



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 28, 2019 – 05:40 PM EST

PDB ID : 5PZV
Title : Human liver fructose-1,6-bisphosphatase 1 (fructose 1,6-bisphosphate 1-phosphatase, E.C.3.1.3.11) complexed with the allosteric inhibitor 1-(4-bromopyridin-2-yl)-3-(4-chlorophenyl)sulfonylurea
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Deposited on : 2017-04-11
Resolution : 2.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.3 (157068), CSD as539be (2018)
Xtriage (Phenix) : 1.13
EDS : rb-20031633
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)
Refmac : 5.8.0158
CCP4 : 7.0 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20031633

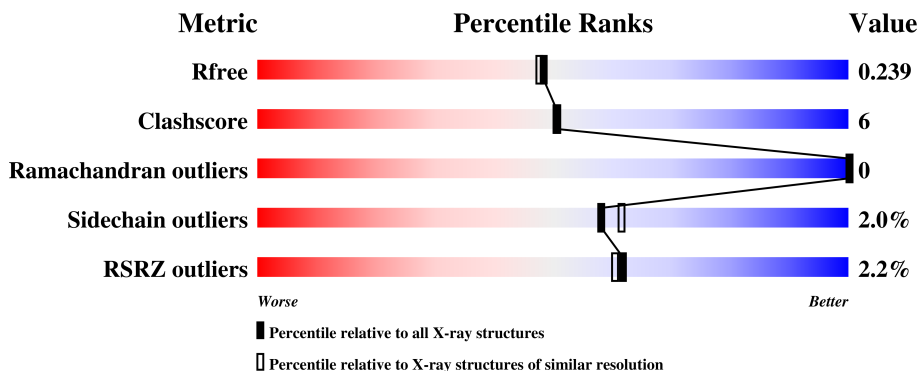
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	111664	7193 (2.00-2.00)
Clashscore	122126	8267 (2.00-2.00)
Ramachandran outliers	120053	8166 (2.00-2.00)
Sidechain outliers	120020	8165 (2.00-2.00)
RSRZ outliers	108989	7011 (2.00-2.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	338	<div style="display: flex; align-items: center;"> <div style="width: 4%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 82%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 12%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 6%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">4% 82% 12% • 6%</p>
1	B	338	<div style="display: flex; align-items: center;"> <div style="width: 2%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 81%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 12%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 6%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">2% 81% 12% 6%</p>
1	C	338	<div style="display: flex; align-items: center;"> <div style="width: 1%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 79%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 15%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 6%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">% 79% 15% • 6%</p>
1	D	338	<div style="display: flex; align-items: center;"> <div style="width: 1%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 84%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 9%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 6%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">% 84% 9% 6%</p>

2 Entry composition [i](#)

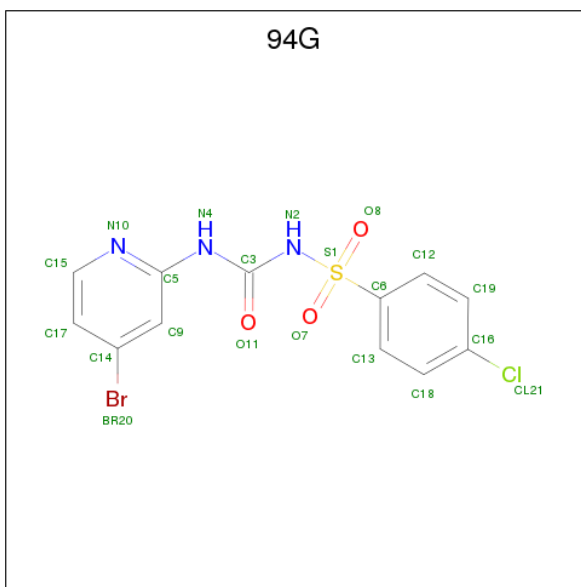
There are 3 unique types of molecules in this entry. The entry contains 10811 atoms, of which 0 are hydrogens and 0 are deuteriums.

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-1,6-bisphosphatase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	319	Total	C	N	O	S	0	0	0
			2441	1556	409	459	17			
1	B	317	Total	C	N	O	S	0	0	0
			2425	1545	406	457	17			
1	C	318	Total	C	N	O	S	0	0	0
			2434	1551	408	458	17			
1	D	317	Total	C	N	O	S	0	0	0
			2425	1545	406	457	17			

- Molecule 2 is N-[(4-bromopyridin-2-yl)carbamoyl]-4-chlorobenzene-1-sulfonamide (three-letter code: 94G) (formula: C₁₂H₉BrClN₃O₃S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	
			Total	Br	C	Cl	N	O			S
2	A	1	Total	Br	C	Cl	N	O	S	0	0
			21	1	12	1	3	3	1		
2	B	1	Total	Br	C	Cl	N	O	S	0	0
			21	1	12	1	3	3	1		

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Mol	Chain	Residues	Atoms							ZeroOcc	AltConf
2	C	1	Total	Br	C	Cl	N	O	S	0	0
			21	1	12	1	3	3	1		
2	D	1	Total	Br	C	Cl	N	O	S	0	0
			21	1	12	1	3	3	1		

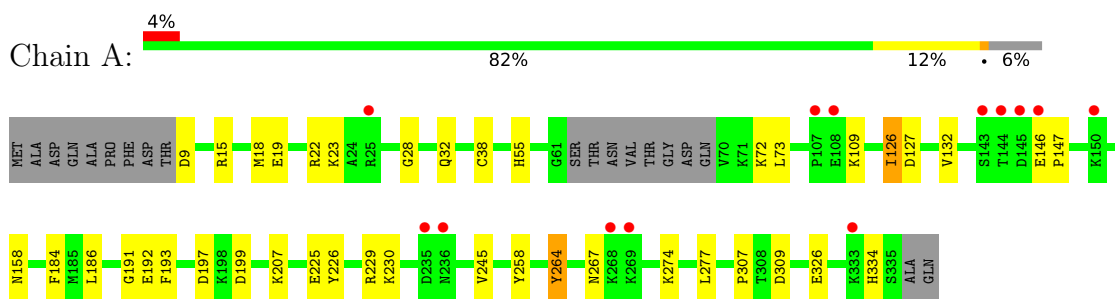
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	219	Total	O	0	0
			219	219		
3	B	226	Total	O	0	0
			226	226		
3	C	282	Total	O	0	0
			282	282		
3	D	275	Total	O	0	0
			275	275		

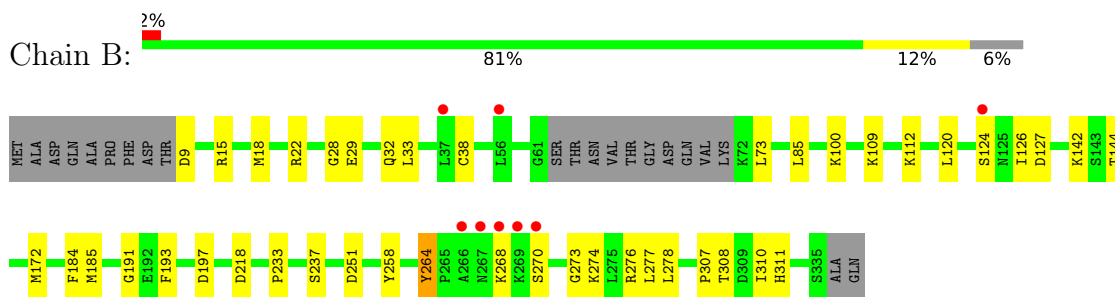
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

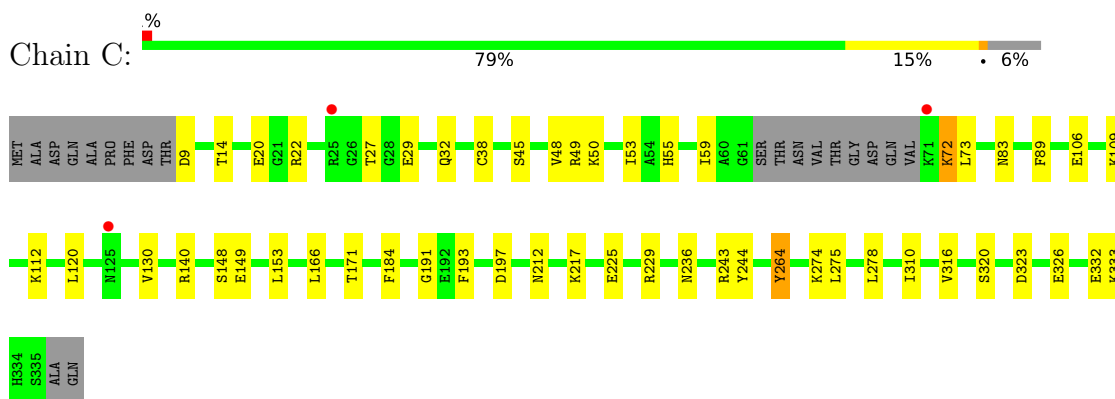
- Molecule 1: Fructose-1,6-bisphosphatase 1



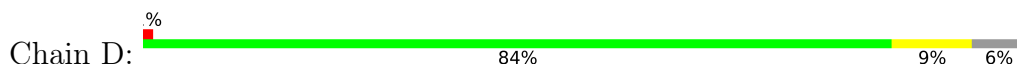
- Molecule 1: Fructose-1,6-bisphosphatase 1

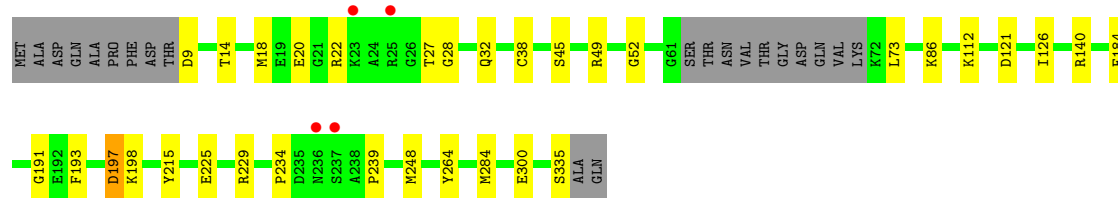


- Molecule 1: Fructose-1,6-bisphosphatase 1



- Molecule 1: Fructose-1,6-bisphosphatase 1





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	67.62Å 83.41Å 277.06Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.67 – 2.00 29.67 – 2.00	Depositor EDS
% Data completeness (in resolution range)	93.8 (29.67-2.00) 93.7 (29.67-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.17 (at 2.00Å)	Xtrriage
Refinement program	PHENIX 1.11.1_2575	Depositor
R, R_{free}	0.191 , 0.239 0.191 , 0.239	Depositor DCC
R_{free} test set	5009 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	23.7	Xtrriage
Anisotropy	0.251	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 64.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.38$, $\langle L^2 \rangle = 0.21$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10811	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.80% 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: 94G

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	0.36	0/2484	0.56	0/3354
1	B	0.38	0/2468	0.57	0/3333
1	C	0.41	0/2477	0.57	0/3344
1	D	0.41	0/2468	0.58	0/3333
All	All	0.39	0/9897	0.57	0/13364

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	2441	0	2491	34	0
1	B	2425	0	2469	32	0
1	C	2434	0	2482	39	0
1	D	2425	0	2469	27	0
2	A	21	0	0	2	0
2	B	21	0	0	2	0
2	C	21	0	0	1	0
2	D	21	0	0	4	0
3	A	219	0	0	13	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	226	0	0	10	0
3	C	282	0	0	15	0
3	D	275	0	0	12	0
All	All	10811	0	9911	123	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 6.

All (123) 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:185:MET:SD	3:B:625:HOH:O	2.21	0.98
1:A:55:HIS:NE2	3:A:502:HOH:O	1.97	0.96
1:C:149:GLU:O	3:C:501:HOH:O	1.89	0.90
1:C:32:GLN:NE2	3:C:504:HOH:O	2.07	0.87
1:A:158:ASN:O	3:A:501:HOH:O	1.95	0.84
1:B:9:ASP:N	3:B:501:HOH:O	2.11	0.81
1:B:32:GLN:OE1	1:D:22:ARG:NH1	2.13	0.80
1:C:229:ARG:NH1	1:C:326:GLU:OE2	2.13	0.80
1:A:9:ASP:N	3:A:506:HOH:O	2.15	0.80
1:C:22:ARG:NH2	3:C:508:HOH:O	2.16	0.78
1:A:22:ARG:NH1	1:C:32:GLN:OE1	2.16	0.78
1:C:83:ASN:OD1	3:C:503:HOH:O	2.03	0.75
1:A:72:LYS:NZ	3:A:507:HOH:O	2.20	0.75
1:B:18:MET:SD	3:B:694:HOH:O	2.48	0.72
1:A:309:ASP:OD2	3:A:504:HOH:O	2.09	0.71
1:D:86:LYS:NZ	3:D:503:HOH:O	2.13	0.71
1:D:121:ASP:OD2	3:D:501:HOH:O	2.09	0.70
1:B:218:ASP:OD1	1:B:268:LYS:NZ	2.22	0.68
1:B:172:MET:HB2	3:B:625:HOH:O	1.94	0.66
1:C:332:GLU:OE2	3:C:506:HOH:O	2.13	0.66
1:A:192:GLU:OE2	3:A:505:HOH:O	2.13	0.66
1:A:277:LEU:HD23	1:A:307:PRO:HB3	1.78	0.66
1:D:284:MET:HE2	3:D:682:HOH:O	1.96	0.66
1:B:308:THR:O	3:B:502:HOH:O	2.15	0.64
1:D:20:GLU:OE1	3:D:504:HOH:O	2.14	0.64
1:D:225:GLU:OE2	1:D:229:ARG:NH2	2.31	0.64
1:A:225:GLU:OE1	1:A:334:HIS:NE2	2.25	0.64
1:B:124:SER:O	3:B:503:HOH:O	2.15	0.63
1:B:22:ARG:NH1	1:D:32:GLN:OE1	2.30	0.63
1:C:275:LEU:HD12	1:C:316:VAL:HG11	1.83	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:333:LYS:NZ	3:C:517:HOH:O	2.35	0.60
1:C:153:LEU:HG	3:C:501:HOH:O	2.02	0.59
1:A:109:LYS:NZ	3:A:517:HOH:O	2.34	0.58
3:C:613:HOH:O	1:D:52:GLY:HA2	2.03	0.58
1:C:9:ASP:N	3:C:519:HOH:O	2.38	0.57
1:A:264:TYR:OH	1:A:274:LYS:HD3	2.04	0.56
1:A:55:HIS:CD2	3:A:502:HOH:O	2.53	0.56
1:D:215:TYR:OH	3:D:505:HOH:O	2.18	0.56
1:B:109:LYS:NZ	1:D:9:ASP:OD2	2.25	0.56
1:B:251:ASP:OD2	3:B:504:HOH:O	2.18	0.56
1:B:73:LEU:HD23	1:B:126:ILE:HD13	1.89	0.55
1:C:184:PHE:HB3	1:C:193:PHE:HB3	1.89	0.55
1:A:22:ARG:HH22	1:C:29:GLU:HG3	1.72	0.53
1:C:9:ASP:N	3:C:521:HOH:O	2.41	0.53
1:D:197:ASP:OD1	3:D:506:HOH:O	2.19	0.53
1:B:28:GLY:HA2	2:B:401:94G:C3	2.40	0.52
1:D:234:PRO:HG3	3:D:688:HOH:O	2.09	0.52
1:B:191:GLY:HA3	1:D:191:GLY:HA3	1.91	0.52
1:B:29:GLU:OE1	1:B:112:LYS:HG2	2.10	0.52
1:A:19:GLU:O	1:A:23:LYS:HG3	2.10	0.52
1:B:28:GLY:HA2	2:B:401:94G:O11	2.10	0.51
1:D:248:MET:SD	3:D:682:HOH:O	2.60	0.51
1:D:73:LEU:HD23	1:D:126:ILE:HD13	1.92	0.50
1:C:50:LYS:HD2	1:C:53:ILE:HD12	1.94	0.50
1:C:55:HIS:HA	1:C:59:ILE:HG22	1.93	0.50
1:D:198:LYS:NZ	3:D:524:HOH:O	2.44	0.50
1:D:300:GLU:HG2	3:D:513:HOH:O	2.11	0.50
1:B:270:SER:HB3	1:B:273:GLY:O	2.12	0.49
1:C:217:LYS:HG2	3:C:646:HOH:O	2.11	0.49
1:A:146:GLU:CG	1:A:147:PRO:HD2	2.42	0.49
1:D:28:GLY:HA2	2:D:401:94G:O11	2.13	0.49
1:B:276:ARG:NH1	1:B:311:HIS:O	2.42	0.49
1:C:112:LYS:HB2	1:C:140:ARG:CZ	2.43	0.49
1:B:270:SER:HB3	1:B:273:GLY:C	2.33	0.49
1:A:184:PHE:HB3	1:A:193:PHE:HB3	1.96	0.48
1:D:9:ASP:N	3:D:526:HOH:O	2.46	0.48
1:A:267:ASN:ND2	3:A:526:HOH:O	2.42	0.48
1:B:184:PHE:HB3	1:B:193:PHE:HB3	1.96	0.48
1:A:264:TYR:CZ	1:A:274:LYS:HD3	2.49	0.47
1:C:264:TYR:OH	1:C:274:LYS:HD3	2.15	0.47
1:B:264:TYR:OH	1:B:274:LYS:HD3	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:191:GLY:HA3	1:C:191:GLY:HA3	1.96	0.47
1:A:127:ASP:HB2	1:B:258:TYR:OH	2.15	0.46
1:A:22:ARG:HG2	2:A:401:94G:C15	2.45	0.46
1:C:148:SER:HB2	3:C:731:HOH:O	2.15	0.46
1:B:233:PRO:HG2	1:B:237:SER:O	2.16	0.46
1:A:199:ASP:OD2	3:A:508:HOH:O	2.20	0.46
1:B:142:LYS:NZ	3:B:516:HOH:O	2.42	0.46
1:B:100:LYS:HE3	1:B:100:LYS:HB2	1.75	0.46
1:D:22:ARG:HG2	2:D:401:94G:C17	2.46	0.46
1:B:33:LEU:CD1	1:B:85:LEU:HD22	2.46	0.46
1:C:243:ARG:NH2	3:C:505:HOH:O	2.13	0.45
1:D:239:PRO:O	3:D:507:HOH:O	2.20	0.45
1:A:207:LYS:NZ	3:A:510:HOH:O	2.24	0.45
1:D:18:MET:HE3	2:D:401:94G:BR20	2.71	0.45
1:A:15:ARG:HG3	3:A:585:HOH:O	2.17	0.45
1:D:184:PHE:HB3	1:D:193:PHE:HB3	1.99	0.45
1:C:120:LEU:HA	1:C:120:LEU:HD12	1.70	0.44
1:C:20:GLU:HB3	2:C:401:94G:CL21	2.54	0.44
1:D:112:LYS:HB2	1:D:140:ARG:NH1	2.33	0.44
1:C:112:LYS:HD2	1:C:140:ARG:NE	2.32	0.44
1:D:112:LYS:HD2	1:D:140:ARG:NE	2.33	0.44
1:C:48:VAL:HA	1:C:73:LEU:HD21	2.00	0.44
1:B:278:LEU:HG	3:B:554:HOH:O	2.17	0.44
1:C:275:LEU:HD12	1:C:316:VAL:CG1	2.47	0.44
1:C:264:TYR:CZ	1:C:274:LYS:HD3	2.53	0.44
1:A:226:TYR:CE1	1:A:230:LYS:HE3	2.53	0.43
1:A:258:TYR:OH	1:B:127:ASP:HB2	2.19	0.43
1:C:45:SER:O	1:C:49:ARG:HD3	2.19	0.43
1:D:45:SER:O	1:D:49:ARG:HD3	2.19	0.43
1:C:212:ASN:HB2	1:C:244:TYR:CE2	2.53	0.42
1:A:229:ARG:NH1	1:A:326:GLU:OE2	2.47	0.42
1:B:120:LEU:HD12	1:B:120:LEU:HA	1.76	0.42
1:C:72:LYS:HB2	1:C:72:LYS:HE3	1.72	0.42
1:A:146:GLU:HG3	1:A:147:PRO:HD2	2.02	0.42
1:C:27:THR:HG23	1:C:29:GLU:H	1.84	0.42
1:A:73:LEU:HD23	1:A:73:LEU:HA	1.72	0.42
1:A:28:GLY:HA2	2:A:401:94G:O11	2.19	0.42
1:B:15:ARG:HD3	3:B:506:HOH:O	2.20	0.42
1:B:278:LEU:HD22	1:B:310:ILE:HA	2.02	0.42
1:C:130:VAL:HA	3:C:614:HOH:O	2.21	0.41
1:D:18:MET:CE	2:D:401:94G:BR20	3.23	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:106:GLU:HG2	3:C:569:HOH:O	2.19	0.41
1:C:225:GLU:O	1:C:229:ARG:HG3	2.20	0.41
1:C:278:LEU:HD22	1:C:310:ILE:HA	2.02	0.41
1:A:18:MET:O	1:A:22:ARG:HG3	2.19	0.41
1:C:320:SER:HB2	1:C:323:ASP:OD2	2.21	0.41
1:C:166:LEU:O	1:C:171:THR:HA	2.21	0.41
1:A:126:ILE:HD13	1:A:132:VAL:HG13	2.03	0.41
1:C:89:PHE:CD1	1:C:109:LYS:HA	2.56	0.41
1:A:186:LEU:HB2	1:A:193:PHE:CE1	2.56	0.40
1:B:277:LEU:HD23	1:B:307:PRO:HB3	2.03	0.40
1:A:245:VAL:HG21	3:A:695:HOH:O	2.21	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	315/338 (93%)	310 (98%)	5 (2%)	0	100	100
1	B	313/338 (93%)	307 (98%)	6 (2%)	0	100	100
1	C	314/338 (93%)	309 (98%)	5 (2%)	0	100	100
1	D	313/338 (93%)	308 (98%)	5 (2%)	0	100	100
All	All	1255/1352 (93%)	1234 (98%)	21 (2%)	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	266/281 (95%)	261 (98%)	5 (2%)	60	64
1	B	264/281 (94%)	260 (98%)	4 (2%)	67	72
1	C	265/281 (94%)	259 (98%)	6 (2%)	53	56
1	D	264/281 (94%)	258 (98%)	6 (2%)	53	56
All	All	1059/1124 (94%)	1038 (98%)	21 (2%)	58	61

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	GLN
1	A	38	CYS
1	A	126	ILE
1	A	197	ASP
1	A	264	TYR
1	B	38	CYS
1	B	144	THR
1	B	197	ASP
1	B	264	TYR
1	C	14	THR
1	C	38	CYS
1	C	72	LYS
1	C	197	ASP
1	C	236	ASN
1	C	264	TYR
1	D	14	THR
1	D	27	THR
1	D	38	CYS
1	D	197	ASP
1	D	264	TYR
1	D	335	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	236	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

4 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	94G	A	401	-	22,22,22	0.23	0	30,31,31	0.43	0
2	94G	B	401	-	22,22,22	0.38	0	30,31,31	0.54	0
2	94G	C	401	-	22,22,22	0.29	0	30,31,31	0.43	0
2	94G	D	401	-	22,22,22	0.29	0	30,31,31	0.49	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	94G	A	401	-	-	0/15/15/15	0/2/2/2
2	94G	B	401	-	-	0/15/15/15	0/2/2/2
2	94G	C	401	-	-	0/15/15/15	0/2/2/2
2	94G	D	401	-	-	0/15/15/15	0/2/2/2

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.

4 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	94G	2	0
2	B	401	94G	2	0
2	C	401	94G	1	0
2	D	401	94G	4	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	319/338 (94%)	0.10	13 (4%) 37 37	15, 29, 64, 92	0
1	B	317/338 (93%)	-0.00	8 (2%) 57 56	12, 27, 57, 92	0
1	C	318/338 (94%)	-0.14	3 (0%) 84 83	12, 22, 47, 74	0
1	D	317/338 (93%)	-0.23	4 (1%) 77 76	11, 21, 43, 72	0
All	All	1271/1352 (94%)	-0.07	28 (2%) 62 60	11, 24, 53, 92	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	71	LYS	3.8
1	A	146	GLU	3.6
1	A	25	ARG	3.6
1	D	25	ARG	3.5
1	A	235	ASP	3.4
1	B	269	LYS	3.4
1	B	268	LYS	3.3
1	A	236	ASN	3.3
1	A	144	THR	3.1
1	A	269	LYS	3.0
1	B	266	ALA	3.0
1	C	125	ASN	3.0
1	B	267	ASN	2.9
1	D	236	ASN	2.9
1	B	270	SER	2.7
1	A	268	LYS	2.7
1	A	333	LYS	2.5
1	B	37	LEU	2.4
1	A	143	SER	2.4
1	A	150	LYS	2.3
1	D	23	LYS	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	107	PRO	2.3
1	A	108	GLU	2.3
1	D	237	SER	2.2
1	B	56	LEU	2.1
1	B	124	SER	2.1
1	A	145	ASP	2.1
1	C	25	ARG	2.1

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 carbohydrates in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	94G	B	401	21/21	0.92	0.14	38,52,79,81	1
2	94G	C	401	21/21	0.93	0.14	21,45,79,84	1
2	94G	A	401	21/21	0.94	0.17	31,46,82,84	1
2	94G	D	401	21/21	0.95	0.18	23,52,104,105	1

6.5 Other polymers [i](#)

There are no such residues in this entry.