Full wwPDB X-ray Structure Validation Report (i)

Nov 7, 2023 – 10:54 AM EST

PDB ID	:	5W1B
Title	:	Tryptophan indole-lyase
Authors	:	Phillips, R.S.; Wood, Z.A.
Deposited on	:	2017-06-02
Resolution	:	2.00 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 (i) 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 (i)) were used in the production of this report:

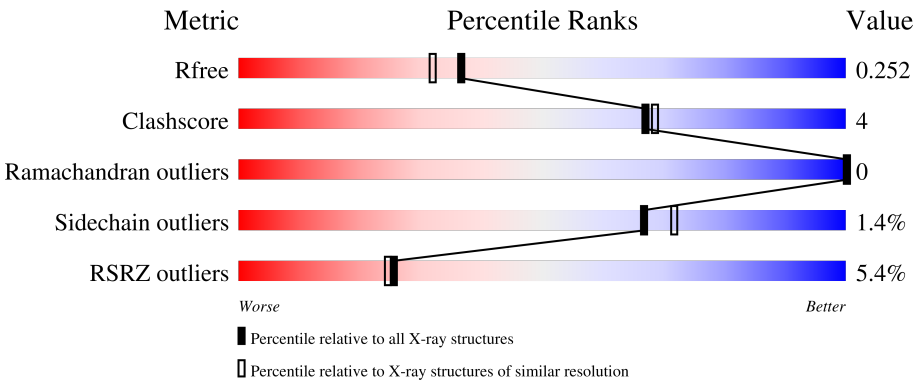
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 2.00 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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 <=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	467	7% 88% 11%
1	B	467	3% 88% 11%
1	C	467	3% 87% 13%
1	D	467	8% 88% 12%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 15641 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 Tryptophanase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	P	S			
1	A	466	3703	2362	628	694	1	18	0	0	0
1	B	466	3703	2362	628	694	1	18	0	0	0
1	C	466	3710	2366	629	696	1	18	0	1	0
1	D	466	3703	2362	628	694	1	18	0	0	0

- Molecule 2 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	K		
2	A	2	2	2	0	0
2	C	2	2	2	0	0

- Molecule 3 is water.

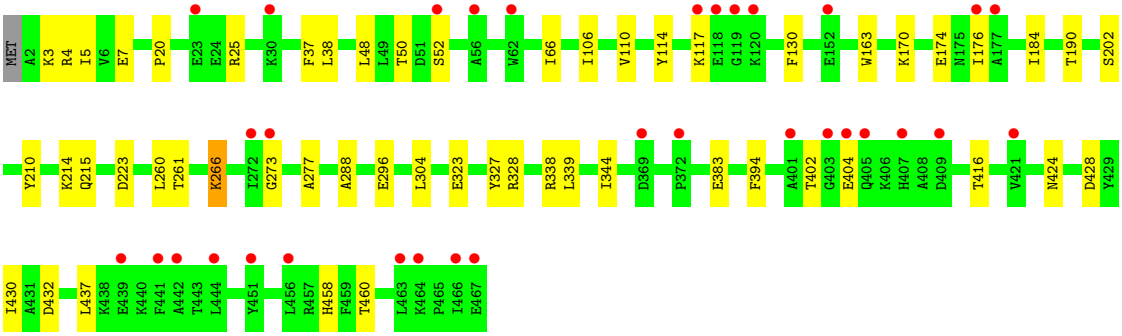
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	190	190	190	0	0
3	B	227	227	227	0	0
3	C	200	200	200	0	0
3	D	201	201	201	0	0

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.

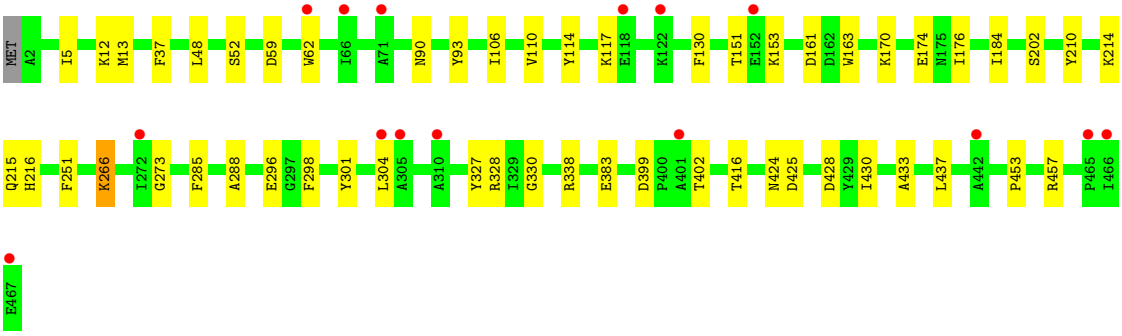
- Molecule 1: Tryptophanase

Chain A:



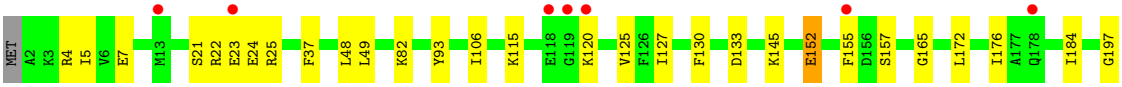
- Molecule 1: Tryptophanase

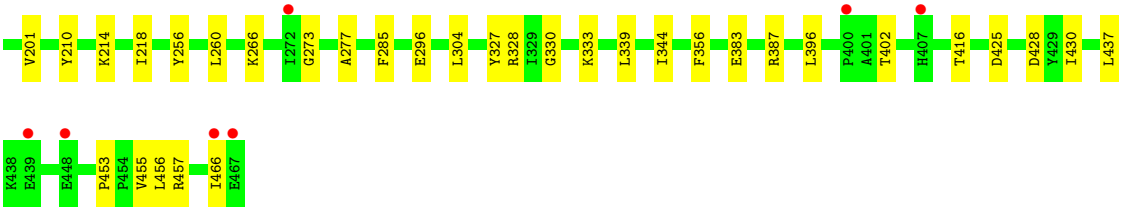
Chain B:



- Molecule 1: Tryptophanase

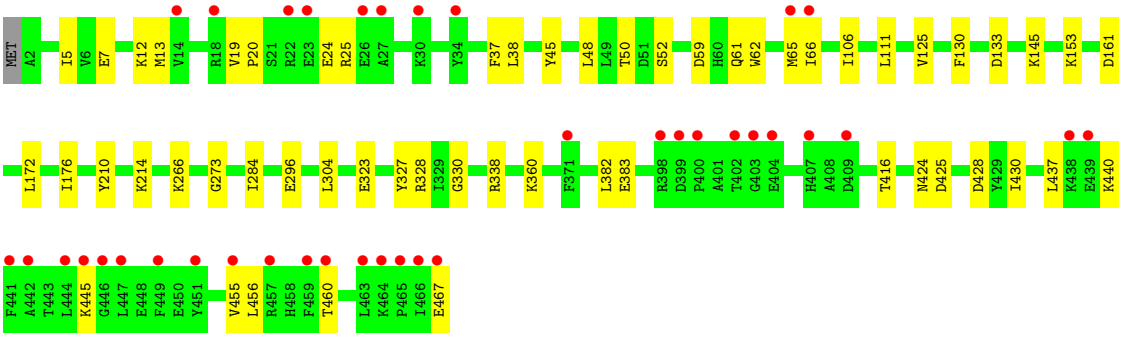
Chain C:





• Molecule 1: Tryptophanase

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, α, β, γ	62.78Å 152.58Å 205.01Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.58 – 2.00 48.59 – 2.00	Depositor EDS
% Data completeness (in resolution range)	96.6 (48.58-2.00) 84.7 (48.59-2.00)	Depositor EDS
R_{merge}	0.23	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	0.13 (at 2.00Å)	Xtriage
Refinement program	PHENIX 1.12_2829	Depositor
R, R_{free}	0.207 , 0.252 0.207 , 0.252	Depositor DCC
R_{free} test set	6628 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å^2)	26.6	Xtriage
Anisotropy	0.616	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 43.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	15641	wwPDB-VP
Average B, all atoms (Å^2)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.90% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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, K

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	0.31	0/3757	0.50	0/5072
1	B	0.30	0/3757	0.50	0/5072
1	C	0.32	0/3764	0.50	0/5082
1	D	0.31	0/3757	0.49	0/5072
All	All	0.31	0/15035	0.50	0/20298

There are no bond length outliers.

There are no bond angle outliers.

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	3703	0	3667	29	0
1	B	3703	0	3667	28	0
1	C	3710	0	3673	32	1
1	D	3703	0	3667	33	0
2	A	2	0	0	0	0
2	C	2	0	0	0	0
3	A	190	0	0	3	0
3	B	227	0	0	0	0
3	C	200	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	201	0	0	2	0
All	All	15641	0	14674	114	1

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 (114) 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:C:115:LYS:HE3	1:C:120:LYS:HE3	1.64	0.79
1:B:383:GLU:HG3	1:B:437:LEU:HD11	1.67	0.74
1:A:4:ARG:NH2	1:A:428:ASP:OD2	2.22	0.73
1:B:62:TRP:CD1	1:D:13:MET:HE2	2.26	0.70
1:A:170:LYS:NZ	1:A:174:GLU:OE2	2.26	0.69
1:B:170:LYS:NZ	1:B:174:GLU:OE2	2.26	0.68
1:C:383:GLU:HG3	1:C:437:LEU:HD21	1.75	0.66
1:A:106:ILE:HD11	1:A:296:GLU:HG3	1.77	0.65
1:A:383:GLU:HG3	1:A:437:LEU:HD21	1.78	0.65
1:A:338:ARG:NH2	1:A:432:ASP:OD1	2.29	0.65
1:C:273:GLY:HA2	1:C:304:LEU:HD21	1.80	0.64
1:D:383:GLU:HG3	1:D:437:LEU:HD21	1.80	0.63
1:B:273:GLY:HA2	1:B:304:LEU:HD21	1.80	0.63
1:B:106:ILE:HD11	1:B:296:GLU:HG3	1.80	0.62
1:B:5:ILE:HD12	1:B:330:GLY:HA3	1.81	0.62
1:C:125:VAL:HG12	1:C:145:LYS:HB3	1.81	0.62
1:A:273:GLY:HA2	1:A:304:LEU:HD21	1.83	0.60
1:A:215:GLN:OE1	3:A:601:HOH:O	2.16	0.59
1:D:455:VAL:HG12	1:D:456:LEU:HG	1.84	0.59
1:A:3:LYS:HE3	1:A:5:ILE:HD11	1.84	0.59
1:C:455:VAL:HG13	1:C:456:LEU:HG	1.85	0.59
1:C:5:ILE:HD12	1:C:330:GLY:HA3	1.84	0.59
1:D:210:TYR:CE1	1:D:214:LYS:HE3	2.39	0.58
1:D:445:LYS:HG3	1:D:467:GLU:HB3	1.85	0.57
1:B:13:MET:HE2	1:D:62:TRP:CD1	2.40	0.57
1:A:176:ILE:HA	1:A:184:ILE:HD11	1.87	0.56
1:D:5:ILE:HD12	1:D:330:GLY:HA3	1.86	0.56
1:D:273:GLY:HA2	1:D:304:LEU:HD21	1.86	0.56
1:D:338:ARG:NH1	1:D:428:ASP:OD1	2.39	0.55
1:C:333:LYS:NZ	3:C:610:HOH:O	2.39	0.55
1:B:13:MET:HE3	1:D:13:MET:HE3	1.88	0.55
1:A:38:LEU:HD21	1:A:460:THR:HA	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:453:PRO:HG2	1:C:457:ARG:HA	1.90	0.54
1:D:48:LEU:HD11	1:D:430:ILE:HD13	1.90	0.54
1:B:453:PRO:HG2	1:B:457:ARG:HA	1.91	0.52
1:A:260:LEU:HG	1:A:277:ALA:HB3	1.91	0.52
1:C:22:ARG:HG2	1:C:25:ARG:NH2	2.25	0.52
1:D:106:ILE:HD11	1:D:296:GLU:HG3	1.91	0.52
1:B:93:TYR:HB3	1:B:285:PHE:CD1	2.44	0.52
1:B:433:ALA:O	1:B:437:LEU:HD13	2.09	0.52
1:B:338:ARG:NH1	1:B:428:ASP:OD1	2.43	0.51
1:A:163:TRP:CD2	1:A:202:SER:HB3	2.46	0.51
1:D:19:VAL:HG22	1:D:45:TYR:CG	2.45	0.51
1:B:90:ASN:HB2	1:B:251:PHE:CE1	2.46	0.50
1:B:110:VAL:HG21	1:B:288:ALA:HA	1.94	0.50
1:D:38:LEU:HD21	1:D:460:THR:HA	1.95	0.49
1:C:4:ARG:NH1	1:C:428:ASP:OD2	2.42	0.48
1:B:424:ASN:ND2	1:D:425:ASP:OD2	2.39	0.48
1:A:394:PHE:CZ	1:A:458:HIS:HA	2.49	0.48
1:C:106:ILE:HD11	1:C:296:GLU:HG3	1.94	0.48
1:B:151:THR:OG1	1:B:153:LYS:HE2	2.13	0.48
1:A:190:THR:HA	1:A:223:ASP:HB3	1.96	0.48
1:D:125:VAL:HG12	1:D:145:LYS:HB3	1.96	0.48
1:D:440:LYS:NZ	3:D:510:HOH:O	2.47	0.48
1:D:7:GLU:HG3	1:D:327:TYR:CE1	2.50	0.47
1:A:424:ASN:ND2	1:C:425:ASP:OD2	2.37	0.46
1:C:214:LYS:NZ	1:C:218:ILE:O	2.40	0.46
1:B:52:SER:HB2	1:B:266:LLP:HE2	1.97	0.46
1:A:7:GLU:HG3	1:A:327:TYR:CZ	2.50	0.46
1:A:50:THR:HG23	1:A:52:SER:H	1.79	0.46
1:C:260:LEU:HG	1:C:277:ALA:HB3	1.97	0.46
1:A:261:THR:OG1	3:A:602:HOH:O	2.19	0.46
1:D:7:GLU:HG3	1:D:327:TYR:CZ	2.52	0.46
1:B:153:LYS:HD2	1:B:161:ASP:CG	2.37	0.45
1:D:25:ARG:NH2	1:D:382:LEU:O	2.40	0.45
1:B:215:GLN:HG3	1:B:216:HIS:ND1	2.32	0.45
1:B:163:TRP:CD2	1:B:202:SER:HB3	2.52	0.45
1:C:339:LEU:HB3	1:C:344:ILE:HB	1.98	0.45
1:A:210:TYR:CE1	1:A:214:LYS:HE3	2.52	0.45
1:B:425:ASP:OD2	1:D:424:ASN:ND2	2.44	0.45
1:D:153:LYS:HD2	1:D:161:ASP:CG	2.37	0.45
1:A:4:ARG:NH1	1:C:4:ARG:NH2	2.65	0.44
1:C:21:SER:OG	1:C:24:GLU:HG3	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:20:PRO:HB2	1:D:24:GLU:HB2	1.98	0.44
1:D:172:LEU:O	1:D:176:ILE:HG12	2.17	0.44
1:C:7:GLU:HG3	1:C:327:TYR:CE1	2.52	0.44
1:A:110:VAL:HG21	1:A:288:ALA:HA	1.99	0.44
1:D:62:TRP:O	1:D:66:ILE:HG12	2.18	0.44
1:B:210:TYR:CE1	1:B:214:LYS:HE3	2.53	0.44
1:C:115:LYS:CE	1:C:120:LYS:HE3	2.41	0.44
1:C:466:ILE:HD11	3:C:661:HOH:O	2.18	0.44
1:C:197:GLY:HA2	1:C:356:PHE:CE1	2.53	0.44
1:D:111:LEU:HD23	1:D:284:ILE:HD12	2.00	0.44
1:A:323:GLU:H	1:A:323:GLU:CD	2.22	0.43
1:D:50:THR:HG23	1:D:52:SER:H	1.84	0.43
1:B:12:LYS:HD3	1:B:59:ASP:HB3	2.01	0.43
1:C:48:LEU:HD11	1:C:430:ILE:HD13	2.00	0.43
1:D:12:LYS:HD3	1:D:59:ASP:HB3	2.01	0.43
1:D:360:LYS:NZ	3:D:514:HOH:O	2.52	0.43
1:C:127:ILE:HD11	1:C:184:ILE:HD12	2.01	0.43
1:B:48:LEU:HD11	1:B:430:ILE:HD13	1.99	0.42
1:B:298:PHE:HB3	1:B:301:TYR:CE2	2.54	0.42
1:C:93:TYR:HB3	1:C:285:PHE:CD1	2.54	0.42
1:A:402:THR:OG1	1:A:404:GLU:HG2	2.20	0.42
1:C:49:LEU:HD13	1:C:387:ARG:HG2	2.01	0.42
1:C:172:LEU:O	1:C:176:ILE:HG13	2.20	0.42
1:A:114:TYR:O	1:A:117:LYS:HG2	2.19	0.42
1:A:66:ILE:HD12	1:D:66:ILE:HG13	2.02	0.42
1:C:165:GLY:O	1:C:201:VAL:HG22	2.20	0.42
1:B:114:TYR:O	1:B:117:LYS:HG2	2.20	0.41
1:D:323:GLU:H	1:D:323:GLU:CD	2.24	0.41
1:C:210:TYR:CE1	1:C:214:LYS:HE3	2.55	0.41
1:C:152:GLU:HA	1:C:155:PHE:HD2	1.86	0.41
1:D:61:GLN:O	1:D:65:MET:HG3	2.21	0.41
1:A:339:LEU:HB3	1:A:344:ILE:HB	2.03	0.41
1:D:20:PRO:HG2	1:D:25:ARG:HG3	2.03	0.41
1:B:176:ILE:HA	1:B:184:ILE:HD11	2.02	0.41
1:B:399:ASP:HB3	1:B:402:THR:OG1	2.21	0.40
1:C:396:LEU:HD12	1:C:396:LEU:HA	1.93	0.40
1:A:48:LEU:HD11	1:A:430:ILE:HD13	2.04	0.40
1:A:266:LLP:H4'1	3:A:765:HOH:O	2.21	0.40
1:C:82:LYS:NZ	3:C:629:HOH:O	2.53	0.40
1:C:115:LYS:HE2	1:C:184:ILE:O	2.21	0.40
1:A:20:PRO:HG2	1:A:25:ARG:CG	2.51	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:23:GLU:OE2	1:C:256:TYR:OH[1_655]	2.14	0.06

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	463/467 (99%)	450 (97%)	13 (3%)	0	100	100
1	B	463/467 (99%)	450 (97%)	13 (3%)	0	100	100
1	C	464/467 (99%)	451 (97%)	13 (3%)	0	100	100
1	D	463/467 (99%)	449 (97%)	14 (3%)	0	100	100
All	All	1853/1868 (99%)	1800 (97%)	53 (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	384/385 (100%)	380 (99%)	4 (1%)	76	81
1	B	384/385 (100%)	379 (99%)	5 (1%)	69	74
1	C	385/385 (100%)	377 (98%)	8 (2%)	53	57
1	D	384/385 (100%)	379 (99%)	5 (1%)	69	74

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	1537/1540 (100%)	1515 (99%)	22 (1%)	67 72

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	PHE
1	A	130	PHE
1	A	328	ARG
1	A	416	THR
1	B	37	PHE
1	B	130	PHE
1	B	327	TYR
1	B	328	ARG
1	B	416	THR
1	C	37	PHE
1	C	130	PHE
1	C	133	ASP
1	C	152	GLU
1	C	157	SER
1	C	328	ARG
1	C	402	THR
1	C	416	THR
1	D	37	PHE
1	D	130	PHE
1	D	133	ASP
1	D	328	ARG
1	D	416	THR

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)

4 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	C	266	1	23,24,25	2.63	6 (26%)	25,32,34	1.52	3 (12%)
1	LLP	D	266	1	23,24,25	2.58	5 (21%)	25,32,34	1.63	5 (20%)
1	LLP	A	266	1	23,24,25	2.56	5 (21%)	25,32,34	1.34	2 (8%)
1	LLP	B	266	1	23,24,25	2.57	6 (26%)	25,32,34	1.43	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	C	266	1	-	5/16/17/19	0/1/1/1
1	LLP	D	266	1	-	8/16/17/19	0/1/1/1
1	LLP	A	266	1	-	6/16/17/19	0/1/1/1
1	LLP	B	266	1	-	6/16/17/19	0/1/1/1

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	266	LLP	C4-C4'	8.18	1.62	1.46
1	B	266	LLP	C4-C4'	8.16	1.62	1.46
1	C	266	LLP	C4-C4'	8.04	1.61	1.46
1	A	266	LLP	C4-C4'	7.97	1.61	1.46
1	A	266	LLP	C4'-NZ	4.98	1.44	1.27
1	B	266	LLP	C4'-NZ	4.94	1.43	1.27
1	D	266	LLP	C4'-NZ	4.92	1.43	1.27
1	C	266	LLP	C4'-NZ	4.88	1.43	1.27
1	C	266	LLP	C4-C5	-4.25	1.36	1.42
1	A	266	LLP	C4-C5	-3.95	1.37	1.42
1	D	266	LLP	C4-C5	-3.87	1.37	1.42
1	C	266	LLP	C2'-C2	3.62	1.56	1.50
1	B	266	LLP	C4-C5	-3.51	1.37	1.42
1	B	266	LLP	C2'-C2	3.45	1.56	1.50
1	D	266	LLP	C2'-C2	3.42	1.56	1.50
1	A	266	LLP	C2'-C2	3.33	1.56	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	266	LLP	C6-N1	3.23	1.41	1.34
1	A	266	LLP	C6-N1	3.03	1.40	1.34
1	B	266	LLP	C6-N1	3.03	1.40	1.34
1	C	266	LLP	C6-N1	2.92	1.40	1.34
1	C	266	LLP	C4-C3	-2.28	1.37	1.40
1	B	266	LLP	C5'-C5	2.06	1.56	1.50

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	D	266	LLP	C4-C4'-NZ	-4.56	103.37	124.31
1	B	266	LLP	C4-C4'-NZ	-4.50	103.65	124.31
1	C	266	LLP	C4-C4'-NZ	-4.41	104.06	124.31
1	A	266	LLP	C4-C4'-NZ	-4.10	105.49	124.31
1	D	266	LLP	OP4-C5'-C5	3.13	115.32	109.35
1	D	266	LLP	OP3-P-OP4	2.58	113.59	106.73
1	B	266	LLP	C5-C6-N1	-2.43	119.77	123.82
1	C	266	LLP	C5-C6-N1	-2.25	120.08	123.82
1	D	266	LLP	C5-C6-N1	-2.21	120.14	123.82
1	C	266	LLP	C3-C4-C4'	-2.14	116.42	120.41
1	D	266	LLP	C2'-C2-N1	2.08	121.73	117.67
1	A	266	LLP	C5-C6-N1	-2.05	120.40	123.82

There are no chirality outliers.

All (25) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	266	LLP	O-C-CA-CB
1	A	266	LLP	CG-CD-CE-NZ
1	B	266	LLP	C-CA-CB-CG
1	B	266	LLP	O-C-CA-CB
1	B	266	LLP	CG-CD-CE-NZ
1	C	266	LLP	C-CA-CB-CG
1	C	266	LLP	O-C-CA-CB
1	C	266	LLP	CG-CD-CE-NZ
1	D	266	LLP	C5'-OP4-P-OP1
1	D	266	LLP	O-C-CA-CB
1	D	266	LLP	CG-CD-CE-NZ
1	A	266	LLP	C-CA-CB-CG
1	D	266	LLP	C-CA-CB-CG
1	A	266	LLP	C3-C4-C4'-NZ
1	B	266	LLP	C3-C4-C4'-NZ

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Mol	Chain	Res	Type	Atoms
1	C	266	LLP	C3-C4-C4'-NZ
1	D	266	LLP	C3-C4-C4'-NZ
1	A	266	LLP	C5-C4-C4'-NZ
1	B	266	LLP	C5-C4-C4'-NZ
1	D	266	LLP	C5'-OP4-P-OP2
1	C	266	LLP	C5-C4-C4'-NZ
1	D	266	LLP	C5-C4-C4'-NZ
1	A	266	LLP	CD-CE-NZ-C4'
1	B	266	LLP	CD-CE-NZ-C4'
1	D	266	LLP	CD-CE-NZ-C4'

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	266	LLP	1	0
1	B	266	LLP	1	0

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	465/467 (99%)	0.51	33 (7%)	16	15	26, 42, 69, 121	0
1	B	465/467 (99%)	0.41	15 (3%)	47	46	27, 41, 63, 103	0
1	C	465/467 (99%)	0.44	14 (3%)	50	49	25, 43, 69, 106	0
1	D	465/467 (99%)	0.53	38 (8%)	11	11	26, 40, 76, 111	0
All	All	1860/1868 (99%)	0.47	100 (5%)	25	24	25, 41, 72, 121	0

All (100) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	407	HIS	7.3
1	D	467	GLU	6.8
1	D	407	HIS	6.0
1	A	467	GLU	5.8
1	B	467	GLU	5.2
1	C	467	GLU	5.1
1	A	119	GLY	5.0
1	D	399	ASP	4.3
1	B	401	ALA	4.2
1	A	401	ALA	3.9
1	D	444	LEU	3.9
1	C	407	HIS	3.9
1	A	409	ASP	3.9
1	D	398	ARG	3.8
1	D	403	GLY	3.7
1	D	464	LYS	3.6
1	D	34	TYR	3.6
1	D	445	LYS	3.5
1	B	272	ILE	3.5
1	D	455	VAL	3.5
1	D	449	PHE	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	120	LYS	3.4
1	C	400	PRO	3.4
1	C	120	LYS	3.3
1	C	23	GLU	3.3
1	C	118	GLU	3.3
1	D	402	THR	3.2
1	D	466	ILE	3.1
1	D	447	LEU	3.1
1	A	441	PHE	3.0
1	B	466	ILE	3.0
1	D	66	ILE	2.9
1	D	400	PRO	2.9
1	D	463	LEU	2.9
1	A	272	ILE	2.9
1	D	23	GLU	2.9
1	D	404	GLU	2.9
1	C	119	GLY	2.8
1	D	439	GLU	2.8
1	A	456	LEU	2.8
1	D	438	LYS	2.8
1	D	26	GLU	2.7
1	B	118	GLU	2.6
1	B	62	TRP	2.6
1	C	466	ILE	2.6
1	A	403	GLY	2.6
1	A	421	VAL	2.6
1	A	369	ASP	2.5
1	A	442	ALA	2.5
1	A	23	GLU	2.5
1	D	27	ALA	2.5
1	A	372	PRO	2.5
1	C	439	GLU	2.5
1	A	444	LEU	2.4
1	D	460	THR	2.4
1	A	463	LEU	2.4
1	A	464	LYS	2.4
1	D	441	PHE	2.4
1	D	457	ARG	2.4
1	C	448	GLU	2.3
1	B	122	LYS	2.3
1	A	404	GLU	2.3
1	A	30	LYS	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	177	ALA	2.3
1	D	22	ARG	2.3
1	A	451	TYR	2.3
1	B	152	GLU	2.3
1	D	442	ALA	2.3
1	D	18	ARG	2.3
1	A	118	GLU	2.3
1	A	152	GLU	2.3
1	B	442	ALA	2.2
1	B	310	ALA	2.2
1	A	176	ILE	2.2
1	B	66	ILE	2.2
1	A	117	LYS	2.2
1	A	56	ALA	2.2
1	A	273	GLY	2.2
1	D	459	PHE	2.2
1	D	446	GLY	2.2
1	C	272	ILE	2.1
1	D	409	ASP	2.1
1	C	155	PHE	2.1
1	D	451	TYR	2.1
1	C	13	MET	2.1
1	A	466	ILE	2.1
1	A	439	GLU	2.1
1	D	371	PHE	2.1
1	B	71	ALA	2.1
1	D	465	PRO	2.1
1	B	465	PRO	2.1
1	A	405	GLN	2.0
1	B	304	LEU	2.0
1	C	178	GLN	2.0
1	D	14	VAL	2.0
1	A	62	TRP	2.0
1	B	305	ALA	2.0
1	A	52	SER	2.0
1	D	30	LYS	2.0
1	D	65	MET	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	D	266	24/25	0.95	0.14	27,35,42,46	0
1	LLP	B	266	24/25	0.96	0.17	29,38,44,46	0
1	LLP	C	266	24/25	0.96	0.17	26,37,43,48	0
1	LLP	A	266	24/25	0.96	0.15	30,37,44,45	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	K	C	502	1/1	0.97	0.10	34,34,34,34	0
2	K	C	501	1/1	0.98	0.07	39,39,39,39	0
2	K	A	501	1/1	0.99	0.08	42,42,42,42	0
2	K	A	502	1/1	0.99	0.03	38,38,38,38	0

6.5 Other polymers (i)

There are no such residues in this entry.