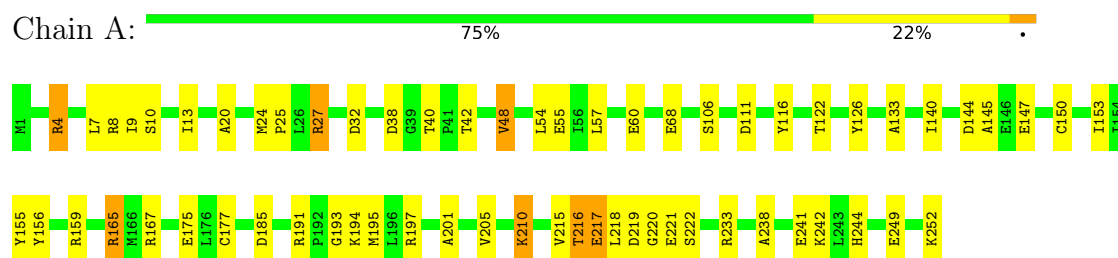
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	35	Total	O	0	0
			35	35		
6	B	36	Total	O	0	0
			36	36		

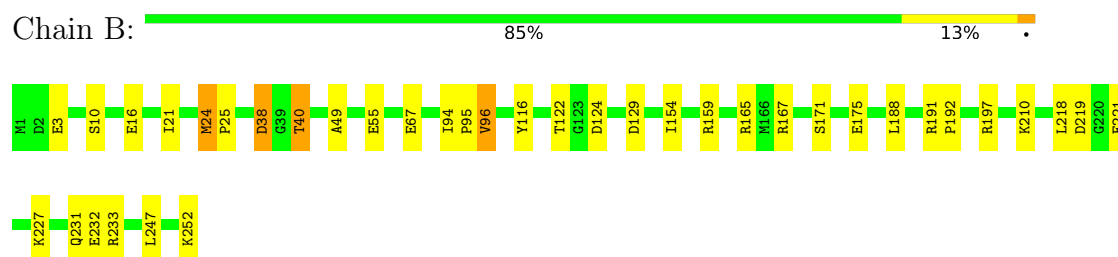
3 Residue-property plots

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: Fructose-1,6-bisphosphatase/inositol-1-monophosphatase



- Molecule 1: Fructose-1,6-bisphosphatase/inositol-1-monophosphatase



4 Data and refinement statistics

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α , β , γ	90.31Å 90.31Å 101.41Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	36.48 – 2.70 36.48 – 2.70	Depositor EDS
% Data completeness (in resolution range)	96.5 (36.48-2.70) 96.6 (36.48-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.11 (at 2.68Å)	Xtriage
Refinement program	PHENIX 1.8_1069	Depositor
R, R_{free}	0.198 , 0.247 0.205 , 0.248	Depositor DCC
R_{free} test set	1261 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	58.2	Xtriage
Anisotropy	0.492	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 30.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Estimated twinning fraction	0.009 for -h,-k,l 0.478 for h,-h-k,-l 0.009 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4060	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.46% 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, MG, IPA, PEG

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	0.64	2/2017 (0.1%)	0.92	3/2713 (0.1%)
1	B	0.63	1/2004 (0.0%)	0.89	6/2698 (0.2%)
All	All	0.63	3/4021 (0.1%)	0.90	9/5411 (0.2%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	8[A]	ARG	N-CA	7.97	1.55	1.46
1	A	8[B]	ARG	N-CA	7.97	1.55	1.46
1	B	96	VAL	C-O	-5.36	1.17	1.24

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	7	LEU	CA-C-N	8.32	131.76	120.44
1	A	7	LEU	C-N-CA	8.32	131.76	120.44
1	B	40	THR	CA-C-N	-5.99	113.77	119.76
1	B	40	THR	C-N-CA	-5.99	113.77	119.76
1	A	217	GLU	N-CA-C	-5.91	102.86	110.19
1	B	154	ILE	N-CA-C	5.64	116.40	108.17
1	B	24	MET	CA-C-N	5.10	125.00	119.85
1	B	24	MET	C-N-CA	5.10	125.00	119.85
1	B	232	GLU	N-CA-C	5.03	117.73	110.28

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	1978	0	1955	51	0
1	B	1968	0	1943	33	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	A	7	0	9	1	0
4	B	14	0	18	1	0
5	B	8	0	16	17	0
6	A	35	0	0	5	0
6	B	36	0	0	2	0
All	All	4060	0	3941	84	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 11.

All (84) 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:217:GLU:HG3	1:A:218:LEU:H	1.08	1.11
1:B:38:ASP:OD1	1:B:40:THR:HG22	1.56	1.05
1:A:165:ARG:HH21	5:B:307:IPA:H12	1.12	1.04
1:B:167:ARG:NH2	5:B:301:IPA:C3	2.23	1.02
1:B:122:THR:HG21	6:B:407:HOH:O	1.60	1.00
1:A:217:GLU:HG3	1:A:218:LEU:N	1.72	0.98
1:A:210:LYS:HE2	6:A:1034:HOH:O	1.63	0.96
1:B:167:ARG:HH21	5:B:301:IPA:C3	1.80	0.92
1:B:167:ARG:HH21	5:B:301:IPA:H31	1.35	0.89
1:B:167:ARG:NH2	5:B:301:IPA:H31	1.89	0.85
1:A:165:ARG:HH21	5:B:307:IPA:C1	1.91	0.83
1:A:165:ARG:NH2	5:B:307:IPA:H12	1.95	0.81
1:A:219:ASP:OD1	1:A:221[B]:GLU:CG	2.33	0.77
1:A:144:ASP:C	1:A:144:ASP:OD1	2.30	0.74
1:B:167:ARG:NH2	5:B:301:IPA:H32	2.02	0.73
1:B:167:ARG:HH22	5:B:301:IPA:H32	1.54	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:217:GLU:CG	1:A:218:LEU:H	1.97	0.71
1:A:194:LYS:HG2	1:A:233:ARG:HB3	1.71	0.71
1:B:165:ARG:HH21	5:B:301:IPA:H12	1.54	0.71
1:B:38:ASP:CG	1:B:40:THR:HG22	2.15	0.71
1:B:210:LYS:NZ	6:B:402:HOH:O	2.25	0.69
1:A:219:ASP:OD1	1:A:221[B]:GLU:HG3	1.94	0.68
1:A:27:ARG:NH1	6:A:1001:HOH:O	2.28	0.66
1:B:231:GLN:OE1	1:B:231:GLN:N	2.30	0.65
1:A:216:THR:OG1	1:A:220:GLY:HA2	1.99	0.63
1:B:124:ASP:OD1	4:B:305:PEG:H32	1.97	0.63
1:B:165:ARG:HE	5:B:301:IPA:H33	1.63	0.62
1:A:194:LYS:HE2	1:A:233:ARG:C	2.27	0.60
1:B:3:GLU:OE1	1:B:129:ASP:HB2	2.03	0.58
1:B:231:GLN:N	1:B:231:GLN:CD	2.65	0.54
1:A:20:ALA:HB1	1:A:48:VAL:HG13	1.91	0.53
1:A:13:ILE:HD11	1:A:57:LEU:HG	1.90	0.52
1:A:216:THR:HG23	1:A:244:HIS:CE1	2.44	0.52
1:B:96:VAL:HG12	1:B:96:VAL:O	2.08	0.52
1:A:249:GLU:O	1:A:252:LYS:HB2	2.10	0.52
1:A:122:THR:HG21	1:B:95:PRO:HG3	1.93	0.51
1:A:10:SER:HB3	1:A:116:TYR:CD2	2.46	0.50
1:A:194:LYS:HE2	1:A:233:ARG:HB2	1.91	0.50
1:A:42:THR:HG23	1:A:68:GLU:HG2	1.92	0.50
1:A:197:ARG:HD3	6:A:1030:HOH:O	2.12	0.49
1:A:197:ARG:HG2	1:A:233:ARG:NH1	2.28	0.48
1:A:144:ASP:OD1	1:A:145:ALA:N	2.46	0.48
1:A:106:SER:HB3	1:A:111:ASP:HB2	1.95	0.48
1:B:197:ARG:HG2	1:B:233:ARG:NH1	2.27	0.48
1:A:167:ARG:NH2	5:B:307:IPA:C3	2.76	0.48
1:B:38:ASP:HB3	1:B:67:GLU:O	2.13	0.48
1:A:42:THR:OG1	1:A:68:GLU:HG3	2.14	0.48
1:A:167:ARG:HH21	5:B:307:IPA:C3	2.27	0.47
1:B:188:LEU:HD11	1:B:247:LEU:HD13	1.96	0.47
1:B:171:SER:O	1:B:175:GLU:HG3	2.15	0.47
1:A:201:ALA:O	1:A:205:VAL:HG23	2.15	0.46
1:B:165:ARG:HE	5:B:301:IPA:C3	2.26	0.46
1:A:215:VAL:HG13	1:A:238:ALA:HB2	1.98	0.46
1:B:10:SER:HB3	1:B:116:TYR:CD2	2.51	0.45
1:A:167:ARG:NH2	5:B:307:IPA:H32	2.31	0.45
1:B:24:MET:HA	1:B:25:PRO:HD3	1.79	0.45
1:A:159:ARG:NH1	6:A:1004:HOH:O	2.36	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:231:GLN:CD	1:B:231:GLN:H	2.19	0.45
1:A:191:ARG:NH2	1:A:195:MET:HB2	2.31	0.45
1:A:159:ARG:HD2	6:A:1004:HOH:O	2.16	0.45
1:B:221:GLU:H	1:B:221:GLU:CD	2.22	0.44
1:A:133:ALA:C	1:A:140:ILE:HG12	2.42	0.44
1:B:227:LYS:HB3	1:B:227:LYS:HE2	1.78	0.44
1:A:9:ILE:HD11	1:A:60:GLU:CD	2.42	0.44
1:A:38:ASP:OD2	1:A:40:THR:OG1	2.36	0.44
1:A:219:ASP:OD1	1:A:221[B]:GLU:OE2	2.36	0.44
1:A:4:ARG:HE	1:A:4:ARG:HB3	1.29	0.43
1:B:165:ARG:NE	5:B:301:IPA:H33	2.30	0.43
1:A:54:LEU:HD23	1:A:54:LEU:HA	1.81	0.43
1:A:193:GLY:HA3	1:A:195:MET:HE2	2.01	0.43
4:A:904:PEG:H42	4:A:904:PEG:H21	1.56	0.43
1:A:155:TYR:HA	1:A:156:TYR:HA	1.86	0.42
1:A:219:ASP:OD1	1:A:221[B]:GLU:HG2	2.15	0.42
1:A:32:ASP:N	1:A:32:ASP:OD1	2.51	0.42
1:A:24:MET:HA	1:A:25:PRO:HD3	1.89	0.42
1:B:159:ARG:HE	1:B:159:ARG:HB2	1.63	0.42
1:B:167:ARG:HH22	5:B:301:IPA:C3	2.09	0.41
1:A:242:LYS:HA	1:A:242:LYS:HD3	1.90	0.41
1:A:126:TYR:CD1	1:A:177:CYS:HB3	2.55	0.41
1:B:21:ILE:HD11	1:B:49:ALA:HB2	2.03	0.41
1:A:150:CYS:HB2	1:A:185:ASP:HB2	2.03	0.40
1:A:153:ILE:HD13	1:A:175:GLU:HB3	2.03	0.40
1:B:191:ARG:HA	1:B:192:PRO:HD3	1.96	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	252/252 (100%)	237 (94%)	15 (6%)	0	100	100
1	B	251/252 (100%)	242 (96%)	9 (4%)	0	100	100
All	All	503/504 (100%)	479 (95%)	24 (5%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	206/204 (101%)	197 (96%)	9 (4%)	25	53
1	B	205/204 (100%)	198 (97%)	7 (3%)	32	62
All	All	411/408 (101%)	395 (96%)	16 (4%)	28	57

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	ARG
1	A	27	ARG
1	A	48	VAL
1	A	55	GLU
1	A	147	GLU
1	A	165	ARG
1	A	210	LYS
1	A	216	THR
1	A	241	GLU
1	B	16	GLU
1	B	38	ASP
1	B	55	GLU
1	B	94	ILE
1	B	218	LEU
1	B	219	ASP
1	B	252	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such

sidechains are listed below:

Mol	Chain	Res	Type
1	B	244	HIS

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 ⓘ

Of 11 ligands modelled in this entry, 4 are monoatomic - leaving 7 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	IPA	B	301	-	3,3,3	0.46	0	3,3,3	0.12	0
2	SO4	A	901	3	4,4,4	0.32	0	6,6,6	0.33	0
2	SO4	B	302	3	4,4,4	0.31	0	6,6,6	0.44	0
4	PEG	A	904	-	6,6,6	1.08	1 (16%)	5,5,5	0.64	0
5	IPA	B	307	-	3,3,3	0.53	0	3,3,3	0.32	0
4	PEG	B	305	-	6,6,6	0.82	0	5,5,5	0.77	0
4	PEG	B	306	-	6,6,6	1.05	1 (16%)	5,5,5	0.62	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
4	PEG	A	904	-	-	4/4/4/4	-
4	PEG	B	305	-	-	2/4/4/4	-
4	PEG	B	306	-	-	3/4/4/4	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	306	PEG	O1-C1	-2.10	1.31	1.42
4	A	904	PEG	O1-C1	-2.03	1.31	1.42

There are no bond angle outliers.

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	904	PEG	C4-C3-O2-C2
4	B	305	PEG	C1-C2-O2-C3
4	A	904	PEG	O1-C1-C2-O2
4	B	305	PEG	O2-C3-C4-O4
4	B	306	PEG	O2-C3-C4-O4
4	B	306	PEG	O1-C1-C2-O2
4	A	904	PEG	O2-C3-C4-O4
4	B	306	PEG	C4-C3-O2-C2
4	A	904	PEG	C1-C2-O2-C3

There are no ring outliers.

4 monomers are involved in 19 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	301	IPA	11	0
4	A	904	PEG	1	0
5	B	307	IPA	6	0
4	B	305	PEG	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 [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	252/252 (100%)	-1.45	0 100 100	26, 40, 60, 76	2 (0%)
1	B	252/252 (100%)	-1.46	0 100 100	29, 41, 59, 81	1 (0%)
All	All	504/504 (100%)	-1.46	0 100 100	26, 40, 60, 81	3 (0%)

There are no RSRZ outliers to report.

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(Å ²)	Q<0.9
5	IPA	B	307	4/4	0.96	0.14	40,42,43,49	0
4	PEG	B	306	7/7	0.97	0.07	51,58,62,71	0
5	IPA	B	301	4/4	0.97	0.11	39,39,44,44	0
4	PEG	B	305	7/7	0.97	0.09	48,55,59,60	0
4	PEG	A	904	7/7	0.98	0.10	47,50,60,61	0
2	SO4	A	901	5/5	0.99	0.06	36,41,50,62	0
2	SO4	B	302	5/5	0.99	0.05	37,43,54,57	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	MG	A	903	1/1	0.99	0.04	45,45,45,45	0
3	MG	B	303	1/1	0.99	0.05	25,25,25,25	0
3	MG	B	304	1/1	0.99	0.07	44,44,44,44	0
3	MG	A	902	1/1	1.00	0.01	24,24,24,24	0

6.5 Other polymers [i](#)

There are no such residues in this entry.