



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:15 PM UTC

PDB ID : 1ZAL / pdb_00001zal
Title : Fructose-1,6-bisphosphate aldolase from rabbit muscle in complex with partially disordered tagatose-1,6-bisphosphate, a weak competitive inhibitor
Authors : St-Jean, M.; Lafrance-Vanasse, J.; Liotard, B.; Sygusch, J.
Deposited on : 2005-04-06
Resolution : 1.89 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

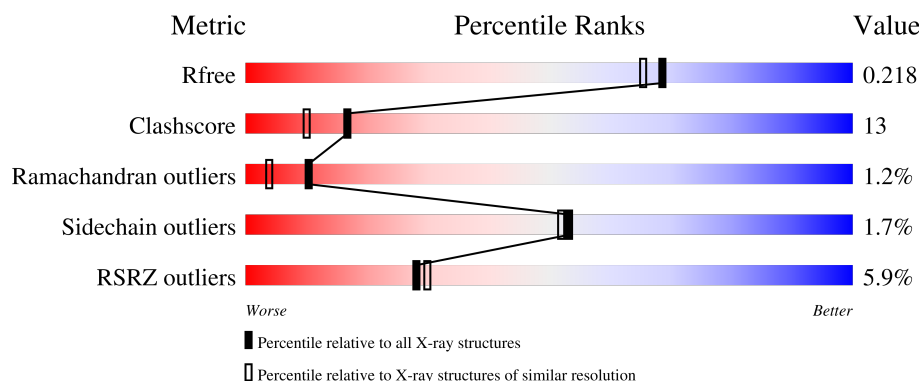
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.89 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	363	6% 72% 26% ..
1	B	363	6% 79% 21% .
1	C	363	6% 70% 28% .
1	D	363	7% 72% 28% .

2 Entry composition [i](#)

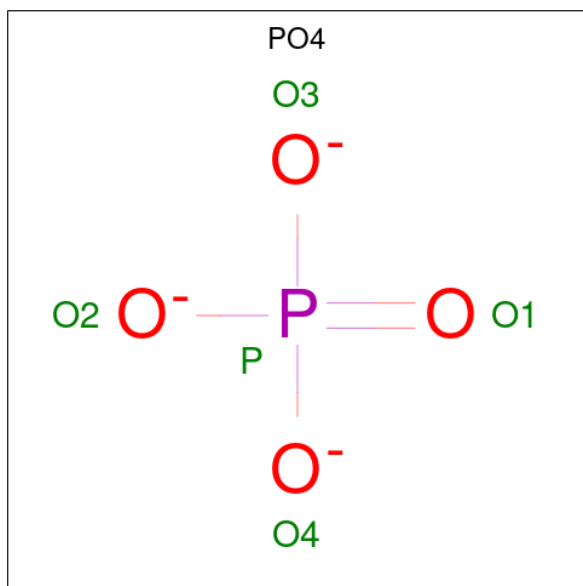
There are 3 unique types of molecules in this entry. The entry contains 13287 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Fructose-bisphosphate aldolase A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	363	Total	C	N	O	S	0	0	0
			2758	1733	489	525	11			
1	B	363	Total	C	N	O	S	0	0	0
			2758	1733	489	525	11			
1	C	363	Total	C	N	O	S	0	0	0
			2758	1733	489	525	11			
1	D	363	Total	C	N	O	S	0	0	0
			2758	1733	489	525	11			

- Molecule 2 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



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

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	O	P	0	1
			5	4	1		
2	B	1	Total	O	P	0	1
			5	4	1		
2	C	1	Total	O	P	0	1
			5	4	1		
2	C	1	Total	O	P	0	1
			5	4	1		
2	D	1	Total	O	P	0	1
			5	4	1		
2	D	1	Total	O	P	0	1
			5	4	1		

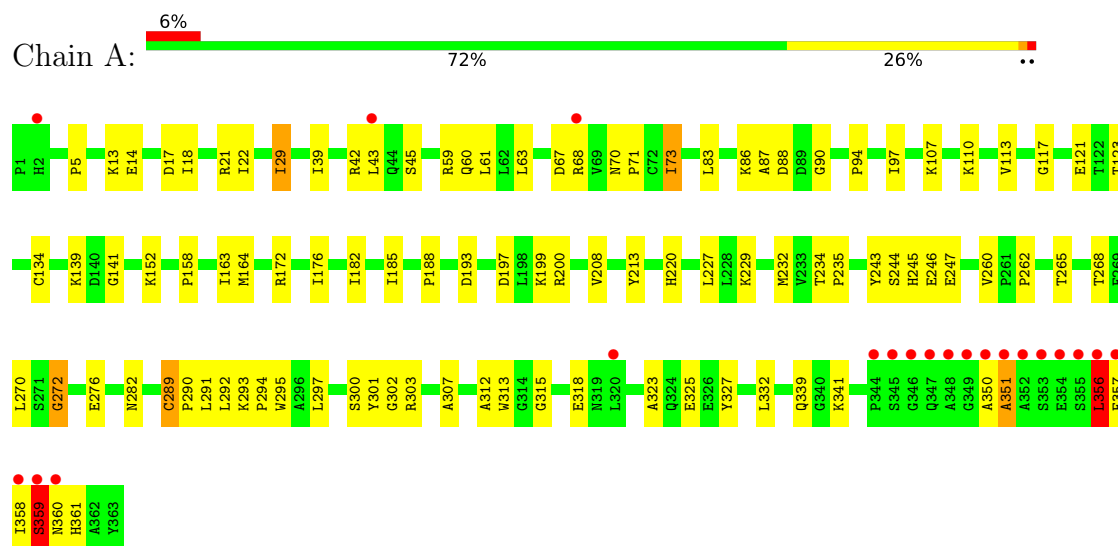
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	496	Total	O	0	7
			496	496		
3	B	538	Total	O	0	8
			538	538		
3	C	592	Total	O	0	8
			592	592		
3	D	589	Total	O	0	8
			589	589		

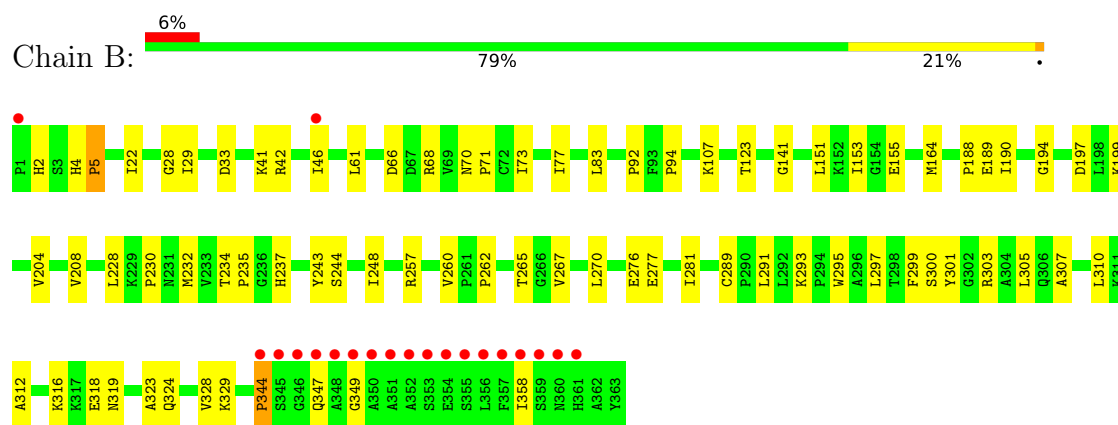
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

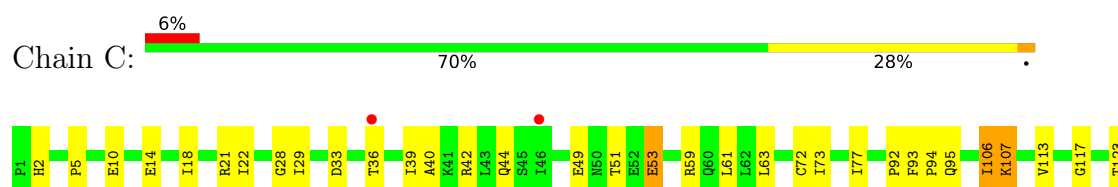
• Molecule 1: Fructose-bisphosphate aldolase A



• Molecule 1: Fructose-bisphosphate aldolase A

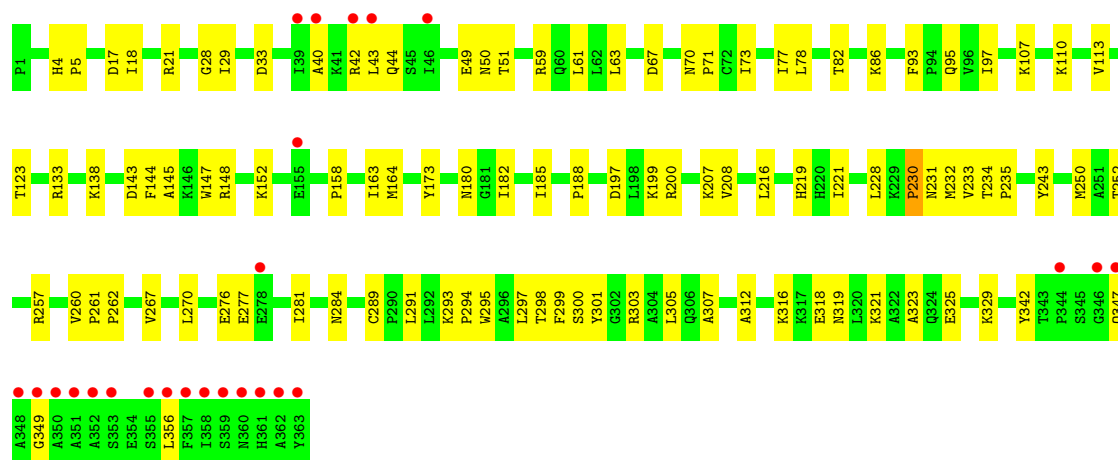
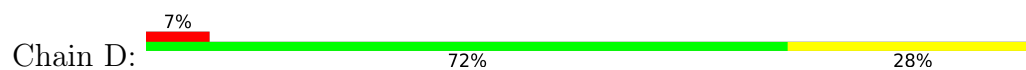


• Molecule 1: Fructose-bisphosphate aldolase A





● Molecule 1: Fructose-bisphosphate aldolase A



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	83.05Å 103.22Å 84.55Å 90.00° 98.56° 90.00°	Depositor
Resolution (Å)	50.00 – 1.89 50.00 – 1.89	Depositor EDS
% Data completeness (in resolution range)	85.4 (50.00-1.89) 92.8 (50.00-1.89)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.89 (at 1.68Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.168 , 0.211 0.176 , 0.218	Depositor DCC
R_{free} test set	6802 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	17.4	Xtriage
Anisotropy	0.655	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 70.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.018 for l,-k,h	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	13287	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.47% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PO4

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	0.39	0/2812	1.00	17/3810 (0.4%)
1	B	0.36	0/2812	0.90	9/3810 (0.2%)
1	C	0.39	0/2812	0.92	7/3810 (0.2%)
1	D	0.36	0/2812	0.90	9/3810 (0.2%)
All	All	0.38	0/11248	0.93	42/15240 (0.3%)

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	A	0	1

There are no bond length outliers.

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	350	ALA	N-CA-C	15.12	127.45	110.97
1	A	260	VAL	N-CA-C	8.91	115.67	107.56
1	A	232	MET	N-CA-C	-8.15	100.02	110.53
1	D	232	MET	N-CA-C	-7.44	100.93	110.53
1	D	260	VAL	N-CA-C	7.43	114.32	107.56
1	C	260	VAL	N-CA-C	7.17	115.00	107.76
1	C	232	MET	N-CA-C	-7.05	101.03	110.55
1	B	232	MET	N-CA-C	-6.95	101.36	110.53
1	B	260	VAL	N-CA-C	6.93	113.87	107.56
1	A	359	SER	N-CA-C	6.76	119.27	111.02
1	A	141	GLY	N-CA-C	6.76	122.89	114.37
1	C	123	THR	N-CA-C	-6.64	99.75	109.18

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	123	THR	N-CA-C	-6.43	100.05	109.18
1	D	289	CYS	CA-C-N	6.33	126.81	119.47
1	D	289	CYS	C-N-CA	6.33	126.81	119.47
1	A	289	CYS	CA-C-N	6.16	126.34	119.32
1	A	289	CYS	C-N-CA	6.16	126.34	119.32
1	B	73	ILE	N-CA-C	6.04	116.31	107.37
1	A	270	LEU	N-CA-C	-6.02	102.43	110.55
1	B	123	THR	N-CA-C	-5.93	100.76	109.18
1	D	73	ILE	N-CA-C	5.82	116.19	107.51
1	A	361	HIS	N-CA-C	-5.76	106.11	114.12
1	C	358	ILE	N-CA-C	-5.74	97.40	109.34
1	C	141	GLY	N-CA-C	5.72	122.65	114.64
1	C	113	VAL	N-CA-C	-5.42	104.75	109.19
1	A	113	VAL	N-CA-C	-5.38	104.77	109.19
1	A	29	ILE	N-CA-C	5.32	116.14	108.48
1	D	148	ARG	N-CA-C	5.26	117.10	108.52
1	B	265	THR	N-CA-C	5.24	116.99	111.28
1	B	289	CYS	CA-C-N	5.21	125.52	119.47
1	B	289	CYS	C-N-CA	5.21	125.52	119.47
1	A	301	TYR	N-CA-C	5.21	118.30	109.76
1	A	302	GLY	N-CA-C	-5.20	105.48	110.21
1	B	141	GLY	N-CA-C	5.18	122.57	115.27
1	D	113	VAL	N-CA-C	-5.17	104.95	109.19
1	C	148	ARG	N-CA-C	5.09	116.81	108.52
1	A	351	ALA	N-CA-C	5.08	118.74	112.54
1	A	123	THR	N-CA-C	-5.07	100.01	110.80
1	D	270	LEU	N-CA-C	-5.04	102.31	110.32
1	A	265	THR	N-CA-C	5.02	116.76	111.28
1	B	189	GLU	N-CA-C	5.01	116.91	109.25
1	A	73	ILE	N-CA-C	5.00	114.77	107.37

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	213	TYR	Sidechain

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen

atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2758	0	2776	80	0
1	B	2758	0	2776	51	0
1	C	2758	0	2776	84	0
1	D	2758	0	2776	77	0
2	A	10	0	0	0	0
2	B	10	0	0	0	0
2	C	10	0	0	0	0
2	D	10	0	0	0	0
3	A	496	0	0	6	0
3	B	538	0	0	6	0
3	C	592	0	0	13	0
3	D	589	0	0	8	0
All	All	13287	0	11104	283	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 13.

All (283) 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:164:MET:HE3	1:C:165:GLU:HA	1.40	1.02
1:A:63:LEU:HD22	1:A:97:ILE:HD11	1.38	1.01
1:A:22:ILE:HG23	1:A:29:ILE:HD11	1.47	0.94
1:C:22:ILE:HG23	1:C:29:ILE:HD11	1.49	0.92
1:A:339:GLN:HB3	1:A:341:LYS:HZ3	1.38	0.88
1:C:22:ILE:CG2	1:C:29:ILE:HD11	2.03	0.86
1:A:22:ILE:CG2	1:A:29:ILE:HD11	2.04	0.86
1:B:22:ILE:HG23	1:B:29:ILE:HD11	1.58	0.84
1:B:22:ILE:CG2	1:B:29:ILE:HD11	2.13	0.78
1:C:250:MET:HE3	1:C:358:ILE:HG13	1.67	0.77
1:B:190:ILE:HD11	1:B:208:VAL:HG21	1.65	0.76
1:A:63:LEU:HD22	1:A:97:ILE:CD1	2.14	0.76
1:C:274:GLN:HB3	1:C:278:GLU:HG3	1.68	0.75
1:B:151:LEU:HB2	1:B:190:ILE:HD13	1.69	0.73
1:C:321:LYS:HE3	1:C:321:LYS:HA	1.70	0.73
1:A:110:LYS:HE3	3:B:3054:HOH:O	1.89	0.73
1:C:164:MET:HE3	1:C:165:GLU:CA	2.17	0.72
1:A:60:GLN:HE22	1:A:88:ASP:H	1.37	0.72
1:D:63:LEU:HD22	1:D:97:ILE:HD11	1.72	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:341:LYS:HB3	1:A:351:ALA:O	1.89	0.71
1:A:39:ILE:HG13	1:A:43:LEU:HD23	1.71	0.70
1:A:60:GLN:NE2	1:A:88:ASP:H	1.90	0.70
1:C:316:LYS:HB2	1:C:319:ASN:ND2	2.07	0.70
1:D:277:GLU:O	1:D:281:ILE:HG12	1.92	0.69
1:C:153:ILE:HD11	1:C:204:VAL:HG11	1.74	0.69
1:A:185:ILE:HD13	1:A:227:LEU:HB2	1.74	0.69
1:B:153:ILE:HD12	1:B:204:VAL:HG21	1.74	0.68
1:A:163:ILE:HD12	1:A:208:VAL:HG22	1.75	0.68
1:B:155:GLU:O	1:C:2:HIS:HE1	1.76	0.68
1:C:95:GLN:HG2	3:C:3212:HOH:O	1.94	0.68
1:D:284:ASN:ND2	1:D:342:TYR:H	1.92	0.67
1:C:153:ILE:CD1	1:C:204:VAL:HG11	2.24	0.66
1:D:321:LYS:HD3	1:D:321:LYS:C	2.21	0.66
1:C:153:ILE:HD12	1:C:204:VAL:HG21	1.78	0.66
1:D:216:LEU:HA	1:D:221:ILE:HD13	1.78	0.65
1:A:200:ARG:HD2	3:D:3366:HOH:O	1.95	0.65
1:D:180:ASN:O	1:D:182:ILE:HD12	1.96	0.65
1:B:66:ASP:OD1	1:B:68:ARG:HD3	1.96	0.65
1:B:197:ASP:HB2	1:B:243:TYR:OH	1.97	0.65
1:D:325:GLU:HG3	1:D:329:LYS:NZ	2.11	0.65
1:C:146:LYS:HE2	1:C:187:GLU:OE1	1.97	0.64
1:C:288:LYS:HD3	3:C:3550:HOH:O	1.97	0.64
1:D:276:GLU:CD	1:D:307:ALA:HB3	2.21	0.64
1:B:293:LYS:HG2	1:B:297:LEU:HD11	1.80	0.64
1:D:70:ASN:HB2	1:D:71:PRO:HD3	1.80	0.64
1:A:293:LYS:HG2	1:A:297:LEU:HD11	1.80	0.64
1:D:163:ILE:HD12	1:D:208:VAL:HG22	1.79	0.64
1:A:134:CYS:HB3	1:A:182:ILE:HD12	1.80	0.63
1:D:293:LYS:HD2	1:D:297:LEU:CD1	2.28	0.63
1:B:83:LEU:HD12	1:B:94:PRO:HG3	1.80	0.63
1:A:68:ARG:HG3	1:A:332:LEU:HD21	1.81	0.62
1:B:228:LEU:HG	1:B:230:PRO:HD3	1.79	0.62
1:A:139:LYS:NZ	1:A:139:LYS:HB3	2.16	0.61
1:A:359:SER:O	1:A:360:ASN:HB3	2.00	0.61
1:B:194:GLY:H	1:B:237:HIS:CE1	2.18	0.61
1:C:106:ILE:HD13	1:C:107:LYS:O	2.01	0.60
1:C:358:ILE:HG21	3:C:3275:HOH:O	2.02	0.60
1:A:185:ILE:HD11	1:A:227:LEU:HD12	1.83	0.60
3:A:3134:HOH:O	1:D:200:ARG:HD2	2.01	0.60
1:A:185:ILE:CD1	1:A:227:LEU:HD12	2.32	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:219:HIS:HB2	1:D:221:ILE:CD1	2.33	0.59
1:C:316:LYS:HD3	3:C:3557:HOH:O	2.02	0.59
1:A:29:ILE:HB	1:A:300:SER:HA	1.83	0.59
1:D:219:HIS:HB2	1:D:221:ILE:HD11	1.85	0.59
1:A:18:ILE:HD13	1:A:21:ARG:NH1	2.18	0.59
1:D:228:LEU:HG	1:D:230:PRO:HD3	1.85	0.58
1:A:290:PRO:CG	1:A:356:LEU:O	2.52	0.58
1:B:153:ILE:CD1	1:B:204:VAL:HG11	2.33	0.58
1:C:250:MET:CE	1:C:358:ILE:HG13	2.34	0.57
1:C:14:GLU:O	1:C:18:ILE:HG12	2.03	0.57
1:C:40:ALA:O	1:C:44:GLN:HG2	2.05	0.57
1:B:277:GLU:O	1:B:281:ILE:HG12	2.05	0.57
1:B:316:LYS:HB2	1:B:319:ASN:ND2	2.20	0.57
1:C:180:ASN:O	1:C:182:ILE:HD12	2.05	0.56
1:D:133:ARG:HD2	3:D:3398:HOH:O	2.04	0.56
1:A:199:LYS:HD3	3:A:3022:HOH:O	2.06	0.56
1:C:244:SER:OG	1:C:247:GLU:HG3	2.06	0.56
1:B:257:ARG:O	1:C:262:PRO:HD2	2.06	0.55
1:A:339:GLN:CB	1:A:341:LYS:HZ3	2.12	0.55
1:A:13:LYS:HE3	1:A:17:ASP:OD1	2.07	0.55
1:C:221:ILE:HD12	1:C:221:ILE:N	2.22	0.55
1:B:291:LEU:O	1:B:293:LYS:HD3	2.06	0.55
1:C:219:HIS:HB2	1:C:221:ILE:CD1	2.37	0.55
1:B:153:ILE:HD11	1:B:204:VAL:HG11	1.89	0.54
1:C:190:ILE:HD11	1:C:208:VAL:HG21	1.88	0.54
1:D:325:GLU:HG3	1:D:329:LYS:HZ2	1.70	0.54
1:A:290:PRO:HG3	1:A:356:LEU:O	2.06	0.54
1:D:356:LEU:HD22	1:D:356:LEU:N	2.23	0.54
1:B:28:GLY:HA3	1:B:299:PHE:CZ	2.42	0.54
1:B:155:GLU:HG3	3:B:3251:HOH:O	2.07	0.54
1:C:176:ILE:HD11	3:D:3208:HOH:O	2.08	0.54
1:B:276:GLU:CD	1:B:307:ALA:HB3	2.33	0.53
1:A:42:ARG:O	1:A:45:SER:HB3	2.08	0.53
1:A:39:ILE:HG13	1:A:43:LEU:CD2	2.38	0.53
1:A:199:LYS:HG3	3:A:3194:HOH:O	2.09	0.53
1:C:29:ILE:HB	1:C:300:SER:HA	1.91	0.53
3:C:3463:HOH:O	1:D:110:LYS:HE3	2.08	0.53
1:C:312:ALA:HB1	3:C:3552:HOH:O	2.09	0.53
1:D:63:LEU:HD22	1:D:97:ILE:CD1	2.39	0.53
1:C:234:THR:HB	1:C:235:PRO:HD2	1.91	0.53
1:C:343:THR:C	1:C:345:SER:H	2.17	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:28:GLY:HA3	1:D:299:PHE:CZ	2.43	0.52
1:C:277:GLU:O	1:C:281:ILE:HG12	2.09	0.52
1:D:221:ILE:HD12	1:D:221:ILE:N	2.23	0.52
1:C:311:LYS:HD3	3:C:3582:HOH:O	2.08	0.52
1:D:298:THR:OG1	1:D:299:PHE:N	2.43	0.52
1:D:250:MET:CE	1:D:291:LEU:HD11	2.39	0.52
1:A:70:ASN:HB2	1:A:71:PRO:HD3	1.92	0.52
1:B:29:ILE:HB	1:B:300:SER:HA	1.92	0.52
1:D:216:LEU:O	1:D:221:ILE:HD13	2.10	0.52
1:B:267:VAL:HB	1:B:297:LEU:HD23	1.92	0.51
1:C:33:ASP:HB3	1:C:77:ILE:HG22	1.93	0.51
1:A:139:LYS:HB3	1:A:139:LYS:HZ3	1.75	0.51
1:C:216:LEU:HA	1:C:221:ILE:HD13	1.93	0.51
1:A:246:GLU:HB3	1:A:358:ILE:CD1	2.41	0.51
1:C:324:GLN:O	1:C:328:VAL:HG23	2.11	0.51
1:A:244:SER:OG	1:A:247:GLU:HG3	2.11	0.50
1:D:29:ILE:HB	1:D:300:SER:HA	1.92	0.50
1:D:312:ALA:HB3	1:D:323:ALA:HA	1.93	0.50
1:B:70:ASN:HB2	1:B:71:PRO:HD3	1.93	0.50
1:C:151:LEU:HB2	1:C:190:ILE:HD13	1.92	0.50
1:C:275:SER:H	1:C:278:GLU:CD	2.20	0.50
1:A:60:GLN:HE22	1:A:88:ASP:N	2.07	0.50
1:C:134:CYS:HB3	1:C:182:ILE:HD13	1.92	0.50
1:A:172:ARG:O	1:A:176:ILE:HG12	2.12	0.50
1:C:291:LEU:O	1:C:293:LYS:HE2	2.12	0.49
1:C:61:LEU:HD23	1:C:61:LEU:C	2.38	0.49
1:D:145:ALA:O	1:D:185:ILE:HD13	2.12	0.49
1:A:61:LEU:C	1:A:61:LEU:HD23	2.38	0.49
1:A:220:HIS:HD2	1:D:207:LYS:NZ	2.11	0.49
1:B:293:LYS:HG2	1:B:297:LEU:CD1	2.40	0.49
1:D:293:LYS:HD2	1:D:297:LEU:HD12	1.95	0.49
1:B:199:LYS:HG3	3:B:3105:HOH:O	2.13	0.49
1:D:63:LEU:HD13	1:D:97:ILE:HD11	1.94	0.49
1:B:324:GLN:O	1:B:328:VAL:HG23	2.13	0.48
1:C:250:MET:HG2	1:C:363:TYR:CE2	2.48	0.48
1:D:93:PHE:O	1:D:97:ILE:HG12	2.13	0.48
1:D:42:ARG:HG3	1:D:42:ARG:HH11	1.77	0.48
1:A:14:GLU:O	1:A:18:ILE:HG12	2.14	0.48
1:A:73:ILE:HD11	1:A:327:TYR:OH	2.14	0.48
1:A:291:LEU:O	1:A:293:LYS:HD3	2.13	0.48
1:B:41:LYS:HD3	3:B:3358:HOH:O	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:216:LEU:O	1:C:221:ILE:HD13	2.14	0.48
1:A:339:GLN:HB3	1:A:341:LYS:NZ	2.20	0.48
1:A:290:PRO:HG2	1:A:356:LEU:O	2.12	0.48
1:C:28:GLY:HA3	1:C:299:PHE:CE1	2.49	0.48
1:D:147:TRP:HB3	1:D:173:TYR:CE2	2.48	0.48
1:A:152:LYS:NZ	1:A:193:ASP:OD1	2.46	0.48
1:B:312:ALA:HB3	1:B:323:ALA:HA	1.96	0.48
1:A:185:ILE:CD1	1:A:227:LEU:HB2	2.42	0.47
1:D:40:ALA:O	1:D:44:GLN:HG3	2.13	0.47
1:A:341:LYS:CB	1:A:351:ALA:O	2.59	0.47
1:D:267:VAL:HB	1:D:297:LEU:HD23	1.96	0.47
1:D:197:ASP:HB2	1:D:243:TYR:OH	2.14	0.47
1:A:245:HIS:HD2	1:A:282:ASN:OD1	1.97	0.47
1:A:117:GLY:HA2	1:B:4:HIS:O	2.15	0.47
1:A:292:LEU:C	1:A:292:LEU:HD23	2.39	0.47
1:C:219:HIS:HB2	1:C:221:ILE:HD11	1.96	0.47
1:D:33:ASP:HB3	1:D:77:ILE:HG22	1.96	0.47
1:D:250:MET:HE2	1:D:291:LEU:HD11	1.96	0.47
1:C:276:GLU:CD	1:C:307:ALA:HB3	2.40	0.47
1:A:73:ILE:CD1	1:A:327:TYR:OH	2.63	0.47
1:A:313:TRP:CE2	1:A:315:GLY:HA2	2.50	0.47
1:B:244:SER:O	1:B:248:ILE:HD13	2.15	0.46
1:D:164:MET:SD	1:D:164:MET:C	2.98	0.46
1:B:28:GLY:HA3	1:B:299:PHE:CE1	2.50	0.46
1:C:28:GLY:HA3	1:C:299:PHE:CZ	2.51	0.46
1:C:53:GLU:HB2	3:C:3112:HOH:O	2.14	0.46
1:A:134:CYS:CB	1:A:182:ILE:HD12	2.45	0.46
1:A:176:ILE:HD11	3:B:3017:HOH:O	2.15	0.46
1:B:92:PRO:HB2	1:B:94:PRO:HD2	1.98	0.46
1:B:164:MET:SD	1:B:164:MET:C	2.98	0.46
1:C:106:ILE:HD13	1:C:106:ILE:C	2.41	0.46
1:C:298:THR:OG1	1:C:299:PHE:N	2.49	0.46
1:A:294:PRO:HG2	1:D:262:PRO:CG	2.46	0.46
1:C:343:THR:C	1:C:345:SER:N	2.73	0.46
1:D:86:LYS:HG3	3:D:3392:HOH:O	2.15	0.46
1:C:358:ILE:HD12	3:C:3275:HOH:O	2.16	0.46
1:D:61:LEU:C	1:D:61:LEU:HD23	2.41	0.46
1:D:49:GLU:HG3	1:D:51:THR:HG23	1.98	0.45
1:A:86:LYS:HD3	1:A:90:GLY:O	2.16	0.45
1:A:272:GLY:HA2	1:A:303:ARG:NH2	2.32	0.45
1:C:293:LYS:HD2	1:C:297:LEU:CD1	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:294:PRO:HG2	1:D:262:PRO:HG3	1.99	0.45
1:D:78:LEU:HD22	1:D:82:THR:HG21	1.98	0.45
1:B:234:THR:HB	1:B:235:PRO:HD2	1.98	0.45
1:C:59:ARG:O	1:C:63:LEU:HG	2.17	0.45
1:D:163:ILE:CD1	1:D:208:VAL:HG22	2.46	0.45
1:D:17:ASP:O	1:D:21:ARG:HG3	2.16	0.45
1:A:59:ARG:O	1:A:63:LEU:HG	2.16	0.45
1:A:234:THR:HB	1:A:235:PRO:HD2	1.99	0.44
1:A:357:PHE:O	3:A:3306:HOH:O	2.21	0.44
1:B:316:LYS:HB2	1:B:319:ASN:HD22	1.80	0.44
1:A:164:MET:SD	1:A:164:MET:C	3.00	0.44
1:B:42:ARG:HH21	1:B:303:ARG:HD3	1.81	0.44
1:C:2:HIS:HB3	3:C:3395:HOH:O	2.16	0.44
1:C:10:GLU:HG2	3:C:3270:HOH:O	2.17	0.44
1:B:316:LYS:HB3	1:B:318:GLU:OE1	2.16	0.44
1:B:66:ASP:CG	1:B:68:ARG:HD3	2.42	0.44
3:A:3442:HOH:O	1:B:2:HIS:HD2	2.00	0.44
1:D:316:LYS:HB2	1:D:319:ASN:ND2	2.33	0.44
1:A:290:PRO:HG2	1:A:357:PHE:HA	1.99	0.44
1:A:325:GLU:HG2	3:A:3351:HOH:O	2.17	0.44
1:B:61:LEU:C	1:B:61:LEU:HD23	2.42	0.44
1:B:262:PRO:CG	1:C:294:PRO:HG3	2.48	0.44
1:C:316:LYS:HD2	1:C:319:ASN:HD21	1.83	0.44
1:A:289:CYS:HA	1:A:290:PRO:HD3	1.88	0.44
1:B:301:TYR:HA	3:B:3009[B]:HOH:O	2.17	0.44
1:A:276:GLU:CD	1:A:307:ALA:HB3	2.43	0.43
1:B:46:ILE:HD13	1:B:310:LEU:HG	2.00	0.43
1:D:28:GLY:HA3	1:D:299:PHE:CE1	2.53	0.43
1:C:228:LEU:HG	1:C:230:PRO:HD3	2.00	0.43
1:C:49:GLU:HG2	1:C:51:THR:HG23	1.99	0.43
1:D:138:LYS:HB2	1:D:182:ILE:HD11	2.00	0.43
1:D:152:LYS:HE2	3:D:3429:HOH:O	2.17	0.43
1:A:262:PRO:HD2	1:D:257:ARG:O	2.19	0.43
1:D:95:GLN:HG3	3:D:3050:HOH:O	2.17	0.43
1:B:301:TYR:HB2	1:B:305:LEU:HG	2.00	0.43
1:C:18:ILE:CD1	1:C:21:ARG:HH22	2.31	0.43
1:D:347:GLN:C	1:D:349:GLY:H	2.25	0.43
1:A:60:GLN:NE2	1:A:88:ASP:N	2.63	0.43
1:C:106:ILE:HD13	1:C:107:LYS:C	2.44	0.43
1:A:139:LYS:NZ	1:A:139:LYS:CB	2.80	0.43
1:D:199:LYS:HG3	3:D:3167:HOH:O	2.17	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:66:ASP:OD2	1:B:68:ARG:NH1	2.52	0.42
1:C:293:LYS:HD2	1:C:297:LEU:HD12	2.00	0.42
1:C:301:TYR:HB3	1:C:304:ALA:HB3	2.00	0.42
1:D:276:GLU:OE2	1:D:307:ALA:HB3	2.17	0.42
1:C:164:MET:HG3	1:C:165:GLU:N	2.34	0.42
1:D:144:PHE:HD2	1:D:185:ILE:HD13	1.83	0.42
1:D:321:LYS:HD3	1:D:321:LYS:O	2.19	0.42
1:D:234:THR:HB	1:D:235:PRO:HD2	2.00	0.42
1:B:329:LYS:HE3	1:B:329:LYS:HB2	1.87	0.42
1:C:337:ALA:C	1:C:339:GLN:H	2.28	0.42
1:D:284:ASN:HD21	1:D:342:TYR:H	1.64	0.42
1:C:207:LYS:NZ	3:C:3471:HOH:O	2.52	0.42
1:D:59:ARG:O	1:D:63:LEU:HG	2.20	0.42
1:A:313:TRP:CZ2	1:A:315:GLY:HA2	2.55	0.42
1:C:313:TRP:CE2	1:C:315:GLY:HA2	2.55	0.42
1:D:158:PRO:CG	1:D:163:ILE:HD11	2.49	0.42
1:C:155:GLU:HA	1:C:155:GLU:OE1	2.20	0.42
1:D:303:ARG:HD2	1:D:356:LEU:HB3	2.02	0.42
1:A:163:ILE:CD1	1:A:208:VAL:HG22	2.47	0.42
1:A:360:ASN:O	1:A:360:ASN:CG	2.62	0.42
1:B:270:LEU:C	1:B:270:LEU:HD12	2.45	0.42
1:A:60:GLN:HE21	1:A:87:ALA:HB1	1.84	0.41
1:B:281:ILE:HD11	1:B:344:PRO:HB3	2.01	0.41
1:C:197:ASP:HB2	1:C:243:TYR:OH	2.20	0.41
1:C:322:ALA:HB1	3:C:3552:HOH:O	2.19	0.41
1:D:301:TYR:HB2	1:D:305:LEU:HG	2.01	0.41
1:A:268:THR:HB	1:A:300:SER:HB2	2.01	0.41
1:A:293:LYS:HE2	1:A:293:LYS:HB2	1.94	0.41
1:D:43:LEU:HD13	1:D:50:ASN:HA	2.01	0.41
1:C:18:ILE:HD13	1:C:21:ARG:HH22	1.85	0.41
1:D:318:GLU:N	1:D:318:GLU:OE1	2.50	0.41
1:B:33:ASP:HB3	1:B:77:ILE:HG22	2.02	0.41
1:C:106:ILE:HD12	1:C:137:TYR:CE2	2.56	0.41
1:D:261:PRO:HA	1:D:262:PRO:HD3	1.93	0.41
1:A:121:GLU:OE2	1:A:158:PRO:HA	2.21	0.41
1:C:93:PHE:N	1:C:94:PRO:CD	2.84	0.41
1:C:176:ILE:CD1	3:D:3208:HOH:O	2.67	0.41
1:A:312:ALA:HB3	1:A:323:ALA:HA	2.01	0.41
1:C:92:PRO:HD2	1:C:95:GLN:HG3	2.02	0.41
1:D:43:LEU:CD1	1:D:50:ASN:HA	2.50	0.41
1:A:83:LEU:HD12	1:A:94:PRO:HG3	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:197:ASP:HB2	1:A:243:TYR:OH	2.20	0.41
1:A:229:LYS:HG3	1:A:268:THR:O	2.21	0.41
1:C:42:ARG:HH21	1:C:303:ARG:HD3	1.86	0.41
1:C:222:TYR:O	1:C:226:THR:HG23	2.21	0.41
1:D:250:MET:CE	1:D:291:LEU:CD1	2.99	0.41
1:C:72:CYS:C	1:C:73:ILE:HD12	2.45	0.41
1:D:219:HIS:CB	1:D:221:ILE:HD11	2.48	0.41
1:C:36:THR:HA	1:C:39:ILE:HG22	2.02	0.40
1:B:194:GLY:H	1:B:237:HIS:HE1	1.66	0.40
1:C:185:ILE:HD12	1:C:185:ILE:N	2.36	0.40
1:D:185:ILE:HD12	1:D:185:ILE:N	2.36	0.40
1:D:231:ASN:H	1:D:231:ASN:ND2	2.18	0.40
1:D:293:LYS:HA	1:D:294:PRO:HD2	1.97	0.40
1:D:18:ILE:HD13	1:D:143:ASP:HB3	2.03	0.40
1:A:67:ASP:HA	1:A:70:ASN:OD1	2.21	0.40
1:D:233:VAL:CG2	1:D:252:THR:HA	2.51	0.40
1:C:117:GLY:HA2	1:D:4:HIS:O	2.22	0.40
1:C:145:ALA:O	1:C:185:ILE:HD13	2.22	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	361/363 (99%)	342 (95%)	15 (4%)	4 (1%)	11	4
1	B	361/363 (99%)	341 (94%)	15 (4%)	5 (1%)	9	3
1	C	361/363 (99%)	337 (93%)	19 (5%)	5 (1%)	9	3
1	D	361/363 (99%)	346 (96%)	12 (3%)	3 (1%)	16	8
All	All	1444/1452 (99%)	1366 (95%)	61 (4%)	17 (1%)	10	4

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	358	ILE
1	C	352	ALA
1	C	358	ILE
1	B	344	PRO
1	B	349	GLY
1	C	351	ALA
1	A	5	PRO
1	B	5	PRO
1	C	188	PRO
1	A	356	LEU
1	B	188	PRO
1	D	67	ASP
1	D	188	PRO
1	A	188	PRO
1	C	5	PRO
1	D	5	PRO
1	A	272	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	291/291 (100%)	286 (98%)	5 (2%)	53	52
1	B	291/291 (100%)	287 (99%)	4 (1%)	59	59
1	C	291/291 (100%)	283 (97%)	8 (3%)	39	34
1	D	291/291 (100%)	288 (99%)	3 (1%)	68	70
All	All	1164/1164 (100%)	1144 (98%)	20 (2%)	53	52

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	107	LYS
1	A	295	TRP
1	A	318	GLU
1	A	356	LEU
1	A	359	SER

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Mol	Chain	Res	Type
1	B	5	PRO
1	B	107	LYS
1	B	295	TRP
1	B	347	GLN
1	C	53	GLU
1	C	106	ILE
1	C	107	LYS
1	C	164	MET
1	C	230	PRO
1	C	295	TRP
1	C	321	LYS
1	C	358	ILE
1	D	107	LYS
1	D	230	PRO
1	D	295	TRP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	60	GLN
1	A	156	HIS
1	A	179	GLN
1	A	180	ASN
1	A	220	HIS
1	A	241	GLN
1	A	245	HIS
1	A	319	ASN
1	A	339	GLN
1	B	95	GLN
1	B	119	ASN
1	B	237	HIS
1	B	287	ASN
1	B	319	ASN
1	C	2	HIS
1	C	4	HIS
1	C	202	GLN
1	C	237	HIS
1	C	241	GLN
1	C	245	HIS
1	C	319	ASN
1	C	339	GLN
1	C	361	HIS

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Mol	Chain	Res	Type
1	D	44	GLN
1	D	54	ASN
1	D	85	GLN
1	D	156	HIS
1	D	231	ASN
1	D	241	GLN
1	D	284	ASN
1	D	319	ASN
1	D	360	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

8 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	$\# Z > 2$
2	PO4	B	3004[A]	-	4,4,4	1.70	0	6,6,6	0.46	0
2	PO4	D	3007[A]	-	4,4,4	1.73	0	6,6,6	0.46	0
2	PO4	A	3002[A]	-	4,4,4	1.75	0	6,6,6	0.47	0
2	PO4	C	3005[A]	-	4,4,4	1.73	0	6,6,6	0.46	0
2	PO4	C	3006[A]	-	4,4,4	1.71	0	6,6,6	0.47	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	PO4	D	3008[A]	-	4,4,4	1.75	0	6,6,6	0.46	0
2	PO4	B	3003[A]	-	4,4,4	1.70	0	6,6,6	0.47	0
2	PO4	A	3001[A]	-	4,4,4	1.72	0	6,6,6	0.47	0

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	363/363 (100%)	0.16	21 (5%)	29 30	11, 22, 44, 64	16 (4%)
1	B	363/363 (100%)	0.17	20 (5%)	30 32	12, 23, 48, 65	16 (4%)
1	C	363/363 (100%)	0.22	20 (5%)	30 32	13, 24, 47, 65	16 (4%)
1	D	363/363 (100%)	0.23	25 (6%)	23 24	12, 24, 62, 103	4 (1%)
All	All	1452/1452 (100%)	0.20	86 (5%)	28 30	11, 23, 50, 103	52 (3%)

All (86) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	356	LEU	8.0
1	A	351	ALA	7.6
1	D	363	TYR	7.6
1	C	350	ALA	7.5
1	A	358	ILE	7.2
1	D	356	LEU	6.9
1	A	357	PHE	6.5
1	B	350	ALA	6.4
1	A	350	ALA	6.3
1	C	346	GLY	6.2
1	C	345	SER	6.1
1	B	358	ILE	6.0
1	B	352	ALA	5.9
1	B	356	LEU	5.9
1	C	356	LEU	5.5
1	B	359	SER	5.5
1	C	358	ILE	5.5
1	B	357	PHE	5.5
1	D	362	ALA	5.4
1	A	355	SER	5.4
1	C	352	ALA	5.3

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Mol	Chain	Res	Type	RSRZ
1	C	351	ALA	5.3
1	B	353	SER	5.1
1	B	351	ALA	4.9
1	C	348	ALA	4.9
1	D	361	HIS	4.8
1	A	352	ALA	4.7
1	D	348	ALA	4.6
1	D	353	SER	4.5
1	D	357	PHE	4.5
1	C	357	PHE	4.5
1	A	359	SER	4.5
1	C	349	GLY	4.5
1	A	348	ALA	4.4
1	B	346	GLY	4.2
1	A	349	GLY	4.1
1	D	358	ILE	4.1
1	B	348	ALA	4.0
1	D	351	ALA	4.0
1	B	345	SER	4.0
1	C	355	SER	3.9
1	B	344	PRO	3.9
1	C	344	PRO	3.9
1	D	350	ALA	3.9
1	A	346	GLY	3.9
1	D	352	ALA	3.8
1	A	345	SER	3.6
1	A	360	ASN	3.6
1	A	353	SER	3.6
1	D	278	GLU	3.6
1	B	349	GLY	3.5
1	C	347	GLN	3.3
1	C	359	SER	3.3
1	C	353	SER	3.3
1	D	349	GLY	3.1
1	A	354	GLU	3.1
1	A	43	LEU	3.1
1	C	354	GLU	3.0
1	A	344	PRO	3.0
1	D	355	SER	2.9
1	B	1	PRO	2.8
1	B	355	SER	2.8
1	B	347	GLN	2.8

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Mol	Chain	Res	Type	RSRZ
1	D	43	LEU	2.7
1	D	346	GLY	2.6
1	D	40	ALA	2.6
1	A	2	HIS	2.6
1	D	46	ILE	2.6
1	A	347	GLN	2.6
1	B	354	GLU	2.6
1	C	36	THR	2.6
1	D	359	SER	2.5
1	B	361	HIS	2.5
1	D	360	ASN	2.4
1	C	46	ILE	2.4
1	A	68	ARG	2.4
1	C	342	TYR	2.4
1	D	39	ILE	2.4
1	D	347	GLN	2.3
1	B	360	ASN	2.3
1	C	360	ASN	2.3
1	D	344	PRO	2.2
1	D	155	GLU	2.2
1	B	46	ILE	2.1
1	A	320	LEU	2.1
1	D	42	ARG	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists the number of atoms with occupancy less than 0.9.

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	PO4	B	3004[A]	5/5	0.84	0.12	52,53,53,54	5
2	PO4	A	3002[A]	5/5	0.86	0.15	45,45,45,45	5
2	PO4	D	3008[A]	5/5	0.86	0.11	41,42,42,42	5
2	PO4	C	3005[A]	5/5	0.87	0.11	37,38,39,40	5
2	PO4	C	3006[A]	5/5	0.88	0.18	41,41,42,43	5
2	PO4	A	3001[A]	5/5	0.88	0.10	33,33,35,35	5
2	PO4	B	3003[A]	5/5	0.90	0.10	37,39,40,40	5
2	PO4	D	3007[A]	5/5	0.97	0.06	24,25,27,27	5

6.5 Other polymers [i](#)

There are no such residues in this entry.