



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 04:06 AM UTC

PDB ID : 3FVX / pdb_00003fvx
Title : Human kynurenine aminotransferase I in complex with tris
Authors : Han, Q.; Robinson, H.; Cai, T.; Tagle, D.A.; Li, J.
Deposited on : 2009-01-16
Resolution : 1.50 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

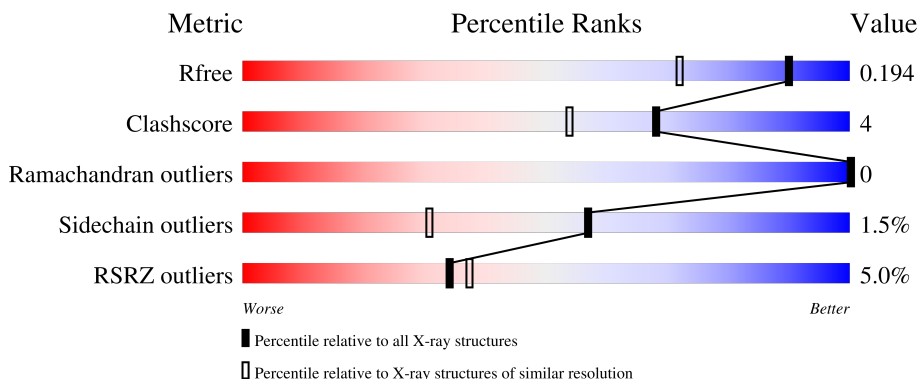
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

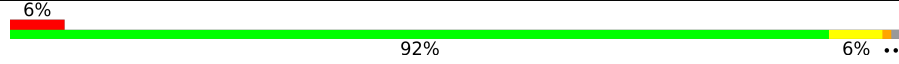
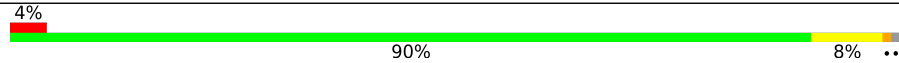
The reported resolution of this entry is 1.50 Å.

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	4037 (1.50-1.50)
Clashscore	190562	4235 (1.50-1.50)
Ramachandran outliers	187476	4153 (1.50-1.50)
Sidechain outliers	187428	4150 (1.50-1.50)
RSRZ outliers	180081	4039 (1.50-1.50)

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	422	 6% 92% 6% ..
1	B	422	 4% 90% 8% ..

2 Entry composition [i](#)

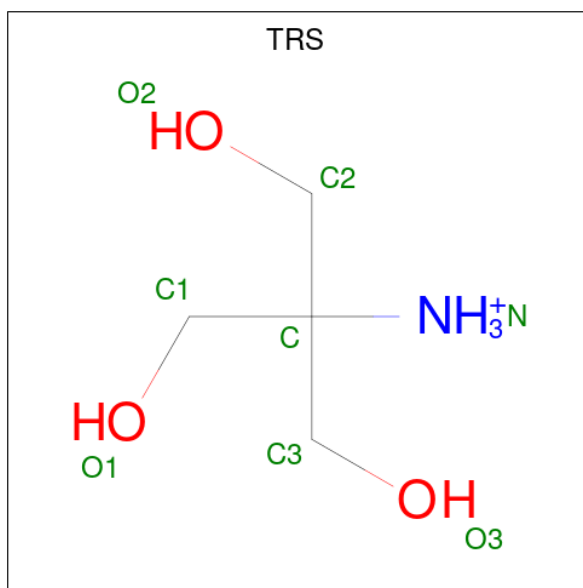
There are 4 unique types of molecules in this entry. The entry contains 7412 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Kynurenine--oxoglutarate transaminase 1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	P	S			
1	A	419	Total 3369	C 2175	N 565	O 609	P 1	S 19	0	0	0
1	B	419	Total 3369	C 2175	N 565	O 609	P 1	S 19	0	0	0

- Molecule 2 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (CCD ID: TRS) (formula: C₄H₁₂NO₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
2	A	1	Total 8	C 4	N 1	O 3	0	0

- Molecule 3 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total 1	Na 1	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Na	0	0
			1	1		

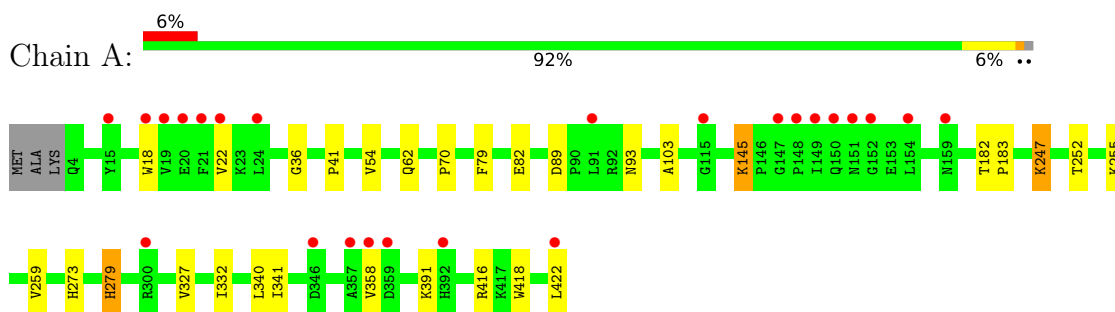
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	342	Total	O	0	0
			342	342		
4	B	322	Total	O	0	0
			322	322		

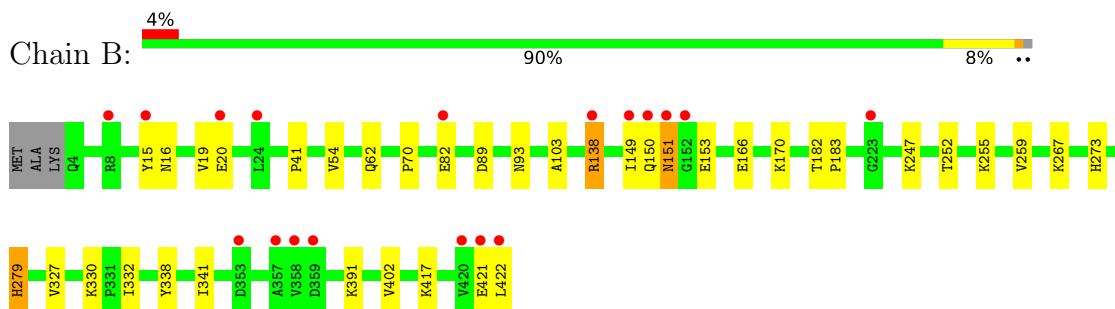
3 Residue-property plots [i](#)

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: Kynurenine--oxoglutarate transaminase 1



- Molecule 1: Kynurenine--oxoglutarate transaminase 1



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	102.88Å 107.44Å 81.69Å 90.00° 112.87° 90.00°	Depositor
Resolution (Å)	29.04 – 1.50 29.04 – 1.50	Depositor EDS
% Data completeness (in resolution range)	98.6 (29.04-1.50) 98.6 (29.04-1.50)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.00 (at 1.50Å)	Xtrriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.178 , 0.200 (Not available) , 0.194	Depositor DCC
R_{free} test set	6430 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	14.3	Xtrriage
Anisotropy	0.056	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 37.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7412	wwPDB-VP
Average B, all atoms (Å ²)	16.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.57% 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: LLP, NA, TRS

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	0.52	0/3441	0.76	4/4671 (0.1%)
1	B	0.53	0/3441	0.75	2/4671 (0.0%)
All	All	0.53	0/6882	0.76	6/9342 (0.1%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	70	PRO	N-CA-C	5.89	117.88	110.70
1	B	70	PRO	N-CA-C	5.59	117.52	110.70
1	B	41	PRO	N-CA-C	5.56	116.53	110.58
1	A	41	PRO	N-CA-C	5.27	116.22	110.58
1	A	145	LYS	CA-C-N	5.20	125.12	119.76
1	A	145	LYS	C-N-CA	5.20	125.12	119.76

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	3369	0	3295	27	0
1	B	3369	0	3295	27	0
2	A	8	0	12	5	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	342	0	0	4	0
4	B	322	0	0	6	0
All	All	7412	0	6602	48	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 (48) 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:A:423:TRS:H21	4:A:661:HOH:O	1.63	0.97
1:A:327:VAL:HG11	1:A:418:TRP:CE2	2.05	0.91
1:B:89:ASP:H	1:B:93:ASN:HD22	1.21	0.87
1:A:89:ASP:H	1:A:93:ASN:HD22	1.19	0.85
1:A:391:LYS:HG3	4:A:656:HOH:O	1.77	0.83
1:B:391:LYS:HG3	4:B:665:HOH:O	1.86	0.76
1:A:327:VAL:HG12	1:A:327:VAL:O	1.93	0.67
1:B:89:ASP:H	1:B:93:ASN:ND2	1.93	0.65
1:A:89:ASP:H	1:A:93:ASN:ND2	1.94	0.65
1:A:327:VAL:HG11	1:A:418:TRP:CD2	2.36	0.59
1:A:247:LLP:C4'	2:A:423:TRS:HN1	2.18	0.56
1:B:15:TYR:HE1	1:B:20:GLU:OE1	1.89	0.55
1:A:273:HIS:HD2	4:A:508:HOH:O	1.91	0.53
1:B:20:GLU:HG3	4:B:723:HOH:O	2.09	0.53
1:A:79:PHE:O	1:A:82:GLU:HG2	2.11	0.51
1:A:327:VAL:CG1	1:A:418:TRP:CE2	2.89	0.50
1:B:150:GLN:O	1:B:153:GLU:HG2	2.12	0.50
1:A:332:ILE:HD13	1:A:341:ILE:HG13	1.94	0.49
1:A:327:VAL:O	1:A:327:VAL:CG1	2.60	0.49
1:B:267:LYS:NZ	4:B:685:HOH:O	2.41	0.49
1:A:252:THR:H	1:B:62:GLN:NE2	2.12	0.48
1:B:332:ILE:HD13	1:B:341:ILE:HG13	1.96	0.48
1:A:54:VAL:HG21	1:B:54:VAL:HG21	1.96	0.48
1:B:273:HIS:HD2	4:B:477:HOH:O	1.97	0.47
1:B:15:TYR:CE1	1:B:20:GLU:OE1	2.66	0.47
1:A:62:GLN:NE2	1:B:252:THR:H	2.13	0.46
1:A:255:LYS:NZ	1:B:279:HIS:HD2	2.14	0.46
1:A:103:ALA:HB1	1:A:259:VAL:HG22	1.98	0.45
1:A:182:THR:HA	1:A:183:PRO:C	2.42	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:279:HIS:HD2	1:B:255:LYS:NZ	2.15	0.45
1:B:103:ALA:HB1	1:B:259:VAL:HG22	1.98	0.45
1:A:36:GLY:HA3	2:A:423:TRS:H31	2.00	0.44
1:B:182:THR:HA	1:B:183:PRO:C	2.43	0.44
1:A:416:ARG:NH2	4:A:556:HOH:O	2.51	0.43
1:B:138:ARG:HA	1:B:138:ARG:HE	1.82	0.43
1:B:330:LYS:NZ	4:B:701:HOH:O	2.51	0.43
1:B:166:GLU:O	1:B:170:LYS:HG2	2.18	0.43
1:A:340:LEU:HD12	1:A:340:LEU:C	2.44	0.43
1:A:18:TRP:O	1:A:22:VAL:HG23	2.19	0.42
1:B:338:TYR:CD1	1:B:402:VAL:HG21	2.55	0.42
1:B:16:ASN:O	1:B:19:VAL:HG22	2.19	0.42
1:A:252:THR:H	1:B:62:GLN:HE22	1.66	0.42
1:A:247:LLP:C4'	2:A:423:TRS:N	2.83	0.42
1:A:62:GLN:HE22	1:B:252:THR:H	1.68	0.41
1:B:151:ASN:HD22	1:B:151:ASN:HA	1.67	0.41
1:B:327:VAL:HA	1:B:422:LEU:HD11	2.02	0.41
1:B:19:VAL:HG23	4:B:723:HOH:O	2.21	0.40
1:A:247:LLP:O3	2:A:423:TRS:N	2.55	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	416/422 (99%)	405 (97%)	11 (3%)	0	100	100
1	B	416/422 (99%)	405 (97%)	11 (3%)	0	100	100
All	All	832/844 (99%)	810 (97%)	22 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	365/367 (100%)	361 (99%)	4 (1%)	65	41
1	B	365/367 (100%)	358 (98%)	7 (2%)	50	22
All	All	730/734 (100%)	719 (98%)	11 (2%)	57	31

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	145	LYS
1	A	279	HIS
1	A	358	VAL
1	A	422	LEU
1	B	82	GLU
1	B	138	ARG
1	B	149	ILE
1	B	151	ASN
1	B	279	HIS
1	B	417	LYS
1	B	421	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	62	GLN
1	A	93	ASN
1	A	109	GLN
1	A	159	ASN
1	A	225	GLN
1	A	273	HIS
1	A	279	HIS
1	A	283	GLN
1	A	296	GLN
1	A	311	GLN
1	A	395	HIS

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Mol	Chain	Res	Type
1	A	409	GLN
1	B	62	GLN
1	B	93	ASN
1	B	109	GLN
1	B	151	ASN
1	B	204	GLN
1	B	273	HIS
1	B	279	HIS
1	B	283	GLN
1	B	395	HIS

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	B	247	1	23,24,25	1.67	6 (26%)	25,32,34	1.49	4 (16%)
1	LLP	A	247	1	23,24,25	1.77	6 (26%)	25,32,34	1.38	2 (8%)

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	B	247	1	-	6/16/17/19	0/1/1/1
1	LLP	A	247	1	-	5/16/17/19	0/1/1/1

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	247	LLP	O3-C3	-5.01	1.25	1.36
1	A	247	LLP	O3-C3	-4.98	1.25	1.36
1	A	247	LLP	CE-NZ	3.29	1.54	1.46
1	A	247	LLP	C4-C4'	3.03	1.53	1.46
1	B	247	LLP	C4-C4'	2.68	1.52	1.46
1	B	247	LLP	CE-NZ	2.64	1.52	1.46
1	A	247	LLP	C4'-NZ	2.60	1.35	1.27
1	B	247	LLP	C4'-NZ	2.47	1.35	1.27
1	B	247	LLP	C6-N1	2.18	1.38	1.34
1	A	247	LLP	C6-N1	2.14	1.38	1.34
1	A	247	LLP	C2-N1	2.03	1.37	1.33
1	B	247	LLP	C2-N1	2.02	1.37	1.33

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	247	LLP	OP4-C5'-C5	4.28	117.39	109.36
1	B	247	LLP	OP4-C5'-C5	4.02	116.90	109.36
1	B	247	LLP	C5-C6-N1	-2.78	119.31	123.83
1	A	247	LLP	C5-C6-N1	-2.48	119.79	123.83
1	B	247	LLP	C4-C3-C2	-2.30	118.85	120.14
1	B	247	LLP	C4-C4'-NZ	-2.13	114.20	124.04

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	247	LLP	C4-C4'-NZ-CE
1	B	247	LLP	C4-C4'-NZ-CE
1	B	247	LLP	C5-C4-C4'-NZ
1	A	247	LLP	C5-C4-C4'-NZ
1	A	247	LLP	C4-C5-C5'-OP4
1	B	247	LLP	C4-C5-C5'-OP4
1	B	247	LLP	CD-CE-NZ-C4'
1	A	247	LLP	C3-C4-C4'-NZ
1	B	247	LLP	C3-C4-C4'-NZ
1	B	247	LLP	C6-C5-C5'-OP4
1	A	247	LLP	CD-CE-NZ-C4'

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	247	LLP	3	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 2 are monoatomic - leaving 1 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
2	TRS	A	423	-	7,7,7	0.46	0	9,9,9	0.70	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
2	TRS	A	423	-	-	3/9/9/9	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	423	TRS	C1-C-C3-O3
2	A	423	TRS	C2-C-C3-O3
2	A	423	TRS	N-C-C3-O3

There are no ring outliers.

1 monomer is involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	423	TRS	5	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

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	418/422 (99%)	0.17	24 (5%) 29 32	8, 14, 28, 44	0
1	B	418/422 (99%)	0.13	18 (4%) 40 44	9, 14, 25, 41	0
All	All	836/844 (99%)	0.15	42 (5%) 34 37	8, 14, 27, 44	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	148	PRO	6.9
1	A	151	ASN	4.9
1	A	152	GLY	4.8
1	B	151	ASN	4.7
1	A	149	ILE	4.4
1	B	359	ASP	4.2
1	A	19	VAL	4.0
1	A	21	PHE	4.0
1	A	358	VAL	3.7
1	A	15	TYR	3.7
1	B	149	ILE	3.6
1	A	422	LEU	3.4
1	B	82	GLU	3.4
1	A	150	GLN	3.2
1	A	147	GLY	3.1
1	A	115	GLY	3.1
1	B	152	GLY	3.1
1	B	358	VAL	3.0
1	B	150	GLN	2.9
1	A	154	LEU	2.7
1	B	422	LEU	2.7
1	B	353	ASP	2.7
1	A	22	VAL	2.7
1	B	20	GLU	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	359	ASP	2.6
1	A	24	LEU	2.5
1	B	15	TYR	2.4
1	B	357	ALA	2.3
1	A	392	HIS	2.3
1	A	18	TRP	2.3
1	A	20	GLU	2.3
1	B	138	ARG	2.3
1	A	91	LEU	2.2
1	B	223	GLY	2.2
1	B	420	VAL	2.2
1	A	300	ARG	2.1
1	A	357	ALA	2.1
1	B	24	LEU	2.1
1	A	346	ASP	2.1
1	B	8	ARG	2.1
1	B	421	GLU	2.1
1	A	159	ASN	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	247	24/25	0.97	0.07	9,11,18,21	0
1	LLP	B	247	24/25	0.97	0.07	8,12,18,20	0

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(\AA^2)	Q<0.9
2	TRS	A	423	8/8	0.81	0.13	23,26,27,27	0
3	NA	A	424	1/1	0.95	0.15	28,28,28,28	0
3	NA	B	423	1/1	0.98	0.05	24,24,24,24	0

6.5 Other polymers [i](#)

There are no such residues in this entry.