



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:00 AM UTC

PDB ID : 3O5B / pdb_00003o5b
Title : Crystal structure of dimeric KIHxk1 in crystal form VII with glucose bound (open state)
Authors : Kuettner, E.B.; Kettner, K.; Keim, A.; Kriegel, T.M.; Strater, N.
Deposited on : 2010-07-28
Resolution : 1.97 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

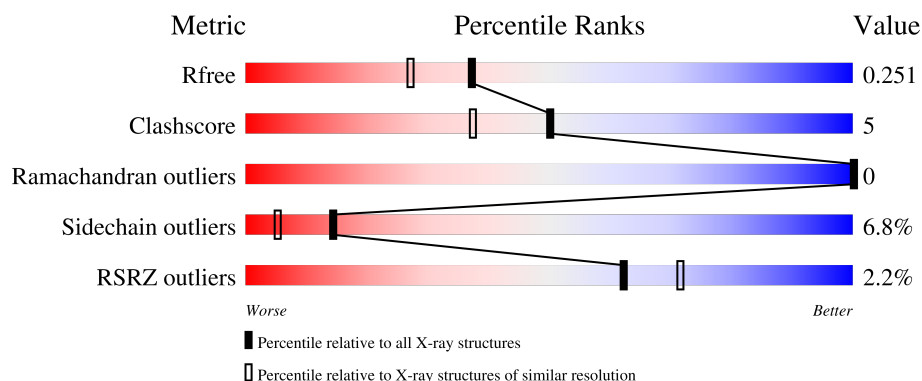
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.97 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	1506 (1.98-1.98)
Clashscore	190562	1534 (1.98-1.98)
Ramachandran outliers	187476	1518 (1.98-1.98)
Sidechain outliers	187428	1518 (1.98-1.98)
RSRZ outliers	180081	1506 (1.98-1.98)

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	485	% 85% 12% ..
1	B	485	4% 81% 15% ..

2 Entry composition [i](#)

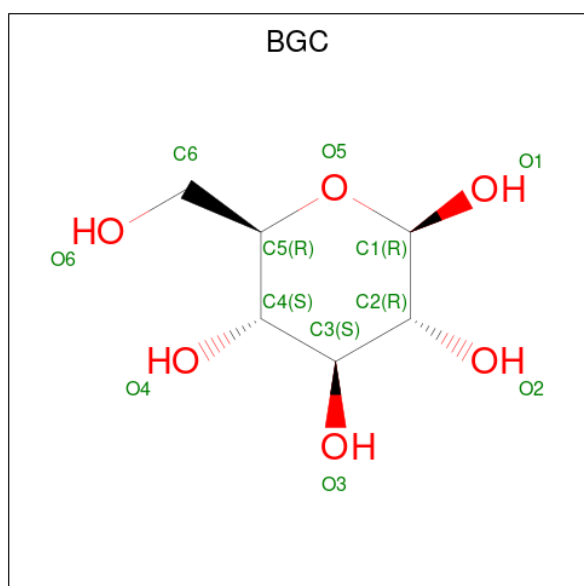
There are 4 unique types of molecules in this entry. The entry contains 7851 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Hexokinase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	479	Total	C	N	O	S	0	2	0
			3736	2377	618	725	16			
1	B	478	Total	C	N	O	S	0	2	0
			3723	2368	616	723	16			

- Molecule 2 is beta-D-glucopyranose (CCD ID: BGC) (formula: C₆H₁₂O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			12	6	6		
2	B	1	Total	C	O	0	0
			12	6	6		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		

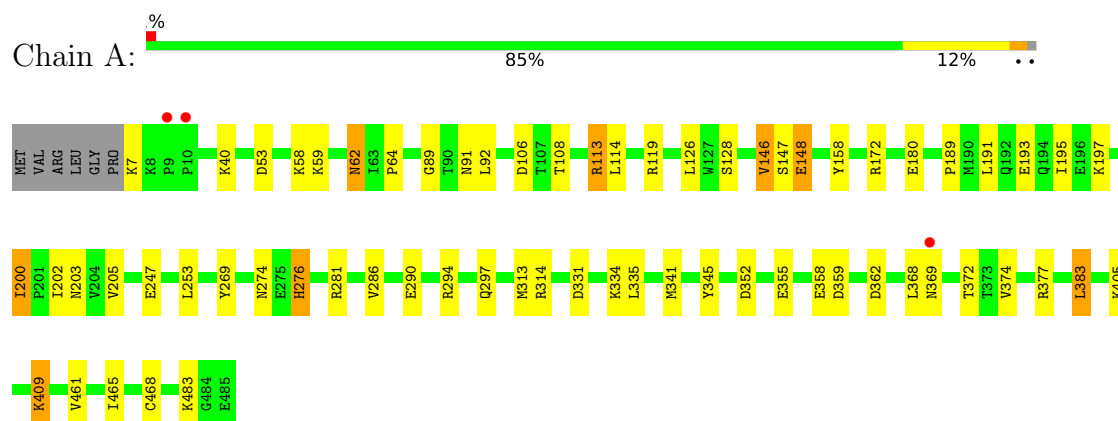
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	211	Total	O	0	0
			211	211		
4	B	142	Total	O	0	0
			142	142		

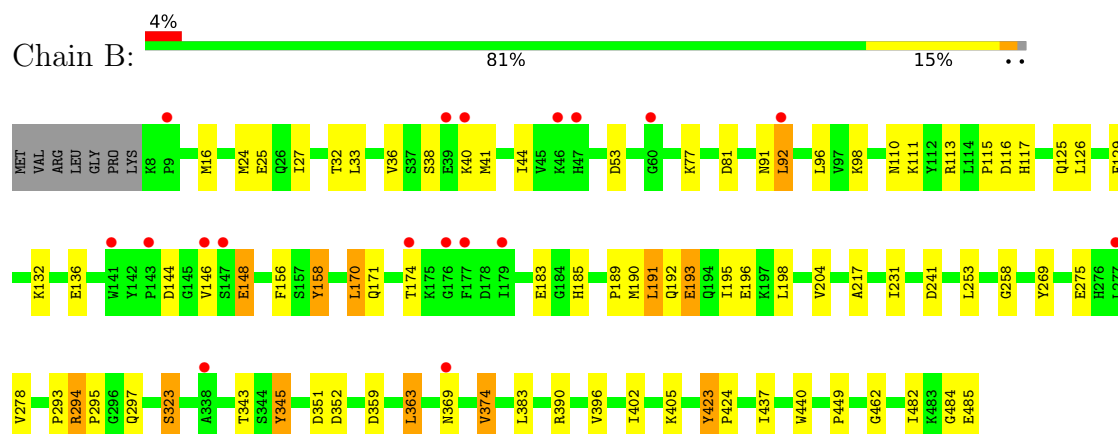
3 Residue-property plots [i](#)

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: Hexokinase



• Molecule 1: Hexokinase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	98.14Å 122.20Å 92.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.90 – 1.97 29.90 – 1.97	Depositor EDS
% Data completeness (in resolution range)	99.6 (29.90-1.97) 99.6 (29.90-1.97)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.11 (at 1.96Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.204 , 0.255 (Not available) , 0.251	Depositor DCC
R_{free} test set	1153 reflections (1.46%)	wwPDB-VP
Wilson B-factor (Å ²)	30.3	Xtriage
Anisotropy	0.619	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 32.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7851	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.39% 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: SO4, BGC

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	1.03	2/3817 (0.1%)	1.10	5/5167 (0.1%)
1	B	0.89	0/3809	1.06	12/5156 (0.2%)
All	All	0.96	2/7626 (0.0%)	1.08	17/10323 (0.2%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	468	CYS	N-CA	5.91	1.53	1.46
1	A	205	VAL	CA-CB	5.22	1.60	1.54

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	374	VAL	CB-CA-C	-7.66	102.16	111.97
1	A	200	ILE	CB-CA-C	-7.24	104.20	110.94
1	B	294	ARG	CA-C-N	-6.63	112.93	119.76
1	B	294	ARG	C-N-CA	-6.63	112.93	119.76
1	B	258	GLY	CA-C-N	6.11	127.47	119.84
1	B	258	GLY	C-N-CA	6.11	127.47	119.84
1	A	148	GLU	CA-C-N	-5.68	113.84	119.92
1	A	148	GLU	C-N-CA	-5.68	113.84	119.92
1	B	345	TYR	CA-C-N	-5.49	113.24	119.28
1	B	345	TYR	C-N-CA	-5.49	113.24	119.28
1	B	374	VAL	N-CA-CB	5.42	116.89	110.55
1	A	158	TYR	N-CA-C	-5.26	99.72	108.23
1	B	323	SER	CB-CA-C	5.21	119.37	110.56
1	B	423	TYR	CA-C-N	5.09	125.33	119.83
1	B	423	TYR	C-N-CA	5.09	125.33	119.83
1	A	108	THR	N-CA-C	-5.02	100.27	108.76
1	B	484	GLY	N-CA-C	-5.01	108.73	114.69

There are no chirality outliers.

There are no planarity outliers.

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	3736	0	3732	37	0
1	B	3723	0	3710	46	0
2	A	12	0	9	0	0
2	B	12	0	12	0	0
3	A	10	0	0	0	0
3	B	5	0	0	0	0
4	A	211	0	0	4	0
4	B	142	0	0	1	0
All	All	7851	0	7463	80	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 5.

All (80) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:191:LEU:HD12	1:B:204:VAL:HG21	1.43	0.96
1:B:345:TYR:CE2	1:B:363:LEU:HD21	2.13	0.83
1:A:313:MET:HG2	1:A:383:LEU:HD22	1.63	0.81
1:B:41:MET:HE2	1:B:437:ILE:HD11	1.66	0.78
1:B:191:LEU:CD1	1:B:204:VAL:HG21	2.15	0.77
1:A:355:GLU:HB2	1:B:110:ASN:HD21	1.47	0.77
1:A:58:LYS:HG3	1:A:247:GLU:HG2	1.71	0.71
1:A:59:LYS:HG3	4:A:584:HOH:O	1.91	0.70
1:A:313:MET:HG2	1:A:383:LEU:CD2	2.23	0.68
1:B:185:HIS:HB2	1:B:190:MET:CE	2.23	0.68
1:B:185:HIS:HB2	1:B:190:MET:HE1	1.77	0.67
1:B:91:ASN:HD22	1:B:113:ARG:HA	1.59	0.66
1:A:358:GLU:HG3	1:B:111:LYS:O	1.98	0.64
1:A:7:LYS:HA	4:A:626:HOH:O	1.98	0.64
1:B:294:ARG:HB3	1:B:297:GLN:HG3	1.82	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:189:PRO:O	1:A:193:GLU:HG2	2.00	0.62
1:B:345:TYR:CE2	1:B:363:LEU:CD2	2.84	0.60
1:A:146:VAL:HG13	1:A:148:GLU:O	2.03	0.58
1:B:192:GLN:O	1:B:196:GLU:HG2	2.04	0.58
1:A:62:ASN:HD21	1:A:274:ASN:ND2	2.02	0.57
1:A:53:ASP:OD1	1:A:405:LYS:HE2	2.05	0.57
1:A:374:VAL:HG23	4:A:619:HOH:O	2.04	0.57
1:B:92:LEU:C	1:B:92:LEU:HD12	2.31	0.56
1:B:53:ASP:OD1	1:B:405:LYS:HE2	2.06	0.55
1:A:286:VAL:O	1:A:290:GLU:HG3	2.06	0.55
1:A:146:VAL:CG1	1:A:148:GLU:O	2.54	0.55
1:B:41:MET:HE2	1:B:437:ILE:CD1	2.36	0.54
1:B:91:ASN:ND2	1:B:113:ARG:HA	2.22	0.54
1:A:383:LEU:C	1:A:383:LEU:HD23	2.33	0.54
1:B:352:ASP:CG	1:B:359:ASP:HB2	2.33	0.53
1:B:36:VAL:CG2	1:B:390:ARG:HG3	2.39	0.53
1:B:241:ASP:OD2	1:B:405:LYS:NZ	2.39	0.52
1:B:191:LEU:C	1:B:191:LEU:HD13	2.35	0.51
1:B:217:ALA:HB2	1:B:462:GLY:HA2	1.92	0.51
1:A:355:GLU:HB2	1:B:110:ASN:ND2	2.22	0.51
1:B:81:ASP:OD1	1:B:98:LYS:HG2	2.10	0.51
1:A:352:ASP:CG	1:A:359:ASP:HB2	2.37	0.50
1:A:383:LEU:HD23	1:A:383:LEU:O	2.11	0.50
1:B:170:LEU:CD1	1:B:183:GLU:HG2	2.42	0.49
1:B:217:ALA:HB2	1:B:462:GLY:CA	2.42	0.49
1:A:91:ASN:HD22	1:A:113:ARG:HA	1.78	0.49
1:A:276:HIS:HB3	1:A:281:ARG:HD3	1.94	0.48
1:A:341:MET:HG2	1:A:345:TYR:CD2	2.49	0.48
1:A:274:ASN:O	1:A:276:HIS:CE1	2.67	0.47
1:B:96:LEU:HD21	1:B:98:LYS:HE3	1.97	0.47
1:A:314:ARG:HD2	1:A:335:LEU:O	2.15	0.47
1:A:409:LYS:HB3	1:A:409:LYS:HE3	1.57	0.47
1:B:189:PRO:O	1:B:193:GLU:HG3	2.16	0.46
1:A:195:ILE:HG23	1:A:200:ILE:HB	1.98	0.46
1:A:200:ILE:HG21	1:A:202:ILE:HD12	1.98	0.46
1:A:294:ARG:NH1	1:A:297:GLN:HG3	2.31	0.46
1:B:36:VAL:HG23	1:B:390:ARG:HG3	1.98	0.45
1:B:38:SER:HB2	4:B:567:HOH:O	2.16	0.45
1:B:146:VAL:HG13	1:B:148:GLU:O	2.16	0.45
1:B:423:TYR:HA	1:B:424:PRO:HD3	1.85	0.44
1:B:482:ILE:HG13	1:B:485:GLU:HG2	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:383:LEU:CD2	1:A:383:LEU:C	2.90	0.44
1:A:89:GLY:O	1:A:119:ARG:HD3	2.18	0.44
1:B:440:TRP:CD2	1:B:449:PRO:HD2	2.53	0.44
1:A:91:ASN:HD21	1:A:113:ARG:HG2	1.83	0.44
1:A:331:ASP:OD2	1:A:331:ASP:C	2.60	0.44
1:B:115:PRO:HB2	1:B:117:HIS:CE1	2.53	0.43
1:B:363:LEU:CD2	1:B:363:LEU:C	2.90	0.43
1:A:334:LYS:HD2	1:A:368:LEU:HA	2.01	0.43
1:B:191:LEU:HD11	1:B:195:ILE:HD11	2.00	0.43
1:B:293:PRO:C	1:B:294:ARG:HD3	2.45	0.42
1:A:334:LYS:HB2	1:A:334:LYS:HE3	1.65	0.42
1:A:372:THR:OG1	1:A:377:ARG:HD3	2.20	0.42
1:B:36:VAL:HG12	1:B:41:MET:HG3	2.02	0.42
1:B:40:LYS:O	1:B:44:ILE:HG13	2.20	0.42
1:A:92:LEU:C	1:A:92:LEU:HD12	2.46	0.41
1:B:156:PHE:HE2	1:B:170:LEU:HD23	1.85	0.41
1:B:44:ILE:HG12	1:B:278:VAL:HG12	2.03	0.41
1:B:16:MET:CE	1:B:27:ILE:HD13	2.51	0.41
1:A:203:ASN:HB2	4:A:566:HOH:O	2.20	0.41
1:B:115:PRO:HD3	1:B:129:PHE:CE1	2.56	0.41
1:B:363:LEU:C	1:B:363:LEU:HD23	2.46	0.40
1:A:461:VAL:O	1:A:465:ILE:HG13	2.20	0.40
1:B:158:TYR:CD1	1:B:158:TYR:N	2.88	0.40
1:B:294:ARG:O	1:B:295:PRO:C	2.63	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	479/485 (99%)	461 (96%)	18 (4%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	478/485 (99%)	459 (96%)	19 (4%)	0	100	100
All	All	957/970 (99%)	920 (96%)	37 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	409/412 (99%)	387 (95%)	22 (5%)	20	9
1	B	408/412 (99%)	375 (92%)	33 (8%)	11	3
All	All	817/824 (99%)	762 (93%)	55 (7%)	14	5

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	LYS
1	A	62	ASN
1	A	64	PRO
1	A	106	ASP
1	A	113	ARG
1	A	114	LEU
1	A	126	LEU
1	A	128	SER
1	A	146	VAL
1	A	147	SER
1	A	172	ARG
1	A	180	GLU
1	A	191	LEU
1	A	197	LYS
1	A	253	LEU
1	A	269	TYR
1	A	276	HIS
1	A	362	ASP

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Mol	Chain	Res	Type
1	A	369	ASN
1	A	383	LEU
1	A	409	LYS
1	A	483	LYS
1	B	24	MET
1	B	25	GLU
1	B	32	THR
1	B	33	LEU
1	B	77	LYS
1	B	92	LEU
1	B	116	ASP
1	B	125	GLN
1	B	126	LEU
1	B	132	LYS
1	B	136	GLU
1	B	144	ASP
1	B	148	GLU
1	B	158	TYR
1	B	170	LEU
1	B	171	GLN
1	B	174	THR
1	B	191	LEU
1	B	193	GLU
1	B	198	LEU
1	B	231	ILE
1	B	253	LEU
1	B	269	TYR
1	B	275	GLU
1	B	323	SER
1	B	343	THR
1	B	351	ASP
1	B	363	LEU
1	B	369	ASN
1	B	374	VAL
1	B	383	LEU
1	B	396	VAL
1	B	402	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	26	GLN

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Mol	Chain	Res	Type
1	A	62	ASN
1	A	91	ASN
1	A	194	GLN
1	A	203	ASN
1	A	209	ASN
1	A	274	ASN
1	A	276	HIS
1	A	369	ASN
1	B	91	ASN
1	B	110	ASN
1	B	117	HIS
1	B	199	ASN
1	B	298	GLN
1	B	367	ASN
1	B	432	GLN
1	B	471	GLN

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](#)

5 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	SO4	A	487	-	4,4,4	0.18	0	6,6,6	0.62	0
3	SO4	A	488	-	4,4,4	0.27	0	6,6,6	0.33	0
2	BGC	A	486	-	12,12,12	0.57	0	17,17,17	2.56	8 (47%)
3	SO4	B	487	-	4,4,4	0.44	0	6,6,6	0.79	0
2	BGC	B	486	-	12,12,12	0.61	0	17,17,17	2.50	7 (41%)

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
2	BGC	A	486	-	-	2/2/22/22	0/1/1/1
2	BGC	B	486	-	-	2/2/22/22	0/1/1/1

There are no bond length outliers.

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	486	BGC	C1-O5-C5	6.11	125.47	113.65
2	B	486	BGC	C1-O5-C5	5.66	124.60	113.65
2	B	486	BGC	C4-C3-C2	-4.90	102.23	110.83
2	A	486	BGC	C1-C2-C3	-3.92	102.36	110.36
2	A	486	BGC	C4-C3-C2	-3.81	104.15	110.83
2	B	486	BGC	O4-C4-C3	3.65	118.97	110.38
2	A	486	BGC	O2-C2-C1	3.18	116.59	109.25
2	B	486	BGC	C1-C2-C3	-3.07	104.09	110.36
2	B	486	BGC	O5-C5-C4	3.00	115.11	109.70
2	A	486	BGC	O3-C3-C4	2.64	116.59	110.38
2	B	486	BGC	O2-C2-C1	2.18	114.29	109.25
2	B	486	BGC	O3-C3-C4	2.17	115.49	110.38
2	A	486	BGC	C6-C5-C4	-2.14	107.75	113.02
2	A	486	BGC	O1-C1-C2	2.13	115.16	108.98
2	A	486	BGC	O3-C3-C2	2.12	115.39	110.38

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	486	BGC	C4-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
2	B	486	BGC	O5-C5-C6-O6
2	A	486	BGC	C4-C5-C6-O6
2	A	486	BGC	O5-C5-C6-O6

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

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	479/485 (98%)	0.12	3 (0%) 85 90	18, 33, 65, 85	2 (0%)
1	B	478/485 (98%)	0.35	18 (3%) 44 54	23, 40, 73, 103	1 (0%)
All	All	957/970 (98%)	0.24	21 (2%) 62 71	18, 36, 69, 103	3 (0%)

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	177	PHE	3.7
1	B	46	LYS	3.6
1	B	174	THR	3.4
1	B	179	ILE	3.3
1	A	10	PRO	3.3
1	B	143	PRO	3.2
1	B	146	VAL	3.1
1	A	9	PRO	2.9
1	B	277	LEU	2.9
1	B	39	GLU	2.7
1	B	9	PRO	2.5
1	A	369	ASN	2.4
1	B	47	HIS	2.4
1	B	369	ASN	2.3
1	B	141	TRP	2.3
1	B	147	SER	2.2
1	B	40	LYS	2.1
1	B	92	LEU	2.1
1	B	60	GLY	2.1
1	B	176	GLY	2.0
1	B	338	ALA	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.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	SO4	A	488	5/5	0.50	0.23	25,29,33,41	5
2	BGC	B	486	12/12	0.78	0.17	27,37,42,46	12
2	BGC	A	486	12/12	0.78	0.15	26,33,40,42	12
3	SO4	B	487	5/5	0.88	0.10	22,26,37,39	5
3	SO4	A	487	5/5	0.92	0.09	22,31,37,37	5

6.5 Other polymers [i](#)

There are no such residues in this entry.