



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 26, 2014 – 04:14 PM GMT

PDB ID : 1AX4
Title : TRYPTOPHANASE FROM PROTEUS VULGARIS
Authors : Isupov, M.N.; Antson, A.A.; Dodson, E.J.; Dodson, G.G.; Dementieva, I.S.; Zakomirdina, L.N.; Wilson, K.S.; Dauter, Z.; Lebedev, A.A.; Harutyunyan, E.H.
Deposited on : 1997-10-28
Resolution : 2.10 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

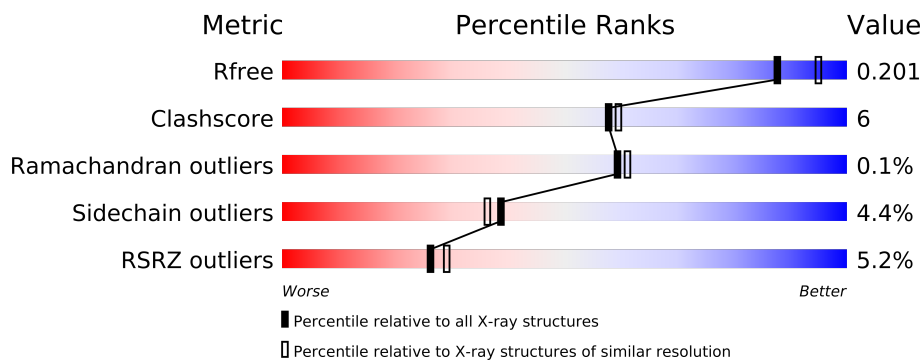
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.10 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}	66092	3012 (2.10-2.10)
Clashscore	79885	3649 (2.10-2.10)
Ramachandran outliers	78287	3610 (2.10-2.10)
Sidechain outliers	78261	3611 (2.10-2.10)
RSRZ outliers	66119	3013 (2.10-2.10)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	467	
1	B	467	
1	C	467	
1	D	467	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 15715 atoms, of which 0 are hydrogen and 0 are deuterium.

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
1	A	465	Total	C	N	O	P	S	0	0	0
			3692	2356	627	690	1	18			
1	B	465	Total	C	N	O	P	S	0	0	0
			3693	2357	627	690	1	18			
1	C	465	Total	C	N	O	P	S	0	0	0
			3693	2357	627	690	1	18			
1	D	465	Total	C	N	O	P	S	0	0	0
			3693	2357	627	690	1	18			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	266	LLP	LYS	MODIFIED RESIDUE	UNP P28796
B	266	LLP	LYS	MODIFIED RESIDUE	UNP P28796
C	266	LLP	LYS	MODIFIED RESIDUE	UNP P28796
D	266	LLP	LYS	MODIFIED RESIDUE	UNP P28796

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	K	0	0
			1	1		
2	A	1	Total	K	0	0
			1	1		
2	D	1	Total	K	0	0
			1	1		
2	C	1	Total	K	0	0
			1	1		

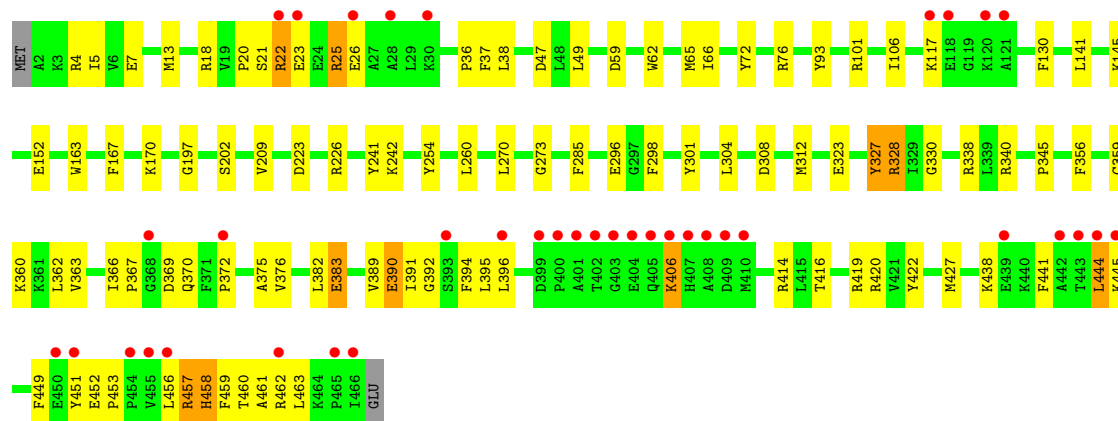
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	225	Total 225	O 225	0	0
3	B	240	Total 240	O 240	0	0
3	C	261	Total 261	O 261	0	0
3	D	214	Total 214	O 214	0	0



● Molecule 1: TRYPTOPHANASE

Chain D:



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	114.99Å 118.23Å 153.68Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	18.00 – 2.10 17.95 – 2.10	Depositor EDS
% Data completeness (in resolution range)	97.6 (18.00-2.10) 97.0 (17.95-2.10)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.28 (at 2.11Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.187 , 0.227 0.168 , 0.201	Depositor DCC
R_{free} test set	2364 reflections (2.05%)	DCC
Wilson B-factor (Å ²)	17.7	Xtriage
Anisotropy	0.544	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 44.3	EDS
Estimated twinning fraction	0.012 for k,h,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 117863 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	15715	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.96% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: K, LLP

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	0.63	0/3747	1.32	34/5060 (0.7%)
1	B	0.62	0/3747	1.24	17/5060 (0.3%)
1	C	0.60	0/3747	1.20	15/5060 (0.3%)
1	D	0.61	0/3747	1.32	27/5060 (0.5%)
All	All	0.62	0/14988	1.27	93/20240 (0.5%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	1

There are no bond length outliers.

All (93) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	327	TYR	CB-CG-CD2	14.96	129.97	121.00
1	A	338	ARG	CD-NE-CZ	14.30	143.61	123.60
1	A	338	ARG	NE-CZ-NH1	11.60	126.10	120.30
1	D	328	ARG	NE-CZ-NH1	11.20	125.90	120.30
1	D	420	ARG	NE-CZ-NH2	-11.01	114.80	120.30
1	A	4	ARG	NE-CZ-NH1	10.66	125.63	120.30
1	B	338	ARG	NE-CZ-NH1	10.62	125.61	120.30
1	D	327	TYR	CB-CG-CD1	-10.42	114.75	121.00
1	A	25	ARG	NE-CZ-NH1	10.23	125.41	120.30
1	D	101	ARG	NE-CZ-NH2	-10.12	115.24	120.30
1	B	419	ARG	NE-CZ-NH1	9.52	125.06	120.30
1	B	340	ARG	NE-CZ-NH1	9.09	124.85	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	327	TYR	CB-CG-CD1	-8.98	115.61	121.00
1	C	419	ARG	NE-CZ-NH1	8.92	124.76	120.30
1	A	338	ARG	NE-CZ-NH2	-8.42	116.09	120.30
1	A	463	LEU	CA-CB-CG	8.35	134.50	115.30
1	B	51	ASP	CB-CG-OD2	8.14	125.63	118.30
1	B	432	ASP	CB-CG-OD1	7.99	125.49	118.30
1	C	296	GLU	OE1-CD-OE2	-7.93	113.79	123.30
1	D	328	ARG	NE-CZ-NH2	-7.91	116.35	120.30
1	D	419	ARG	NE-CZ-NH1	7.91	124.25	120.30
1	C	398	ARG	CD-NE-CZ	7.85	134.59	123.60
1	A	59	ASP	CB-CG-OD1	7.61	125.14	118.30
1	C	10	ARG	NE-CZ-NH1	7.56	124.08	120.30
1	A	296	GLU	OE1-CD-OE2	-7.51	114.29	123.30
1	D	226	ARG	NE-CZ-NH1	7.45	124.03	120.30
1	C	10	ARG	NE-CZ-NH2	-7.44	116.58	120.30
1	A	4	ARG	CD-NE-CZ	7.37	133.92	123.60
1	D	422	TYR	CB-CG-CD2	-7.36	116.58	121.00
1	D	427	MET	CA-CB-CG	7.34	125.77	113.30
1	B	338	ARG	NE-CZ-NH2	-7.32	116.64	120.30
1	C	59	ASP	CB-CG-OD1	7.30	124.87	118.30
1	A	51	ASP	CB-CG-OD2	7.27	124.84	118.30
1	B	398	ARG	CD-NE-CZ	7.20	133.68	123.60
1	A	328	ARG	NE-CZ-NH1	7.11	123.86	120.30
1	B	340	ARG	NE-CZ-NH2	-7.03	116.78	120.30
1	A	248	GLU	OE1-CD-OE2	-7.02	114.87	123.30
1	A	72	TYR	CB-CG-CD1	-6.99	116.81	121.00
1	D	59	ASP	CB-CG-OD1	6.98	124.58	118.30
1	D	328	ARG	CD-NE-CZ	6.90	133.26	123.60
1	C	72	TYR	CB-CG-CD1	-6.84	116.90	121.00
1	B	419	ARG	NE-CZ-NH2	-6.83	116.88	120.30
1	A	419	ARG	NE-CZ-NH2	-6.83	116.89	120.30
1	C	328	ARG	NE-CZ-NH1	6.80	123.70	120.30
1	A	420	ARG	NE-CZ-NH2	-6.76	116.92	120.30
1	A	59	ASP	CB-CG-OD2	-6.57	112.39	118.30
1	D	4	ARG	NE-CZ-NH1	6.46	123.53	120.30
1	D	25	ARG	NE-CZ-NH1	6.44	123.52	120.30
1	D	308	ASP	CB-CG-OD1	-6.41	112.53	118.30
1	D	422	TYR	CB-CG-CD1	6.39	124.83	121.00
1	A	10	ARG	NE-CZ-NH2	-6.37	117.12	120.30
1	D	47	ASP	CB-CG-OD1	6.30	123.97	118.30
1	A	414	ARG	NE-CZ-NH1	-6.30	117.15	120.30
1	A	428	ASP	CB-CG-OD1	6.28	123.95	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	308	ASP	CB-CG-OD1	-6.25	112.68	118.30
1	A	427	MET	CA-CB-CG	6.18	123.81	113.30
1	B	337	ASP	CB-CG-OD1	-6.17	112.75	118.30
1	B	267	ASP	CB-CG-OD1	6.14	123.83	118.30
1	D	338	ARG	NE-CZ-NH1	6.11	123.35	120.30
1	A	101	ARG	NE-CZ-NH2	-6.10	117.25	120.30
1	A	237	ARG	NE-CZ-NH2	-6.08	117.26	120.30
1	A	248	GLU	CG-CD-OE1	6.03	130.36	118.30
1	D	72	TYR	CB-CG-CD1	-5.86	117.48	121.00
1	C	327	TYR	CB-CG-CD2	5.75	124.45	121.00
1	B	296	GLU	OE1-CD-OE2	-5.73	116.43	123.30
1	D	226	ARG	NH1-CZ-NH2	-5.59	113.25	119.40
1	D	65	MET	CG-SD-CE	5.55	109.09	100.20
1	B	162	ASP	CB-CG-OD2	-5.52	113.33	118.30
1	A	420	ARG	NE-CZ-NH1	5.52	123.06	120.30
1	C	291	ARG	NE-CZ-NH2	-5.51	117.55	120.30
1	C	328	ARG	CD-NE-CZ	5.49	131.29	123.60
1	A	4	ARG	NE-CZ-NH2	-5.46	117.57	120.30
1	A	340	ARG	NE-CZ-NH1	-5.45	117.58	120.30
1	B	334	TYR	CB-CG-CD2	-5.42	117.75	121.00
1	D	254	TYR	CA-CB-CG	-5.38	103.17	113.40
1	B	429	TYR	CB-CG-CD2	-5.37	117.78	121.00
1	D	414	ARG	NE-CZ-NH2	5.36	122.98	120.30
1	D	457	ARG	NE-CZ-NH1	-5.32	117.64	120.30
1	A	210	TYR	CB-CG-CD1	-5.30	117.82	121.00
1	D	241	TYR	CB-CG-CD2	5.27	124.16	121.00
1	D	76	ARG	CA-CB-CG	5.23	124.91	113.40
1	A	51	ASP	CB-CG-OD1	-5.21	113.61	118.30
1	C	462	ARG	NE-CZ-NH2	-5.21	117.69	120.30
1	B	4	ARG	CD-NE-CZ	5.21	130.90	123.60
1	A	307	ARG	NE-CZ-NH2	-5.21	117.70	120.30
1	A	327	TYR	CB-CG-CD2	5.21	124.12	121.00
1	D	223	ASP	CB-CG-OD2	5.16	122.95	118.30
1	A	327	TYR	CB-CG-CD1	-5.15	117.91	121.00
1	A	308	ASP	CB-CG-OD1	-5.15	113.66	118.30
1	C	4	ARG	CD-NE-CZ	5.14	130.80	123.60
1	A	168	ASP	CB-CG-OD2	-5.12	113.69	118.30
1	B	327	TYR	CB-CG-CD2	5.06	124.04	121.00
1	A	47	ASP	CB-CG-OD1	5.02	122.81	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	162	ASP	Mainchain

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3692	0	3659	66	0
1	B	3693	0	3661	37	0
1	C	3693	0	3662	43	0
1	D	3693	0	3661	56	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	225	0	0	4	0
3	B	240	0	0	3	0
3	C	261	0	0	4	0
3	D	214	0	0	4	0
All	All	15715	0	14643	184	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 6.

All (184) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:4:ARG:HH21	1:C:4:ARG:NH2	1.11	1.48
1:A:4:ARG:NH2	1:C:4:ARG:NH2	1.67	1.41
1:A:4:ARG:NH2	1:C:4:ARG:HH21	1.16	1.36
1:A:266:LLP:C4	1:A:266:LLP:NZ	2.34	0.90
1:A:26:GLU:HG3	1:A:382:LEU:HD21	1.52	0.89
1:C:383:GLU:HG2	1:C:437:LEU:HD21	1.56	0.85
1:A:4:ARG:NH2	1:C:4:ARG:CZ	2.43	0.81
1:B:453:PRO:HG2	1:B:457:ARG:HA	1.64	0.78
1:D:145:LYS:HE3	3:D:615:HOH:O	1.85	0.77
1:D:141:LEU:HG	3:D:692:HOH:O	1.85	0.76
1:C:266:LLP:H4'1	3:C:589:HOH:O	1.84	0.75

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:4:ARG:CZ	1:C:4:ARG:NH2	2.49	0.75
1:A:399:ASP:HB3	1:A:402:THR:OG1	1.94	0.68
1:A:453:PRO:HG2	1:A:460:THR:HB	1.74	0.68
1:B:383:GLU:HG3	1:B:437:LEU:HD21	1.74	0.67
1:A:410:MET:HB2	3:A:723:HOH:O	1.94	0.67
1:D:453:PRO:HG2	1:D:457:ARG:HA	1.76	0.67
1:D:375:ALA:HB1	1:D:444:LEU:HD22	1.76	0.65
1:A:197:GLY:HA2	1:A:356:PHE:CE1	2.32	0.65
1:B:197:GLY:HA2	1:B:356:PHE:CE1	2.32	0.64
1:D:367:PRO:HG2	1:D:370:GLN:HG3	1.80	0.63
1:D:390:GLU:O	1:D:391:ILE:HD13	1.98	0.63
1:C:145:LYS:HE3	3:C:698:HOH:O	1.99	0.62
1:A:87:GLU:HB2	3:A:643:HOH:O	2.00	0.62
1:D:197:GLY:HA2	1:D:356:PHE:CE1	2.35	0.62
1:D:366:ILE:HD11	1:D:444:LEU:HB3	1.83	0.61
1:B:383:GLU:HG2	3:B:668:HOH:O	2.00	0.61
1:D:445:LYS:HD2	1:D:445:LYS:H	1.65	0.61
1:C:197:GLY:HA2	1:C:356:PHE:CE1	2.37	0.60
1:A:399:ASP:OD1	1:A:400:PRO:HD2	2.02	0.60
1:D:20:PRO:HG2	1:D:25:ARG:HG2	1.82	0.60
1:B:402:THR:HB	1:B:404:GLU:HG2	1.84	0.59
1:A:359:CYS:HA	1:A:362:LEU:HG	1.84	0.59
1:D:458:HIS:H	1:D:458:HIS:CD2	2.20	0.59
1:C:448:GLU:HG3	1:C:466:ILE:HG13	1.84	0.59
1:B:410:MET:SD	3:B:705:HOH:O	2.57	0.59
1:B:370:GLN:NE2	1:B:466:ILE:HD13	2.18	0.58
1:D:21:SER:O	1:D:25:ARG:HG3	2.04	0.58
1:A:4:ARG:CZ	1:C:4:ARG:CZ	2.82	0.57
1:C:152:GLU:HG2	3:C:646:HOH:O	2.05	0.56
1:A:32:ALA:HB2	1:A:39:LEU:HD23	1.88	0.56
1:D:26:GLU:HG2	1:D:382:LEU:HD22	1.89	0.55
1:A:394:PHE:CZ	1:A:463:LEU:HD21	2.42	0.55
1:A:21:SER:OG	1:A:24:GLU:HG3	2.07	0.55
1:D:396:LEU:HD23	1:D:406:LYS:HB3	1.89	0.54
1:D:438:LYS:HA	1:D:441:PHE:CD2	2.42	0.54
1:C:340:ARG:HD3	3:C:645:HOH:O	2.06	0.53
1:D:345:PRO:HD2	1:D:362:LEU:HD21	1.90	0.53
1:A:4:ARG:HH21	1:C:4:ARG:HH21	0.75	0.53
1:A:4:ARG:NH1	1:A:428:ASP:OD2	2.41	0.53
1:C:22:ARG:O	1:C:26:GLU:HG3	2.09	0.53
1:D:451:TYR:HB3	1:D:462:ARG:HG3	1.91	0.53
1:D:93:TYR:HB3	1:D:285:PHE:CD1	2.44	0.53

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:362:LEU:O	1:D:441:PHE:HB3	2.09	0.52
1:A:5:ILE:HD12	1:A:330:GLY:HA3	1.91	0.52
1:A:438:LYS:HA	1:A:441:PHE:CD2	2.43	0.52
1:C:448:GLU:HG3	1:C:466:ILE:CG1	2.39	0.52
1:D:340:ARG:HD3	3:D:676:HOH:O	2.10	0.52
1:A:397:GLY:HA2	1:A:457:ARG:HH11	1.74	0.52
1:A:390:GLU:O	1:A:391:ILE:HD13	2.09	0.52
1:D:5:ILE:HD12	1:D:330:GLY:HA3	1.92	0.51
1:A:456:LEU:HD23	1:A:460:THR:OG1	2.10	0.51
1:A:323:GLU:H	1:A:323:GLU:CD	2.14	0.51
1:A:145:LYS:HE3	3:A:671:HOH:O	2.11	0.51
1:A:447:LEU:HD23	1:A:465:PRO:HA	1.93	0.50
1:C:398:ARG:HG2	1:C:398:ARG:HH11	1.76	0.50
1:D:38:LEU:HD21	1:D:460:THR:HA	1.93	0.50
1:B:13:MET:HE2	1:D:62:TRP:CG	2.47	0.49
1:C:410:MET:HE2	1:C:410:MET:HA	1.94	0.49
1:B:7:GLU:HG3	1:B:327:TYR:CZ	2.47	0.49
1:A:366:ILE:HG22	1:A:372:PRO:HD3	1.94	0.49
1:B:398:ARG:HH11	1:B:398:ARG:HG2	1.77	0.49
1:A:359:CYS:HB3	1:A:376:VAL:HG21	1.94	0.49
1:A:449:PHE:CD1	1:A:463:LEU:HD22	2.46	0.49
1:A:22:ARG:HH21	1:A:382:LEU:HD22	1.77	0.49
1:D:395:LEU:HG	1:D:395:LEU:O	2.12	0.49
1:C:93:TYR:HB3	1:C:285:PHE:CD1	2.47	0.49
1:B:93:TYR:HB3	1:B:285:PHE:CD1	2.48	0.49
1:B:21:SER:O	1:B:25:ARG:HG3	2.13	0.49
1:B:197:GLY:HA2	1:B:356:PHE:CZ	2.48	0.48
1:B:323:GLU:CD	1:B:323:GLU:H	2.16	0.48
1:C:7:GLU:HG3	1:C:327:TYR:CZ	2.49	0.48
1:B:451:TYR:OH	1:B:453:PRO:HB3	2.13	0.48
1:A:345:PRO:HD2	1:A:362:LEU:HD21	1.95	0.48
1:C:5:ILE:HD12	1:C:330:GLY:HA3	1.95	0.48
1:D:7:GLU:HG3	1:D:327:TYR:CZ	2.48	0.48
1:D:106:ILE:HD11	1:D:296:GLU:HG3	1.96	0.48
1:A:369:ASP:HA	1:A:405:GLN:HE21	1.78	0.48
1:A:453:PRO:CG	1:A:460:THR:HB	2.43	0.48
1:D:21:SER:HG	1:D:23:GLU:HG2	1.79	0.48
1:D:451:TYR:HB3	1:D:462:ARG:HB2	1.96	0.48
1:B:5:ILE:HD12	1:B:330:GLY:HA3	1.95	0.47
1:A:210:TYR:CE1	1:A:214:LYS:HE3	2.49	0.47
1:D:390:GLU:OE1	1:D:392:GLY:N	2.45	0.47
1:A:62:TRP:CG	1:C:13:MET:HE2	2.49	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:360:LYS:HD2	1:B:411:GLU:HG2	1.97	0.47
1:D:449:PHE:CD1	1:D:452:GLU:HB2	2.49	0.47
1:B:163:TRP:CD2	1:B:202:SER:HB3	2.49	0.47
1:A:421:VAL:CG1	1:C:10:ARG:HD2	2.44	0.47
1:A:62:TRP:CD1	1:C:13:MET:HE2	2.50	0.47
1:C:106:ILE:HD11	1:C:296:GLU:HG3	1.96	0.46
1:B:260:LEU:HG	1:B:277:ALA:HB3	1.96	0.46
1:B:266:LLP:H4'1	3:B:583:HOH:O	2.15	0.46
1:D:22:ARG:HH21	1:D:382:LEU:HD13	1.81	0.46
1:D:167:PHE:CD1	1:D:209:VAL:HG21	2.50	0.46
1:C:383:GLU:HG2	1:C:437:LEU:CD2	2.38	0.46
1:A:19:VAL:HG22	1:A:45:TYR:CD2	2.50	0.46
1:C:16:LYS:HG2	1:C:17:ILE:N	2.30	0.46
1:A:273:GLY:HA2	1:A:304:LEU:HD21	1.97	0.46
1:D:359:CYS:HB3	1:D:376:VAL:HG21	1.97	0.45
1:A:197:GLY:HA2	1:A:356:PHE:CZ	2.52	0.45
1:D:163:TRP:CD2	1:D:202:SER:HB3	2.50	0.45
1:D:36:PRO:HB2	3:D:610:HOH:O	2.16	0.45
1:B:371:PHE:CE1	1:B:394:PHE:HA	2.51	0.45
1:A:106:ILE:HD11	1:A:296:GLU:HG3	1.98	0.45
1:A:132:PHE:HD1	1:A:194:ASN:HD22	1.65	0.45
1:B:62:TRP:CG	1:D:13:MET:HE2	2.52	0.45
1:C:21:SER:OG	1:C:24:GLU:HG3	2.17	0.45
1:A:167:PHE:CD1	1:A:209:VAL:HG21	2.52	0.44
1:D:394:PHE:CZ	1:D:463:LEU:HD21	2.52	0.44
1:B:366:ILE:HD11	1:B:444:LEU:O	2.16	0.44
1:D:21:SER:OG	1:D:23:GLU:HG2	2.18	0.44
1:B:4:ARG:NH1	1:B:428:ASP:OD2	2.49	0.44
1:A:372:PRO:HG3	1:A:411:GLU:HG3	1.99	0.44
1:A:448:GLU:O	1:A:463:LEU:HB3	2.18	0.44
1:A:396:LEU:HA	1:A:396:LEU:HD12	1.83	0.44
1:B:399:ASP:OD1	1:B:400:PRO:HD2	2.18	0.44
1:C:197:GLY:HA2	1:C:356:PHE:CZ	2.52	0.44
1:B:13:MET:HE2	1:D:62:TRP:CD1	2.52	0.44
1:C:163:TRP:CD2	1:C:202:SER:HB3	2.53	0.44
1:C:49:LEU:HD12	1:C:389:VAL:HB	1.99	0.44
1:D:372:PRO:O	1:D:376:VAL:HG23	2.18	0.43
1:C:323:GLU:CD	1:C:323:GLU:H	2.22	0.43
1:D:273:GLY:HA2	1:D:304:LEU:HD21	1.99	0.43
1:A:293:VAL:HG11	1:B:459:PHE:CE2	2.54	0.43
1:D:323:GLU:H	1:D:323:GLU:CD	2.21	0.43
1:A:358:ASP:O	1:A:362:LEU:HG	2.19	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:132:PHE:HD1	1:B:194:ASN:HD22	1.65	0.43
1:A:13:MET:HE2	1:C:62:TRP:CG	2.54	0.43
1:B:449:PHE:CZ	1:B:457:ARG:NH1	2.87	0.42
1:B:449:PHE:CZ	1:B:457:ARG:HD2	2.54	0.42
1:D:22:ARG:NH2	1:D:382:LEU:HB3	2.34	0.42
1:A:298:PHE:HB3	1:A:301:TYR:CE2	2.54	0.42
1:A:13:MET:HE2	1:C:62:TRP:CD1	2.54	0.42
1:B:273:GLY:HA2	1:B:304:LEU:HD21	2.01	0.42
1:D:457:ARG:O	1:D:461:ALA:N	2.39	0.42
1:D:340:ARG:HD3	1:D:340:ARG:HH11	1.64	0.42
1:C:273:GLY:HA2	1:C:304:LEU:HD21	2.00	0.42
1:A:93:TYR:HB3	1:A:285:PHE:CD1	2.54	0.42
1:D:456:LEU:HD22	1:D:459:PHE:HD2	1.84	0.42
1:A:232:TYR:CD2	1:A:329:ILE:HD13	2.54	0.42
1:C:298:PHE:HB3	1:C:301:TYR:CE2	2.54	0.42
1:B:449:PHE:CD1	1:B:452:GLU:HB2	2.54	0.42
1:A:397:GLY:HA2	1:A:457:ARG:NH1	2.34	0.42
1:A:393:SER:HB2	1:A:405:GLN:HE22	1.85	0.42
1:A:449:PHE:CD1	1:A:452:GLU:HB2	2.55	0.42
1:D:22:ARG:NH1	1:D:383:GLU:OE2	2.49	0.41
1:D:49:LEU:HD12	1:D:389:VAL:HB	2.02	0.41
1:D:298:PHE:HB3	1:D:301:TYR:CE2	2.55	0.41
1:A:367:PRO:O	1:A:372:PRO:HD3	2.21	0.41
1:A:252:ASP:O	1:A:255:LYS:HB2	2.21	0.41
1:A:369:ASP:HA	1:A:405:GLN:NE2	2.35	0.41
1:B:453:PRO:HA	1:B:454:PRO:HD3	1.88	0.41
1:B:260:LEU:C	1:B:260:LEU:HD12	2.41	0.41
1:D:260:LEU:HD12	1:D:260:LEU:C	2.41	0.41
1:A:367:PRO:HG2	1:A:370:GLN:HG3	2.02	0.41
1:C:132:PHE:HD1	1:C:194:ASN:HD22	1.69	0.41
1:C:4:ARG:NH1	1:C:428:ASP:OD2	2.53	0.41
1:C:223:ASP:OD1	1:C:266:LLP:H2'2	2.21	0.41
1:D:396:LEU:HA	1:D:396:LEU:HD12	1.96	0.41
1:D:394:PHE:CE1	1:D:463:LEU:HD21	2.56	0.41
1:B:298:PHE:HB3	1:B:301:TYR:CE1	2.56	0.41
1:C:62:TRP:CZ3	1:D:66:ILE:HD13	2.56	0.41
1:C:167:PHE:CD1	1:C:209:VAL:HG21	2.56	0.41
1:A:129:ASN:O	1:A:150:VAL:HB	2.22	0.40
1:A:270:LEU:HD11	1:A:312:MET:HA	2.02	0.40
1:D:270:LEU:HD11	1:D:312:MET:HA	2.03	0.40
1:B:62:TRP:CD1	1:D:13:MET:HE2	2.55	0.40
1:A:340:ARG:HD3	1:A:340:ARG:HH11	1.68	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:3:LYS:CE	3:A:651:HOH:O	2.69	0.40
1:B:76:ARG:HH11	1:B:76:ARG:HG3	1.85	0.40
1:C:362:LEU:HD13	1:C:441:PHE:CZ	2.57	0.40
1:D:367:PRO:HG2	1:D:369:ASP:OD1	2.22	0.40
1:C:260:LEU:HD12	1:C:260:LEU:C	2.42	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

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	462/467 (99%)	440 (95%)	22 (5%)	0	100	100
1	B	462/467 (99%)	439 (95%)	21 (4%)	2 (0%)	43	39
1	C	462/467 (99%)	446 (96%)	16 (4%)	0	100	100
1	D	462/467 (99%)	442 (96%)	20 (4%)	0	100	100
All	All	1848/1868 (99%)	1767 (96%)	79 (4%)	2 (0%)	59	61

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	409	ASP
1	B	368	GLY

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	383/385 (100%)	364 (95%)	19 (5%)	34	30

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	383/385 (100%)	366 (96%)	17 (4%)	39	36
1	C	383/385 (100%)	369 (96%)	14 (4%)	45	45
1	D	383/385 (100%)	366 (96%)	17 (4%)	39	36
All	All	1532/1540 (100%)	1465 (96%)	67 (4%)	39	36

All (67) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	PHE
1	A	117	LYS
1	A	130	PHE
1	A	152	GLU
1	A	170	LYS
1	A	242	LYS
1	A	328	ARG
1	A	361	LYS
1	A	363	VAL
1	A	382	LEU
1	A	383	GLU
1	A	390	GLU
1	A	405	GLN
1	A	409	ASP
1	A	412	PHE
1	A	416	THR
1	A	460	THR
1	A	462	ARG
1	A	463	LEU
1	B	18	ARG
1	B	37	PHE
1	B	117	LYS
1	B	130	PHE
1	B	152	GLU
1	B	170	LYS
1	B	242	LYS
1	B	328	ARG
1	B	338	ARG
1	B	340	ARG
1	B	390	GLU
1	B	404	GLU
1	B	406	LYS
1	B	410	MET

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Mol	Chain	Res	Type
1	B	416	THR
1	B	458	HIS
1	B	466	ILE
1	C	37	PHE
1	C	117	LYS
1	C	130	PHE
1	C	152	GLU
1	C	170	LYS
1	C	328	ARG
1	C	338	ARG
1	C	360	LYS
1	C	398	ARG
1	C	410	MET
1	C	416	THR
1	C	443	THR
1	C	457	ARG
1	C	464	LYS
1	D	18	ARG
1	D	22	ARG
1	D	37	PHE
1	D	117	LYS
1	D	130	PHE
1	D	152	GLU
1	D	170	LYS
1	D	242	LYS
1	D	328	ARG
1	D	360	LYS
1	D	363	VAL
1	D	383	GLU
1	D	390	GLU
1	D	406	LYS
1	D	416	THR
1	D	444	LEU
1	D	458	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	194	ASN
1	A	374	GLN
1	A	378	ASN
1	B	194	ASN

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Mol	Chain	Res	Type
1	C	194	ASN
1	D	194	ASN
1	D	458	HIS

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

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	A	266	1	21,22,25	275.18	8 (38%)	21,28,34	1.79	4 (19%)
1	LLP	B	266	1	24,24,25	4.57	6 (25%)	30,32,34	2.07	9 (30%)
1	LLP	C	266	1	24,24,25	4.24	8 (33%)	30,32,34	2.26	10 (33%)
1	LLP	D	266	1	24,24,25	4.34	7 (29%)	30,32,34	2.04	11 (36%)

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	266	1	-	0/11/13/19	0/1/1/1
1	LLP	B	266	1	-	0/15/17/19	0/1/1/1
1	LLP	C	266	1	-	0/15/17/19	0/1/1/1
1	LLP	D	266	1	-	0/15/17/19	0/1/1/1

All (29) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	266	LLP	C4-C3	977.89	1.44	1.39
1	A	266	LLP	C4-C5	795.90	1.43	1.39
1	B	266	LLP	O-C	21.05	1.25	1.11
1	A	266	LLP	O-C	20.12	1.25	1.11
1	D	266	LLP	O-C	19.89	1.25	1.11
1	C	266	LLP	O-C	19.11	1.24	1.11
1	C	266	LLP	O3-C3	-3.97	1.27	1.37
1	B	266	LLP	O3-C3	-3.41	1.28	1.37
1	C	266	LLP	C4'-C4	-3.37	1.47	1.51
1	B	266	LLP	C2'-C2	3.20	1.56	1.50
1	B	266	LLP	C4'-NZ	-2.89	1.30	1.45
1	C	266	LLP	C4'-NZ	-2.78	1.31	1.45
1	D	266	LLP	C4'-NZ	-2.76	1.31	1.45
1	C	266	LLP	C2-N1	2.72	1.38	1.33
1	D	266	LLP	O3-C3	-2.68	1.30	1.37
1	A	266	LLP	C2'-C2	2.63	1.55	1.50
1	B	266	LLP	C4'-C4	-2.57	1.48	1.51
1	A	266	LLP	CA-C	2.50	1.53	1.48
1	D	266	LLP	CA-C	2.48	1.52	1.48
1	D	266	LLP	C2-N1	2.41	1.38	1.33
1	D	266	LLP	C2'-C2	2.33	1.55	1.50
1	D	266	LLP	C4'-C4	-2.31	1.49	1.51
1	A	266	LLP	P-OP3	-2.29	1.46	1.54
1	A	266	LLP	C2-N1	2.27	1.38	1.33
1	C	266	LLP	C3-C4	2.17	1.43	1.40
1	C	266	LLP	C2'-C2	2.14	1.54	1.50
1	A	266	LLP	O3-C3	-2.11	1.30	1.35
1	B	266	LLP	C5-C4	2.09	1.43	1.40
1	C	266	LLP	CA-C	2.01	1.52	1.48

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	266	LLP	C-CA-N	-6.45	107.39	113.83
1	B	266	LLP	C-CA-N	-6.19	107.65	113.83
1	D	266	LLP	C-CA-N	-5.63	108.20	113.83
1	C	266	LLP	C4'-NZ-CE	5.44	133.47	113.47
1	B	266	LLP	C4'-NZ-CE	4.89	131.45	113.47
1	A	266	LLP	C-CA-N	-4.63	109.21	113.83
1	C	266	LLP	C3-C4-C5	-4.34	114.13	118.74
1	D	266	LLP	C4'-NZ-CE	3.82	127.51	113.47
1	A	266	LLP	O3-C3-C2	3.56	120.52	116.44
1	B	266	LLP	C5-C6-N1	-3.22	118.04	123.86

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	266	LLP	C6-C5-C4	3.14	120.48	118.10
1	B	266	LLP	C6-N1-C2	3.07	125.86	119.28
1	C	266	LLP	OP3-P-OP4	3.00	114.94	106.65
1	D	266	LLP	C6-C5-C4	2.99	120.37	118.10
1	D	266	LLP	C5-C6-N1	-2.96	118.52	123.86
1	A	266	LLP	OP3-P-OP2	2.91	118.96	107.61
1	C	266	LLP	C4'-C4-C5	2.90	122.36	119.70
1	D	266	LLP	C3-C2-N1	-2.62	117.07	120.66
1	D	266	LLP	C6-N1-C2	2.50	124.63	119.28
1	B	266	LLP	C3-C2-N1	-2.49	117.25	120.66
1	C	266	LLP	C4'-C4-C3	2.48	123.46	120.31
1	D	266	LLP	C4'-C4-C5	2.48	121.97	119.70
1	C	266	LLP	C4-C3-C2	2.45	124.05	119.82
1	D	266	LLP	OP3-P-OP2	2.39	116.93	107.61
1	D	266	LLP	OP3-P-OP1	-2.35	102.75	110.44
1	B	266	LLP	P-OP4-C5'	-2.33	112.81	121.22
1	D	266	LLP	O3-C3-C2	2.31	121.71	117.61
1	C	266	LLP	C3-C2-N1	-2.29	117.52	120.66
1	B	266	LLP	C6-C5-C4	2.29	119.84	118.10
1	A	266	LLP	OP3-P-OP4	2.26	112.89	106.65
1	B	266	LLP	OP3-P-OP2	2.16	116.03	107.61
1	B	266	LLP	OP2-P-OP4	-2.12	100.79	106.65
1	D	266	LLP	C4-C4'-NZ	2.08	115.40	111.52
1	C	266	LLP	C4-C4'-NZ	2.01	115.27	111.52

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

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.

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 ⓘ

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	465/467 (99%)	-0.01	38 (8%) 12 13	8, 21, 64, 89	0
1	B	465/467 (99%)	-0.35	18 (3%) 37 42	8, 21, 64, 89	0
1	C	465/467 (99%)	-0.47	3 (0%) 86 90	8, 21, 64, 88	0
1	D	465/467 (99%)	-0.13	38 (8%) 12 13	7, 21, 64, 89	0
All	All	1860/1868 (99%)	-0.24	97 (5%) 26 29	7, 21, 65, 89	0

All (97) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	401	ALA	10.6
1	A	400	PRO	7.9
1	B	403	GLY	7.6
1	A	401	ALA	7.2
1	D	407	HIS	7.2
1	D	456	LEU	7.2
1	A	399	ASP	6.2
1	D	402	THR	5.9
1	D	400	PRO	5.7
1	A	456	LEU	5.6
1	B	401	ALA	5.6
1	B	407	HIS	5.5
1	D	405	GLN	5.4
1	A	407	HIS	5.4
1	D	404	GLU	5.2
1	B	402	THR	5.2
1	A	404	GLU	5.0
1	A	451	TYR	5.0
1	A	403	GLY	4.8
1	D	442	ALA	4.8
1	D	118	GLU	4.1

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Mol	Chain	Res	Type	RSRZ
1	B	365	GLN	3.9
1	A	455	VAL	3.9
1	A	402	THR	3.9
1	A	23	GLU	3.8
1	A	393	SER	3.8
1	A	445	LYS	3.7
1	A	396	LEU	3.7
1	A	405	GLN	3.7
1	B	404	GLU	3.7
1	A	372	PRO	3.6
1	B	456	LEU	3.5
1	D	454	PRO	3.5
1	B	399	ASP	3.3
1	A	439	GLU	3.3
1	D	410	MET	3.3
1	D	393	SER	3.2
1	A	366	ILE	3.2
1	A	398	ARG	3.2
1	B	396	LEU	3.2
1	C	445	LYS	3.2
1	A	34	TYR	3.2
1	D	408	ALA	3.1
1	B	119	GLY	3.1
1	A	466	ILE	3.1
1	D	399	ASP	3.1
1	D	451	TYR	3.1
1	D	121	ALA	3.0
1	D	439	GLU	3.0
1	A	408	ALA	3.0
1	B	410	MET	2.9
1	A	30	LYS	2.9
1	D	120	LYS	2.9
1	B	400	PRO	2.9
1	A	397	GLY	2.8
1	D	23	GLU	2.8
1	A	454	PRO	2.8
1	D	465	PRO	2.7
1	A	370	GLN	2.7
1	A	26	GLU	2.6
1	A	449	PHE	2.6
1	D	444	LEU	2.6
1	D	372	PRO	2.5

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Mol	Chain	Res	Type	RSRZ
1	D	450	GLU	2.5
1	B	451	TYR	2.5
1	D	409	ASP	2.5
1	B	442	ALA	2.5
1	D	396	LEU	2.5
1	A	410	MET	2.5
1	A	457	ARG	2.4
1	B	364	PRO	2.4
1	C	407	HIS	2.4
1	D	368	GLY	2.4
1	D	455	VAL	2.3
1	A	465	PRO	2.3
1	A	409	ASP	2.3
1	A	442	ALA	2.3
1	B	405	GLN	2.3
1	C	155	PHE	2.2
1	D	403	GLY	2.2
1	D	26	GLU	2.2
1	D	30	LYS	2.2
1	D	466	ILE	2.2
1	A	27	ALA	2.2
1	B	408	ALA	2.2
1	A	21	SER	2.2
1	A	360	LYS	2.2
1	D	22	ARG	2.2
1	B	454	PRO	2.2
1	D	406	LYS	2.2
1	A	368	GLY	2.2
1	D	462	ARG	2.1
1	A	367	PRO	2.1
1	D	28	ALA	2.0
1	D	443	THR	2.0
1	D	445	LYS	2.0
1	D	117	LYS	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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	RSR	LLDF	B-factors(\AA^2)	Q<0.9
1	LLP	C	266	24/25	0.08	0.49	7,12,18,21	0
1	LLP	A	266	23/25	0.07	-0.09	7,12,17,19	0
1	LLP	D	266	24/25	0.06	-0.27	7,12,18,22	0
1	LLP	B	266	24/25	0.05	-1.10	7,12,18,21	0

6.3 Carbohydrates

There are no carbohydrates in this entry.

6.4 Ligands

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	K	C	500	1/1	0.03	-2.20	10,10,10,10	0
2	K	B	500	1/1	0.02	-2.65	10,10,10,10	0
2	K	A	500	1/1	0.01	-3.96	12,12,12,12	0
2	K	D	500	1/1	0.03	-5.30	12,12,12,12	0

6.5 Other polymers

There are no such residues in this entry.