



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 24, 2023 – 01:52 PM EST

PDB ID : 3JZL
Title : CRYSTAL STRUCTURE OF A PUTATIVE CYSTATHIONINE BETA-LYASE INVOLVED IN ALUMINUM RESISTANCE (LMOF2365_1314) FROM LISTERIA MONOCYTOGENES STR. 4B F2365 AT 1.91 Å RESOLUTION
Authors : Joint Center for Structural Genomics (JCSG)
Deposited on : 2009-09-23
Resolution : 1.91 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.31.2
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.31.2

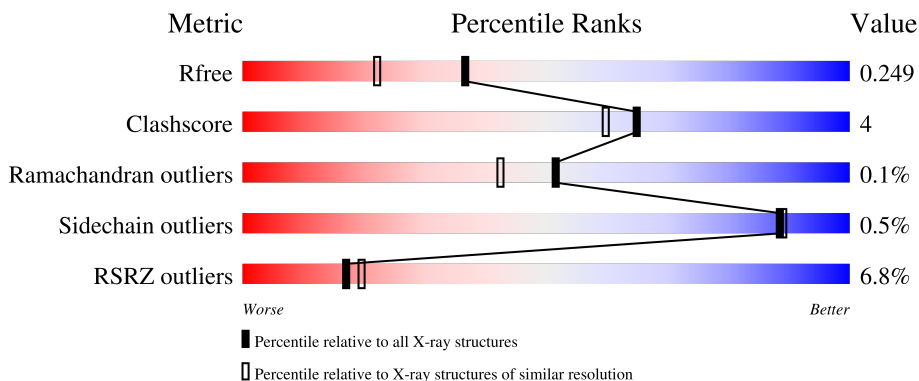
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.91 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}	130704	7937 (1.94-1.90)
Clashscore	141614	8644 (1.94-1.90)
Ramachandran outliers	138981	8530 (1.94-1.90)
Sidechain outliers	138945	8530 (1.94-1.90)
RSRZ outliers	127900	7793 (1.94-1.90)

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	409	
1	B	409	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6397 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 Putative cystathionine beta-lyase involved in aluminum resistance.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace	
			Total	C	N	O	P	S				Se
1	A	405	3109	1997	509	592	1	2	8	0	6	0
1	B	394	2990	1916	494	571	1	2	6	0	10	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	expression tag	UNP Q720C5
B	0	GLY	-	expression tag	UNP Q720C5

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



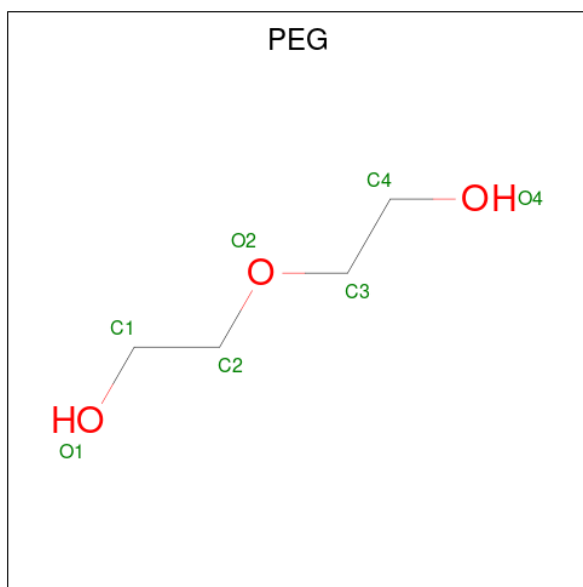
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	6	3	3	0	0

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

- Molecule 3 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C₄H₁₀O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	4	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	154	Total	O	0	0
			154	154		
4	B	120	Total	O	0	0
			120	120		

4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	54.41Å 153.17Å 183.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.44 – 1.91 39.43 – 1.91	Depositor EDS
% Data completeness (in resolution range)	97.9 (39.44-1.91) 97.9 (39.43-1.91)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.89 (at 1.91Å)	Xtrriage
Refinement program	REFMAC 5.5.0072, PHENIX	Depositor
R, R_{free}	0.210 , 0.241 0.221 , 0.249	Depositor DCC
R_{free} test set	2967 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	23.3	Xtrriage
Anisotropy	0.927	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 57.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6397	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.89% 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: GOL, LLP, 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	0.65	0/3160	0.79	1/4284 (0.0%)
1	B	0.66	0/3044	0.83	3/4123 (0.1%)
All	All	0.65	0/6204	0.81	4/8407 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	291	ARG	NE-CZ-NH1	6.71	123.66	120.30
1	A	291	ARG	NE-CZ-NH1	6.66	123.63	120.30
1	B	252	ARG	NE-CZ-NH2	-6.13	117.24	120.30
1	B	305	ASP	CB-CG-OD1	5.18	122.97	118.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3109	0	3029	21	0
1	B	2990	0	2890	29	0
2	A	6	0	8	0	0
2	B	12	0	16	1	0
3	A	6	0	7	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	154	0	0	1	0
4	B	120	0	0	1	0
All	All	6397	0	5950	50	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 4.

All (50) 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:331:ALA:HB3	4:B:447:HOH:O	1.78	0.82
1:B:114[B]:LEU:HD23	1:B:117:ILE:HD12	1.68	0.74
1:B:331:ALA:HB2	1:B:385:LEU:CD1	2.19	0.72
1:B:331:ALA:HB2	1:B:385:LEU:HD11	1.73	0.71
1:B:329:ALA:O	1:B:333:ALA:HB2	1.99	0.61
1:B:268:SER:HB3	1:B:272:MSE:HE2	1.86	0.57
1:A:189[A]:LYS:NZ	1:A:197:VAL:HG23	2.23	0.53
1:B:93:LEU:HD22	1:B:103:LEU:HD22	1.92	0.52
1:B:82:ILE:O	1:B:236:GLY:HA2	2.10	0.52
1:B:331:ALA:CB	1:B:385:LEU:HD11	2.39	0.52
1:B:111:TYR:CE1	1:B:114[B]:LEU:HG	2.46	0.51
1:B:90:SER:HB2	1:B:117:ILE:HD13	1.93	0.50
1:B:330:PHE:HE1	1:B:387:VAL:HG21	1.76	0.50
1:A:114:LEU:HD23	1:A:117:ILE:HD12	1.93	0.49
1:A:234:LYS:HA	1:A:366[A]:ILE:HD13	1.95	0.49
1:A:227:ASN:HA	1:A:371:LEU:HD22	1.94	0.49
1:A:268:SER:HB3	1:A:272:MSE:HE2	1.95	0.49
1:B:227:ASN:HA	1:B:371:LEU:HD22	1.94	0.49
1:B:11:VAL:HG13	1:B:295:ALA:HB1	1.95	0.49
1:B:331:ALA:HB2	1:B:385:LEU:HD13	1.96	0.48
1:B:297[A]:LEU:HD22	1:B:404:ILE:HD11	1.95	0.48
1:A:4:ILE:HG23	1:A:7:ILE:H	1.79	0.47
1:A:83:ILE:HG22	4:A:491:HOH:O	2.14	0.47
1:B:47:PHE:CD2	1:B:270:LEU:HD13	2.49	0.47
1:A:189[A]:LYS:HZ2	1:A:197:VAL:HG23	1.79	0.47
1:A:64:GLU:HG2	1:A:78:VAL:HG23	1.97	0.47
1:A:338:SER:O	1:A:342:ALA:HB2	2.15	0.46
1:B:120:ILE:HD11	1:B:135:TYR:HE2	1.81	0.46
1:A:90:SER:HB2	1:A:117:ILE:HD13	1.99	0.45
1:A:224:LEU:O	1:A:229:GLY:HA3	2.16	0.45
1:B:289:GLY:HA3	1:B:396:ILE:HG21	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:199:VAL:HB	1:B:220:ILE:HG22	1.98	0.44
1:B:114[B]:LEU:HD23	1:B:114[B]:LEU:HA	1.74	0.44
1:A:268:SER:CB	1:A:272:MSE:HE2	2.47	0.43
1:B:330:PHE:CD2	1:B:403:ALA:HB1	2.53	0.43
1:A:375:ALA:HA	1:A:386:TYR:O	2.19	0.43
1:A:199:VAL:HB	1:A:220:ILE:HG22	2.01	0.43
1:A:185:ILE:HG12	1:A:197:VAL:HG11	2.01	0.42
1:A:111:TYR:CZ	1:A:226:LLP:H6	2.55	0.42
1:B:114[B]:LEU:CD2	1:B:117:ILE:HD12	2.43	0.42
1:B:185:ILE:HG12	1:B:197:VAL:HG11	2.02	0.41
1:B:289:GLY:HA3	1:B:396:ILE:HD13	2.03	0.41
1:A:279:ALA:HB3	1:A:280:PRO:HD3	2.01	0.41
1:A:111:TYR:CE2	1:A:113:THR:OG1	2.70	0.41
1:A:289:GLY:HA3	1:A:396:ILE:HD13	2.02	0.41
1:A:335[B]:GLN:OE1	1:A:363:GLY:N	2.54	0.41
1:B:226:LLP:OP2	2:B:410:GOL:O3	2.37	0.41
1:B:385:LEU:HD23	1:B:387:VAL:CG2	2.51	0.41
1:B:300:PHE:CE1	1:B:404:ILE:HG21	2.57	0.40
1:B:114[A]:LEU:HD12	1:B:117:ILE:HB	2.04	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	408/409 (100%)	395 (97%)	13 (3%)	0	100	100
1	B	395/409 (97%)	374 (95%)	20 (5%)	1 (0%)	41	31
All	All	803/818 (98%)	769 (96%)	33 (4%)	1 (0%)	51	42

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	326	LYS

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	314/323 (97%)	313 (100%)	1 (0%)	92	93
1	B	296/323 (92%)	294 (99%)	2 (1%)	84	83
All	All	610/646 (94%)	607 (100%)	3 (0%)	88	89

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	228	PRO
1	B	105	TYR
1	B	227	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

2 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	LLP	A	226	1	23,24,25	1.75	3 (13%)	25,32,34	1.54	5 (20%)
1	LLP	B	226	1	23,24,25	1.78	5 (21%)	25,32,34	1.61	4 (16%)

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	LLP	A	226	1	-	1/16/17/19	0/1/1/1
1	LLP	B	226	1	-	2/16/17/19	0/1/1/1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	226	LLP	O3-C3	-5.98	1.23	1.37
1	A	226	LLP	O3-C3	-5.78	1.23	1.37
1	A	226	LLP	C4-C4'	2.89	1.52	1.46
1	B	226	LLP	C4-C4'	2.80	1.52	1.46
1	B	226	LLP	C4'-NZ	2.14	1.34	1.27
1	B	226	LLP	P-OP3	-2.11	1.46	1.54
1	A	226	LLP	C4-C5	-2.11	1.39	1.42
1	B	226	LLP	C4-C5	-2.07	1.39	1.42

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	226	LLP	OP4-C5'-C5	4.62	118.15	109.35
1	A	226	LLP	C5-C4-C4'	-3.13	116.42	121.56
1	B	226	LLP	C5-C4-C4'	-2.90	116.79	121.56
1	A	226	LLP	CE-NZ-C4'	-2.87	110.09	118.90
1	A	226	LLP	OP4-P-OP1	-2.75	98.77	106.47
1	A	226	LLP	OP4-C5'-C5	2.68	114.46	109.35
1	A	226	LLP	C3-C4-C4'	2.50	125.07	120.41
1	B	226	LLP	C3-C4-C4'	2.47	125.02	120.41
1	B	226	LLP	CE-NZ-C4'	-2.04	112.63	118.90

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	B	226	LLP	C4-C4'-NZ-CE
1	A	226	LLP	C4-C4'-NZ-CE
1	B	226	LLP	C5'-OP4-P-OP3

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	226	LLP	1	0
1	B	226	LLP	1	0

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 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
3	PEG	A	410	-	5,5,6	0.53	0	4,4,5	0.45	0
2	GOL	B	409	-	5,5,5	0.47	0	5,5,5	0.30	0
2	GOL	B	410	-	5,5,5	0.50	0	5,5,5	0.45	0
2	GOL	A	409	-	5,5,5	0.44	0	5,5,5	0.30	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. '2' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	PEG	A	410	-	-	1/3/3/4	-
2	GOL	B	409	-	-	1/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GOL	B	410	-	-	4/4/4/4	-
2	GOL	A	409	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	410	GOL	O1-C1-C2-C3
2	B	410	GOL	C1-C2-C3-O3
2	B	409	GOL	C1-C2-C3-O3
2	B	410	GOL	O2-C2-C3-O3
3	A	410	PEG	O1-C1-C2-O2
2	B	410	GOL	O1-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	410	GOL	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	396/409 (96%)	0.36	13 (3%) 46 49	13, 22, 35, 48	0
1	B	385/409 (94%)	0.69	40 (10%) 6 7	13, 23, 35, 49	0
All	All	781/818 (95%)	0.52	53 (6%) 17 19	13, 22, 35, 49	0

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	352	PRO	6.9
1	B	329	ALA	5.2
1	B	330	PHE	4.9
1	B	385	LEU	4.7
1	B	352	PRO	4.7
1	B	350	TYR	4.5
1	B	322	HIS	4.4
1	B	321	PHE	4.2
1	B	344	VAL	4.1
1	B	359	ILE	4.0
1	A	408	VAL	3.9
1	B	348	GLY	3.9
1	B	349	ALA	3.7
1	A	348	GLY	3.7
1	B	353	GLY	3.6
1	B	354	TYR	3.5
1	A	349	ALA	3.5
1	B	328	VAL	3.5
1	B	170	TYR	3.5
1	B	331	ALA	3.4
1	A	278	LEU	3.3
1	B	347	ILE	3.2
1	B	326	LYS	3.2
1	B	371	LEU	3.0

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Mol	Chain	Res	Type	RSRZ
1	B	377	GLY	3.0
1	B	6	ALA	3.0
1	A	350	TYR	3.0
1	B	356	ASP	3.0
1	B	171	ALA	2.9
1	B	282	VAL	2.9
1	A	45	PHE	2.8
1	B	346	PRO	2.7
1	B	278	LEU	2.7
1	B	190	ASN	2.7
1	B	332	GLN	2.6
1	B	83	ILE	2.6
1	B	114[A]	LEU	2.5
1	B	122	LYS	2.5
1	B	403	ALA	2.5
1	B	56	ASP	2.5
1	A	371	LEU	2.4
1	A	391	LEU	2.3
1	A	56	ASP	2.3
1	A	347	ILE	2.3
1	B	141	LEU	2.2
1	B	320	SER	2.2
1	B	230	GLY	2.2
1	A	231	GLY	2.1
1	B	232	LEU	2.1
1	B	375	ALA	2.1
1	B	407	ILE	2.1
1	A	345	LEU	2.0
1	B	224	LEU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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	LLP	A	226	24/25	0.94	0.13	16,27,31,32	0
1	LLP	B	226	24/25	0.95	0.15	17,26,30,31	0

6.3 Carbohydrates [i](#)

There are no monosaccharides 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
2	GOL	B	409	6/6	0.84	0.16	40,47,48,53	0
3	PEG	A	410	6/7	0.84	0.18	55,56,57,59	0
2	GOL	B	410	6/6	0.88	0.22	67,68,69,69	0
2	GOL	A	409	6/6	0.93	0.12	26,32,34,34	0

6.5 Other polymers [i](#)

There are no such residues in this entry.