



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 03:40 PM GMT

PDB ID : 4CZU
Title : Crystal structure of the kinase domain of SnRK3.23 At1g30270 T190D mutant
Authors : Chaves-Sanjuan, A.; Sanchez-Barrena, M.J.; Albert, A.
Deposited on : 2014-04-22
Resolution : 1.90 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

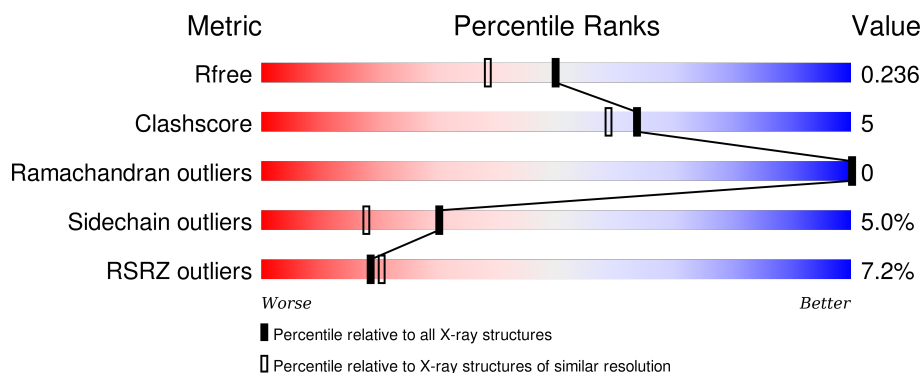
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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}	91344	4755 (1.90-1.90)
Clashscore	102246	5398 (1.90-1.90)
Ramachandran outliers	100387	5338 (1.90-1.90)
Sidechain outliers	100360	5339 (1.90-1.90)
RSRZ outliers	91569	4766 (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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	464	<div> <div>6%</div> <div> <div></div> <div>52%</div> <div>11%</div> <div>•</div> <div>36%</div> </div> </div>
1	B	464	<div> <div>2%</div> <div> <div></div> <div>54%</div> <div>7%</div> <div></div> <div>38%</div> </div> </div>
1	C	464	<div> <div>5%</div> <div> <div></div> <div>55%</div> <div>6%</div> <div></div> <div>39%</div> </div> </div>
1	D	464	<div> <div>5%</div> <div> <div></div> <div>57%</div> <div>7%</div> <div>•</div> <div>35%</div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	CPS	A	1321	-	-	-	X
2	CPS	B	1315	-	-	-	X
2	CPS	C	1314	X	-	-	X
2	CPS	D	1325	-	-	-	X
2	CPS	D	1326	-	-	-	X
3	SO4	B	1319	-	-	-	X
3	SO4	C	1317	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10595 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 CBL-INTERACTING SERINE/THREONINE-PROTEIN KINASE 23.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	297	Total	C	N	O	S	0	5	0
			2384	1532	405	438	9			
1	B	287	Total	C	N	O	S	0	6	0
			2317	1492	393	422	10			
1	C	281	Total	C	N	O	S	0	4	0
			2268	1462	385	412	9			
1	D	302	Total	C	N	O	S	0	3	0
			2425	1554	418	444	9			

There are 28 discrepancies between the modelled and reference sequences:

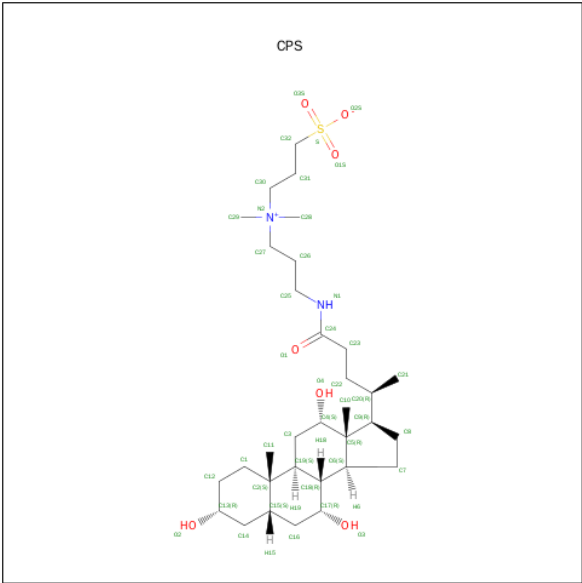
Chain	Residue	Modelled	Actual	Comment	Reference
A	19	PRO	-	EXPRESSION TAG	UNP Q93VD3
A	20	GLY	-	EXPRESSION TAG	UNP Q93VD3
A	21	ILE	-	EXPRESSION TAG	UNP Q93VD3
A	22	HIS	-	EXPRESSION TAG	UNP Q93VD3
A	23	SER	-	EXPRESSION TAG	UNP Q93VD3
A	24	GLY	-	EXPRESSION TAG	UNP Q93VD3
A	190	ASP	THR	ENGINEERED MUTATION	UNP Q93VD3
B	19	PRO	-	EXPRESSION TAG	UNP Q93VD3
B	20	GLY	-	EXPRESSION TAG	UNP Q93VD3
B	21	ILE	-	EXPRESSION TAG	UNP Q93VD3
B	22	HIS	-	EXPRESSION TAG	UNP Q93VD3
B	23	SER	-	EXPRESSION TAG	UNP Q93VD3
B	24	GLY	-	EXPRESSION TAG	UNP Q93VD3
B	190	ASP	THR	ENGINEERED MUTATION	UNP Q93VD3
C	19	PRO	-	EXPRESSION TAG	UNP Q93VD3
C	20	GLY	-	EXPRESSION TAG	UNP Q93VD3
C	21	ILE	-	EXPRESSION TAG	UNP Q93VD3
C	22	HIS	-	EXPRESSION TAG	UNP Q93VD3
C	23	SER	-	EXPRESSION TAG	UNP Q93VD3
C	24	GLY	-	EXPRESSION TAG	UNP Q93VD3

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Chain	Residue	Modelled	Actual	Comment	Reference
C	190	ASP	THR	ENGINEERED MUTATION	UNP Q93VD3
D	19	PRO	-	EXPRESSION TAG	UNP Q93VD3
D	20	GLY	-	EXPRESSION TAG	UNP Q93VD3
D	21	ILE	-	EXPRESSION TAG	UNP Q93VD3
D	22	HIS	-	EXPRESSION TAG	UNP Q93VD3
D	23	SER	-	EXPRESSION TAG	UNP Q93VD3
D	24	GLY	-	EXPRESSION TAG	UNP Q93VD3
D	190	ASP	THR	ENGINEERED MUTATION	UNP Q93VD3

- Molecule 2 is 3-[(3-CHOLAMIDOPROPYL)DIMETHYLAMMONIO]-1-PROPANESULFO NATE (three-letter code: CPS) (formula: C₃₂H₅₈N₂O₇S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			42	32	2	7	1		
2	B	1	Total	C	N	O	S	0	0
			42	32	2	7	1		
2	B	1	Total	C	N	O	S	0	0
			42	32	2	7	1		
2	C	1	Total	C	N	O	S	0	0
			42	32	2	7	1		
2	C	1	Total	C	N	O	S	0	0
			42	32	2	7	1		
2	D	1	Total	C	N	O	S	0	0
			42	32	2	7	1		
2	D	1	Total	C	N	O	S	0	0
			42	32	2	7	1		

- Molecule 3 is SULFATE ION (three-letter code: 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	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		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		

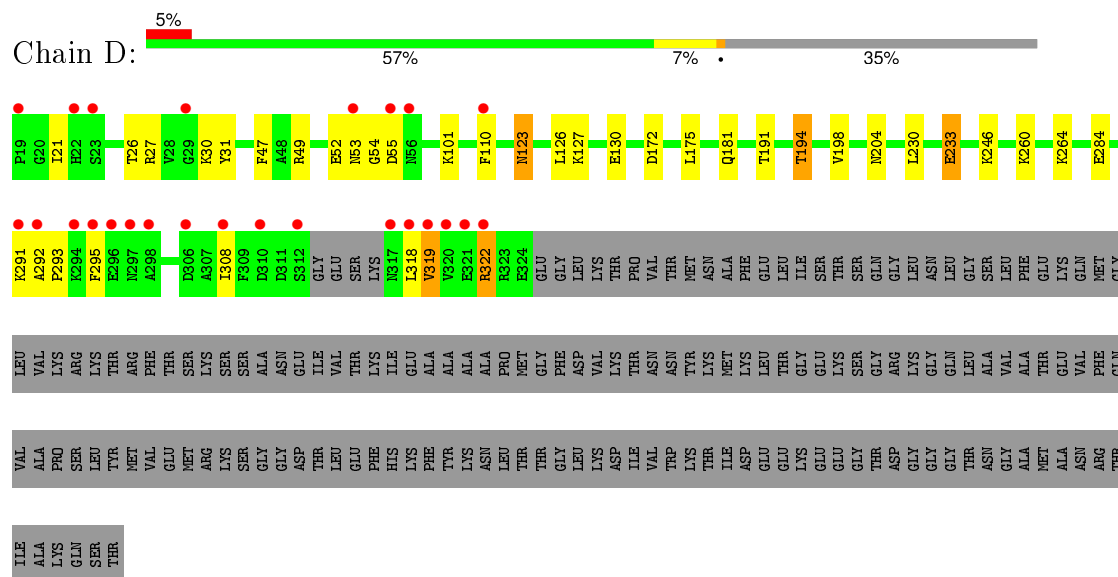
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	246	Total	O	0	0
			246	246		
4	B	234	Total	O	0	0
			234	234		
4	C	166	Total	O	0	0
			166	166		
4	D	176	Total	O	0	0
			176	176		



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	72.70 Å 91.73 Å 207.99 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.97 – 1.90 49.97 – 1.90	Depositor EDS
% Data completeness (in resolution range)	100.0 (49.97-1.90) 96.2 (49.97-1.90)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.18 (at 1.90 Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.192 , 0.233 0.194 , 0.236	Depositor DCC
R_{free} test set	5512 reflections (5.47%)	DCC
Wilson B-factor (Å ²)	24.9	Xtriage
Anisotropy	0.426	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 53.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 110291 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	10595	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.33% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: CPS, SO4

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	0.38	0/2447	0.56	0/3293
1	B	0.40	0/2382	0.55	0/3205
1	C	0.37	0/2326	0.55	0/3131
1	D	0.38	0/2483	0.56	0/3344
All	All	0.38	0/9638	0.56	0/12973

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	2384	0	2426	38	0
1	B	2317	0	2361	26	0
1	C	2268	0	2310	13	0
1	D	2425	0	2456	23	0
2	A	42	0	58	3	0
2	B	84	0	116	2	0
2	C	84	0	116	7	0
2	D	84	0	116	6	0
3	A	20	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	20	0	0	1	0
3	C	25	0	0	0	0
3	D	20	0	0	1	0
4	A	246	0	0	5	0
4	B	234	0	0	6	0
4	C	166	0	0	2	0
4	D	176	0	0	4	0
All	All	10595	0	9959	103	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 (103) 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:A:27:ARG:NH1	1:A:29:GLY:O	2.01	0.94
1:A:123:ASN:O	1:C:35:ARG:NH2	2.22	0.73
1:B:246:LYS:NZ	1:B:248:GLU:OE1	2.22	0.72
1:A:312:SER:HB3	1:A:314:GLU:HG2	1.73	0.70
1:C:27:ARG:NH1	1:C:29:GLY:O	2.26	0.67
1:B:230:LEU:HD12	1:D:191:THR:HG21	1.76	0.66
1:D:246:LYS:HD3	2:D:1326:CPS:H30	1.81	0.62
1:A:140[A]:ASN:OD1	1:A:290:TYR:OH	2.17	0.62
1:B:165:ASN:N	1:B:165:ASN:OD1	2.34	0.60
1:B:119:LYS:NZ	4:B:2094:HOH:O	2.33	0.60
2:C:1314:CPS:H19	2:C:1314:CPS:H10B	1.84	0.60
1:D:126:LEU:HD13	1:D:130:GLU:HG2	1.85	0.58
2:C:1314:CPS:H10A	2:C:1314:CPS:H21A	1.83	0.58
1:A:42:PHE:CD1	1:A:66:LYS:HG3	2.41	0.56
1:B:193:GLY:HA3	1:B:230:LEU:HD11	1.89	0.54
1:B:46:LYS:NZ	4:B:2023:HOH:O	2.40	0.54
1:A:127:LYS:HG2	4:A:2117:HOH:O	2.08	0.54
1:C:66:LYS:NZ	1:C:181:GLN:OE1	2.32	0.53
1:A:112:THR:HA	2:A:1321:CPS:H28B	1.91	0.52
1:B:62:ILE:HD12	1:B:106:PHE:CE2	2.44	0.52
1:D:246:LYS:HE3	2:D:1326:CPS:H32A	1.92	0.51
1:A:312:SER:CB	1:A:314:GLU:HG2	2.38	0.51
3:D:1330:SO4:O1	4:D:2175:HOH:O	2.18	0.51
1:A:29:GLY:H	1:A:314:GLU:HB3	1.75	0.51
1:C:206:LYS:HB2	2:C:1314:CPS:H30A	1.93	0.51
1:A:144:TYR:OH	1:A:148:ARG:NH1	2.43	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:26:THR:HG22	1:D:322:ARG:HG3	1.92	0.51
1:D:204:ASN:ND2	4:D:2110:HOH:O	2.32	0.50
1:D:194:THR:HG23	1:D:198:VAL:HG21	1.93	0.50
1:C:237:LEU:HB3	2:C:1314:CPS:H11A	1.94	0.50
1:D:26:THR:HG22	1:D:322:ARG:HA	1.94	0.50
1:A:238:THR:OG1	2:C:1314:CPS:N1	2.45	0.50
1:B:130:GLU:HG2	4:B:2105:HOH:O	2.11	0.49
1:A:37:LEU:HD21	4:A:2243:HOH:O	2.12	0.49
1:A:236:ASN:HD21	2:C:1314:CPS:H271	1.77	0.49
1:A:69[C]:LYS:HD2	1:B:208:TYR:CD2	2.47	0.49
1:D:318:LEU:HD12	1:D:319:VAL:N	2.28	0.48
1:B:265:ARG:NH1	4:B:2212:HOH:O	2.41	0.48
1:A:65:GLU:HG2	4:A:2052:HOH:O	2.14	0.48
1:A:19:PRO:HB3	1:A:49:ARG:HH12	1.79	0.48
1:D:123:ASN:ND2	4:D:2066:HOH:O	2.46	0.48
1:D:246:LYS:HD3	2:D:1326:CPS:H271	1.96	0.47
1:A:69[A]:LYS:HG3	1:B:208:TYR:CE2	2.50	0.47
1:C:311:ASP:OD1	1:C:312:SER:N	2.48	0.47
1:A:22:HIS:CG	1:A:25:ARG:HG3	2.50	0.47
1:B:93:ARG:NE	3:B:1317:SO4:O3	2.47	0.47
1:A:178:LEU:HB2	1:A:183:ARG:HG3	1.97	0.47
1:A:111:VAL:HG11	1:A:169:LYS:HG3	1.96	0.47
1:B:238:THR:HG22	1:B:242:LYS:HE2	1.98	0.46
1:B:193:GLY:HA3	1:B:230:LEU:CD1	2.45	0.46
1:D:26:THR:CG2	1:D:322:ARG:HG3	2.45	0.46
1:D:181:GLN:H	1:D:181:GLN:CD	2.19	0.46
2:D:1326:CPS:H272	2:D:1326:CPS:C24	2.45	0.45
2:C:1314:CPS:H28	2:C:1314:CPS:H31A	1.64	0.45
1:A:35:ARG:HH11	1:A:35:ARG:HG2	1.80	0.45
1:D:127:LYS:HD3	1:D:127:LYS:HA	1.76	0.45
1:A:60:LYS:HE3	1:A:185:ASP:HB3	1.99	0.45
2:B:1314:CPS:H261	2:B:1314:CPS:H29B	1.71	0.44
2:D:1326:CPS:H31	2:D:1326:CPS:H28	1.73	0.44
1:A:132:ARG:HG3	1:A:285:TRP:CH2	2.53	0.44
1:B:130:GLU:OE1	4:B:2103:HOH:O	2.21	0.44
1:B:81:SER:OG	1:B:308:ILE:HD12	2.18	0.44
1:A:27:ARG:HH12	1:A:30:LYS:HA	1.83	0.43
1:A:78:ARG:NH1	1:B:75:GLN:HG3	2.33	0.43
1:C:194:THR:HB	1:C:195:PRO:HD3	1.99	0.43
1:D:127:LYS:HE3	4:D:2023:HOH:O	2.18	0.43
2:A:1321:CPS:H29A	2:A:1321:CPS:H32	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:148:ARG:HA	1:A:148:ARG:HD3	1.88	0.43
1:B:239:SER:HA	1:B:242:LYS:HE3	2.00	0.43
1:C:144:TYR:OH	1:C:148:ARG:NH1	2.49	0.43
1:C:275:ILE:HG12	1:C:279:GLU:HB2	2.01	0.43
1:D:318:LEU:HD12	1:D:319:VAL:H	1.83	0.43
1:B:153:ARG:NH2	1:B:207:GLY:O	2.51	0.43
1:C:183[C]:ARG:NH2	4:C:2098:HOH:O	2.52	0.43
1:D:260:LYS:HE3	1:D:264:LYS:HE3	2.01	0.42
1:C:59:ILE:HD12	1:C:107:VAL:HG22	2.01	0.42
1:A:93:ARG:HG2	1:A:95:PHE:CE1	2.53	0.42
1:B:246:LYS:HZ2	1:B:248:GLU:HB2	1.84	0.42
1:A:75:GLN:NE2	4:A:2065:HOH:O	2.47	0.42
1:A:146:HIS:CE1	1:A:211:ALA:HA	2.55	0.42
1:C:156:LYS:HE3	4:C:2088:HOH:O	2.19	0.42
1:B:37[B]:LEU:HD21	1:B:47:PHE:CZ	2.55	0.42
1:B:144:TYR:CZ	1:B:148:ARG:CZ	3.02	0.42
1:C:178:LEU:O	1:C:183[A]:ARG:NH1	2.52	0.42
1:A:31:TYR:OH	4:A:2009:HOH:O	2.21	0.42
1:A:206:LYS:HB3	1:A:206:LYS:HE2	1.87	0.42
1:D:292:ALA:HA	1:D:293:PRO:HD3	1.76	0.41
2:B:1315:CPS:H31	2:B:1315:CPS:H271	1.65	0.41
1:B:128:GLU:HB2	4:B:2104:HOH:O	2.20	0.41
1:A:112:THR:HG22	2:A:1321:CPS:H28	2.03	0.41
1:A:33:LEU:HD23	1:A:48:ALA:HB2	2.03	0.41
1:D:27:ARG:HG2	1:D:31:TYR:O	2.21	0.41
1:D:230:LEU:HB2	1:D:233:GLU:HG2	2.03	0.41
1:B:212:LYS:HE2	1:B:274:ARG:HB3	2.03	0.41
1:D:21:ILE:HD11	1:D:47:PHE:HB2	2.03	0.41
1:A:188:LEU:HA	1:A:188:LEU:HD23	1.96	0.41
1:A:69[C]:LYS:HD2	1:B:208:TYR:HD2	1.84	0.41
1:B:133:LYS:HE3	1:B:133:LYS:HB2	1.90	0.41
1:A:303:ASP:HA	1:A:304:ASP:HA	1.65	0.41
1:A:85:LEU:HA	1:A:85:LEU:HD23	1.87	0.40
1:D:246:LYS:CD	2:D:1326:CPS:H271	2.51	0.40
1:D:49:ARG:HE	1:D:54:GLY:HA2	1.87	0.40
1:A:101:LYS:HE3	1:A:101:LYS:HB3	1.91	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	298/464 (64%)	291 (98%)	7 (2%)	0	100	100
1	B	289/464 (62%)	282 (98%)	7 (2%)	0	100	100
1	C	281/464 (61%)	268 (95%)	13 (5%)	0	100	100
1	D	301/464 (65%)	293 (97%)	8 (3%)	0	100	100
All	All	1169/1856 (63%)	1134 (97%)	35 (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	260/393 (66%)	245 (94%)	15 (6%)	25	13
1	B	252/393 (64%)	243 (96%)	9 (4%)	42	30
1	C	245/393 (62%)	235 (96%)	10 (4%)	37	25
1	D	263/393 (67%)	246 (94%)	17 (6%)	21	10
All	All	1020/1572 (65%)	969 (95%)	51 (5%)	30	18

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	21	ILE
1	A	23	SER

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Mol	Chain	Res	Type
1	A	26	THR
1	A	35	ARG
1	A	69[A]	LYS
1	A	69[C]	LYS
1	A	101	LYS
1	A	125	ARG
1	A	172	ASP
1	A	204	ASN
1	A	237	LEU
1	A	294	LYS
1	A	296	GLU
1	A	303	ASP
1	A	308	ILE
1	B	23	SER
1	B	25	ARG
1	B	84	LYS
1	B	110	PHE
1	B	130	GLU
1	B	165	ASN
1	B	205	ASN
1	B	291	LYS
1	B	303	ASP
1	C	27	ARG
1	C	69	LYS
1	C	71	LYS
1	C	72	MET
1	C	155	LEU
1	C	182	VAL
1	C	206	LYS
1	C	294	LYS
1	C	297	ASN
1	C	304	ASP
1	D	30	LYS
1	D	52	GLU
1	D	53	ASN
1	D	55	ASP
1	D	101	LYS
1	D	110	PHE
1	D	123	ASN
1	D	172	ASP
1	D	175	LEU
1	D	194	THR

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Mol	Chain	Res	Type
1	D	233	GLU
1	D	284	GLU
1	D	291	LYS
1	D	295	PHE
1	D	308	ILE
1	D	319	VAL
1	D	322	ARG

Some 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](#)

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

24 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	CPS	A	1321	-	44,45,45	0.20	0	67,70,70	0.59	1 (1%)
3	SO4	A	1322	-	4,4,4	0.27	0	6,6,6	0.13	0
3	SO4	A	1323	-	4,4,4	0.44	0	6,6,6	0.13	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	SO4	A	1324	-	4,4,4	0.22	0	6,6,6	0.16	0
3	SO4	A	1325	-	4,4,4	0.20	0	6,6,6	0.08	0
2	CPS	B	1314	-	44,45,45	0.16	0	67,70,70	0.56	1 (1%)
2	CPS	B	1315	-	44,45,45	0.14	0	67,70,70	0.56	1 (1%)
3	SO4	B	1316	-	4,4,4	0.39	0	6,6,6	0.13	0
3	SO4	B	1317	-	4,4,4	0.43	0	6,6,6	0.12	0
3	SO4	B	1318	-	4,4,4	0.07	0	6,6,6	0.06	0
3	SO4	B	1319	-	4,4,4	0.23	0	6,6,6	0.08	0
2	CPS	C	1313	-	44,45,45	0.17	0	67,70,70	0.50	0
2	CPS	C	1314	-	44,45,45	0.47	0	67,70,70	1.15	5 (7%)
3	SO4	C	1315	-	4,4,4	0.55	0	6,6,6	0.09	0
3	SO4	C	1316	-	4,4,4	0.50	0	6,6,6	0.18	0
3	SO4	C	1317	-	4,4,4	0.30	0	6,6,6	0.07	0
3	SO4	C	1318	-	4,4,4	0.22	0	6,6,6	0.09	0
3	SO4	C	1319	-	4,4,4	0.39	0	6,6,6	0.15	0
2	CPS	D	1325	-	44,45,45	0.18	0	67,70,70	0.54	1 (1%)
2	CPS	D	1326	-	44,45,45	0.19	0	67,70,70	0.47	0
3	SO4	D	1327	-	4,4,4	0.70	0	6,6,6	0.16	0
3	SO4	D	1328	-	4,4,4	0.28	0	6,6,6	0.21	0
3	SO4	D	1329	-	4,4,4	0.48	0	6,6,6	0.17	0
3	SO4	D	1330	-	4,4,4	0.13	0	6,6,6	0.07	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	CPS	A	1321	-	-	1/25/90/90	0/4/4/4
3	SO4	A	1322	-	-	0/0/0/0	0/0/0/0
3	SO4	A	1323	-	-	0/0/0/0	0/0/0/0
3	SO4	A	1324	-	-	0/0/0/0	0/0/0/0
3	SO4	A	1325	-	-	0/0/0/0	0/0/0/0
2	CPS	B	1314	-	-	0/25/90/90	0/4/4/4
2	CPS	B	1315	-	-	0/25/90/90	0/4/4/4
3	SO4	B	1316	-	-	0/0/0/0	0/0/0/0
3	SO4	B	1317	-	-	0/0/0/0	0/0/0/0
3	SO4	B	1318	-	-	0/0/0/0	0/0/0/0
3	SO4	B	1319	-	-	0/0/0/0	0/0/0/0
2	CPS	C	1313	-	-	1/25/90/90	0/4/4/4
2	CPS	C	1314	-	1/1/14/15	0/25/90/90	0/4/4/4
3	SO4	C	1315	-	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	SO4	C	1316	-	-	0/0/0/0	0/0/0/0
3	SO4	C	1317	-	-	0/0/0/0	0/0/0/0
3	SO4	C	1318	-	-	0/0/0/0	0/0/0/0
3	SO4	C	1319	-	-	0/0/0/0	0/0/0/0
2	CPS	D	1325	-	-	0/25/90/90	0/4/4/4
2	CPS	D	1326	-	-	0/25/90/90	0/4/4/4
3	SO4	D	1327	-	-	0/0/0/0	0/0/0/0
3	SO4	D	1328	-	-	0/0/0/0	0/0/0/0
3	SO4	D	1329	-	-	0/0/0/0	0/0/0/0
3	SO4	D	1330	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1315	CPS	C22-C23-C24	-2.36	107.66	113.27
2	C	1314	CPS	C10-C5-C4	-2.05	107.09	109.09
2	C	1314	CPS	C3-C4-C5	2.16	113.39	111.20
2	A	1321	CPS	C27-C26-C25	2.19	115.84	111.34
2	B	1314	CPS	C27-C26-C25	2.75	116.99	111.34
2	C	1314	CPS	C5-C9-C20	2.75	122.85	119.50
2	C	1314	CPS	C10-C5-C6	2.78	115.60	111.22
2	D	1325	CPS	C27-C26-C25	2.85	117.19	111.34
2	C	1314	CPS	C27-C26-C25	4.67	120.93	111.34

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	C	1314	CPS	C5

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	1313	CPS	C32-C31-C30-N2
2	A	1321	CPS	C32-C31-C30-N2

There are no ring outliers.

7 monomers are involved in 20 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1321	CPS	3	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1314	CPS	1	0
2	B	1315	CPS	1	0
3	B	1317	SO4	1	0
2	C	1314	CPS	7	0
2	D	1326	CPS	6	0
3	D	1330	SO4	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	297/464 (64%)	0.43	26 (8%) 12 14	13, 24, 62, 92	0
1	B	287/464 (61%)	0.20	11 (3%) 44 48	9, 22, 47, 95	0
1	C	281/464 (60%)	0.30	22 (7%) 16 17	11, 25, 55, 83	0
1	D	302/464 (65%)	0.51	25 (8%) 14 15	8, 25, 57, 95	0
All	All	1167/1856 (62%)	0.36	84 (7%) 18 20	8, 24, 59, 95	0

All (84) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	21	ILE	11.9
1	D	319	VAL	6.4
1	D	295	PHE	6.2
1	A	319	VAL	6.2
1	D	294	LYS	5.7
1	C	101	LYS	5.0
1	A	317	ASN	4.9
1	C	28	VAL	4.8
1	A	318	LEU	4.8
1	A	297	ASN	4.8
1	A	295	PHE	4.7
1	D	318	LEU	4.7
1	A	292	ALA	4.6
1	A	315	SER	4.6
1	A	303	ASP	4.6
1	C	292	ALA	4.5
1	A	21	ILE	4.2
1	A	294	LYS	4.2
1	D	317	ASN	4.0
1	D	292	ALA	4.0
1	D	321	GLU	3.9

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Mol	Chain	Res	Type	RSRZ
1	B	22	HIS	3.9
1	C	35	ARG	3.9
1	C	290	TYR	3.8
1	D	298	ALA	3.6
1	A	314	GLU	3.6
1	A	24	GLY	3.6
1	A	291	LYS	3.6
1	D	19	PRO	3.6
1	C	294	LYS	3.5
1	D	110	PHE	3.4
1	A	293	PRO	3.4
1	B	24	GLY	3.4
1	D	320	VAL	3.4
1	C	291	LYS	3.3
1	B	296	GLU	3.3
1	A	313	GLY	3.3
1	D	291	LYS	3.2
1	C	310	ASP	3.1
1	D	56	ASN	3.1
1	A	290	TYR	3.0
1	A	320	VAL	2.9
1	D	297	ASN	2.9
1	D	55	ASP	2.8
1	A	23	SER	2.8
1	A	110	PHE	2.8
1	D	322	ARG	2.8
1	D	296	GLU	2.8
1	C	297	ASN	2.7
1	C	293	PRO	2.7
1	D	310	ASP	2.7
1	B	23	SER	2.6
1	B	295	PHE	2.5
1	B	165	ASN	2.5
1	A	22	HIS	2.5
1	A	296	GLU	2.4
1	D	306	ASP	2.4
1	A	284	GLU	2.4
1	C	307	ALA	2.4
1	B	294	LYS	2.3
1	B	303	ASP	2.3
1	C	27	ARG	2.3
1	C	69	LYS	2.3

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Mol	Chain	Res	Type	RSRZ
1	D	312	SER	2.3
1	C	296	GLU	2.3
1	D	22[A]	HIS	2.2
1	C	298	ALA	2.2
1	C	202	VAL	2.2
1	C	29	GLY	2.2
1	A	316	LYS	2.2
1	A	282	GLU	2.2
1	B	25	ARG	2.2
1	C	284	GLU	2.1
1	D	29	GLY	2.1
1	A	27	ARG	2.1
1	C	102	THR	2.1
1	A	123	ASN	2.1
1	C	282	GLU	2.1
1	C	288	LYS	2.1
1	D	53	ASN	2.0
1	C	295	PHE	2.0
1	D	23	SER	2.0
1	B	292	ALA	2.0
1	D	308	ILE	2.0

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

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

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	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	CPS	C	1314	42/42	0.73	0.31	7.49	56,70,114,137	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	CPS	D	1326	42/42	0.79	0.26	4.79	31,42,83,105	0
2	CPS	A	1321	42/42	0.75	0.26	4.68	33,45,76,89	0
2	CPS	B	1315	42/42	0.85	0.17	4.60	25,35,81,113	0
3	SO4	B	1319	5/5	0.86	0.27	4.00	61,64,76,84	0
3	SO4	C	1317	5/5	0.82	0.43	3.67	75,83,98,110	0
2	CPS	D	1325	42/42	0.79	0.24	2.74	32,52,106,131	0
3	SO4	B	1318	5/5	0.72	0.20	1.72	62,64,83,93	0
3	SO4	A	1324	5/5	0.86	0.16	1.62	61,70,76,86	0
2	CPS	C	1313	42/42	0.91	0.16	1.37	21,31,128,140	0
3	SO4	A	1323	5/5	0.90	0.18	1.28	70,77,79,82	0
2	CPS	B	1314	42/42	0.80	0.22	0.96	24,37,99,117	0
3	SO4	C	1319	5/5	0.88	0.15	0.68	80,83,99,104	0
3	SO4	D	1328	5/5	0.75	0.20	0.59	82,93,106,111	0
3	SO4	D	1327	5/5	0.86	0.17	0.21	65,70,73,73	0
3	SO4	C	1318	5/5	0.83	0.26	0.19	86,90,106,106	0
3	SO4	D	1329	5/5	0.86	0.20	0.17	80,90,103,117	0
3	SO4	C	1315	5/5	0.97	0.10	-0.51	40,43,56,56	0
3	SO4	B	1316	5/5	0.96	0.08	-0.81	48,58,62,70	0
3	SO4	B	1317	5/5	0.89	0.15	-0.94	75,78,86,101	0
3	SO4	A	1325	5/5	0.71	0.27	-	79,92,103,111	0
3	SO4	A	1322	5/5	0.90	0.20	-	72,73,97,100	0
3	SO4	C	1316	5/5	0.96	0.18	-	40,47,53,55	0
3	SO4	D	1330	5/5	0.62	0.26	-	70,78,102,104	0

6.5 Other polymers [i](#)

There are no such residues in this entry.